

The nuclear Cooper pair

Structure and Reactions

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November 5, 2020

Preface

The elementary modes of nuclear excitation are vibrations and rotations, single-particle motion, and pairing vibrations and rotations. The specific reactions probing these modes are inelastic scattering and Coulomb excitation, single- and two-particle transfer processes respectively.

The interweaving of the elementary modes of excitation leads to the renormalization of energy, radial wavefunction and particle content of the single-particles. Also to the energy, width and collectivity of vibrations and rotations. This implies renormalization of the formfactors and transition densities, Q -value and effective deformation parameters both in 3D- and in gauge-space, and state and mass number dependence of the optical potential. As a consequence, the emergence of a variety of properties, like generalized rigidity and long range correlations. Also of resonant phenomena as a function of the bombarding energy of the projectile inducing the anelastic and/or transfer processes, implying the need to go beyond lowest order distorted wave Born approximation, namely to second order DWBA.

Within this context one can posit that nuclear structure (bound) and reactions (continuum) are but two aspects of the same physics. Even more so concerning the study of light exotic halo nuclei, in which case the distinction between bound and continuum states is almost completely blurred. This is also the reason why these two aspects of nuclear physics are treated in the present monograph on equal footing, within the framework of a unified nuclear field theory of structure and reactions (NFT)_(s+r).

This theory provides the (graphical) rules to diagonalize in a compact and economic way the nuclear Hamiltonian for both bound and continuum states. It does so in terms of Feynman diagrams which describe the coupling of elementary modes of excitation, correcting for the overcompleteness of the basis (structure) and for the non-orthogonality of the scattering states (reaction), as well as for Pauli principle violation. The outcome connects directly with observables: absolute reaction cross sections and decay probabilities.

In other words (NFT)_(s+r) focuses on the scattering amplitudes which determine the absolute cross sections for the variety of physical processes, involving also those in which (quasi) bosons and fermions are created or annihilated, connecting such processes to formfactors and transition densities. Processes where one set of particles with given energies, momenta, angular momenta, etc. go in and another group (or the same), comes out. That is, as it happens in the laboratory, let alone

in nature.

Pairing vibrations and rotations, closely connected with nuclear superfluidity are paradigms of quantal nuclear phenomena. They thus play an important role within the field of nuclear structure. It is only natural that two-nucleon transfer plays a similar role concerning the probing of the nucleus.

At the basis of fermionic pairing phenomena one finds Cooper pairs, weakly bound, largely extended, strongly overlapping (quasi-) bosonic entities, made out of pairs of nucleons dressed by collective vibrations and interacting mainly through the exchange of these vibrations as well as through the bare NN -interaction, eventually corrected by $3N$ contributions. Cooper pairs not only change the statistics of the nuclear stuff around the Fermi surface and, condensing, the properties of nuclei close to their ground state. They also display a rather remarkable mechanism of tunneling between target and projectile in direct two-nucleon transfer reaction.

Cooper pair partners being weakly bound ($\ll \epsilon_F$, Fermi energy), are correlated over distances (correlation length) much larger than nuclear dimensions ($\gg R$, nuclear radius). On the other hand, Cooper pairs –building blocks of the so-called abnormal (pair) density– are forced to be confined within regions in which normal, single-particle density is present and thus, within nuclear dimensions. Said it differently, the mean field acts on Cooper pairs, as a strong external field, distorting their spatial structure.

The correlation length paradigm comes into evidence, for example, when two nuclei are set into weak contact in a direct reaction. In such a case, the partner nucleons of a Cooper pair have a finite probability to be confined each within the mean field of a different nucleus, equally well pairing correlated than when both nucleons are in the same nucleus. It is then natural that a Cooper pair can also tunnel between target and projectile, equally well correlated, through simultaneous than through successive transfer processes. Because of the weak binding of the Cooper pair, let alone the fact that tunneling probability falls off with increasing mass, successive is the dominant transfer process.

Although one does not expects supercurrents in nuclei, one can study long-range pairing correlations in terms of individual quantal states and of the tunneling of single Cooper pairs. Such weak coupling Cooper pair transfer reminds the tunneling mechanism of electronic Cooper pairs across a barrier (e.g. a dioxide layer of dimensions much smaller than the correlation length) separating two superconductors, known as a Josephson junction¹. The main difference is that, as a rule, in the nuclear time dependent junction transiently established in direct two-nucleon transfer process, only one or even none of the two weakly interacting nuclei are, as a rule, superfluid. On the other hand in nuclei, paradigmatic example of fermionic quantum finite many-body system, zero point fluctuations (ZPF) in general, and those associated with pair addition and pair subtraction modes known as pairing vibrations in particular, are much stronger than in condensed matter. Thus, pairing correlations based on even a single Cooper pair can lead to distinct pairing

¹Josephson (1962); Anderson (1964b).

correlation effects in two-nucleon transfer processes.

Nucleonic Cooper pair tunneling has played and is playing a central role in the probing of these subtle quantal phenomena, both in the case of light exotic nuclei as well as of medium and heavy nuclei lying along the stability valley. They have been instrumental in shedding light on the subject of pairing in nuclei at large, and on nuclear superfluidity in particular. Consequently, and as already said, the subject of two-nucleon transfer occupies a central place in the present monograph. Both concerning the conceptual and the computational aspects of the description of nuclear pairing, as well as regarding the quantitative confrontation of the theoretical results with the experimental findings, in terms of absolute differential cross sections.

Concerning exotic nuclei, recent experiments carried out at TRIUMF (Canada) have provided, through the magnifying glass of $(NFT)_{(s+r)}$, a microscopic view of what can be considered a unique embodiment of Copper's pair model²: a pair of fermions (neutrons) moving in time reversal states on top of a Fermi surface and interacting mainly through the exchange of a long wavelength vibration (phonon)³, leading to a barely bound system. The two neutrons give rise to an isotropic halo. Because the vibration (phonon) results from the sloshing back and forth of the neutron halo against the nucleons of the core moving in phase, one is in presence of a realization of Nambu's tumbling⁴ or, more precisely, symbiotic mechanism of spontaneously broken symmetry in gauge space.

Regarding the case of medium heavy nuclei lying along the stability valley, recent studies of heavy ion reactions between superfluid nuclei carried out at energies around and below the Coulomb barrier at the National Laboratory of Legnaro (Italy) have provided a measure of the neutron Cooper pair correlation length. Within this context, in the present monograph interdisciplinarity⁵ is used as a tool to attack concrete nuclear problems. But also, making use of the unique laboratory provided by the finite quantum many-body system of which the atomic nucleus is a paradigmatic example⁶, to shed light on condensed matter results, in terms of analogies involving individual, quantal single-particle states, let alone tunneling of single Cooper pairs.

Because of the central role the interweaving of the variety of elementary modes of nuclear excitation, namely single-particle motion and collective vibrations and rotations play in nuclear superfluidity, the study of Cooper pair tunneling in nuclei requires a consistent description of nuclear structure in terms of dressed quasiparticles and, making use of the resulting renormalized wavefunctions (formfactors), of one-nucleon transfer processes⁷. This is similar to the situation encountered in su-

²Cooper (1956).

³Fröhlich (1952); Bardeen and Pines (1955); Bardeen et al. (1957a,b).

⁴Nambu (1991).

⁵Quoting de Gennes (1974): "...what a theorist can and should systematically introduce is comparison with other fields".

⁶Broglia (2019) (Overview).

⁷Within this context one recognizes the difficulties of extracting spectroscopic factors from ex-

perconductors, in connection with strongly coupled systems⁸, and experimentally studied through one-electron tunneling experiments⁹

Thus, in the present monograph the general physical arguments and technical computational details concerning the calculation of absolute one-and two nucleon transfer differential cross sections, making use of state of the art NFT structure input, are discussed in detail.

As a result of this approach, theoretical and experimental nuclear practitioners can use the present monograph at profit. To help this use, the basic nuclear structure formalism, in particular that associated with single-particle and collective motion in both normal and superfluid nuclei, is economically introduced through general physical arguments. This is also in keeping with the availability in the current literature, of detailed discussions of the corresponding material. Within this context, the monographs *Nuclear Superfluidity* by Brink and Broglia and *Oscillations in Finite Quantum Systems* by Bertsch and Broglia, published also by Cambridge University Press, can be considered companion volumes to the present one. Volume which shares with those a similar aim: to provide a broad physical view of central issues in the study of finite quantal many-body nuclear systems accessible to motivated students and practitioners. However, neither the present one, nor the other two are introductory texts. In particular the present one in which an attempt at unifying structure and reactions as it happens in nature, is made. On the other hand, unifying discrete (mainly structure) and continuum (mainly reactions) configuration spaces, implies that one will be dealing with those structure results which can be tested by means of experiment. A fact which makes the subject of the present monograph a chapter of quantum mechanics, and thus opened to a wide range of practitioners¹⁰.

Concerning the notation, we have divided each chapter into sections. Each section may, in turn, be broken down into subsections. Equations and Figures

periment, in terms of single-particle transfer cross sections calculated making use of mean field wavefunctions.

⁸Eliashberg (1960).

⁹Giaever (1973).

¹⁰Within this context let us mention the intimately correlated subjects of Random Phase Approximation (RPA) and Particle Vibration Coupling (PVC) not found in a fourth year curriculum. They are explained and refer to in a number of places throughout the present monograph, starting from a pedestrian level and for both surface (particle-hole) and pairing (particle-particle and hole-hole) vibrations (Sects. 1.2 and 1.3), and then extended to include further details and facets (see Sects. 1.5, 1.6 and 2.3 Fig. (2.3.1), Sect. 2.3.1 (Fig. 2.3.9), App. 2.B (Fig. 1.2.3 (inset)), Sect. 3.1 (Fig. 3.1.6). Furthermore, in the case of RPA of pairing vibrations around the closed shell system ^{208}Pb , one provides in Sect. 3.5, detailed documentation of the numerical calculation of the associated wavefunctions (X - and Y -amplitudes) at the level of an exercise in a fourth year course. A similar situation is encountered in connection with the subject of the Distorted Wave Born Approximation (DWBA), again a subject not found in fourth year curricula. It is treated at the pedestrian level in Apps. 5.6 and 6.4 in connection with one-particle and two-particle transfer reactions respectively. And, once more in full detail, without eschewing complexities, but again in the style of an example-exercise (“Sometimes one has to say difficult things, but one ought to say them as simply as one know-how” (G. H. Hardy, *A mathematician’s apology*, Cambridge University Press, Cambridge (1969))) in Sects. 5.1 and 6.2.

are identified by the number of the chapter and that of the section. Thus (5.1.33) labels the thirtythird equation of section 1 of chapter 5. Similarly, Fig. 5.1.2 labels the second figure of section 1 of chapter 5. Concerning the Appendices, they are labelled by the chapter number and by a Latin letter in alphabetical order, e.g. App. 2.A, App. 2.B, etc. Concerning equations and Figures, a sequential number is added. Thus (2.B.2) labels the second equation of Appendix B of chapter 2, while Fig. 2.B.1 labels the first figure of Appendix B of Chapter 2. References are called in terms of the author's surname and publication year and are found in alphabetic order in the bibliography at the end of the monograph.

A methodological approach used in the present monograph concerns a certain degree of repetition. Similar, but not the same issues are dealt with more than once using different but equatable terminologies. This approach reflects the fact that useful concepts like reaction channels, or correlation length, let alone elementary modes of excitation, are easy to understand but difficult to define¹¹. This is because their validity is not exhausted by a single perspective. But even more important, because their power in helping at connecting¹² seemingly unrelated results and phenomena is difficult to be fully appreciated the first time around, spontaneous symmetry breaking and associated emergent properties providing an example of this fact.

Throughout, a number of footnotes are found. This is in keeping with the fact that footnotes can play a special role within the framework of an elaborated presentation. In particular, they are useful to emphasize relevant issues in an economic way. Being outside the main text, they give the possibility of stating eventual important results, without the need of elaborating on the proof, but referring to the corresponding sources. Within this context, and keeping the natural distances, one can mention that in the paper in which Born¹³ introduces the probabilistic interpretation of Schrödinger's wavefunction, the fact that this probability is connected with its [modulus] squared and not with the wavefunction itself, is only referred to in a footnote.

A large fraction of the material contained in this monograph have been the subject of the lectures of the fourth year course "Nuclear Structure Theory" which RAB delivered throughout the years at the Department of Physics of the University of Milan, as well as at the Niels Bohr Institute (Copenhagen) and at Stony Brook (State University of New York). Part of it was also presented by the authors in the

¹¹"Gentagelsen den er virkeligheden, og Tilværelsens Alvor" (Repetition is the reality and the seriousness of life: S. Kierkegaard Gjentagelsen (1843)).

¹²"The concepts and propositions get "meaning" viz. "content", only through their connection with sense-experiences... The degree of certainty with which this connection, viz., intuitive combination, can be undertaken, and nothing else, differentiates empty fantasy from scientific "truth"... A correct proposition borrows its "truth" from the truth-content of the system to which it belongs" (A. Einstein, Autobiographical notes, in Albert Einstein, Ed. P. A. Schilpp, Vol I, Harper, New York (1951) p. 13.).

¹³Born (1926a). Within this context, it is of notice that the extension of Born probabilistic interpretation to the case of many-particle systems is also found in a footnote (Pauli (1927), footnote on p. 83 of the paper). Source of information Pais (1986).

course Nuclear Reactions held at the PhD School of Physics of the University of Milan.

GP wants to thank the tutoring of Ben Bayman concerning specific aspects of two-particle transfer reactions. Discussions with Ian Thompson and Filomena Nunes on a variety of reaction subjects are gratefully acknowledged. RAB acknowledges the essential role the collaboration with Francisco Barranco and Enrico Vigezzi has played concerning the variety of nuclear structure aspects of the present monograph. His debt with the late Aage Winther regarding the reaction aspects of it is difficult to express in words. The overall contributions of Daniel Bès, Ben Bayman and Pier Francesco Bortignon¹⁴ are only too explicitly evident throughout the text and constitute a daily source of inspiration. G. P. and R. A. B. have received important suggestions and comments regarding concrete points and the overall presentation of the material discussed below from Ben Bayman, Pier Francesco Bortignon, David Brink, Willem Dickhoff and Vladimir Zelevinsky and are here gratefully acknowledged. We are specially beholden to Elena Litvinova and Horst Lenske for much constructive criticism and suggestions.

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Chapter 1

Introduction

Toute théorie physique est fondée sur l'analogie qu'on établi entre des choses malconnues et des choses simples.

Simone Weil

1.1 Views of the nucleus

In the atom, the nucleus provides the Coulomb field in which negatively charged electrons ($-e$) move independently of each other in single-particle orbitals. The filling of these orbitals explains Mendeleev's periodic table. Thus the valence of the chemical elements as well as the particular stability of the noble gases associated with the closing of shells (2(He), 10(Ne), 18(Ar), 36(Kr), 54(Xe), 86(Ra)). The dimension of the atom is measured in angstroms ($\text{\AA}=10^{-8}\text{cm}$), and typical energies in eV, the electron mass being $m_e \approx 0.511 \text{ MeV}$ ($\text{MeV}=10^6\text{eV}$).

The atomic nucleus is made out of positively charged protons ($+e$) and of (uncharged) neutrons, nucleons, of mass $\approx 10^3 \text{ MeV}$ ($m_p = 938.3 \text{ MeV}$, $m_n = 939.6 \text{ MeV}$). Nuclear dimensions are of the order of few fermis ($\text{fm}= 10^{-13} \text{ cm}$). While the stability of the atom is provided by a source external to the electrons, namely the atomic nucleus, this system is self-bound as a result of the strong interaction of range $a_0 \approx 0.9 \text{ fm}$ and strength $v_0 \approx -100 \text{ MeV}$ acting among nucleons.

1.1.1 The liquid drop and the shell models

While most of the atom is empty space, the density of the atomic nucleus is conspicuous ($\rho = 0.17 \text{ nucleon/fm}^3$). The “closed packed” nature of this system implies a short mean free path as compared to nuclear dimensions. This can be estimated from classical kinetic theory $\lambda \approx (\rho\sigma)^{-1} \approx 1 \text{ fm}$, where $\sigma \approx 2\pi a_0^2$ is the nucleon–nucleon cross section. It seems then natural to liken the atomic nucleus to

a liquid drop¹. This picture of the nucleus provided the framework to describe the basic features of the fission process².

The leptodermic properties of the atomic nucleus are closely connected with the semi-empirical mass formula³.

$$m(N, Z) = (Nm_n + Zm_p) - \frac{1}{c^2}B(N, Z), \quad (1.1.1)$$

the binding energy being

$$B(N, Z) = \left(b_{vol}A - b_{surf}A^{2/3} - \frac{1}{2}b_{sym}\frac{(N-Z)^2}{A} - \frac{3}{5}\frac{Z^2e^2}{R_c} \right). \quad (1.1.2)$$

The first term is the volume energy representing the binding energy in the limit of large A , for $N = Z$ and in the absence of the Coulomb interaction ($b_{vol} \approx 15.6$ MeV). The second term represents the surface energy, where

$$b_{surf} = 4\pi r_0^2 \gamma. \quad (1.1.3)$$

The nuclear radius is written as $R = r_0 A^{1/3}$, with $r_0 = 1.2$ fm, the surface tension energy being $\gamma \approx 0.95$ MeV/fm². The third term in (1.1.2) is the symmetry term which reflects the tendency towards stability for $N = Z$, with $b_{sym} \approx 50$ MeV. The symmetry energy can be divided into a kinetic and a potential energy part. A simple estimate of the kinetic energy part can be obtained by making use of the Fermi gas model which gives $(b_{sym})_{kin} \approx (2/3)\epsilon_F \approx 25$ MeV ($\epsilon_F \approx 36$ MeV). Consequently,

$$V_1 = (b_{sym})_{pot} = b_{sym} - (b_{sym})_{kin} \approx 25 \text{ MeV}. \quad (1.1.4)$$

The last term of (1.1.2) is the Coulomb energy corresponding to a uniformly charged sphere of radius $R_c = 1.25 A^{1/3}$ fm.

When in a heavy-ion reaction two nuclei come within the range of the nuclear forces, the Coulomb trajectory of relative motion will be changed by the attraction which will act between the nuclear surfaces. This surface interaction is a fundamental quantity in heavy ion reactions. Assuming two spherical nuclei at a relative distance $r_{AA} = R_a + R_A$, where R_a and R_A are the corresponding half-density radii, the (maximum) force acting between the two surfaces is

$$\left(\frac{\partial U_{AA}^N}{\partial r} \right)_{r_{AA}} = 4\pi\gamma \frac{R_a R_A}{R_a + R_A} \quad (1.1.5)$$

This result allows for the calculation of the ion-ion (proximity) potential which, supplemented with a position dependent absorption, can be used to accurately describe heavy ion reactions⁴.

¹Bohr and Kalckar (1937).

²Meitner and Frisch (1939); Bohr and Wheeler (1939).

³Weizsäcker (1935).

⁴See e.g. Broglia and Winther (2004) and refs. therein.

In such reactions, not only elastic processes are observed, but also anelastic reactions in which one, or both surfaces of the interacting nuclei are set into vibration (Fig. 1.1.1). The restoring force parameter associated with oscillations of multipolarity λ is

$$C_\lambda = (\lambda - 1)(\lambda + 2)R_0^2\gamma - \frac{3}{2\pi} \frac{\lambda - 1}{2\lambda + 1} \frac{Z^2 e^2}{R_c}, \quad (1.1.6)$$

where the second term corresponds to the contribution of the Coulomb energy to C_λ . Assuming the flow associated with surface vibrations to be irrotational, the associated inertia for small amplitude oscillations is,

$$D_\lambda = \frac{3}{4\pi} \frac{1}{\lambda} A M R^2, \quad (1.1.7)$$

the energy of the corresponding mode being

$$\hbar\omega_\lambda = \hbar \sqrt{\frac{C_\lambda}{D_\lambda}}. \quad (1.1.8)$$

The label λ stands for the angular momentum of the vibrational mode. Furthermore, the vibrations can be characterized by the parity quantum number $\pi = (-1)^\lambda$ and the third component of λ , denoted μ . Aside from λ, μ , surface vibrations can also be characterized by an integer $n (= 1, 2, \dots)$, an ordering number indicating increasing energy. For simplicity, a single common label α will also be used.

A picture apparently antithetic to that of the liquid drop, the shell model, emerged from the study of experimental data, plotting them against either the number of protons (atomic number), or the number of neutrons in nuclei, rather than against the mass number. One of the main nuclear features which led to the development of the shell model was the study of the stability and abundance of nuclear species and the discovery of what are usually called magic numbers⁵. What makes a number magic is that a configuration of a magic number of neutrons, or of protons, is unusually stable whatever the associated number of other nucleons is⁶.

The strong binding of a magic number of nucleons and weak binding for one more, reminds the results concerning the atomic stability of rare gases. In the nuclear case, the spin-orbit coupling plays an important role, as can be seen from the level scheme shown in Fig. 1.1.2, obtained by assuming that nucleons move independently of each other in an average potential of spherical symmetry.

A closed shell, or a filled level, has angular momentum zero. Thus, nuclei with one nucleon outside (missing from) closed shell, should have the spin and parity of the orbital associated with the odd nucleon (-hole), a prediction confirmed by the data (available at that time) throughout the mass table. Such a picture implies that the nucleon mean free path is large compared to nuclear dimensions.

⁵Elsasser (1933); Mayer (1948); Haxel et al. (1949).

⁶Mayer (1949); Mayer and Teller (1949).

The systematic studies of the binding energies leading to the shell model found also, that the relation (1.1.2) had to be supplemented to take into account the fact that nuclei with A odd, i.e. with an odd number of either protons or neutrons, are energetically unfavored compared with the neighbors even-even ones by a quantity of the order of $\delta \approx 33 \text{ MeV}/A^{3/4}$ called the pairing energy⁷ and found at the basis of the odd-even staggering effect.

1.1.2 Nuclear excitations

In addition to the quantum numbers λ , μ and π , one can characterize nuclear excitations by additional quantum numbers such as isospin τ and spin σ . Furthermore one can assign a particle (baryon or transfer) quantum number β . For a nucleon moving above the Fermi surface one has $\beta = +1$, while for a hole in the Fermi sea $\beta = -1$. For (quasi-) bosonic excitations, $\beta = 0$ for a mode associated with e.g. surface oscillations, which can also be viewed as a correlated particle-hole excitation (within this context see Fig. 1.2.3). In particular, the low-lying quadrupole and octupole vibrations of even-even nuclei (see e.g. Fig. 1.1.3) have quantum numbers $\beta = 0$, $\lambda^\pi = 2^+, 3^-$, $\tau = 0$ (protons and neutrons oscillate in phase) and $\sigma = 0$ (no spin-flips in the excitation).

For modes which involve the addition or subtraction of two correlated nucleons to the nucleus, $\beta = +2$ (Fig. 1.3.1) and $\beta = -2$ respectively. The excitation which, around closed shells, connects the ground state of an even nucleus, to the ground state of the next even nucleus, a monopole pairing vibration (i.e. $\lambda^\pi = 0^+$, $\beta = +2$) is of this type (Fig. 1.3.2). Multipole pairing vibrations with quantum numbers $\beta = \pm 2$ and $\lambda^\pi = 2^+, 4^+ \dots$, have also been observed throughout the mass table⁸.

The low-lying excited state of closed shell nuclei can be interpreted as a rule, as a harmonic quadrupole or octupole collective surface vibration (Fig. 1.1.3) described by the collective Hamiltonian⁹

$$H_{coll} = \sum_{\lambda\mu} \left(\frac{1}{2D_\lambda} |\hat{\Pi}_{\lambda\mu}|^2 + \frac{C_\lambda}{2} |\hat{\alpha}_{\lambda\mu}|^2 \right) \quad (1.1.9)$$

Following Dirac (1930) one can describe the oscillatory motion introducing boson creation (annihilation) operator $\Gamma_{\lambda\mu}^\dagger$ ($\Gamma_{\lambda\mu}$) obeying

$$\left[\Gamma_\alpha, \Gamma_{\alpha'}^\dagger \right] = \delta(\alpha, \alpha'), \quad (1.1.10)$$

⁷ Mayer and Jensen (1955) p.9. Connecting with further developments associated with the BCS theory of superconductivity (Bardeen et al. (1957a,b)) and its extension to the atomic nucleus (Bohr et al. (1958)), the quantity δ is identified with the pairing gap Δ parametrized according to $\Delta = 12 \text{ MeV}/\sqrt{A}$ (Bohr and Mottelson (1969)). It is of notice that for typical superfluid nuclei like ^{120}Sn , the expression of δ leads to a numerical value which can be parametrized as $\delta \approx 33 \text{ MeV}/(A^{1/4} \times A^{1/2}) \approx 10 \text{ MeV}/\sqrt{A}$.

⁸ See e.g. Flynn et al. (1971); Brink and Broglia (2005) Ch. 5.

⁹ Classically $\Pi_{\lambda\mu} = D_\lambda \dot{\alpha}_{\lambda\mu}$.

and leading to,

$$\hat{a}_{\lambda\mu} = \sqrt{\frac{\hbar\omega_\lambda}{2C_\lambda}} (\Gamma_{\lambda\mu}^\dagger + (-1)^\mu \Gamma_{\lambda-\mu}), \quad (1.1.11)$$

a similar expression being valid for the conjugate momentum variable $\hat{\Pi}_{\lambda\mu}$, resulting in

$$\hat{H}_{coll} = \sum_{\lambda\mu} \hbar\omega_\lambda ((-1)^\mu \Gamma_{\lambda\mu}^\dagger \Gamma_{\lambda-\mu} + 1/2). \quad (1.1.12)$$

The frequency of the mode is $\omega_\lambda = (C_\lambda/D_\lambda)^{1/2}$, while $(\hbar\omega_\lambda/2C_\lambda)^{1/2}$ is the amplitude of the zero-point fluctuation of the vacuum state $|0\rangle_B, |n_{\lambda\mu} = 1\rangle = \Gamma_{\lambda\mu}^\dagger |0\rangle_B$ being the one-phonon state. To simplify the notation, in many cases one writes $|n_\alpha = 1\rangle$.

The ground and low-lying states of nuclei with one nucleon outside closed shell can be described by the single-particle Hamiltonian

$$H_{sp} = \sum_\nu \epsilon_\nu a_\nu^\dagger a_\nu, \quad (1.1.13)$$

where $a_\nu^\dagger(a_\nu)$ is the single-particle creation (annihilation) operator,

$$|\nu\rangle = a_\nu^\dagger |0\rangle_F, \quad (1.1.14)$$

being the single-particle state of quantum numbers $\nu (\equiv nljm$, namely number of nodes, orbital and total angular momentum, and its third component) and energy ϵ_ν , $|0\rangle_F$ being the Fermion vacuum. It is of notice that

$$[H_{coll}, \Gamma_{\lambda'\mu'}^\dagger] = \hbar\omega_{\lambda'} \Gamma_{\lambda'\mu'}^\dagger \quad (1.1.15)$$

and

$$[H_{sp}, a_{\nu'}^\dagger] = \epsilon_{\nu'} a_{\nu'}^\dagger. \quad (1.1.16)$$

This outcome results from the bosonic

$$[\Gamma_\alpha, \Gamma_{\alpha'}^\dagger] = \delta(\alpha, \alpha'), \quad (1.1.17)$$

and fermionic

$$\{a_\nu, a_{\nu'}^\dagger\} = \delta(\nu, \nu'), \quad (1.1.18)$$

commutation (anti-commutation) relations.

The existence of drops of nuclear matter displaying both collective surface vibrations and independent-particle motion are emergent properties not contained in the particles forming the system, neither in the forces acting among them.

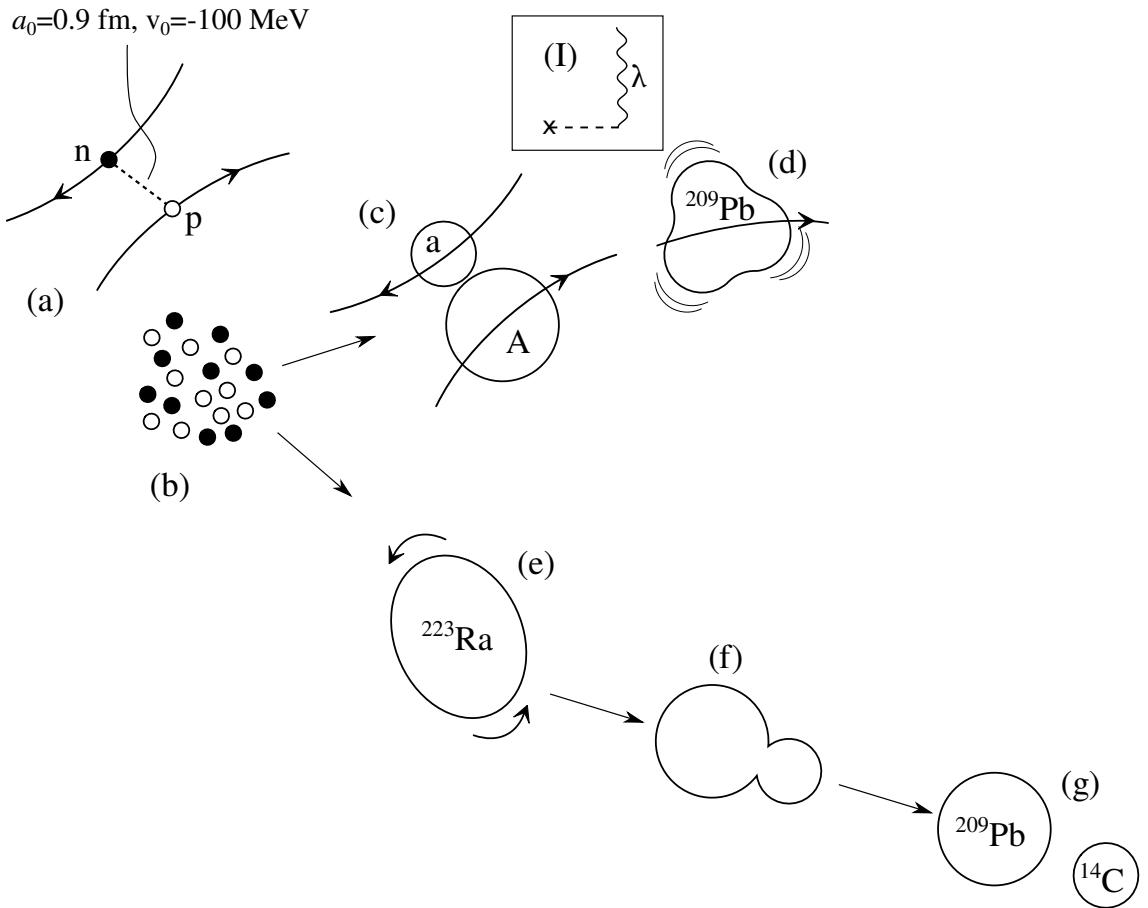


Figure 1.1.1: Emergent properties (collective modes) (a) Nucleon-Nucleon (NN) interaction in a scattering experiment; emergent properties (collective nuclear modes) (b) assembly of a swarm of nucleons condensing into drops of nuclear matter, examples shown in (c) and (e); (c) anelastic heavy ion reaction $a + A \rightarrow a + A^*$ setting the nucleus A into an octupole surface oscillations (d); in inset (I) the time-dependent nuclear plus Coulomb fields associated with the reaction (c) is represented by a cross followed by a dashed line, while the wavy line labeled λ describes the propagation of the $\lambda^\pi = 3^-$ surface vibration schematically shown in (d), time running upwards; (e) another possible outcome of nucleon condensation: the (weakly) quadrupole deformed nucleus ^{223}Ra which can rotate as a whole with a moment of inertia considerably smaller than the rigid moment of inertia, a fact intimately connected with the role played by pairing in nuclei, effect which becomes overwhelming in the phenomenon of exotic decay displayed in (f) in which the nuclear surface zero point fluctuations (quadrupole ($\lambda = 2$), octupole ($\lambda = 3$), etc.) can get, with a small but finite probability ($P \approx 10^{-10}$), spontaneously in phase and produce a neck-in (saddle conformation) leading eventually to the (exotic) decay mode $^{223}\text{Ra} \rightarrow ^{209}\text{Pb} + ^{14}\text{C}$, as experimentally observed (g) (Rose and Jones (1984), see Brink and Broglia (2005), Ch. 7 and refs. therein). As correctly explained in Matsuyanagi et al. (2013) in general, and valid also in the case of the ZPF leading to the saddle (neck-in) configuration, such fluctuations are associated with genuine quantum vibrations essentially different in character from surface oscillations of a classical liquid drop. The intimate connection between pairing and collective vibrations reveals itself through the inertial masses governing the collective kinetic energies.

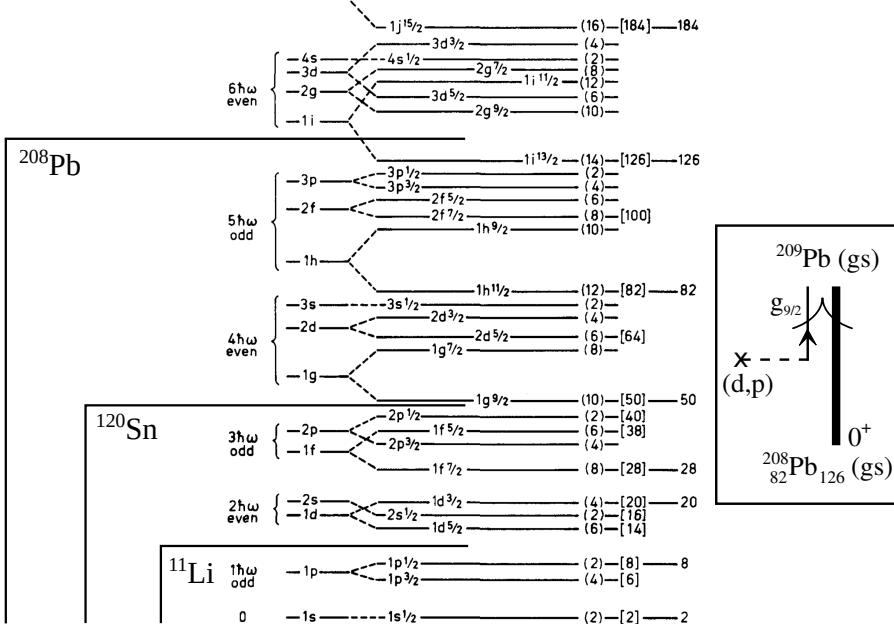


Figure 1.1.2: Sequence of levels of the harmonic oscillator potential labeled with the principal oscillator quantum number ($N(\hbar\omega) = 0(\hbar\omega), 1(\hbar\omega), 2(\hbar\omega), \dots$) the parity being $\pi = (-1)^N$). The next column shows the splitting of major shell degeneracies obtained using a more realistic potential (Woods-Saxon), the quantum number being the number of radial nodes of the associated single-particle s, p, d , etc., states. The levels shown at the center result when a spin-orbit term is also considered, the quantum numbers nlj characterizing the states of degeneracy $(2j+1)$ ($j = l \pm 1/2$). To the left we schematically (in particular in the case of Li which displays non Meyer and Jensen sequence) indicate the Fermi energy associated with a light (exotic), medium, and heavy nucleus, namely ^{11}Li , ^{120}Sn and ^{208}Pb . In the inset, a schematic graphical representation of the reaction $^{208}\text{Pb}(d, p)^{209}\text{Pb}(\text{gs})$ is shown. A cross followed by a horizontal dashed line represents, in the present case, the (d, p) field, while a single arrowed line describes the odd nucleon moving in the $g_{9/2}$ orbital above the $N = 126$ shell closure drawn as a bold line labeled 0^+ .

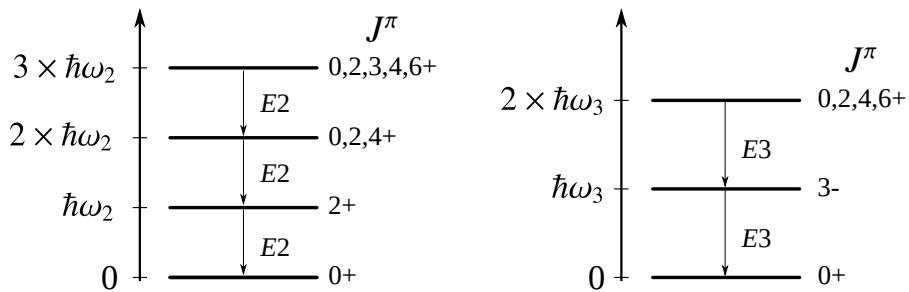


Figure 1.1.3: Schematic representation of harmonic quadrupole and octupole liquid drop collective surface vibrational modes.

Expressed it differently, generalized rigidity closely connected to the inertial parameter D_λ implies that acting on a nucleus with an external $\beta = 0$, time-dependent (nuclear/Coulomb) field, the system reacts as a whole (collective vibrations; also rotations see Sect. 1.4), while acting with fields which change particle number by one ($\beta = \pm 1$; e.g. (d, p) and (p, d) reactions) the system reacts in terms of independent particle motion, feeling the pushings and pullings of the other nucleons only when trying to leave the nucleus. Such a behaviour can hardly be inferred from the study of the NN -forces in free space, being truly emergent properties of the finite, quantum many-body nuclear system.

Collective surface vibrations and independent particle motion are examples of what are called elementary modes of excitation in finite many-body physics, and collective variables in soft-matter physics.

1.2 The particle-vibration coupling

The oscillation of the nucleus under the influence of the surface tension implies that the potential $U(r, R)$ in which nucleons move independently of each other change with time. For low-energy collective vibrations this change is slow as compared with single-particle motion. Within this scenario the nuclear radius can be written as

$$R = R_0 \left(1 + \sum_{\lambda\mu} \alpha_{\lambda\mu} Y_{\lambda\mu}^*(\hat{r}) \right) \quad (1.2.1)$$

Assuming small amplitude motion,

$$U(r, R) = U(r, R_0) + \delta U(r), \quad (1.2.2)$$

where

$$\delta U = \kappa \hat{\alpha} \hat{F} = \Lambda_\alpha \left(\Gamma_{\lambda\mu}^\dagger + (-1)^\mu \Gamma_{\lambda-\mu} \right) \hat{F} = H_c, \quad (1.2.3)$$

with

$$\Lambda_\alpha = \kappa \sqrt{\frac{\hbar\omega_\lambda}{2C_\lambda}}, \quad (1.2.4)$$

is the particle-vibration coupling (PVC) strength (Fig. 1.2.1), product of the dynamic deformation

$$\beta_\lambda = \sqrt{2\lambda + 1} \sqrt{\frac{\hbar\omega_\lambda}{2C_\lambda}}, \quad (1.2.5)$$

and of the strength κ , while

$$\hat{F} = \sum_{\nu_1\nu_2} \langle \nu_1 | F | \nu_2 \rangle a_{\nu_1}^\dagger a_{\nu_2}, \quad (1.2.6)$$

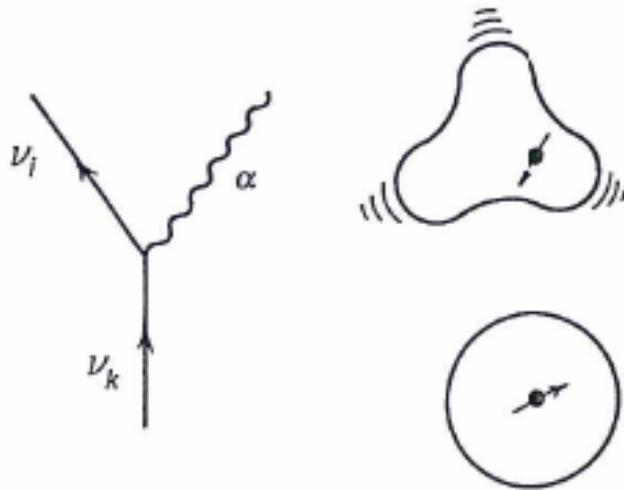


Figure 1.2.1: Graphical representation of a process by which a fermion, bouncing inelastically off the surface, sets it into vibration. Particles are represented by an arrowed line pointing upwards which is also the direction of time, while the vibration is represented by a wavy line. In the cartoon to the right, the black dot represents a nucleon moving in a spherical mean field of which it excites, through the PVC vertex, an octupole vibration after bouncing inelastically off the surface.

is a single-particle field with (dimensionless) form factor,

$$F = -\frac{R_0}{\kappa} \frac{\partial U}{\partial r} Y_{\lambda\mu}^*(\hat{r}). \quad (1.2.7)$$

An estimate of κ is given below (Eq. (1.2.13)).

Diagonalizing δU making use of the graphical (Feynman) rules of nuclear field theory (NFT) to be discussed in the following chapter, one obtains structure results which can be used in the calculation of absolute transition probabilities and differential reaction cross sections, quantities which can be compared with the experimental findings.

Within the framework of NFT, single-particles are to be calculated as the Hartree-Fock solution of the NN -interaction $v(|\mathbf{r} - \mathbf{r}'|)$ (Fig. 1.2.2, diagrams (a)–(c)) – e.g. a regularized NN -bare interaction in terms of renormalization group methods or alternative techniques (v_{low-k}), taking eventually also $3N$ terms into account¹⁰ – leading, in particular to

$$U(r) = \int d\mathbf{r}' \rho(r') v(|\mathbf{r} - \mathbf{r}'|), \quad (1.2.8)$$

¹⁰See e.g. Duguet (2013, 2004); Duguet and Lesinski (2008); Lesinski et al. (2009); Hebeler et al. (2009); Baroni et al. (2010); Duguet et al. (2010); Lesinski et al. (2011); Hergert and Roth (2009), and references therein.

that is, the Hartree field¹¹ expressing the selfconsistency between density ρ and potential U (Fig. 1.2.2 (b)), while vibrations are to be calculated in the Random Phase Approximation (RPA) making use of the same interaction¹² (Fig. 1.2.3), extending the selfconsistency to fluctuations $\delta\rho$ of the density and δU of the mean field, that is,

$$\delta U(r) = \int d\mathbf{r}' \delta\rho(\mathbf{r}') v(|\mathbf{r} - \mathbf{r}'|). \quad (1.2.9)$$

Making use of the solution to this relation one obtains the transition density $\delta\rho$. The matrix elements $\langle n_\lambda = 1, \nu_i | \delta\rho | \nu_k \rangle$ provide the particle-vibration coupling matrix elements to work out the variety of coupling processes between single-particle and collective motion¹³ (Fig. 1.2.1). That is, the matrix elements of the PVC Hamiltonian¹⁴ H_c . However, the role of the NN -interaction v is not exhausted neither by H_{HF} nor by $H_{RPA} + H_c$. Diagonalizing

$$H = H_{HF} + H_{RPA} + H_c + v, \quad (1.2.10)$$

by applying, in the basis of single-particle and collective modes, that is solutions of H_{HF} and of H_{RPA} respectively, the graphical NFT rules (see next chapter) one obtains an exact solution of the total Hamiltonian, to the order of $1/\Omega$ of the Feynman diagrams calculated.

The quantity Ω is the effective degeneracy in which the nucleonic excitations are allowed to correlate through v to give rise to the collective modes, $1/\Omega$ being the small parameter of the NFT diagrams¹⁵ (see Sect. 2.5.4, Eq. (2.5.85)). Concerning the rules of NFT (Sect. 2.5), they codify the way in which H_c (three-point vertices)

¹¹To this potential one has to add the Fock potential resulting from the fact that nucleons are fermions (Fig. 1.2.2 (c)).

¹²Bohm and Pines (1951, 1953). The sum of the so called bubble (ring) diagrams are taken into account to infinite order in RPA. This is the reason why bubble contributions in the diagonalization of Eq. (1.2.10) are not allowed in NFT, being already contained in the basis states (see next chapter, Sect. 2.5).

¹³It is of notice that the, so-called, scattering vertex shown in Fig. 1.2.1 is not operative in RPA. Being an harmonic approximation, either two fermion lines (particle-hole) enter a vertex and a (quasi) boson line comes out (forwards going process), or two-fermion lines and a boson line come in or go out from the vertex (backwards going process). See inset Fig. 1.2.3, graphs (b) and (c) respectively, as well as Fig. 1.2.4 (a).

¹⁴Mottelson (1968), Mottelson (1969); Hamamoto (1969, 1970a,b, 1977); Bès and Broglia (1971a); Broglia et al. (1971a,b); Flynn et al. (1971), and references therein.

¹⁵According to renormalized NFT v is the NN -interaction which eventually combined with a k -mass, is used to calculate the bare single-particle states (HF-approximation), collective vibrations (RPA) and particle-vibration coupling vertices so that once the corresponding renormalization diagrams (including also four-point vertices, i.e. v) have been worked out, the resulting dressed elementary modes of excitation reproduce the experimental findings. In the case in which one is only interested in the collective vibrations for the purpose to dress the single-particle degrees of freedom, one can take them from experiment (empirical renormalization). In other words, determine the Λ -values by making use of the experimental dynamical deformation parameter $\beta_\lambda = \sqrt{\frac{\hbar\omega_\lambda}{2C_\lambda}} \frac{1}{\sqrt{2\lambda+1}}$ and energies $\hbar\omega_\lambda$ (Broglia et al. (2016)), in conjunction with the expression of the RPA amplitudes X and Y and dispersion relation displayed in caption to Fig. 1.2.3.

and v (four-point vertices) are to be treated to all orders of perturbation theory. Also which processes (diagrams) are not allowed because they will imply overcounting of correlations already included in the basis states.

As will be shown in the following chapters, NFT allows in an economic fashion to sum to infinite order weakly convergent processes as well as particularly important ones, without at the same time being forced to do the same in connection with other processes which either are rapidly convergent, or which lead to small contributions that can be neglected¹⁶.

Before proceeding, a simple estimate of the coupling strength κ is worked out, making use of a schematic separable interaction¹⁷

$$v = \kappa \hat{F} \hat{F}^\dagger \quad (1.2.11)$$

and of an expansion of the nuclear density $\rho(r, R)$ similar to (1.2.2), that is,

$$\delta\rho(\mathbf{r}) = -R_0 \frac{\partial\rho(r)}{\partial r} Y_{\lambda\mu}^*(\hat{r}). \quad (1.2.12)$$

With the help of the dynamical selfconsistent relation (1.2.9) one obtains,

$$\kappa = \int r^2 dr R_0 \frac{\partial\rho(r)}{\partial r} R_0 \frac{\partial U(r)}{\partial r}. \quad (1.2.13)$$

For attractive fields, both U and κ are negative.

Because of quantal zero point fluctuations, a nucleon propagating in the nuclear medium moves through a cloud of (quasi) bosonic virtual excitations to which it couples becoming dressed and acquiring an effective mass, charge, etc. (Fig. 1.2.4; see also Sects. 5.4 and 5.8).

Furthermore, the sharp transition expected to take place, in the independent-particle motion, between occupied ($V_i^2 = 1$, Fig. 1.2.2 (f)) and empty states ($V_k^2 = 0$) becomes blurred due to the processes (b) and (c) (Fig. 1.2.4) which dress the nucleons. This is illustrated in Fig. 1.2.5, where a schematic representation of a detailed (NFT) calculation¹⁸ of the occupation probability of neutron orbits in the closed shell system ^{208}Pb is displayed¹⁹. Similar results have been obtained in other nuclei, for example the neutron open shell nucleus ^{120}Sn (Fig. 1.1.2). Within this context let us consider the example of the $1i_{11/2}$ neutron state which, in the independent particle model of ^{208}Pb , would be unoccupied. Taking the particle vibration coupling mechanism into account, which we limit for simplicity to the consideration of $\beta = 0$, $p-h$ collective surface vibration, all those ZPF of the ^{208}Pb

¹⁶The reason for this flexibility is to be partially found in the fact that bare elementary modes of excitation used as basis states in NFT, contain an important fraction of the many-body nuclear correlations, making the diagonalization of the nuclear Hamiltonian a low-dimensional problem (see also Sect. 2.1).

¹⁷See e.g. Bohr and Mottelson (1975) Eq. (6-37).

¹⁸See Bortignon et al. (1998) pp. 77,78.

¹⁹See Sect. 5.8 as well as Mahaux et al. (1985), in particular Sect. 4.7 of this reference.

“vacuum” (see Fig. 1.2.4 (a)) which involves the $1i_{11/2}$ orbital in the structure of the collective mode would contribute to $n_{1i_{11/2}}$.

Vibrational modes can also become renormalized through the coupling to dressed nucleons which, in intermediate virtual states, can exchange the vibrations which produce their clothing, with the second fermion (hole state). Such a process leads to a renormalization of the PVC vertex²⁰ (Fig. 1.2.6), as well as of the bare NN -interaction, in particular 1S_0 component (bare pairing interaction)²¹.

The analytic procedures equivalent to the diagrammatic ones to obtain the HF and RPA solutions associated with the bare NN -interaction v are provided by the relations (1.1.16) and (1.1.15) respectively, replacing the corresponding Hamiltonians by $(T + v)$, where T is the kinetic energy operator. The phonon operator associated with surface vibrations is defined as,

$$\Gamma_\alpha^\dagger = \sum_{ki} X_{ki}^\alpha \Gamma_{ki}^\dagger + Y_{ki}^\alpha \Gamma_{ki}, \quad (1.2.14)$$

the normalization condition being,

$$[\Gamma_\alpha, \Gamma_\alpha^\dagger] = \sum_{ki} (X_{ki}^{\alpha 2} - Y_{ki}^{\alpha 2}) = 1. \quad (1.2.15)$$

The operator $\Gamma_{ki}^\dagger = a_k^\dagger a_i (\epsilon_k > \epsilon_F, \epsilon_i \leq \epsilon_F)$ creates a particle-hole excitation acting on the HF vacuum state $|0\rangle_F$. It is assumed that

$$[\Gamma_{ki}, \Gamma_{k'i'}^\dagger] = \delta(k, k')\delta(i, i'). \quad (1.2.16)$$

Within this context, RPA is a harmonic, quasi-boson approximation.

From being antithetic views of the nuclear structure, a proper analysis of the experimental data testifies to the fact that the collective and the independent particle pictures of the nuclear structure require and support each other²². To obtain a quantitative description of nucleon motion and nuclear phonons (vibrations), one needs a proper description of the k - and ω -dependent “dielectric” function of the nuclear medium, in a similar way in which a proper description of the reaction processes used as probes of the nuclear structure requires the use of the optical potential (continuum “dielectric” function).

The NFT solutions which diagonalize (1.2.10) provide all the elements to calculate the structure properties of nuclei, and also of the optical potential needed

²⁰ Bertsch et al. (1983); Barranco et al. (2004) and refs. therein. It is to be noted that in the case in which the renormalized vibrational modes, i.e. the initial and final wavy lines in Fig. 1.2.6 have angular momentum and parity $\lambda^\pi = 0^+$, and one uses a model in which there is symmetry between the particle and the hole subspaces, the four diagrams sum to zero, because of particle (gauge) conservation. A fact closely connected with Wick’s theorem of QED and generalized Ward’s identity (Ward (1950)).

²¹ See e.g. Brink and Broglia (2005) Ch. 10 and references therein.

²² Bohr and Mottelson (1975).

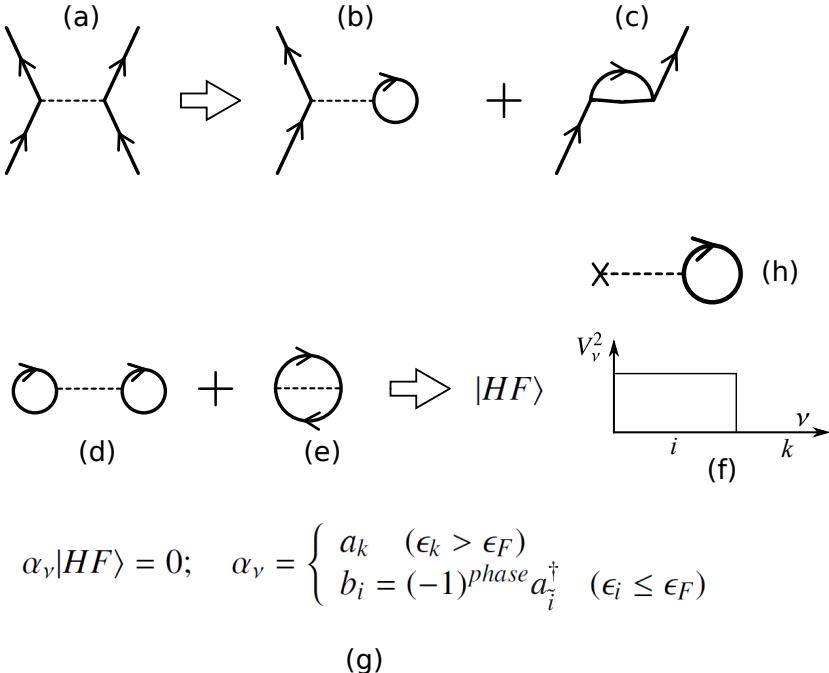


Figure 1.2.2: Schematic representation of the processes characterizing the Hartree-Fock ground state (single-particle vacuum), in terms of Feynman diagrams. (a) nucleon-nucleon interaction through the bare (instantaneous) NN -potential. (b) Hartree mean field contribution. (c) Fock mean field contribution. (d,e) ground state correlations (ZPF) associated with the Hartree and Fock fields. (f) There is, in HF (mean field) theory, a complete decoupling between occupied and empty states, labeled i and k respectively, and thus a sharp discontinuity at the Fermi energy of the occupation probability, from the value of 1 to 0. (g) This decoupling allows for the definition of two annihilation operators: $a_k(b_i)$ particle (hole) annihilation operators, implying the existence of hole (antiparticle) states ($b_i^\dagger|HF\rangle$) with quantum numbers time reversed to that of particle states, (for details see e.g. Brink and Broglia (2005) App. A). In other words, the $|HF\rangle$ ground (vacuum) state is filled up to the Fermi energy (ϵ_F) with N nucleons. The system with $(N - 1)$ nucleons can, within the language of (Feynman's) field theory, be described in terms of the degrees of freedom of that of the missing nucleon (hole-, antiparticle state). Such a description is considerably more economic than that corresponding to an antisymmetric wavefunction with $(N - 1)$ spatial and spin coordinates (\mathbf{r}_i, σ_i). Within the above scenario, a stripping reaction $N(d, p)(N + 1)$ can be viewed as the creation of a particle state ($a_k^\dagger|HF\rangle = |k\rangle$) and that of a pickup reaction $N(p, d)(N - 1)$ as that of a hole state ($b_i^\dagger|HF\rangle \equiv |\tilde{i}\rangle$). (h) Hartree, mean field contribution to the nuclear density, the density operator being represented by a cross followed by a dashed horizontal line (see also Fig. 1.8.1).

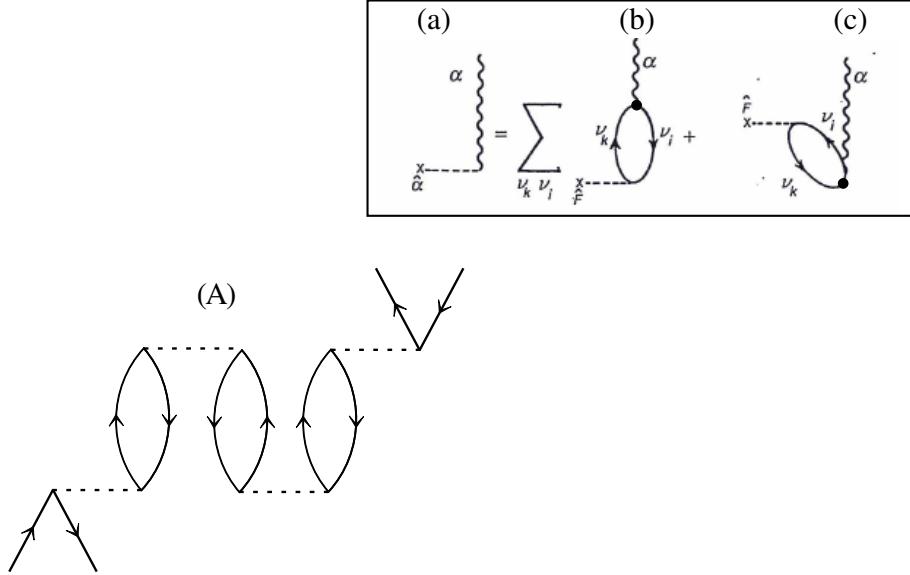


Figure 1.2.3: (A) typical Feynman diagram diagonalizing, in the harmonic approximation (RPA), the NN -interaction $v(|\mathbf{r} - \mathbf{r}'|)$ (horizontal dashed line) in a particle-hole basis provided by the Hartree-Fock solution of v . Bubbles going forward in time (inset (b); lead to the amplitudes $X_{ki}^\alpha = \frac{\Lambda_\alpha \langle \tilde{i}|F|k \rangle}{(\epsilon_k - \epsilon_i) - \hbar\omega_\alpha}$) are associated with configuration mixing of particle-hole states ($\langle \tilde{i} | = \tau |i \rangle$, τ being the time reversal operator). It is of notice that $\epsilon_k > \epsilon_F$ and $\epsilon_i \lesssim \epsilon_F$, and $\epsilon_j = \epsilon_{\tilde{j}}$ (Kramers degeneracy). Bubbles going backwards in time (inset (c), leading to the amplitudes $Y_{ki}^\alpha = -\frac{\Lambda_\alpha \langle \tilde{i}|F|k \rangle}{(\epsilon_k - \epsilon_i) + \hbar\omega_\alpha}$) are associated with zero point fluctuations (ZPF) of the ground state (term $(1/2)\hbar\omega$ for each degree of freedom in Eq. (1.1.12)). The self consistent solutions of (A), eigenstates of the dispersion relation $\sum_{ki} \frac{2(\epsilon_k - \epsilon_i)|\langle \tilde{i}|F|k \rangle|^2}{(\epsilon_k - \epsilon_i)^2 - (\hbar\omega_\alpha)^2} = 1/\kappa_\alpha$ (see inset) are represented by a wavy line (inset), that is a collective mode which can be viewed as a correlated particle (arrowed line going upward)- hole (arrowed line going downward) excitation. The quantity $\Lambda_\alpha = \kappa_\alpha \sqrt{\frac{\hbar\omega_\lambda}{2C_\lambda}}$ is the particle-vibration coupling vertex (normalization constant of the RPA amplitudes $\sum_{ki} (X_{ki}^\alpha)^2 - (Y_{ki}^\alpha)^2 = 1$). Bohm and Pines (1951, 1953), see also Brink and Broglia (2005) Sect. 8.3.

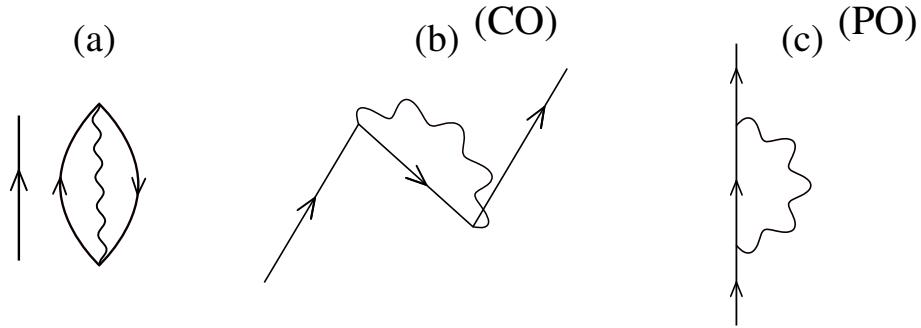


Figure 1.2.4: (a) a nucleon (single arrowed line pointing upward) moving in presence of the zero point fluctuation of the nuclear ground state associated with a collective surface vibration; (b) Pauli principle leads to a dressing process of the nucleon (so called correlation (CO) diagram) resulting from the exchange of the fermion line considered explicitly and that participating in the correlated particle-hole excitation of the ZPF; (c) time ordering gives rise to the second possible lowest order dressing process (so called polarization (PO) diagram). The above are phenomena closely related to the Lamb shift of Quantum Electrodynamics (see e.g. Weinberg (1996a) Sect. 14.3 of this reference).

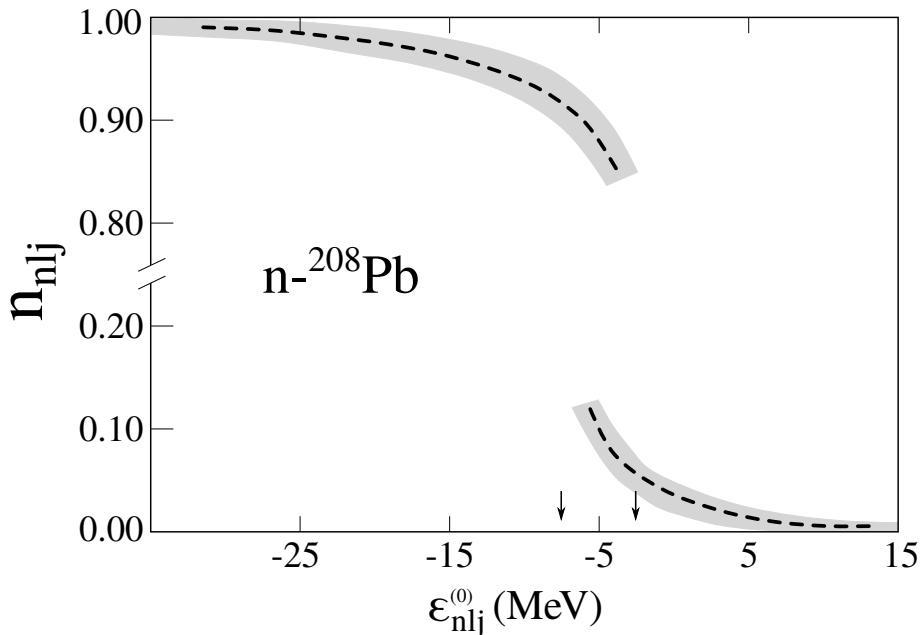


Figure 1.2.5: Schematic representation of the occupation probability (see Sect. 5.8) of ^{208}Pb neutron orbits: $1f_{7/2}, 2p_{1/2}, 1g_{7/2}, 1h_{11/2}, 1h_{9/2}, 2f_{7/2}, 2f_{5/2}, 1i_{13/2}, 3p_{3/2}, 3p_{1/2}$ hole states and $2g_{9/2}, 1i_{11/2}, 1j_{15/2}, 3d_{5/2}, 2g_{7/2}, 42_{1/2}, 2d_{3/2}, 2h_{11/2}, 2h_{9/2}$ particle states. The arrows show the location of $\epsilon_F^- = \epsilon_{3p_{1/2}}$ and $\epsilon_F^+ = \epsilon_{2g_{9/2}}$.

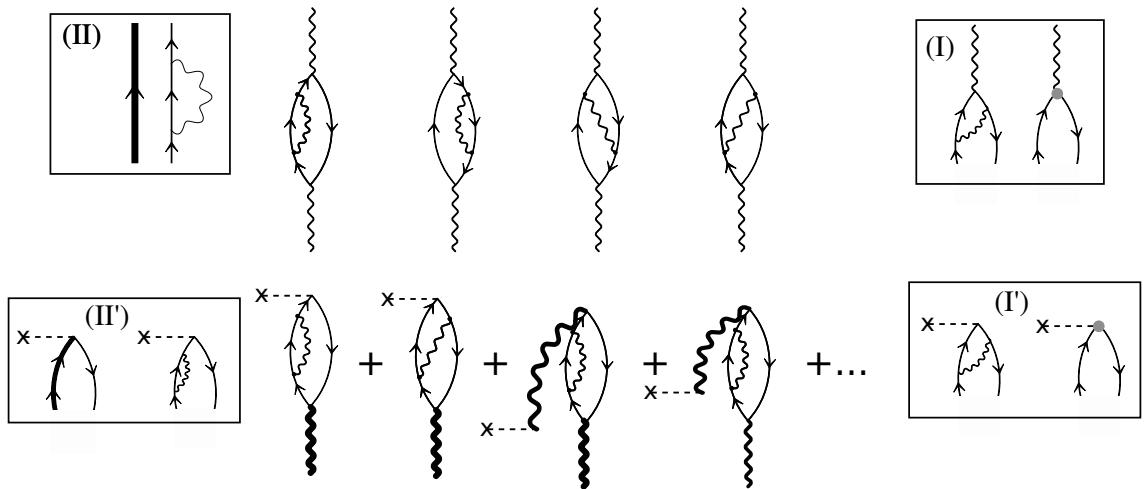


Figure 1.2.6: (Upper part) Examples of renormalization processes dressing a surface collective vibrational state. (Lower part) intervening the process with an external electromagnetic field ($E\lambda$: cross followed by dashed horizontal line; bold wavy lines, renormalized vibration of multipolarity λ) the $B(E\lambda)$ transition strength can be determined. In insets (I) and (I'), the hatched circle in the diagram to the right stands for the renormalized PVC strength resulting from the processes described by the corresponding diagrams to the left (vertex corrections). In (II') the bold face arrowed line (left diagram) represents the motion of a nucleon of effective mass m^* in a potential $(m/m^*)U(r)$, generated by the self-energy process shown to the right (see also (II) diagram to the right), $U(r)$ being the potential describing the motion of nucleons (drawn as a thin arrowed line) of bare mass m_k .

to describe nucleon-nucleus as well as nucleus-nucleus elastic scattering and reaction processes²³. Furthermore, the NFT solutions of (1.2.10) show that both single-particle (fermionic) and collective (bosonic) elementary modes of excitation emerge from the same properties of the NN -interaction. For example, a bunching of levels of the same or opposite parity just below and above the Fermi energy, imply low-lying quadrupole (e.g. $^{40}_{50}\text{Sn}_N$ -isotopes), dipole ($^{11}_3\text{Li}_8$) or octupole ($^{208}_{82}\text{Pb}_{126}$) collective vibrational modes. And, as a bonus, one has the building blocks to construct the nuclear spectrum, and bring quantitative simplicity into the experimental findings. Within this scenario, the NFT solutions of (1.2.10) also indicate the minimum set of experimental probes needed to have a “complete” picture of the nuclear structure associated with a given energy region. This is a consequence of the central role played by the quantal many-body renormalization process which interweave the variety of elementary modes of excitation. Renormalization processes which acts on par on the radial dependence of the wavefunctions (formfactors) and on the single-particle content of the orbitals involved in the reaction process under discussion (see e.g. Sect. 5.2.2). In other words, structure ad reactions are to be treated on equal footing²⁴.

The development of experimental techniques and associated hardware has allowed for the identification of a rich variety of elementary modes of excitation aside from collective surface vibrations and independent particle motion: quadrupole and octupole rotational bands, giant resonances of varied multipolarity and isospin, as well as pairing vibrations and rotation, together with giant pairing vibrations of transfer quantum number $\beta = \pm 2$. Modes which can be specifically excited in inelastic and Coulomb excitation processes, and one- and two-particle transfer reactions.

1.3 Pairing vibrations

Let us introduce this new type of elementary mode of excitation by making a parallel with quadrupole surface vibrations within the framework of RPA, namely

$$\left[(H_{sp} + H_i), \Gamma_\alpha^\dagger \right] = \hbar\omega_\alpha \Gamma_\alpha^\dagger, \quad (1.3.1)$$

²³ It is of notice that in the present monograph the subject of the optical potential is not treated. There exists a vast literature concerning this subject. See e.g. Feshbach (1958, 1962); Jackson (1970); Jeukenne et al. (1976); Sartor and Mahaux (1980); Pollarolo et al. (1983); Satchler (1983); Mahaux et al. (1985); Broglia and Winther (2004); Dickhoff and Van Neck (2005); Montanari et al. (2014); Dickhoff et al. (2017); Dickhoff and Charity (2019); Fernández-García et al. (2010a); Fernández-García et al. (2010b); Jenning (2011); Barbieri and Jennings (2005); Rotureau et al. (2017); Broglia et al. (1981) and references therein.

²⁴ Within this context, and referring to one-particle transfer reactions for concreteness, the prescription of using the ratio of the absolute experimental cross section and of the theoretical one (calculated in the Distorted Wave Born Approximation (DWBA) making use of Saxon-Woods single-particle wavefunctions as formfactors) to extract the single-particle content of the orbital under study may not be completely appropriate.

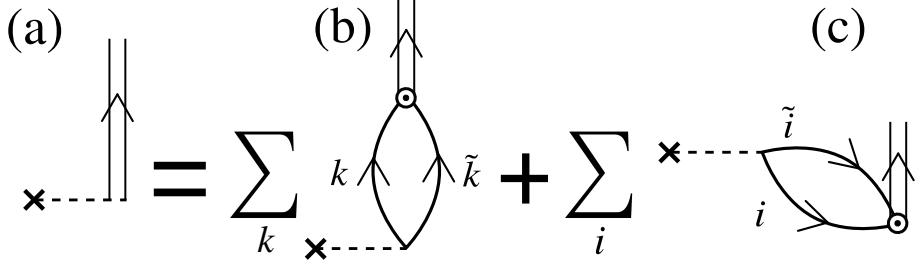


Figure 1.3.1: Graphical representation of the RPA dispersion relation describing the pair addition vibrational mode, drawn as a double arrowed line. A cross followed by a dashed horizontal line stands for (see Eq. (1.3.7)): (a) the collective operator Γ_α^\dagger , (b) the operator Γ_k^\dagger creating *a pair of nucleons moving in time reversal states* (k, \tilde{k}) *above the Fermi energy* ($\epsilon_k > \epsilon_F$); (c) The operator Γ_i annihilating *a pair of time reversal nucleon holes* (i, \tilde{i}) associated with ground state correlations ($\epsilon_i \leq \epsilon_F$).

where for simplicity we use, instead of v , a quadrupole-quadrupole separable interaction ($i = QQ$) defined as²⁵

$$H_{QQ} = -\kappa Q^\dagger Q, \quad (1.3.2)$$

with

$$Q^\dagger = \sum_{ki} \langle k | r^2 Y_{2\mu} | i \rangle a_k^\dagger a_i, \quad (1.3.3)$$

while H_{sp} and Γ_α^\dagger are defined in (1.1.13) and (1.2.14) supplemented by (1.2.15).

In connection with the pairing energy mentioned above (odd-even staggering), it is a consequence of correlation of pairs of like nucleons moving in time reversed states (v, \tilde{v}). A similar phenomenon to that found in metals at low temperatures and giving rise to superconductivity²⁶. The pairing interaction ($i = P$) can be written, within the approximation (1.3.2) used in the case of the quadrupole-quadrupole force, as

$$H_P = -GP^\dagger P, \quad (1.3.4)$$

where

$$P^\dagger = \sum_{v>0} P_v^\dagger, \quad (1.3.5)$$

²⁵It is of notice that in this case, and at variance with (1.2.11)–(1.2.13), the fact that the force is attractive is explicitly expressed through the minus sign, assuming $\kappa > 0$. This is done for didactical purposes, to connect with the standard notation for the pairing interaction given in Eq. (1.3.4) below.

²⁶Bardeen et al. (1957a,b). With regards to the extension of BCS to nuclei see Bohr et al. (1958). See also Broglia and Zelevinsky (2013).

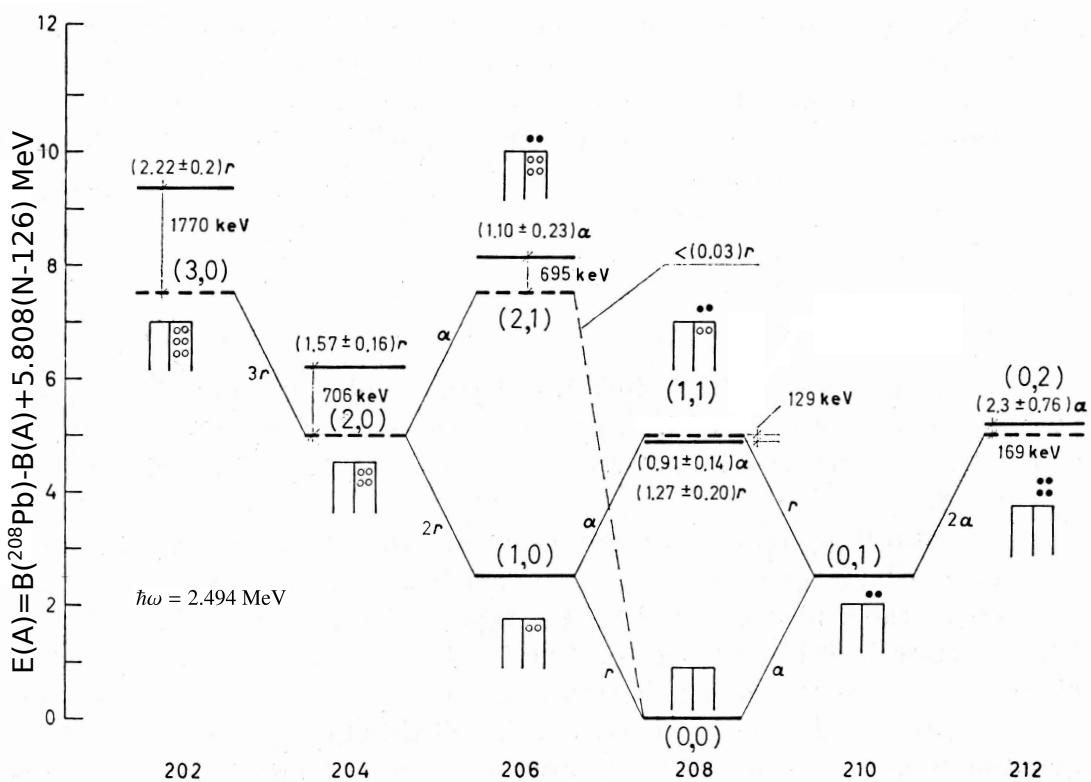


Figure 1.3.2: The many-phonon pairing vibrational spectrum around the “vacuum” $|gs(^{208}\text{Pb})\rangle$. The energies predicted by the pairing vibrational model (harmonic approximation) are displayed as dashed horizontal lines. The harmonic quantum numbers (n_r, n_a) indicating the number of pair addition and pair removal modes are displayed for each level. A schematic representation of the many-particle many-hole structure of the state is also given. The transitions predicted by the model (harmonic approximation) are indicated in units of r and a (see equation (1.3.10) and subsequent text). The corresponding experimental numbers are also given together with their errors. The dashed line between the states (0,0) and (2,1) indicates that the $^{208}\text{Pb}(p, t)^{206}\text{Pb}$ reaction to the three-phonon state in ^{206}Pb was carried out, and an upper limit of $0.03r$ for the corresponding cross-section determined (see Flynn et al. (1972a) Broglia et al. (1973), see also Lanford and McGrory (1973)). The quantities $B(A)$ (y -axis) stand for the binding energy of the corresponding isotope of Pb (see x -axis), N being the corresponding neutron number. The linear term in N has been introduced so as to make $\hbar\omega_a = \hbar\omega_r (= 2.494 \text{ MeV})$.

and

$$P_\nu^\dagger = a_\nu^\dagger a_{\tilde{\nu}}^\dagger, \quad (1.3.6)$$

the quantity G being the pairing coupling constant (attractive $G > 0$). Concerning the states $(\nu, \tilde{\nu})$ see caption to Fig. 1.3.1. Consequently, in this case the concept of independent particle field \hat{Q} (see also (1.2.6) and (1.2.7)) associated with particle-hole (ph) excitations and carrying transfer quantum number $\beta = 0$ has to be generalized to include fields describing independent pair motion of different multipolarity and (normal) parity, in particular 0^+ , in which case $\alpha \equiv (\beta = +2, J^\pi = 0^+)$. The associated RPA creation operator can be written as,

$$\Gamma_a^\dagger(n) = \sum_k X_n^a(k) \Gamma_k^\dagger + \sum_i Y_n^a(i) \Gamma_i, \quad (1.3.7)$$

with

$$\Gamma_k^\dagger = a_k^\dagger a_{\bar{k}}^\dagger (\epsilon_k > \epsilon_F), \quad \Gamma_i^\dagger = a_i^\dagger a_i (\epsilon_i \leq \epsilon_F), \quad (1.3.8)$$

and

$$\sum_k (X_n^a(k))^2 - \sum_i (Y_n^a(i))^2 = 1, \quad (1.3.9)$$

for the pair addition ($(pp), \beta = +2$) mode, and a similar expression for the pair removal ($(hh), \beta = -2$) mode.

In Fig. 1.3.1 the NFT graphical representation of the RPA equations for the pair addition mode is given. The state $\Gamma_a^\dagger(\beta = +2)|\tilde{0}\rangle$, where $|\tilde{0}\rangle$ is the correlated ground state of a closed shell nucleus, can be viewed as the nuclear parallel of a Cooper pair found at the basis of the microscopic theory of superconductivity²⁷. While surface vibrations are associated with the normal ($\beta = 0$) nuclear density, pairing vibrations are connected with the so called abnormal ($\beta = \pm 2$) nuclear density (density of Cooper pairs), both static and dynamic.

Similar to the quadrupole and octupole vibrational bands built out of n_α phonons of quantum numbers $\alpha \equiv (\beta = 0, \lambda^\pi = 2^+, 3^-)$ schematically shown in Fig. 1.1.3 and experimentally observed in inelastic and Coulomb excitation and associated γ -decay processes, pairing vibrational bands build of n_α phonons of quantum numbers $\alpha \equiv (\beta = \pm 2, \lambda^\pi = 0^+, 2^+)$ have been identified around closed shells in terms of two-nucleon transfer reactions throughout the mass table²⁸. In Fig. 1.3.2, the monopole pairing vibrational band built around the vacuum state $|gs(^{208}\text{Pb})\rangle$ in terms of the pair addition and subtraction modes

$$|a\rangle = |n_a = 1\rangle = |gs(^{210}\text{Pb})\rangle, \quad \text{and} \quad |r\rangle = |n_r = 1\rangle = |gs(^{206}\text{Pb})\rangle, \quad (1.3.10)$$

²⁷ The solution of the Cooper pair problem, being of TD-type does not take into account ground state correlations. Within this context see Ambegaokar (1969) p. 285.

²⁸ See e.g. Flynn et al. (1971), Flynn et al. (1972a) and references therein.

is displayed. The indices $n_a (= 1, 2, \dots)$ and $n_r (= 1, 2, \dots)$ indicate the number of pair addition and/or of pair removal modes of which the states are made. The label a and r indicate in Fig. 1.3.2, the transition probability, i.e. two-nucleon (Cooper pair) transfer cross section which connects the ground (“vacuum”) state $|gs(^{208}\text{Pb})\rangle$ with $|a\rangle (\equiv (0, 1))$ and $|r\rangle (\equiv (1, 0))$ respectively.

1.4 Spontaneously broken symmetry

Because empty space is homogeneous and isotropic, and to the degree of accuracy relevant in many-body problems, also invariant under time-reversal, the nuclear Hamiltonian is translational, rotational and time-reversal invariant. According to quantum mechanics, the corresponding wavefunctions transform in an irreducible way under the corresponding groups of transformation. When the solution of the Hamiltonian does not have some of these symmetries, for example defines a privileged direction in space violating rotational invariance, one is confronted with the phenomenon of spontaneously broken symmetry. Strictly speaking, this can take place only for idealized systems that are infinitely large. But when one sees similar phenomena in atomic nuclei, although not so clear or regular, one recognizes that this system is after all a finite quantum many-body system (FQMBS)²⁹.

1.4.1 Quadrupole deformations in 3D-space

A nuclear embodiment of the spontaneous symmetry breaking phenomenon is provided by a quadrupole deformed mean field. A situation one is confronted with, when the value of the lowest quadrupole frequency ω_2 of the RPA solution (1.1.15) tends to zero ($C_2 \rightarrow 0$, D_2 finite). A phenomenon resulting from the interplay of the interaction v (H_{QQ} in (1.3.2)), and of the nucleons outside closed shell, leading to tidal-like polarization of the spherical core.

Coordinate and linear momentum ((x, p_x) single-particle motion) as well as Euler angles and angular momentum (ω, \mathbf{I}) associated with rotation in three-dimensional (3D)-space, are conjugate variables. Similarly, the gauge angle and the number of particles ((ϕ, N) rotation in gauge space), fulfill $[\phi, N] = i$. The operators $e^{-ip_x x}$, $e^{-i\omega \cdot \mathbf{I}}$ and $e^{-iN\phi}$ induce Galilean transformation in 1D-space and rotations in 3D- and in gauge (2D-) space respectively.

Making again use, for didactical purposes, of H_{QQ} instead of v , and calling $|N\rangle$, for Nilsson³⁰, the eventual mean field solution of the Hamiltonian $T + H_{QQ}$, one can write,

$$\langle N | \hat{Q} | N \rangle = Q_0, \quad (1.4.1)$$

where, for simplicity, we assumed axial symmetry ($\lambda = 2, \mu = 0$). One is then confronted with the emergence of a static quadrupole deformation. Rewriting H_{QQ}

²⁹Anderson (1972, 1984).

³⁰Nilsson (1955).

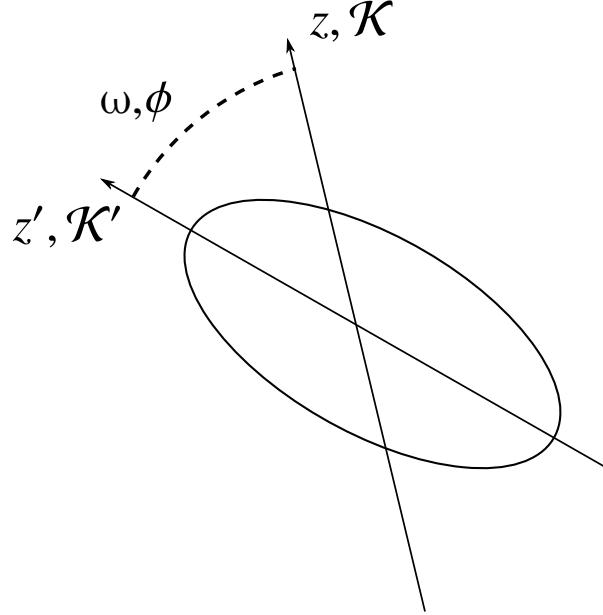


Figure 1.4.1: Schematic representation of deformation in 3D- (gauge-) space, where the laboratory (\mathcal{K}) and the intrinsic (\mathcal{K}' , body fixed) frames of reference make a relative relative angle ((ω , Euler angles); (ϕ , gauge angle)) are indicated (see Eqs. (1.4.7) and (1.4.50)).

in terms of $((\hat{Q}^\dagger - Q_0) + Q_0)$ and its Hermitian conjugate leads to,

$$H = H_{sp} + H_{QQ} = H_{MF} + H_{fluct}, \quad (1.4.2)$$

where

$$H_{MF} = H_{sp} - \kappa(\hat{Q}^\dagger + \hat{Q}), \quad (1.4.3)$$

is the mean field, and

$$H_{fluct} = -\kappa(\hat{Q}^\dagger - Q_0)(\hat{Q} - Q_0) \quad (1.4.4)$$

the residual interaction inducing fluctuations around Q_0 .

Let us concentrate on the mean filed Hamiltonian H_{MF} . It describes the motion of nucleons in a single-particle potential of radius $R_0(1 + \beta_2 Y_{20}(\hat{r}))$, with β_2 proportional to the intrinsic quadrupole moment³¹ Q_0 ($\beta_2 \approx Q_0/(ZR_0^2)$). The reflection invariance and axial symmetry of the Nilsson Hamiltonian implies that parity π and projection Ω of the total angular momentum along the symmetry axis are constants of motion for the one-particle Nilsson states. These states, are two-fold degenerate, since two orbits that differ only in the sign of Ω represent the same motion,

³¹Mottelson and Nilsson (1959), where use has been made of $\beta_2 \approx 0.95\delta$, see Bohr and Mottelson (1975) p. 47 and Eq. (4-191).

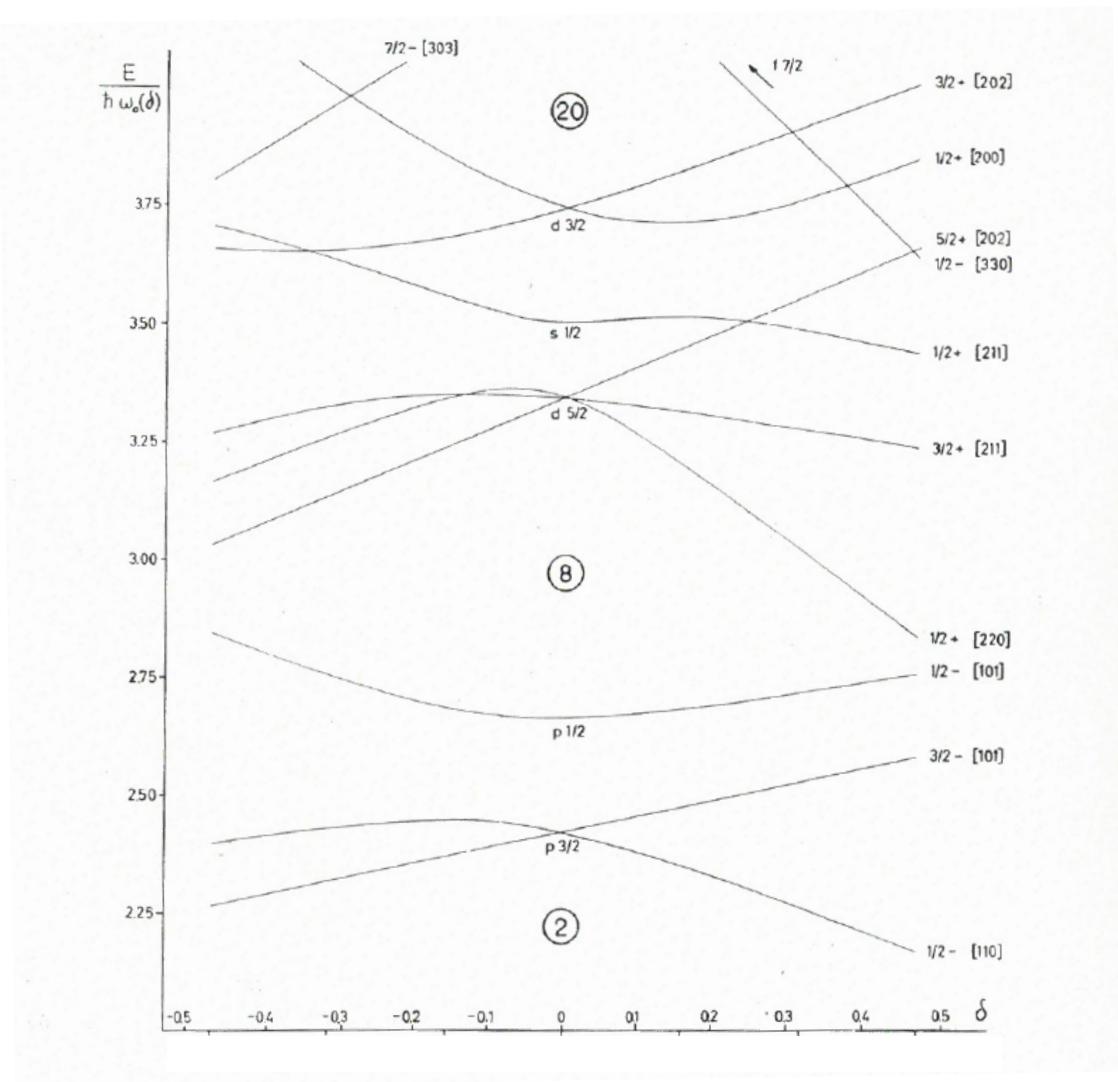


Figure 1.4.2: Nilsson single-particle levels in a quadrupole deformed potential in the regions $2 < Z < 20$ and $2 < N < 20$.

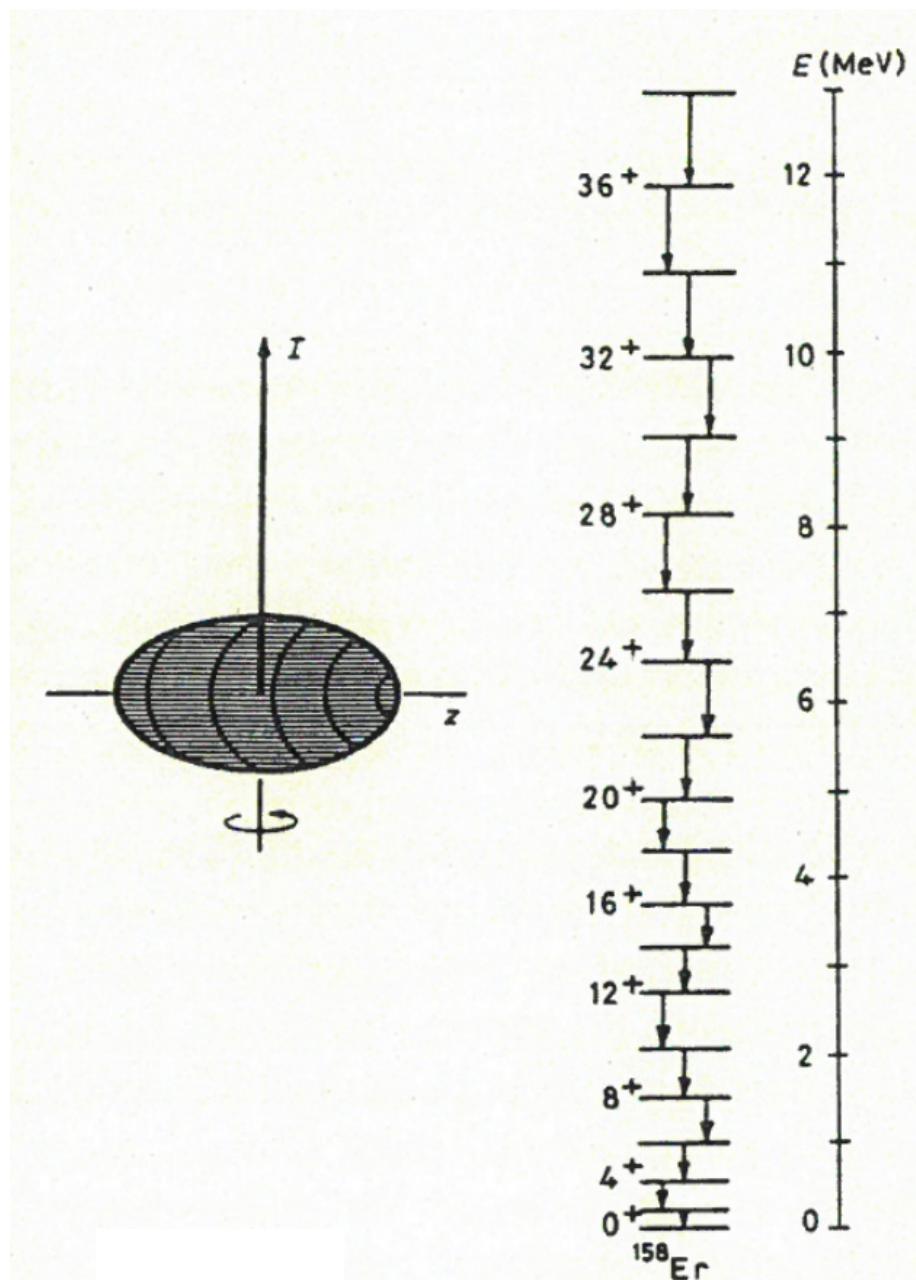


Figure 1.4.3: Ground state rotational band of ^{158}Er .

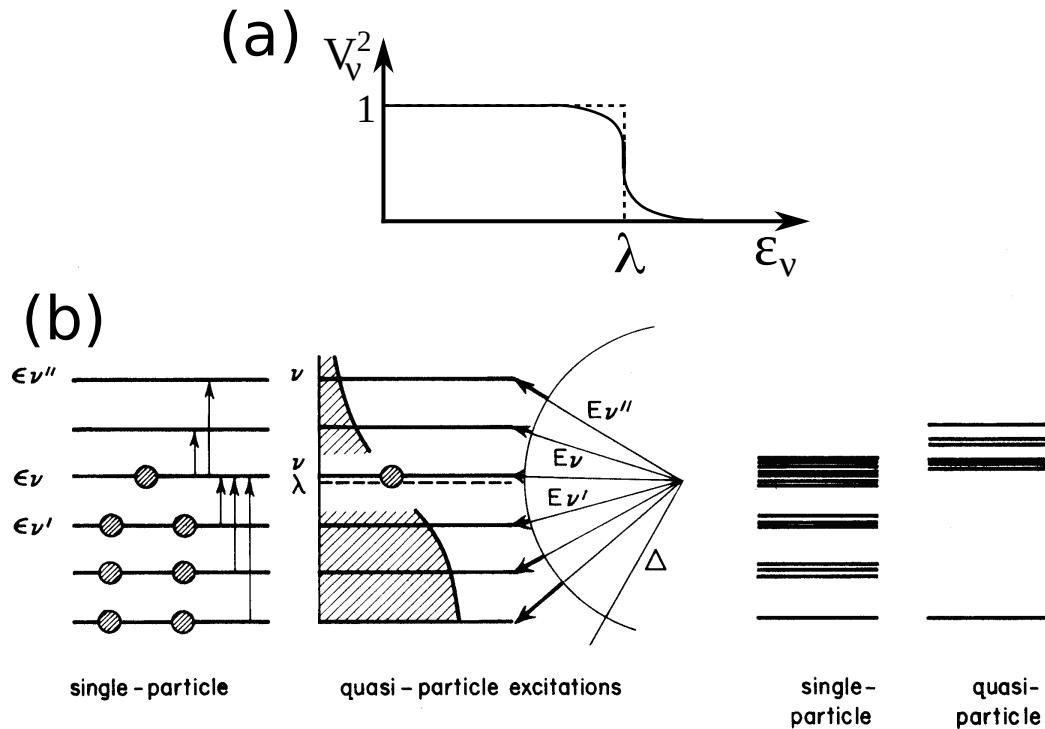


Figure 1.4.4: (a) Independent (dashed line) and BCS occupation numbers; (b) ground state and excited states in the extreme single-particle model and in the pairing-correlated, superfluid model in the case of a system with an odd number of particles. In the first case, the energy of the ground state of the odd system differs from that of the even with one particle fewer by the energy difference $\epsilon_{\nu} - \epsilon_{\nu''}$ while in the second case by the energy $E_{\nu} = \sqrt{(\epsilon_{\nu} - \lambda)^2 + \Delta^2} \approx \Delta$, associated with the fact the odd particle has no partner. Excited states can be obtained in the independent particle case, where it is assumed that levels are two-fold degenerate (Kramers degeneracy) by promoting the odd particle to states above the level ϵ_{ν} , or one particle of the two-fold occupied levels to states above ν' (arrows). To the left only a selected number of these excitations are shown. In the superfluid case excited states can be obtained by breaking of pairs in any orbit. The associated quasiparticle energy is drawn also here by an arrow of which the thin part indicates the contribution of the pairing gap and the thick part indicates the kinetic energy contribution, i.e. the contribution arising from the single-particle motion. Note the very different density of levels emerging from these two pictures, which are shown at the far right of the figure.

apart from the clockwise and anticlockwise sense of revolution around the symmetry axis. One can thus write the Nilsson creation operators in terms of a linear combination of creation operators carrying good total angular momentum j ,

$$\gamma_{a\Omega}^\dagger = \sum_j A_j^a a_{aj\Omega}^\dagger, \quad (1.4.5)$$

where the label a stands for the quantum numbers which aside from Ω , specify the orbital.

Expressing (1.4.5) in the intrinsic, body-fixed, system of coordinates \mathcal{K}' (Fig. 1.4.1) where the 3 (z') axis lies along the symmetry axis and the 1 and 2 (x', y') axis lie in a plane perpendicular to it,

$$\gamma'_{a\Omega}^\dagger = \sum_j A_j^a \sum_{\Omega'} \mathcal{D}_{\Omega\Omega'}^2(\omega) a_{aj\Omega'}^\dagger, \quad (1.4.6)$$

and one can write the Nilsson state as

$$|N(\omega)\rangle_{\mathcal{K}'} = \prod_{a\Omega>0} \gamma'_{a\Omega}^\dagger \gamma'_{a\Omega}^\dagger |0\rangle_F, \quad (1.4.7)$$

where ω represent the Euler angles, $|0\rangle_F$ is the particle vacuum, and $\gamma'_{a\Omega}^\dagger |0\rangle_F$ is the state time-reversed to $\gamma_{a\Omega}^\dagger |0\rangle_F$. For well deformed nuclei, a conventional description of the one-particle motion is based on the similarity of the nuclear potential to that of an anisotropic oscillator potential,

$$V = \frac{1}{2}M \left(\omega_3^2 x_3^2 + \omega_\perp^2 (x_1^2 + x_2^2) \right) = \frac{1}{2}M\omega_0 r^2 \left(1 - \frac{4}{3}\delta \times P_2(\cos\theta) \right), \quad (1.4.8)$$

with $\omega_3 \omega_\perp^2 = \omega_0^3$. That is a volume which is independent of the deformation $\delta \approx 0.95\beta_2$. The corresponding single-particle states have energy

$$\epsilon(n_3 n_\perp) = (n_3 + \frac{1}{2})\hbar\omega_3 + (n_\perp + 1)\hbar\omega_\perp, \quad (1.4.9)$$

where n_3 and $n_\perp = n_1 + n_2$ are the number of quanta along and perpendicular to the symmetry axis, $N = n_3 + n_\perp$ being the total oscillator quantum number. The degenerate states with the same value of n_\perp can be specified by the component Λ of the orbital angular momentum along the symmetry axis,

$$\Lambda = \pm n_\perp, \pm(n_\perp - 2), \dots, \pm 1 \text{ or } 0. \quad (1.4.10)$$

The complete expression of the Nilsson potential includes, aside from the central term discussed above, a spin-orbit term ($\Sigma = \pm 1/2$, $\Omega = \Lambda + \Sigma$), and a term proportional to the orbital angular momentum quantity squared, so as to make the shape of the oscillator to resemble more that of a Saxon-Woods potential. One can label the Nilsson levels in terms of the asymptotic quantum numbers $\Omega[Nn_3\Lambda]$.

The resulting states provide an overall account of the experimental findings. An example of relevance for light nuclei (N and $Z < 20$) is given in³² Fig. 1.4.2.

The Nilsson intrinsic state (1.4.7) does not have a definite angular momentum but is rather a superposition of such states,

$$|N(\omega)\rangle_{\mathcal{K}'} = \sum c_I |I\rangle. \quad (1.4.11)$$

Because there is no restoring force associated with different orientations of $|N(\omega)\rangle_{\mathcal{K}'}$, fluctuations in the Euler angle diverge in the right way so as to restore rotational invariance, leading to a rotational band whose members are

$$|IKM\rangle \sim \int d\omega \mathcal{D}_{MK}^I(\omega) |N(\omega)\rangle_{\mathcal{K}'}, \quad (1.4.12)$$

with energy

$$E_I = \frac{\hbar^2}{2I} I(I+1). \quad (1.4.13)$$

The quantum numbers I, M, K are the total angular momentum I , and its third component M and K along the laboratory (z) and intrinsic (z') frame references respectively. Rotational bands have been observed up to rather high angular momenta in terms of individual transitions. In particular, up to $I = 60\hbar$ in the case of³³ the nucleus ^{152}Dy . An example of quadrupole rotational bands is given in Fig. 1.4.3.

1.4.2 Deformation in gauge space

Let us now return to the Hamiltonian $H = H_{sp} + H_P$ (Eqs. (1.1.13) and (1.3.4)), in connection with pairing vibrations. In the case in which $\hbar\omega_1(\beta = -2) = \hbar\omega_1(\beta = +2) = 0$, the system deforms, this time in gauge space. The mean field solution of the pairing Hamiltonian³⁴ called $|BCS\rangle$, leads to the finite expectation value

$$\alpha_0 = \langle BCS | P^\dagger | BCS \rangle, \quad (1.4.14)$$

of the pair creation operator P^\dagger , quantity which can be viewed as the order parameter of the new deformed phase of the system in gauge space. The pairing Hamiltonian can be written as

$$H = H_{MF} + H_{fluct}, \quad (1.4.15)$$

where

$$H_{MF} = H_{sp} - \Delta(P^\dagger + P) + \frac{\Delta^2}{G} \quad (1.4.16)$$

³²Mottelson and Nilsson (1959).

³³Nolan and Twin (1988) and refs. therein.

³⁴Bardeen et al. (1957a,b).

and

$$H_{fluct} = -G(P^\dagger - \alpha_0)(P - \alpha_0). \quad (1.4.17)$$

The quantity

$$\Delta = G\alpha_0, \quad (1.4.18)$$

is the so called pairing gap (Fig. 1.4.4), 2Δ measuring the binding energy of Cooper pairs, the quantity α_0 being the number of them (see Eq. (1.4.42)).

The mean field pairing Hamiltonian

$$H_{MF} = \sum_{\nu>0} (\epsilon_\nu - \lambda) (a_\nu^\dagger a_\nu + a_{\tilde{\nu}}^\dagger a_{\tilde{\nu}}) - \Delta \sum_{\nu>0} (\epsilon_\nu - \lambda) (a_\nu^\dagger a_{\tilde{\nu}}^\dagger + a_{\tilde{\nu}} a_\nu) + \frac{\Delta^2}{G} \quad (1.4.19)$$

is a bilinear expression in the creation and annihilation operators, ν labeling the quantum numbers of the single-particle orbitals where nucleons are allowed to correlate (e.g. (nlj, m)), while $\tilde{\nu}$ denotes the time reversal state which in this case is degenerate with ν and has quantum numbers $(nlj, -m)$. The condition $\nu > 0$ in Eq. (1.4.19) implies that one sums over $m > 0$. It is of notice that

$$\hat{N} = \sum_{\nu>0} (a_\nu^\dagger a_\nu + a_{\tilde{\nu}}^\dagger a_{\tilde{\nu}}), \quad (1.4.20)$$

is the number operator, and $\lambda \hat{N}$ in Eq. (1.4.19) acts as the Coriolis force in the body-fixed frame of reference in gauge space.

One can diagonalize H_{MF} by a rotation in the (a^\dagger, a) -space. This can be accomplished through the Bogoliubov-Valatin (quasiparticle) transformation³⁵

$$\alpha_\nu^\dagger = U_\nu a_\nu^\dagger - V_\nu a_{\tilde{\nu}}, \quad (1.4.21)$$

where the occupation numbers (U_ν, V_ν) fulfill $|U_\nu|^2 + |V_\nu|^2 = 1$. The BCS solution does not change the energies ϵ_ν of the single-particle levels (measured in (1.4.19) from the Fermi energy λ), nor the associated wavefunctions $\varphi_\nu(\mathbf{r})$, but the occupation probabilities of levels around the Fermi energy within an energy range 2Δ ($2\Delta/\epsilon_F \approx 3$ MeV/36 MeV ≈ 0.06)³⁶. The quasiparticle operator α_ν^\dagger creates a particle in the single-particle state ν with probability U_ν^2 , while it creates a hole (annihilates a particle) with probability V_ν^2 . To be able to create a particle, the state ν should be empty, while to annihilate a particle it has to be filled, so U_ν^2 and V_ν^2 are the probabilities that the state ν is empty and is occupied respectively. Within this context, the one-quasiparticle states

$$|\nu\rangle = \alpha_\nu^\dagger |BCS\rangle, \quad (1.4.22)$$

³⁵Bogoliubov (1958a,b); Valatin (1958).

³⁶See also discussion starting just before Eq. (1.4.53) and ending just before Eq. (1.4.55).

are orthonormal. In particular

$$\langle \nu | \nu \rangle = 1 = \langle BCS | \alpha_\nu \alpha_\nu^\dagger | BCS \rangle = \langle BCS | \{ \alpha_\nu, \alpha_\nu^\dagger \} | BCS \rangle = U_\nu^2 + V_\nu^2, \quad (1.4.23)$$

where the anticommutators

$$\{ \alpha_\nu, \alpha_{\nu'}^\dagger \} = \delta(\nu, \nu'), \quad (1.4.24)$$

and

$$\{ \alpha_\nu, \alpha_{\nu'} \} = \{ \alpha_\nu^\dagger, \alpha_{\nu'}^\dagger \} = 0, \quad (1.4.25)$$

have been used. Note that the $|BCS\rangle$ state is the quasiparticle vacuum

$$\alpha_\nu |BCS\rangle = 0, \quad (1.4.26)$$

in a similar way in which $|0\rangle_F$ is the particle vacuum. Inverting the quasiparticle transformation (1.4.21) and its complex conjugate, i.e. expressing α_ν^\dagger and α_ν (and their time reversals (tr)) as function of α_ν^\dagger and α_ν (and tr) one can rewrite (1.4.19) in terms of quasiparticles.

Minimizing the energy $E_0 = \langle BCS | H | BCS \rangle$ in terms of³⁷ V_ν

$$\frac{\partial E_0}{\partial V_\nu} = 0, \quad (1.4.27)$$

making use of the expression for the average number of particles

$$N_0 = \langle BCS | \hat{N} | BCS \rangle = 2 \sum_{\nu > 0} V_\nu^2, \quad (1.4.28)$$

of the number of Cooper pairs

$$\alpha_0 = \langle BCS | P^\dagger | BCS \rangle = \sum_{\nu > 0} U_\nu V_\nu, \quad (1.4.29)$$

and of the pairing gap

$$\Delta = G \sum_{\nu > 0} U_\nu V_\nu, \quad (1.4.30)$$

one obtains,

$$H_{MF} = H_{11} + U, \quad (1.4.31)$$

where

$$H_{11} = \sum_\nu E_\nu \alpha_\nu^\dagger \alpha_\nu, \quad (1.4.32)$$

³⁷For details see e.g. Ragnarsson and Nilsson (2005). See also Nathan and Nilsson (1965).

and

$$U = 2 \sum_{\nu>0} (\epsilon_\nu - \lambda) V_\nu^2 - \frac{\Delta^2}{G}. \quad (1.4.33)$$

The quantity

$$E_\nu = \sqrt{(\epsilon_\nu - \lambda)^2 + \Delta^2}, \quad (1.4.34)$$

is the quasiparticle energy, while the occupation (emptiness) probability amplitudes are

$$V_\nu = \frac{1}{\sqrt{2}} \left(1 - \frac{\epsilon_\nu - \lambda}{E_\nu} \right)^{1/2}, \quad (1.4.35)$$

$$U_\nu = \frac{1}{\sqrt{2}} \left(1 + \frac{\epsilon_\nu - \lambda}{E_\nu} \right)^{1/2}, \quad (1.4.36)$$

respectively. From the relations (1.4.28), (1.4.30) and (1.4.34) one obtains

$$N_0 = 2 \sum_{\nu>0} V_\nu^2, \quad (1.4.37)$$

and

$$\frac{1}{G} = \sum_{\nu>0} \frac{1}{2E_\nu}. \quad (1.4.38)$$

These equations, known as the BCS number and gap equations, allow one to calculate the parameters λ and Δ from the knowledge of N_0 , G and ϵ_ν , parameters which completely determine E_ν , V_ν and U_ν and thus the BCS mean field solution (Fig. 1.4.4). The validity of the BCS description of superfluid open shell nuclei have been extensively confirmed³⁸.

The relation (1.4.26) implies, as a consequence of Pauli principle, that

$$\begin{aligned} |BCS\rangle &= \frac{1}{\text{Norm}} \prod_{\nu>0} \alpha_\nu \alpha_{\bar{\nu}} |0\rangle_F = \prod_{\nu>0} (U_\nu + V_\nu P_\nu^\dagger) |0\rangle_F \\ &= \left(\prod_{\nu>0} U_\nu \right) \sum_{N \text{ even}} \frac{(\sum_{\nu>0} c_\nu P_\nu^\dagger)^{N/2}}{(N/2)!} |0\rangle_F \\ &= \left(\prod_{\nu>0} U_\nu \right) \sum_{n=0,1,2,\dots} \frac{(\sum_{\nu>0} c_\nu P_\nu^\dagger)^n}{n!} |0\rangle_F \\ &= \left(\prod_{\nu>0} U_\nu \right) \exp \left(\sum_{\nu>0} c_\nu P_\nu^\dagger \right) |0\rangle_F, \end{aligned} \quad (1.4.39)$$

³⁸See e.g. Broglia and Zelevinsky (2013) and refs therein.

where

$$c_\nu = V_\nu / U_\nu, \quad (1.4.40)$$

while n denotes the number of pairs of particles ($2n = N$). *The last expression of (1.4.39) implies that $|BCS\rangle$ is a coherent state.*

In the above discussion of BCS we have treated in a rather cavalier fashion the fact that the amplitudes U_ν and V_ν are in fact complex quantities. A possible choice of phasing is³⁹

$$U_\nu = U'_\nu; \quad V_\nu = V'_\nu e^{-2i\phi}, \quad (1.4.41)$$

and, as a consequence,

$$\alpha_0 = e^{-2i\phi} \alpha'_0 \quad (\alpha'_0 = \sum_{\nu>0} U'_\nu V'_\nu) \text{ and } \Delta = e^{-2i\phi} \Delta' \quad (\Delta' = G \alpha'_0), \quad (1.4.42)$$

U'_ν and V'_ν being real quantities, while ϕ is the gauge angle, conjugate variable to the number of particles operator. Then⁴⁰

$$\hat{\phi} = i\partial/\partial\mathcal{N}, \quad \mathcal{N}, \quad (1.4.43)$$

and

$$[\hat{\phi}, \mathcal{N}] = i, \quad (1.4.44)$$

where $\mathcal{N} \equiv \hat{N}$ (Eq. (1.4.20)), the gauge transformations being induced by the operator

$$\mathcal{G}(\phi) = e^{-i\mathcal{N}\phi}. \quad (1.4.45)$$

Let us replace the amplitudes (1.4.41) in (1.4.39),

$$|BCS\rangle = \left(\prod_{\nu>0} U'_\nu \right) \sum_{N \text{ even}} e^{-iN\phi} |\Phi_N\rangle = \left(\prod_{\nu>0} U'_\nu \right) \sum_{N \text{ even}} e^{-iN\phi} |\Phi_N\rangle, \quad (1.4.46)$$

where

$$|\Phi_N\rangle = \frac{\left(\sum_{\nu>0} c'_\nu P_\nu^\dagger \right)^{N/2}}{(N/2)!} |0\rangle_F, \quad (1.4.47)$$

with $c'_\nu = V'_\nu / U'_\nu$. It is of notice that

$$|\Phi_0\rangle = \sum_{\nu>0} c'_\nu P_\nu^\dagger |0\rangle_F \quad (1.4.48)$$

³⁹Schrieffer (1973). The same results as those which will be derived are obtained with the alternative choice $U_\nu = U'_\nu e^{i\phi}$, $V_\nu = V'_\nu e^{-i\phi}$ (see e.g. Potel et al. (2013b)).

⁴⁰See e.g. Brink and Broglia (2005) App. H and refs. therein.

is the single Cooper pair state. The $|BCS\rangle$ state does not have a definite number of particles, but only in average, being a wavepacket in N . In fact, (1.4.46) defines a privileged direction in gauge space, being an eigenstate of $\hat{\phi}$

$$\hat{\phi}|BCS\rangle = i \frac{\partial}{\partial N} \left(\prod_{\nu>0} U'_\nu \right) \sum_{N \text{ even}} e^{-iN\phi} |\Phi_N\rangle = \phi |BCS\rangle. \quad (1.4.49)$$

Expressing it differently (1.4.46) can be viewed as an axially symmetric deformed system in gauge space, whose symmetry axis coincides with the z' component of the body-fixed frame of reference \mathcal{K}' , which makes an angle ϕ with the laboratory z -axis (Fig. 1.4.1).

With the help of Eq. (1.4.39) (first line) one can write

$$\begin{aligned} |BCS(\phi=0)\rangle_{\mathcal{K}'} &= \prod_{\nu>0} (U'_\nu + V'_\nu P_\nu^\dagger) |0\rangle_F = \prod_{\nu>0} (U_\nu + V_\nu P_\nu^\dagger) |0\rangle_F \\ &= \prod_{\nu>0} (U'_\nu + e^{-2i\phi} V'_\nu P_\nu^\dagger) |0\rangle_F, \end{aligned} \quad (1.4.50)$$

where use was made of the relations⁴¹

$$\mathcal{G}(\phi) a_\nu^\dagger \mathcal{G}^{-1}(\phi) = e^{-i\phi} a_\nu^\dagger = a_\nu'^\dagger, \quad (1.4.51)$$

and

$$\mathcal{G}(\phi) P_\nu^\dagger \mathcal{G}^{-1}(\phi) = e^{-2i\phi} P_\nu^\dagger = P_\nu'^\dagger. \quad (1.4.52)$$

Spontaneously broken symmetry in nuclei is, as a rule, associated with the presence of rotational bands, as already found in the case of quadrupole deformed nuclei. *Consequently, one expects in nuclei with $\alpha_0 \neq 0$ rotational bands in which particle number plays the role of angular momentum. That is pairing rotational bands (see Eq. (1.4.65)).*

Before proceeding with the discussion of the term H_{fluct} let us return to the relation (1.4.26). That is the fact that one can view the state $|BCS\rangle$ as the quasiparticle vacuum. Thus implying that it can be written as

$$|BCS\rangle \sim \prod_\nu \alpha_\nu |0\rangle_F, \quad (1.4.53)$$

⁴¹It is to be noted that \mathcal{G} induces a counter clockwise rotation, $\mathcal{G}(\chi) \hat{\phi} \mathcal{G}^{-1}(\chi) = \hat{\phi} - \chi$,

$$\begin{aligned} [\mathcal{N}, \phi] &= -i, \quad \mathcal{G}(\chi) = e^{-i\mathcal{N}\chi} \quad \chi : \text{c-number}, \quad [e^{-i\mathcal{N}\chi}, \phi] = e^{-i\mathcal{N}\chi} \phi - \phi e^{-i\mathcal{N}\chi}, \\ [e^{-i\mathcal{N}\chi}, \phi] e^{i\mathcal{N}\chi} &= e^{-i\mathcal{N}\chi} \phi e^{i\mathcal{N}\chi} - \phi, \quad e^{-i\mathcal{N}\chi} = 1 - i\mathcal{N}\chi \quad \chi^2 \ll \chi \\ [1 - i\mathcal{N}\chi, \phi] &= -i[\chi, \phi] = -i(\mathcal{N}[\chi, \phi] + [\mathcal{N}, \phi]\chi) = -i(0 - i\chi) = -\chi \\ [e^{-i\mathcal{N}\chi}, \phi] &= -\chi e^{-i\mathcal{N}\chi} \\ [e^{-i\mathcal{N}\chi}, \phi] e^{i\mathcal{N}\chi} &= -\chi e^{-i\mathcal{N}\chi} e^{i\mathcal{N}\chi} = -\chi, \quad -\chi = \mathcal{G}(\chi) \phi \mathcal{G}^{-1}(\chi) - \phi \\ \mathcal{G}(\chi) \phi \mathcal{G}^{-1}(\chi) &= \phi - \chi \end{aligned}$$

as well as

$$|BCS\rangle \sim \prod_{\nu>0} \alpha_\nu \alpha_{\tilde{\nu}} |0\rangle_F, \quad (1.4.54)$$

The state $|\nu'\rangle \sim \prod_{\nu \neq \nu'} \alpha_\nu |0\rangle_F$ corresponds to a one quasiparticle state, while $|\nu' \tilde{\nu}'\rangle \sim \prod_{\nu>0, \nu \neq \nu'} \alpha_\nu^\dagger \alpha_{\tilde{\nu}}^\dagger |0\rangle_F$ is proportional to a two-quasiparticle state. Both states display an energy (pairing) gap with respect to the ground state. In the first case of the order of Δ , in the second of 2Δ . A property closely connected with the fact that in BCS condensation as described by $|BCS\rangle$, a macroscopic number of Cooper pairs are phase coherent, each of them being made out of pairs of fermions entangled in time reversal states. Implying an energy gap for both pair dissociation and single pair breaking. Property referred to as off-diagonal-long-range-order (ODLRO)⁴² (see Sect. 4.9.3).

1.4.3 The normal and the superconducting state

Within this scenario it is useful in discussing BCS condensation in nuclei to go back to its origin⁴³, namely superconductivity in metals. It is found that for superconductors based on metals displaying a strong electron-phonon coupling⁴⁴, the behaviour of the BCS occupation probability V_k^2 resembles very much that of the same quantity but for the metal at temperature close but above the critical temperature, i.e. the metal in the normal state. A consequence of processes like those shown in graphs (b) and (c) of Fig. 1.2.4 where, in that case, the fermions are electrons (or electron holes), and the bosons are lattice phonons⁴⁵. Such similitude is also observed in the nuclear case (Fig. 1.4.4, and Fig. 1.2.5).

Concerning superconductors, it is found that the changes observed in the behaviour of the systems upon cooling below the critical temperature, cannot be described at profit just in terms of changes in the occupation probability of one-electron momentum eigenstate (from the sharp line to the smooth curve around λ ($\approx \epsilon_F$), Fig. 1.4.4 (a)), in keeping with the fact that in the process, no gap opens up in E_k -space. In fact, the partial occupation V_ν^2 of the normal state carrying random (gauge) phases is, by lowering the temperature through T_c , progressively replaced by a single quantum state. In it, about the same set of many-body states, in which fermions are dressed and interact pairwise by exchanging bosons (aside than through the bare interaction; screened Coulomb potential in the case of superconductors, NN -potential in the case of superfluid nuclei), become coherently superposed with a fixed phase relation leading to ODLRO. It is then not so much, or better not only, the superconducting (superfluid) state that is special, but also the so called normal state.

⁴² Penrose (1951); Penrose and Onsager (1956); Yang (1962); see also Anderson (1996).

⁴³ Bohr et al. (1958).

⁴⁴ These systems display, in the normal state a consistent electric resistance. In other words, bad conductors are, as a rule, good superconductors, like for example Pb. On the other hand Au does not became superconducting at any measured temperature.

⁴⁵ See e.g. Tinkham (1996) p.28.

In what follows we discuss the structure of H_{fluct} and single out the term responsible for restoring gauge invariance to the mean field solution, giving thus rise to pairing rotational bands for discrete values of the angular momentum in gauge space, namely N , differing in two units from each other. In terms of quasiparticles, H_{fluct} can be expressed as

$$H_{fluct} = H'_p + H''_p + C \quad (1.4.55)$$

where

$$H'_p = -\frac{G}{4} \left(\sum_{\nu>0} (U_\nu^2 - V_\nu^2) (\Gamma_\nu^\dagger + \Gamma_\nu) \right)^2 \quad (1.4.56)$$

and

$$H''_p = \frac{G}{4} \left(\sum_\nu (\Gamma_\nu^\dagger - \Gamma_\nu) \right)^2, \quad (1.4.57)$$

with

$$\Gamma_\nu^\dagger = \alpha_\nu^\dagger \alpha_{\bar{\nu}}^\dagger. \quad (1.4.58)$$

The term C stands for constant terms, as well as for terms proportional to the number of quasiparticles, which consequently vanish when acting on $|BCS\rangle$. The term H'_p gives rise to two-quasiparticle pairing vibrations with energies $\gtrsim 2\Delta$. It can be shown that it is the term H''_p which restores gauge invariance⁴⁶. In fact,

$$[H_{MF} + H''_p, \tilde{N}] = 0. \quad (1.4.59)$$

We now diagonalize $H_{MF} + H''_p$ in the quasiparticle RPA (QRPA),

$$[H_{MF} + H''_p, \Gamma_n^\dagger] = \hbar\omega_n \Gamma_n^\dagger, \quad [\Gamma_n, \Gamma_{n'}^\dagger] = \delta(n, n'), \quad (1.4.60)$$

where

$$\Gamma_n^\dagger = \sum_\nu (a_{n\nu} \Gamma_\nu^\dagger + b_{n\nu} \Gamma_\nu), \quad \Gamma_\nu^\dagger = \alpha_\nu^\dagger \alpha_{\bar{\nu}}^\dagger, \quad (1.4.61)$$

is the creation operator of the n th vibrational mode. In the case of the $n = 1$, lowest energy root, it can be written as

$$|1''\rangle = \Gamma_1^\dagger |0''\rangle = \frac{\Lambda_1''}{2\Delta} (\tilde{N} - N_0) |0''\rangle, \quad (1.4.62)$$

where \tilde{N} is the particle number operator written in terms of Γ_ν^\dagger and Γ_ν , and Λ_1'' is the strength of the quasiparticle-mode coupling.

⁴⁶Högaasen-Feldman (1961); Bès and Broglia (1966). For details see Brink and Broglia (2005), App. I (Sect. I.4) and App. J.

It can be shown that the dimensionless prefactor ($\Lambda_1''/2\Delta$), is the zero point amplitude of the mode⁴⁷, that is

$$\sqrt{\frac{\hbar\omega_1''}{2C_1''}} = \sqrt{\frac{\hbar^2}{2D_1''\hbar\omega_1''}}. \quad (1.4.63)$$

Because the lowest frequency is $\omega_1'' = 0$, the associated ZPF diverge ($\Lambda_1'' \sim (\hbar\omega_1'')^{-1/2}$). It can be seen that this is because $C_1'' \rightarrow 0$, while D_1'' remains finite. In fact,

$$\frac{D_1''}{\hbar^2} = \frac{2\Delta^2}{\Lambda_1''^2\hbar\omega_1''} = 4 \sum_{\nu>0} \frac{U_\nu^2 V_\nu^2}{E_\nu}. \quad (1.4.64)$$

In keeping with the fact that a rigid rotation in gauge space can be generated by a series of infinitesimal operations of type $\mathcal{G}(\delta\phi) = e^{i(\hat{N}-N_0)\delta\phi}$, the one phonon state $|1''\rangle = \Gamma_1^\dagger |0''\rangle$ results from rotations of the deformed (correlated) intrinsic state of divergent amplitude. That is, fluctuations of ϕ over the whole $0 - 2\pi$ range. By proper inclusion of these fluctuations one can restore gauge invariance violated by $|BCS\rangle_{\mathcal{K}}$. The resulting states⁴⁸

$$|N\rangle \sim \int_0^{2\pi} d\phi e^{iN\phi} |BCS(\phi)\rangle_{\mathcal{K}} \sim \left(\sum_{\nu>0} c'_\nu P_\nu^\dagger \right)^{N/2} |0\rangle_F, \quad (1.4.65)$$

have fixed particle number and constitute the members of a pairing rotational band. Making use of a simplified model⁴⁹ (single j -shell) it can be shown that the energies of these states can be written as,

$$E_N = \lambda(N - N_0) + \frac{G}{4} (N - N_0)^2, \quad (1.4.66)$$

where

$$\frac{G}{4} = \frac{\hbar^2}{2D_1''}. \quad (1.4.67)$$

An example of pairing rotational bands is provided by the ground state of the single open closed shell superfluid $^{50}_{50}\text{Sn}_N$ -isotopes (Fig. 1.4.5), $N_0 = 68$ having been found to minimize $E_0 = \langle BCS | H | BCS \rangle$ (see Eq. (1.4.27)), and thus has been used in the solution of the BCS number equation (1.4.37).

Making use of the BCS pair transfer amplitudes,

$$B_\nu \sim \langle BCS | P_\nu^\dagger | BCS \rangle = U_\nu V_\nu, \quad (1.4.68)$$

⁴⁷Amplitude which in the case of surface vibrations was defined in Eq. (1.1.11).

⁴⁸Where use has been made of Eqs. (1.4.50), (1.4.46) and (4.7.27).

⁴⁹See Brink and Broglia (2005) App. H. and refs. therein.

in combination with a computer code (see App. 7.A) and of global optical parameters, one can account for the absolute value of the Cooper pair transfer differential cross section, within experimental errors (see e.g. Figs. 3.1.7 and 7.2.1). The fact that projecting out the ground state of the different Sn-isotopes from the intrinsic BCS state describing ^{118}Sn one obtains a quantitative description of observations carried out with the help of the specific probe of pairing correlations (Cooper pair transfer), testifies to the fact that pairing rotational bands can be considered elementary modes of nuclear excitation, emergent properties of spontaneous symmetry breaking of gauge invariance⁵⁰.

The fact that this insight follows from the use of QRPA⁵¹ in the calculation of the ZPF of the collective solutions of the pairing Hamiltonian underscores the importance of conserving approximations⁵² to describe quantum many-body problems in general, and finite size quantum many-body problems (FSQMB), of which the nucleus represents a paradigmatic example, in particular.

1.5 Giant dipole resonance

Aside from low-lying collective states, that is rotations and low-energy vibrations, nuclei also display high-lying collective modes known as giant resonances.

If one shines a beam of photons on a nucleus, it is observed that the system absorbs energy resonantly essentially at a single frequency, of the order of⁵³ $\nu = 5 \times 10^{21}$ Hz, corresponding to an energy of $h\nu \approx 20$ MeV.

It is not difficult to understand how γ -rays excite a nuclear dipole vibration. A photon carries with it an oscillating electric field. Although the wavelength of a 20 MeV γ -ray is smaller than that of other forms of electromagnetic radiation such as visible light, it is still large ($\lambda \approx 63$ fm) compared to the dimensions of e.g. ^{40}Ca ($R_0 \approx 4.1$ fm). As a result the photon electric field is nearly uniform across the nucleus at any time. This field exerts a force on the positively charged protons. Consequently, it can set the center of mass into an antenna like, dipole oscillation (Thompson scattering), in which case no photon is absorbed. Another possibility is that it leads to an internal excitation of the system. In this case, because the center of mass of the system does not move, the neutrons have to oscillate against the protons, again in an antenna-like motion. The restoring force of the vibration, known as the giant dipole resonance (GDR), is provided by the attractive force between protons and neutrons.

The connotation of giant is in keeping with the fact that it essentially carries

⁵⁰A finding related to the fact that $|BCS\rangle$ is a coherent state.

⁵¹Using the Tamm-Dancoff approximation, i.e. setting $Y \equiv 0$ (and thus $\sum X^2 = 1$) in the QRPA solution, does not lead to particle number conservation, in keeping with the fact that the amplitudes Y are closely connected to ZPF.

⁵²That is, approximations respecting sum rules.

⁵³Making use of $h = 4.1357 \times 10^{-15}$ eV·s one obtains for $h\nu = 1$ eV the frequency $\nu = 2.42 \times 10^{14}$ Hz and thus $\nu = 4.8 \times 10^{21}$ Hz for $h\nu = 20$ MeV. The wavelength of a photon of energy E is $\lambda = hc/E \approx 2\pi \times 200$ MeV fm/ E , which for $E = 20$ MeV leads to $\lambda \approx 63$ fm.

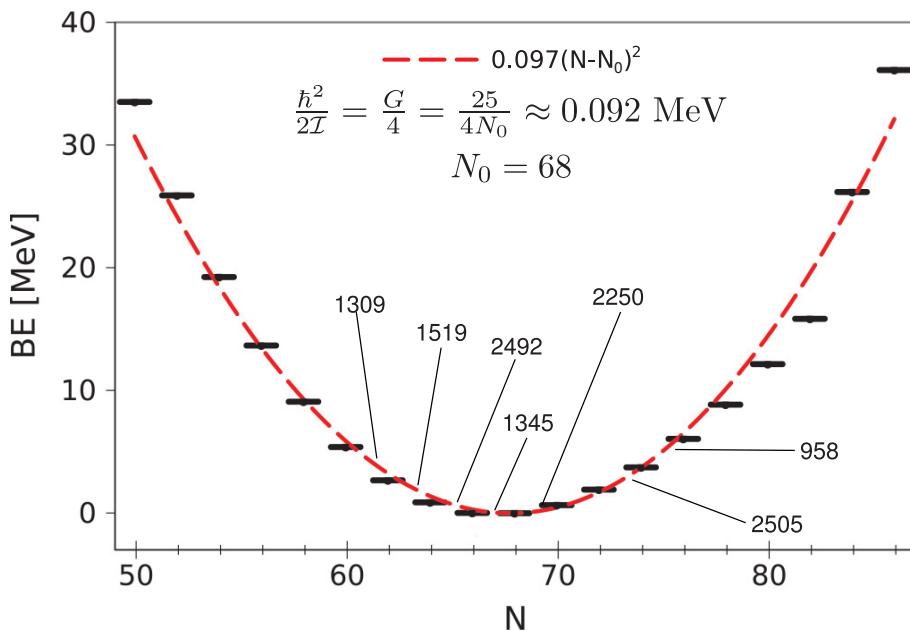


Figure 1.4.5: Pairing rotational band associated with the ground states of the Sn-isotopes. The lines represent the energies calculated according to the expression $BE = B(^{50+N}\text{Sn}_N) - 8.124N + 46.33$ (Brink and Broglia (2005)), subtracting the contribution of the single nucleon addition to the nuclear binding energy obtained by a linear fitting of the binding energies of the whole Sn chain. The estimate of $\hbar^2/2I$ was obtained using the single j -shell model Eq. (1.4.67) (see also Brink and Broglia (2005), Appendix H and references therein). The numbers indicated with thin lines are the absolute value of the experimental $gs \rightarrow gs$ cross section (in units of μb).

the full photo absorption cross section (energy weighted sum rule, see below), and resonance because it displays a Lorentzian-like shape with a full width at half maximum of few MeV ($\lesssim 5$ MeV), considerably smaller than the energy centroid⁵⁴ $\hbar\omega_{GDR} \approx 80/A^{1/3}$ MeV. Microscopically, the GDR can be viewed as a correlated particle-hole excitation, that is a state made out of a coherent linear combination of proton and neutron particle-hole excitations with essentially $\Delta N = 1$ character, as well as small $\Delta N = 3, 5, \dots$ components (Fig. 1.2.3). Because the difference in energy between major shells is $\hbar\omega_0 \approx 41A^{1/3}$ MeV, one expects that about half of the contribution to the energy of the GDR arises from the neutron-proton interaction. More precisely, from the so called symmetry potential V_1 (see Eq. (1.1.4)), which measures the energy price the system has to pay to separate protons from neutrons. Theoretical estimates lead to⁵⁵

$$(\hbar\omega_{GDR})^2 = (\hbar\omega_0)^2 + \frac{3\hbar^2 V_1}{m\langle r^2 \rangle} = \frac{1}{A^{2/3}} [(41)^2 + (60)^2] \text{ MeV}^2, \quad (1.5.1)$$

resulting in

$$\hbar\omega_{GDR} \approx \frac{73}{A^{1/3}} \approx \frac{87}{R} \text{ MeV}, \quad (1.5.2)$$

where $R = 1.2A^{1/3}$ is the numerical value of the nuclear radius when measured in fm. The above quantity is to be compared with the empirical value $\hbar\omega_{GDR} \approx (80/A^{1/3})$ MeV $\approx (95/R)$ MeV.

It is of notice that the elastic vibrational frequency of a spherical solid made out of particles of mass m can be written as $\omega_{el}^2 \sim \mu/(m\rho R^2) \sim v_t^2/R^2$, where R is the radius, ρ the density and v_t the transverse sound velocity proportional to the Lamé shear modulus of elasticity μ . In other words, giant resonances in general, and the GDR in particular, are embodiments of the elastic response of the nucleus to an impulsive external field, like that provided by a photon. The nuclear rigidity to sudden solicitations is provided by the shell structure, quantitatively measured by the energy separation between major shells.

1.6 Giant pairing vibrations

Due to spatial quantization, in particular to the existence of major shells of pair degeneracy $\Omega(\equiv (2j+1)/2)$ separated by an energy $\hbar\omega_0 \approx 41/A^{1/3}$ MeV, the Cooper

⁵⁴Within this context we refer to the discussion concerning the renormalization of collective modes carried out in the text before Eq. (1.2.14), in particular regarding the cancellation between self-energy and vertex corrections (see Fig. 1.2.6 and footnote 20). This is a basic result of NFT –as discussed in later chapters– being connected with sum rule arguments in general, and particle conservation in particular. Argument which has been extended to the case of finite temperature as well as to include relativistic effects Ward (1950); Nambu (1960); Bortignon and Broglia (1981); Bertsch et al. (1983); Bortignon et al. (1998); Litvinova and Wibowo (2018); Wibowo and Litvinova (2019); Lalazissis and Ring (2019) and references therein.

⁵⁵See e.g. Bertsch and Broglia (2005) Ch. 5, and references therein.

pair model can be extended to encompass pair addition and pair subtraction modes across major shells⁵⁶. Assuming that both the single-particle energies ϵ_k and ϵ_i appearing in Fig. 1.3.1 are both at an energy $\hbar\omega_0$ away from the Fermi energy, one obtains the dispersion relation

$$-\frac{1}{G} = \frac{\Omega}{W - 2\hbar\omega_0} - \frac{\Omega}{W + 2\hbar\omega_0}, \quad (1.6.1)$$

leading to

$$(2\hbar\omega_0)^2 - W^2 = 4\hbar\omega_0 G \Omega, \quad (1.6.2)$$

and implying a high lying pair addition mode of energy

$$W = 2\hbar\omega_0 \left(1 - \frac{G\Omega}{\hbar\omega_0}\right)^{1/2}. \quad (1.6.3)$$

The forwards- (backwards)-going RPA amplitudes are, in the present case

$$X = \frac{\Lambda_0 \Omega^{1/2}}{2\hbar\omega_0 - W} \quad \text{and} \quad Y = \frac{\Lambda_0 \Omega^{1/2}}{2\hbar\omega_0 + W} \quad (1.6.4)$$

normalized according to the relation

$$1 = X^2 - Y^2 = \Lambda_0^2 \Omega \frac{8\hbar\omega_0 W}{((2\hbar\omega_0)^2 - W^2)^2}, \quad (1.6.5)$$

where Λ_0 stands for the particle-pair vibration coupling vertex. Making use of (1.6.2) one obtains

$$\left(\frac{\Lambda_0}{G}\right)^2 = \Omega \left(1 - \frac{G\Omega}{\hbar\omega_0}\right)^{-1/2} \quad (1.6.6)$$

quantity corresponding, within the framework of the simplified model used, to the two-nucleon transfer cross section. Summing up, the monopole giant pairing vibration has an energy close to $2\hbar\omega_0$, and is expected to be populated in two-particle transfer processes with a cross section of the order of that associated with the low-lying pair addition mode, being this one of the order of Ω . Simple estimates of (1.6.3) and (1.6.6) can be obtained making use of $\Omega \approx \frac{2}{3}A^{2/3}$ and $G \approx 17/A$ MeV, namely

$$W = 0.85 \times 2\hbar\omega_0, \quad \left(\frac{\Lambda_0}{G}\right)^2 \approx 1.2\Omega. \quad (1.6.7)$$

Experimental evidence of GPV in light nuclei have been reported⁵⁷.

⁵⁶Broglia and Bes (1977).

⁵⁷Cappuzzello et al. (2015); Cavallaro et al. (2019), Bortignon and Broglia (2016). See also Laskin et al. (2016); Id Betan et al. (2002); Dussel et al.; Mouginot et al. (2011); Khan et al. (2004); Avez et al. (2008); Khan et al. (2009); Dasso et al. (2015); Fortunato et al. (2002); Herzog et al. (1986); Id Betan et al. (2002).

1.7 Sum rules

There are important operator identities which restrict the possible matrix elements in a physical system. Let us calculate the double commutator of the Hamiltonian describing the system and a single-particle operator F . That is

$$[\hat{F}, [\hat{H}, \hat{F}]] = (2\hat{F}\hat{H}\hat{F} - \hat{F}^2\hat{H} - H\hat{F}^2). \quad (1.7.1)$$

Let us assume that $\hat{F} = \sum_k F(\mathbf{r}_k)$ and $H = T + v(\mathbf{r}, \mathbf{r}')$, where $v(\mathbf{r}, \mathbf{r}') = -\kappa_1 \hat{F}(\mathbf{r})\hat{F}(\mathbf{r}')$. Thus

$$[\hat{F}, [\hat{H}, \hat{F}]] = \sum_k \frac{\hbar^2}{m} (\nabla_k F(\mathbf{r}_k))^2. \quad (1.7.2)$$

Let us take the average value on the correlated ground state

$$S(F) = \sum_{\alpha'} |\langle \alpha' | F | \tilde{0} \rangle|^2 (E_{\alpha'} - E_0) = \frac{\hbar^2}{2m} \int d\mathbf{r} |\nabla F|^2 \rho(\mathbf{r}), \quad (1.7.3)$$

where we have used $H|\alpha\rangle = E_\alpha|\alpha\rangle$ and $H|\tilde{0}\rangle = E_0|\tilde{0}\rangle$, and the sum $\sum_{\alpha'}$ is over the complete set of eigenstates of the system. The above result describes the reaction of a system at equilibrium to which one applies an impulsive field, giving to the particles a momentum ∇F . On the average, the particles started at rest so their mean energy after the sudden impulse is $\hbar^2|\nabla F|^2/2m$, a result which does not depend on the interaction acting among the nucleons, the energy being absorbed from the (instantaneous) external field before the system is disturbed from equilibrium. The relation (1.7.3) is known as the energy weighted sum rule (EWSR).

An important application of (1.7.3) implies a situation where F has a constant gradient. Inserting $\mathbf{F} = z \times \mathbf{z}$ in (1.7.3), the integral simplifies because $\nabla F = 1$, and the integral leads just to the number of particles,

$$\sum_{\alpha} |\langle \alpha | F | \tilde{0} \rangle|^2 (E_{\alpha} - E_0) = \frac{\hbar^2 N}{2m}. \quad (1.7.4)$$

The electric field of a photon is of this form in the dipole approximation, which is valid when the size of the system is small compared to the wavelength of the photon, the single-particle field being

$$F(\mathbf{r}) = e \sum_k \left[\frac{N-Z}{A} - t_z(k) \right] r_k Y_{1\mu}(\hat{r}_k), \quad (1.7.5)$$

with $t_z = -1/2$ for protons and $+1/2$ for neutrons, t_z being the z -component of the isospin operator. For the dipole operator referred to the nuclear center of mass one obtains

$$\sum_{\alpha'} |\langle \alpha' | F | \tilde{0} \rangle|^2 (E_{\alpha'} - E_0) = \frac{9}{4\pi} \frac{\hbar^2 e^2}{2m} \frac{NZ}{A}. \quad (1.7.6)$$

The above relation is known as the Thomas-Reiche-Kuhn (TRK) sum rule, and is equal to the maximum energy a system can absorb from the dipole field. The RPA solution respects the EWSR.

1.8 Ground state correlations

The zero point fluctuations associated with collective nuclear vibrations affect, among other thing, the nuclear mean field properties⁵⁸. Within the harmonic approximation discussed in Sections 1.1.1 and 1.1.2 in connection with ($\beta = 0$) surface vibrational modes, the ground state Ψ_0 of (1.1.9) has an energy $\frac{1}{2}\hbar\left(\frac{C_\lambda}{D_\lambda}\right)^{1/2} = \hbar\omega_\lambda/2$ for each degree of freedom (e.g. $(2\lambda + 1)$ in the case of a vibration of multipolarity λ). The associated coordinate-momentum indeterminacy relation assumes its minimum value⁵⁹

$$\Delta\alpha_{\lambda\mu}(n)\Delta\pi_{\lambda\mu}(n) = \frac{\hbar}{2}, \quad (1.8.1)$$

as

$$\Delta\alpha_{\lambda\mu}(n) = \sqrt{\frac{\hbar\omega_\lambda(n)}{2C_\lambda(n)}}, \quad (1.8.2)$$

and $\Delta\pi_{\lambda\mu} = \sqrt{\hbar D\omega_\lambda/2}$, and in keeping with the fact that $|\Psi_0|^2$ is mathematically a Poisson distribution. The above expression implies that the mean square radius will be modified from its mean field value R_0 and thus also the nuclear density distribution. The value of $\hbar\omega_\lambda(n)/2C_\lambda(n)$ can be determined by working out the collective mode $|n_\lambda(n) = 1\rangle = \Gamma_{\lambda\mu}^\dagger(n)|\tilde{0}\rangle$ in RPA, the zero point fluctuations being closely connected with the Y -amplitudes of the mode.

Let us start by calculating the effect of the zero point fluctuations on the nuclear density distribution. The corresponding operator can be written as

$$\hat{\rho}(\mathbf{r}) = a^\dagger(\mathbf{r})a(\mathbf{r}), \quad (1.8.3)$$

⁵⁸Gogny (1978); Esbensen and Bertsch (1983); Reinhard and Drechsel (1979); Khodel et al. (1982); Barranco and Broglia (1987, 1985). See also Brown and Jacob (1963); Anderson and Thouless (1962) (it is likely a coincidence in connection with this inaugural issue of Phys. Lett. that short of hundred pages after, one finds the paper of B. D. Josephson, Possible new effects in superconducting tunneling, Phys. Lett. **1**, 251 (1962)).

⁵⁹The same result is found for Ψ_n describing a state with n -quanta, being at the basis that solutions with $n \gg 1$ behaves as “quasiclassical” or “coherent” states of the harmonic oscillator (Glauber (2007)) in keeping with the fact that the contribution of the zero point energy is negligible in such case ($(n + 1/2)\hbar\omega \approx n\hbar\omega$) and that the many quanta wavepacket always attain the lower limit of (1.8.1) (Basdevant and Dalibard (2005) pp. 153,465) (discussions with Pier Francesco Bortignon in March 2018 concerning coherent states are gratefully acknowledged). Schrödinger used this result in a paper (Schrödinger (1926b)) to suggest that waves (material waves) described by his wave function were the only reality, particles being only derivative things. In support of his view he considered a superposition of linear harmonic oscillator wavefunctions and showed that the wave group holds permanently together in the course of time. And he adds that the same will be true for the electron as it moves in high orbits of the hydrogen atom, hoping that wave mechanics would turn out to be a branch of classical physics (Pais (2000)). It was Born who first provided the correct interpretation of Schrödinger’s wavefunction (modulus square) in his paper “Quantum mechanical collision phenomena” (Born (1926a)). In it it is stated that the result of solving with wave mechanics the process of elastic scattering of a beam of particles by a static potential is not what the state after the collision is, but how probable is a given effect of the collision.

where $a^\dagger(\mathbf{r})$ is the creation operator of a nucleon at point \mathbf{r} . It can be expressed in terms of the phase space creation operators $a_\nu^\dagger(\nu \equiv n, l, j, m)$ as

$$a^\dagger(\mathbf{r}) = \sum_\nu \varphi_\nu^*(\mathbf{r}) a_\nu^\dagger, \quad (1.8.4)$$

where $\varphi_\nu(\mathbf{r})$ are the single-particle wavefunctions. Thus

$$\hat{\rho}(\mathbf{r}) = \sum_{\nu\nu'} \varphi_\nu^*(\mathbf{r}) \varphi_{\nu'}(\mathbf{r}) a_\nu^\dagger a_{\nu'}. \quad (1.8.5)$$

The matrix element in the HF ground state is (Fig. 1.2.2 (h))

$$\rho_0(\mathbf{r}) =_F \langle 0 | \hat{\rho}(\mathbf{r}) | 0 \rangle_F = \sum_{i, (\epsilon_i \leq \epsilon_F)} |\varphi_i(\mathbf{r})|^2. \quad (1.8.6)$$

To lowest order of perturbation theory in the particle-vibration coupling vertex, the NFT diagrams associated with the change of ρ_0 due to ZPF are shown in Fig. 1.8.1. Graphs (a), (b) and (c), (d) describe the changes in the density operator and in the single-particle potential respectively⁶⁰. This can be seen from the insets (I) and (II). The dashed horizontal line starting with a cross and ending at a hatched circle represents the renormalized density operator. This phenomenon is similar to that encountered in connection with vertex renormalization in Fig. 1.2.6, that is the renormalization of the particle-vibration coupling (insets (I) and (I')). Concerning potential renormalization, the bold face arrowed line shown in inset (II) of Fig. 1.8.1 represents the motion of a renormalized nucleon due to the self-energy process induced by the coupling to vibrational modes. A phenomenon which can be described at profit through an effective mass, the so called ω -mass m_ω , in which case particle motion is described by the Hamiltonian⁶¹ $(\hbar^2/2m_\omega) \nabla^2 + \left(\frac{m}{m_\omega}\right) U(r)$. The ω -mass can be written as $m_\omega = (1 + \lambda)m$, where m is the nucleon mass and λ is the so called mass enhancement factor $\lambda = N(0)\Lambda$, where $N(0)$ is the density of levels at the Fermi energy, and Λ the PVC vertex strength, typical values being $\lambda = 0.4$.

The fact that in calculating $\delta\rho$, that is, the correction to the nuclear density distribution (renormalization of the density operator), one finds to the same order of perturbation a correction to the potential, is in keeping with the self consistency existing between the two quantities (Eq. (1.2.8)). *Now, what changes is not only the single-particle energy, but also the single-particle content measured by $Z_\omega = m/m_\omega$, as well as the radial dependence of the wavefunctions of the states.* It is of notice that the effective mass approximation, although being quite useful, cannot take care of the energy dependence of the renormalization process which leads, in the case of single-particle motion to renormalized energies, spectroscopic amplitudes and radial wavefunctions. The analytic expressions associated with

⁶⁰Barranco (1985), Barranco and Broglia (1985).

⁶¹The ratio (m/m_ω) gives a measure of the single-particle energy content (Mahaux et al. (1985), Eq. (3.5.18)), while m_ω is proportional to the (energy-) slope of the (single-particle) self-energy dispersion relation (Sect. 5.4.1; see also Brink and Broglia (2005) and refs. therein).

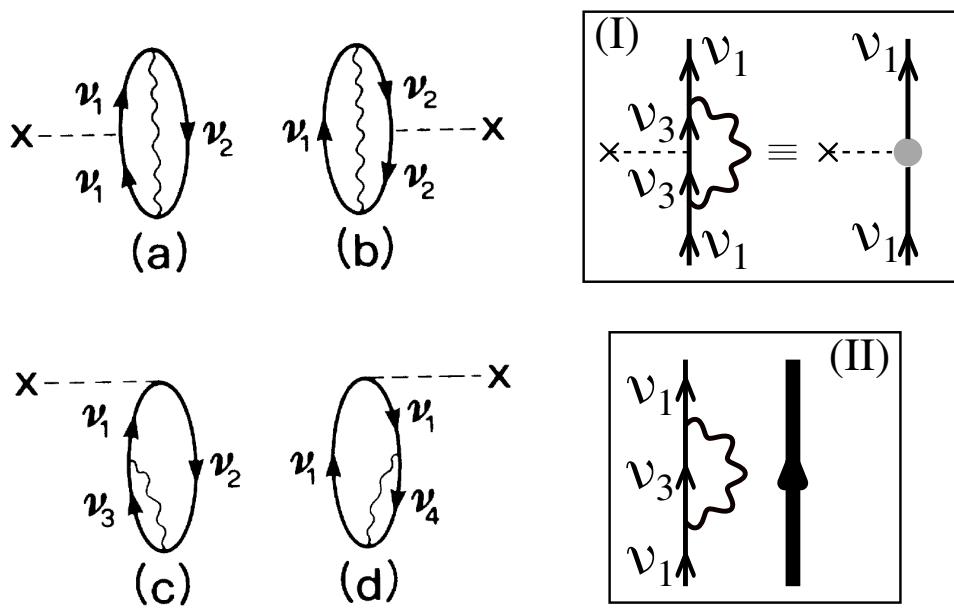


Figure 1.8.1: Lowest-order corrections in the particle-vibration coupling vertex of the nuclear density due to the presence of zero-point fluctuations associated with density vibrations. An arrowed line pointing upwards denotes a particle, while one pointing downward a hole. A wavy line represents a surface phonon. The density operator is described through a dotted horizontal line starting with a cross. Graphs (a) and (b) are typical examples of density contributions to $\delta\rho$ (see inset (I); the dashed horizontal line starting with a cross and ending at a hatched circle in the diagram to the right, represents the renormalized density operator, resulting from the processes displayed to the left); (c) and (d) are examples of potential contributions (see inset (II); the bold face arrowed line represents the renormalized single-particle state due to the coupling to the vibrations leading to the self energy process shown to the left).

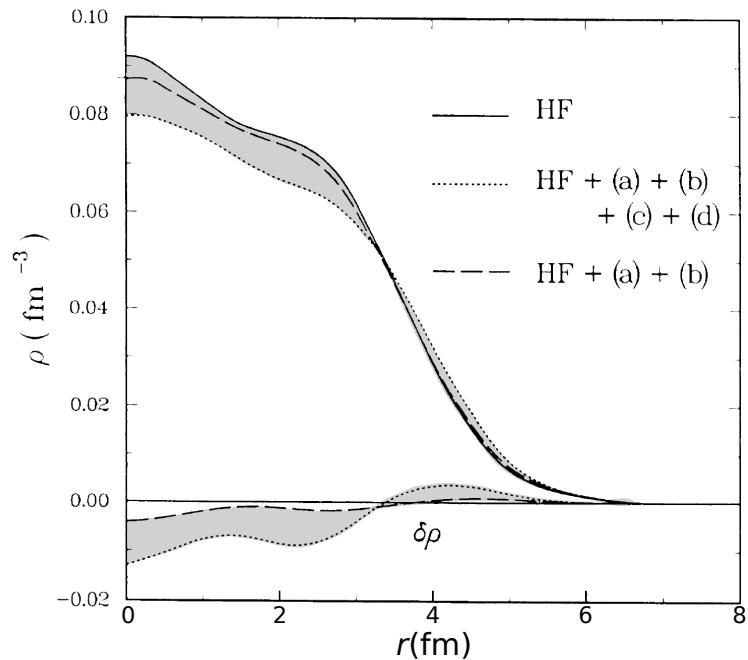


Figure 1.8.2: Modification in the charge density of ^{40}Ca induced by the zero-point fluctuations associated with vibrations of the surface modes. The results labeled HF, HF+(a)+(b), and HF+(a)+(b)+(c)+(d) are the Hartree-Fock density, and that resulting from adding to it the corrections $\delta\rho$ associated with the processes (a)+(b) and (a)+(b)+(c)+(d) displayed in Fig. 1.8.1, respectively. In the lower part of the figure the quantities $\delta\rho$ are displayed.

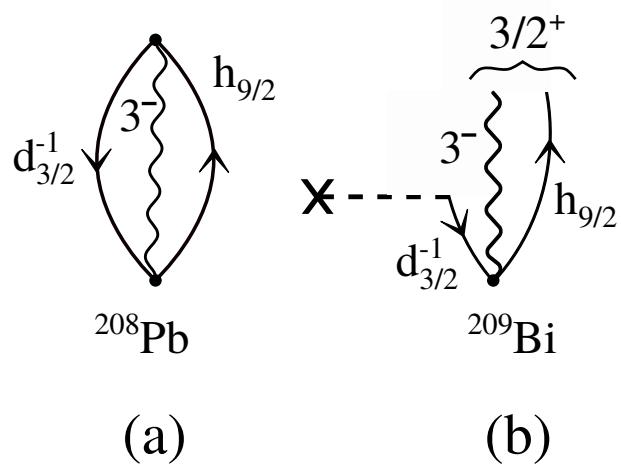


Figure 1.8.3: (a) Example of zero point fluctuation of the ground state of the double-magic nucleus $^{208}_{82}\text{Pb}_{126}$ associated with the low-lying octupole vibration of this system, observed at an energy of 2.615 MeV and displaying a collective electromagnetic decay to the ground state of $32 B_{sp}$. The **proton** particle-hole component $(h_{9/2}, d_{3/2}^{-1})_{3^-}$ displayed carries a large amplitude in the octupole vibration wavefunction. (b) Diagram representing the transfer of one proton to ^{208}Pb , which fills the $d_{3/2}^{-1}$ hole state leading to a $3/2^+$ in $^{209}_{83}\text{Bi}_{126}$, member of the septuplet of states $|(3^- \otimes h_{9/2} J^\pi)\rangle$ with $J^\pi = 3/2^+, 5/2^+, \dots, 15/2^+$. The horizontal dashed line starting with a cross stands for the stripping process ($^3\text{He}, d$).

diagrams (a) and (c) of Fig. 1.8.1 are

$$\delta\rho(r)_{(a)} = \frac{(2\lambda+1)}{4\pi} \sum_{\nu_1\nu_2n} [Y_n(\nu_1\nu_2; \lambda)]^2 R_{\nu_1}(r)R_{\nu_2}(r), \quad (1.8.7)$$

and

$$\begin{aligned} \delta\rho(r)_{(c)} &= (2\lambda+1)\Lambda_n(\lambda) \sum_{\nu_1\nu_2\nu_3} \frac{M(\nu_1, \nu_3; \lambda)}{\epsilon_{\nu_1} - \epsilon_{\nu_2}} (2j_1 + 1)^{-1/2} \\ &\times Y_n(\nu_3\nu_2; \lambda) \times R_{\nu_1}(r)R_{\nu_2}(r), \end{aligned} \quad (1.8.8)$$

where M is the matrix element of $-\frac{R_0}{\kappa} \frac{\partial U}{\partial r} Y_{\lambda\mu}(\hat{r})$, $n = 1, 2, \dots$ labeling the first, second, etc vibrational modes as a function of increasing energy, and Λ_n is the strength of the particle-vibration coupling associated with the n -mode of multipolarity λ . The functions $R(r)$ are the radial wavefunctions associated with the states ν_1 and ν_2 . While $\delta\rho_{(a)}$ can be written in terms of the RPA Y -amplitudes which are directly associated with the zero point fluctuations of harmonic motion, $\delta\rho_{(c)}$ contains a scattering vertex not found in RPA – that is going beyond the harmonic approximation – and essential to describe renormalization processes of the different degrees of freedom, namely single-particle (energy, single-particle content and radial dependence of the wavefunction) and collective motion, as well as medium polarization interactions. In particular the induced pairing interaction (see Sect. 4.8.1).

In Fig. 1.8.2 results of calculations of $\delta\rho$ carried out for the closed shell nucleus ^{40}Ca are shown. The vibrations were determined by diagonalizing multipole-multipole separable interactions in the RPA. All the roots of multipolarity and parity $\lambda^\pi = 2^+, 3^-, 4^+$ and 5^- which exhaust the EWSR were included in the calculations. Both isoscalar and isovector degrees of freedom were considered, as well as low-lying and giant resonances.

From the point of view of single-particle motion, the vibrations associated with low-lying modes display rather low frequency ($\hbar\omega_\lambda/\epsilon_F \approx 0.1$) and lead to an ensemble of essentially adiabatic deformed shapes. Nucleons can thus reach to distances from the nuclear center which are considerably larger than the radius R of the static spherical potential. Because the frequency of the giant resonances are of similar magnitude to those corresponding to the single-particle motion, the associated surface deformations average out. Thus low-lying vibrational modes account for most of the contributions to the changes in the density distribution⁶².

Making use of $(\delta\rho)_{low-lying}$, the change in the mean square radius was calculated, leading⁶³ to $\delta\langle r^2 \rangle = 0.494 \text{ fm}^2$. It implies a relative correction $\delta\langle r^2 \rangle / ((3/5)R_0^2) \approx$

⁶² Another example of the recurrent central role played by low-frequency modes in determining the properties and behavior of systems at all levels of organization, from the atomic nucleus to the Casimir effect in QED (see Fig. 7.E.3), to superconductivity in metals (Sect. 4.8.2) as well as to the folding of proteins (Micheletti et al. (2004)) in connection with metabolism and of brain activity ($\nu < 0.1 \text{ Hz}$) (Mitra et al. (2018); Vyazovskiy and Harris (2013) in different phases of sleep).

⁶³ $\delta\langle r^2 \rangle = (0.371 + 0.123)\text{fm}^2 = 0.494\text{fm}^2$, see fourth column (Micr. (a)+(b)+(c)+(d)) of Table 1 of Barranco and Broglia (1987).

5%, where $R_0 = r_0 A^{1/3} = 4.1$ fm is the value of ${}^{40}\text{Ca}$ radius obtained from systematics ($r_0 = 1.2$ fm), $(3/5)R_0^2 \approx 10.11\text{fm}^2$ being the associated value of the mean square radius.

Similar calculations to the ones discussed above, but in this case taking into account only the contributions of the low-lying octupole vibration⁶⁴ indicate that nucleons are to be found a reasonable part of the time in higher shells than those assigned to them by the shell model. The average number of “excited” particles being ≈ 2.4 . Correspondingly, hole states will be left behind. Because of their presence in the closed shell nucleus, one can transfer a nucleon to states below the Fermi energy in, for example, (d, p) or $({}^3\text{He}, d)$ one-neutron or one-proton stripping reactions respectively, leaving the final nucleus with one-nucleon above closed shell coupled to a vibration.

Systematic studies of such multiplets have been carried out throughout the mass table. In particular around the closed shell nucleus ${}_{82}^{208}\text{Pb}_{126}$ (Fig. 1.8.3). Within this context it is not only quite natural but also necessary, to deal with structure and reactions on equal footing. This is one of the main goals of the present monograph.

1.9 Interactions

A number of subjects are not touched upon in the present monograph. In particular, the role temperature plays in the structure and decay of nuclei and that of bare (e.g. the Argonne NN -potential) and/or effective forces (e.g. Gogny, Skyrme, etc.). In what follows we briefly elaborate on this last point, referring to Bortignon et al. (1998) and refs. therein concerning the first one.

The Coulomb interaction resulting from the exchange of photons between charged particles defines the domain area of quantum electrodynamics (QED). Being the electron and the photon fields in interaction, what we call an electron is only partially to be associated with the electron field alone. It is also partially to be associated with the photon field which dresses the electron field. And conversely what is called the photon field can materialize itself in space in terms of an electron and a positron⁶⁵.

The Coulomb interaction is the best known of all physical interactions, and QED constitutes the paradigm of theories which can be considered correct. In natural units in which the magnetic moment of the electron emerging from Dirac equation has the value of 1, the results of QED corrections agree –once divergent contributions have been renormalized by properly adjusting the value of the bare electron mass and charge⁶⁶–, within experimental errors, namely down to the

⁶⁴Brown and Jacob (1963). See also Anderson and Thouless (1962).

⁶⁵Schwinger (2001).

⁶⁶Changing indirectly the Coulomb interaction. Not that observed in the laboratory and acting between physical (dressed) electrons, but the “bare” Coulomb interaction acting between the “naked” electron fields, and used in the calculation of the variety of processes through which the photon field dresses the electrons.

eleventh decimal figure, with observation. That is $1.00115965221 \pm 0.00000000003$ (exp.); $1.00115965246 \pm 0.00000000020$ (QED)⁶⁷.

The energy difference between the $2S_{1/2}$ and the $2P_{1/2}$ states of the hydrogen atom which according to Dirac's theory should be degenerate, emerges naturally, according to QED, from the dressing of the hydrogen's electron by the photon associated with the zero point fluctuation of the electromagnetic vacuum. The measured value of the Lamb shift⁶⁸ is 1057.845 (9) MHz, an experimental value⁶⁹ whose accuracy is limited by the ≈ 100 MHz natural line width of the $2P_{1/2}$ state. The best QED value, limited by uncertainties in the radius of the proton, is 1057.865 MHz⁷⁰

The above two numbers (electron magnetic moment and Lamb shift) are results of *ab initio* calculations within the framework of an effective field theory (QED). *Ab initio* to the extent that quantum mechanics in general⁷¹ and the Dirac equation⁷² in particular can be considered, on grounds of their universal validity, to be so, including QED with its bare mass and charge parameters used to cure infinities (renormalization)⁷³.

Nuclear field theory (NFT) was developed following the (graphical) approach adopted by Feynman in his formulation of QED, making use of the correspondence electron→nucleon, photons→collective vibrations, fine structure constant→particle-vibration coupling. In keeping with the fact that in Feynman's QED *nothing is really free*⁷⁴ (*bare*), a fact that transfers to NFT, and of the very large value of the induced (retarded) nuclear interaction –as large or larger than that of the bare NN -interaction in e.g. the case of the 1S_0 (pairing channel)⁷⁵– arising from the exchange of collective vibrations between nucleons, the role which specific (*bare*) nuclear forces (four-point vertices v) play in the results of NFT calculations, becomes somewhat blurred.

Starting from the mean field (see Eq. (1.2.8)) at the basis of the non-observable bare single-particle energies, one employs a generic shape (Woods-Saxon), adjusting the depth, radius and diffusivity of the central potential, and the depth of the spin-orbit one in connection with an r -dependent effective k -mass (exchange poten-

⁶⁷Kinoshita (1990).

⁶⁸Lamb Jr (1964).

⁶⁹Lundeen and Pipkin (1981, 1986); Pipkin (1990).

⁷⁰Sapirstein and Yennie (1990); Grotch (1994).

⁷¹Heisenberg (1925); Born and Jordan (1925); Born et al. (1926); Dirac (1925); Schrödinger (1926b); Born (1926a); Heisenberg (1927); Pauli (1925).

⁷²Dirac (1928a,b).

⁷³Feynman (1949); Schwinger (1948); Tomonaga (1946); see also Dyson (1949); Schwinger (1958); Schweber (1994).

⁷⁴Feynman (1975).

⁷⁵For example, concerning the contribution to the pairing gap of ^{120}Sn ($\Delta \approx 1.45$ MeV) arising from the bare 1S_0 (pairing) interaction v_{14} (Argonne NN -potential) and that associated with the exchange of collective vibrational modes between nucleons moving in time reversal states close to the Fermi energy (Idini et al. (2015)). In the case of ^{11}Li , essentially all of the pairing interaction which binds the two halo neutrons to the core ^9Li , arise from the exchange between them, of the soft $E1$ -mode (pygmy dipole resonance of ^{11}Li). (Barranco et al. (2001); Broglia et al. (2019a)).

tial) so that the dressed single-particle states reproduce the experimental findings⁷⁶. Parallel to the bare mass parameter entering renormalized QED. In connection with the coupling between nucleons and surface vibrations, namely the strength Λ_α , one adjusts κ so as to reproduce the properties of the collective modes (see inset Fig. 1.2.3), a procedure which parallels the tuning of the bare charge, and thus the strength of the coupling between electrons and photons (and of the bare Coulomb interaction) in renormalized QED. In other words, renormalized nuclear field theory (NFT)_{ren}

The Mayer and Jensen sequence of levels around the $N = 8$ magic nuclei is $1p_{1/2}, 1d_{5/2}, 2s_{1/2}$ (Fig. 1.1.2), while experimentally $^{11}_4\text{Be}_7$ displays the sequence⁷⁷ $1/2^+, 1/2^-$ (bound), $5/2^+$ (resonance). A consequence of the dressing of the bare $1p_{1/2}, 1d_{5/2}$ and $2s_{1/2}$ states by the quadrupole vibration⁷⁸ of ^{10}Be ($\beta_2 \approx 0.9$). This extremely large value of the dynamical deformation parameter implies a Lamb-shift like mechanism which shifts the bare $1p_{1/2}$ and $2s_{1/2}$ orbitals by more than 3 MeV with respect to each other, the resulting dressed levels reproducing the experimental ones ($\epsilon_{\widetilde{1/2}^+} = -0.5$ MeV, $\epsilon_{\widetilde{1/2}^-} = -0.18$ MeV, i.e. $\Delta\epsilon = (\epsilon_{\widetilde{1/2}^-} - \epsilon_{\widetilde{1/2}^+}) = 0.32$ MeV, see Fig. 5.2.4).

The predicted $E1$ -decay between the parity inverted levels is $B(E1; \widetilde{1/2}^- \rightarrow \widetilde{1/2}^+) = 0.11e^2 \text{ fm}^2$, a quantity to be compared with the experimental value $B(E1) = 0.102 \pm 0.002e^2 \text{ fm}^2$. The two (three) numbers ($\epsilon_{\widetilde{1/2}^+}, \epsilon_{\widetilde{1/2}^-}, \Delta\epsilon$), and $B(E1)$) can, arguably, be used to assess the accuracy of renormalized NFT results on something which one can call equal footing, with the two numbers (electron magnetic moment and Lamb shift) discussed above in connection with the assessment of the level of accuracy of QED results.

It would be computationally useful and conceptually satisfying to have a bare NN -potential, eventually derived from QCD, reproducing the NN -phase shifts and which employed in nuclear structure calculations, lead to bare static (Eq. (1.2.8)) and time-dependent (dynamic, Eq. (1.2.9)) mean fields whose solutions (bare single-particle states and bare vibrational modes), interweaved according to the rules of a field theory, e.g. NFT (Sect. 2.5.2), accounted for the value of the observables resulting from a “complete” set of probes (see e.g. Figs. 2.8.1 and 5.2.6). Concerning single-particle motion: energies, absolute one-particle transfer differential cross sections and thus insight into both single-particle content and renormalized single-particle wavefunctions (form factors), γ -decay and associated effective charges, etc. Concerning collective vibrations: energies, γ -decay, absolute differential cross sections for inelastic and Coulomb excitations (surface modes) and for two-nucleon transfer processes (pairing vibrations). Until such desideratum becomes reality⁷⁹, the aim at shedding light onto the physics at the basis of new experimental results and that of providing guidance in the quest to achieve a

⁷⁶See e.g. Barranco et al. (2017) and Barranco et al. (2020).

⁷⁷Kwan et al. (2014).

⁷⁸Barranco et al. (2017).

⁷⁹See however footnote 10 of this Chapter.

deeper and unified picture of nuclear structure and reactions, can be carried out in terms of empirical renormalization, i.e. (NFT)_{ren}.

In keeping with the fact that the domain area of QED is all of chemistry and much of biology⁸⁰, the important *renormalization effects* undergone by *bare forces* when used to describe even one of the simplest (real) many-body system –i.e. the hydrogen atom in the case of the Coulomb interaction– is quite universal. Because of the variety of facets and the importance this issue has in connection with the unified nuclear field theory of structure and reactions we discuss in the present monograph, tailored after Feynman’s diagrammatic version of QED, we attempt at shedding light on bare force-renormalization process phenomena through a number of interdisciplinary examples. A crown one is provided by dispersive (retarded) forces like the van der Waals interaction (App. 3.B). Also by the Casimir effect and the hydrophobic force (App. 7.E), let alone by the renormalization the Coulomb force undergoes in metals (Sect. 4.8). Dressing which in this case changes not only the value, but also the sign of the interaction resulting eventually in the phenomenon of superconductivity. Hallmarks of which are perfect conductivity and perfect diamagnetism, and the special effects found in terms of the renormalized Coulomb force acting across a Josephson junction (Sect. 4.6).

Within the above scenario one can mention that at the basis of the mechanism which inverts, in the odd $N = 7$ isotones ${}^3_3\text{Li}$ and ${}^11_4\text{Be}$, the standard Mayer-Jensen sequence of levels $1p_{1/2}$ and $2s_{1/2}$ and comes close to do so also concerning the $1d_{5/2}$ and $1p_{1/2}$ single-particle levels constitutes, arguably, the largest Lamb shift-like phenomenon displayed by a physical system.

1.10 Hindsight

In order for a nucleon moving in a level close to the Fermi energy to display a mean free path larger than nuclear dimensions and to be reflected at the nuclear surface through an elastic process, all other nucleons must move in a rather ordered, correlated fashion. Within this context to posit that single-particle motion is the most collective of all nuclear motions⁸¹ seems natural. Associated with mean field and single-particle motion one finds the typical bunching of the corresponding levels closely connected with the major shells lying above and below the Fermi energy. A fact which determines the parity (and angular momentum) of the lowest excitations associated with promoting a nucleon across the Fermi surface. By correlating these excitations through the same components of the NN -interaction leading to the single-particle bunching within the mean field approximation, one obtains collective multipole vibrations of particle-hole type. Similar arguments result in the presence of multipole pairing vibrations in the low-energy spectrum, and in their coupling to single-particle motion.

It is then natural to consider single-particle motion and collective states on

⁸⁰Feynman (2006).

⁸¹Mottelson (1962),

equal footing, and as basis states of a physical description of the atomic nucleus in which to diagonalize both three point (PVC) and four-point (v) vertices , eventually taking in this second case also $3N$ terms into account⁸². Such an approach provides also direct indication of the minimum set of experiments needed to obtain a “complete” picture (test) of the low-energy properties of the atomic nucleus. They constitute the specific probes of each of the basis states (elementary modes of excitation). Namely inelastic scattering and Coulomb excitation (particle-hole collective vibrations and rotations), one-particle transfer processes (independent-particle motion) and two-particle transfer reactions (pairing vibrations and rotations).

Summing up, because of the interweaving existing between the variety of elementary modes of excitation, experimental probes associated with fields which carry transfer quantum number $\beta = 0, \pm 1$ and ± 2 , and different multipolarities, spins and isospins are needed to characterize the structure of nuclei. This is what we attempt at explaining and formulating in the following chapters, setting special emphasis on transfer processes, in which case the relative motion of the reacting nuclei and the intrinsic motion of the nucleons in target and projectile cannot be separated, and one is forced to treat structure and reactions in an unified way.

⁸²See footnote 10 of this Chapter.

Chapter 2

Structure and reactions

To introduce the quanta of nuclear physics, let us consider, as an example the spectrum of elementary excitations based on the ground state of ^{208}Pb . This “vacuum” state has specially simple properties (has a minimum of degeneracy) on account of the closed-shell configuration of 82 protons and 126 neutrons

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In what follows, the connection between the concept of elementary modes of excitation and associated specific probes is discussed in terms of selected experiments. Connection which finds its ultimate test in terms of predicted and measured absolute transition probabilities and differential cross sections.

2.1 Elementary modes of excitation and specific probes

Subject to external probes which couple weakly to the nucleus, that is, in such a way that the system can be expressed in terms of the properties of the excitation in the absence of probes¹, the nucleus reacts in terms of single-particle (-hole) motion (one-particle transfer), vibrations (surface, spin, etc.) and rotations (Coulomb excitation and inelastic scattering) and pairing vibrations and rotations (two-nucleon transfer reactions), that is, in terms of elementary modes of excitation.

Collective vibrations in nuclei can be characterized by a variety of quantum numbers. In particular angular momentum (J), parity (π) and transfer quantum number (β). Let us consider the doubly magic nucleus $^{208}\text{Pb}_{126}$ to illustrate some

¹ See e.g. Pines and Nozières (1966), Bohr and Mottelson (1975) and refs. therein. Within the context of linear response see also Sect. 7.4, text after Eq. (7.4.1).

aspects of the transfer quantum number β . The ground state of this nucleus ^{208}Pb , namely $|\text{gs}({}^{208}\text{Pb})\rangle$ has angular momentum 0 and positive parity ($J^\pi = 0^+$). It can be viewed as the vacuum state of the variety of elementary modes of excitation at the basis of the nuclear field theory (NFT) description of this system. The lowest lying collective vibration ($E_x \approx 2.6$ MeV) is an octupole surface vibration. Microscopically, it can be viewed as a correlated particle-hole (p, h) excitation. Consequently, this state is characterized by the quantum numbers $J^\pi = 3^-, \beta = 0$. A neutron moving around ^{208}Pb in levels above the Fermi energy can have quantum numbers $j^\pi = 9/2^+, 11/2^+, 5/2^+, \dots (2g_{9/2}, 1i_{11/2}, 3d_{5/2}, \dots)$, all of them having $\beta = 1$.

In the case of quantum electrodynamics (Feynman graphical formulation²), these two modes (octupole and single-particle modes) parallel the photon and the electron. The nucleus ^{208}Pb , aside from having a rich variety of $\beta = 0$ collective (p, h) correlated modes (“photons”) like the $J^\pi = 2^+$ (4.1 MeV), 5^- (3.2 MeV) and the octupole vibration, it also displays collective $\beta = \pm 2$ modes. These pairing vibrations can be viewed as correlated two-particle ($\beta = +2$) or two-hole ($\beta = -2$) vibrations. For example, pair addition modes ($\beta = +2$) with quantum numbers $J^\pi = 0^+, 2^+, 4^+, \dots$ correspond to the ground state and to the lowest $2^+, 4^+, \dots$ states of ^{210}Pb . In particular, the $|J^\pi = 0_1^+, \beta = 2\rangle \equiv |\text{gs}({}^{210}\text{Pb})\rangle$ can be viewed as a nuclear Cooper pair³: a weakly correlated pair of fermions moving in time reversal states lying close to the Fermi energy. Cooper pairs are the building blocks of the microscopic theory of superconductivity developed by Bardeen, Cooper, Schrieffer (BCS)⁴, and of its extension to the nuclear case⁵.

In Figs. 2.1.1–2.1.3, examples of specific reactions which have identified different elementary modes of excitation of ^{208}Pb mentioned above, are given. In particular, a cartoon representation of elastic, inelastic, one- and two-particle direct transfer reactions induced by alpha, deuteron and triton projectiles impinging on ^{208}Pb are shown. In all cases a standard setup is used, in which a light projectile is aimed at a fixed target (thin foil made out of ^{208}Pb). The outgoing particles carrying the corresponding physical information, i.e. momentum, angular momentum, energy, etc. transferred to or from the target, are deflected by the electromagnetic fields of a spectrometer and eventually recorded at a given angle by particle detectors (points a,b, ..., 0⁺, 2⁺, ... in a magnet). These events provide structural information as shown in the two dimensional strength function displayed below the cartoon laboratory setup.

The associated strength functions, recorded at a wide range of angles provide the absolute differential cross sections associated with each of the nuclear states populated in the process. They are typically measured in⁶ millibarns per steradian

²“The practical usefulness of the Feynman rules and diagrams made them one of the most essential elements of the scientific training of every theoretical physicist” (Mehra, 1996).

³See footnote 27 Ch. 1.

⁴Bardeen et al. (1957a,b).

⁵Bohr et al. (1958).

⁶A barn is defined as $1\text{b}=10^{-28}\text{m}^2 = 100\text{fm}^2$. In three dimensions, the solid angle Ω is related

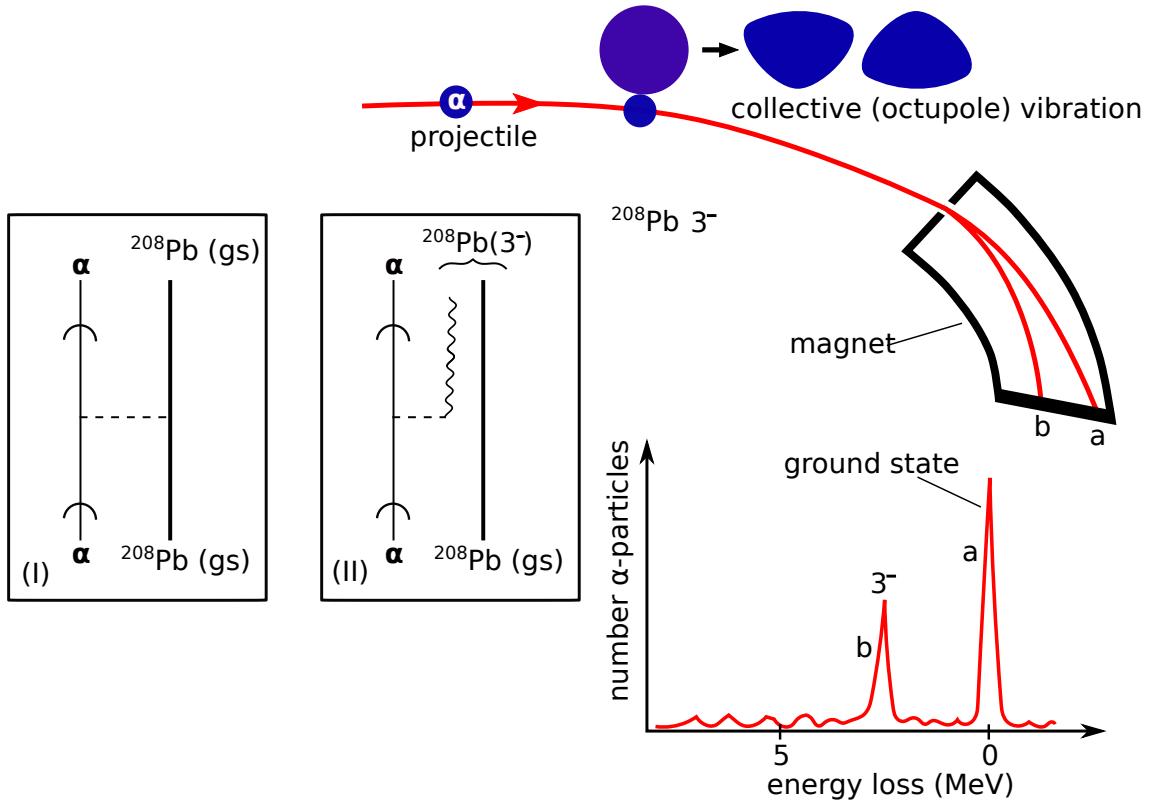


Figure 2.1.1: Schematic representation of: an **elastic** process labeled a in both the spectrograph and in the strength function plotted below it (population of the ground state), and **inelastic** b (population of the lowest octupole vibration lying at 2.62 MeV) processes associated with the reaction $^{208}\text{Pb}(\alpha, \alpha')^{208}\text{Pb}^*$. In the inset (I) a schematic nuclear field theory reaction plus structure, NFT_(s+r), diagram describing the elastic process (potential scattering, dashed horizontal line) is displayed. The α -projectile moving in the continuum is represented by an arrowed (curved) line. From the measurement of the elastic differential cross section one can deduce the partial wave phase shifts (Sect. 4.2). In the inset (II) a schematic NFT_(s+r) diagram describing the inelastic excitation (see Fig. 1.1.1 (inset (I))) of the low-lying octupole vibration (wavy line) of ^{208}Pb by the action of the transient field created by the α -particle on the target (horizontal dashed line) is given. Outgoing α particles are deflected in a magnetic spectograph and recorded in a detector. The corresponding excitation function is given in the lowest part of the figure.

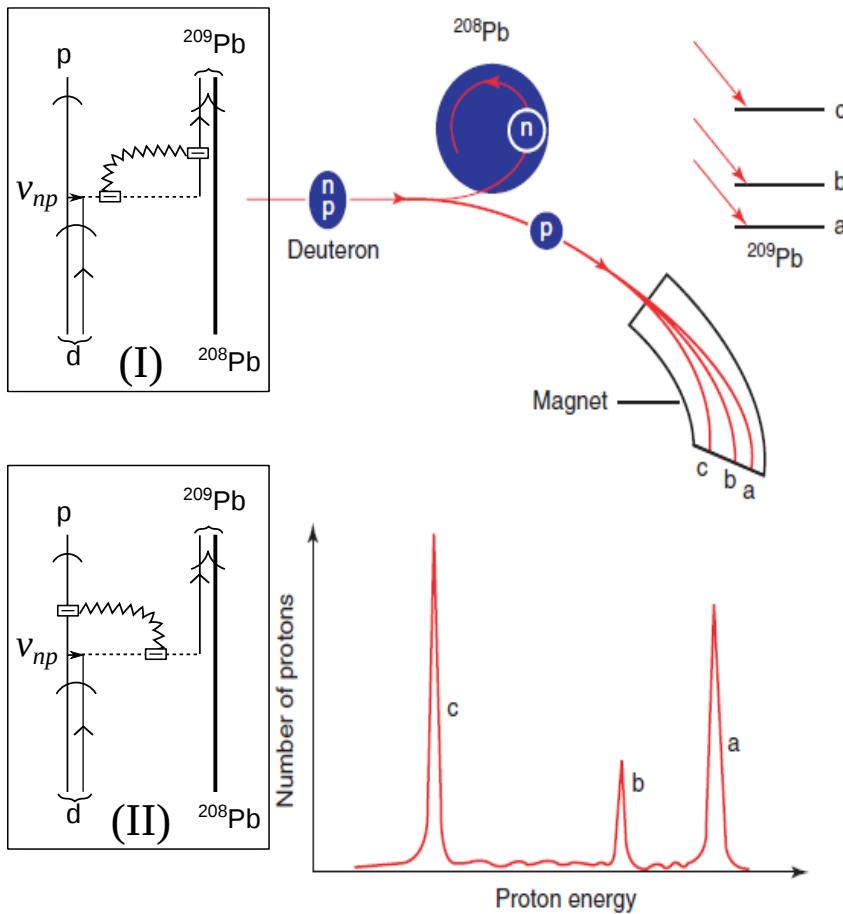


Figure 2.1.2: Schematic representation of the one-nucleon transfer reaction $^{208}\text{Pb}(d, p)^{209}\text{Pb}$ populating valence single-particle states of ^{209}Pb . In the inset a schematic NFT_(s+r) diagram describing the process is shown. Curved arrowed lines describe the projectile d (deuteron) and outgoing particle p (proton) moving in the continuum. The short horizontal arrowed line labeled v_{np} represents the proton–neutron interaction inducing the transfer process (dashed horizontal line) while the open dashed rectangle indicates the Particle Recoil Coupling (PRC) vertex. That is, the coupling of the relative motion to the recoil process described in terms of a jagged line (App. 2.B). This information is carried out in the center of mass system by the outgoing particles in the final channel. Within this context the jagged line is involved in a virtual process (insets (I) and (II)). The energy and momentum of the outgoing proton reflects the recoil, the Q -value of the reaction and the angular momentum and excitation energy of the final state as analyzed in the magnet and recorded in the particle detector (a,b,c).

to the area of the spherical surface A it spans ($\Omega = A/R^2$ sr), in a similar way in which in two

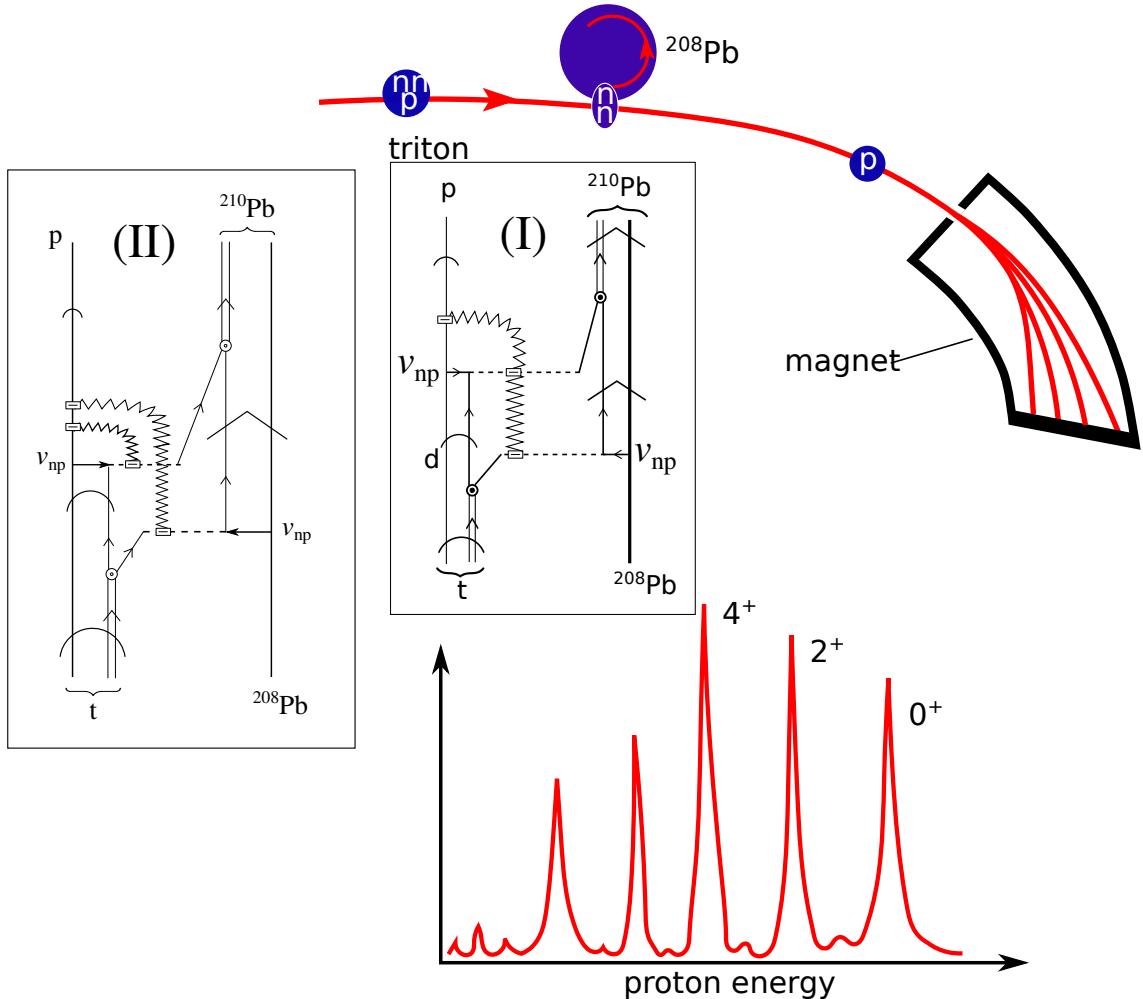


Figure 2.1.3: Schematic representation of the two-nucleon transfer reaction $^{208}\text{Pb}(t, p)^{210}\text{Pb}$ populating the ground state 0^+ , and lowest lying excited states 2^+ and 4^+ . That is, monopole, quadrupole and hexadecapole pair addition modes (multipole pairing vibrations) of ^{208}Pb (App. 7.D; see also Brink and Broglia (2005) Sect. 5.3.1 p. 108). In the inset (I) a $\text{NFT}_{(\text{s+r})}$ diagram associated with the (successive) transfer process is displayed. The jagged line brings information to the outgoing particle in the final channel, of the change in scaling in the asymptotic outgoing waves with respect to the incoming ones, concerning the different mass partitions (recoil) of summed value $2m$ (App. 2.B; concerning the apparent non-linearity that is the direct coupling of two recoil modes, this can be avoided drawing the process as shown in inset (II); see also App. 2.A).

(mb/sr). To translate these quantities into nuclear structure information, a model of structure and of reactions is needed to calculate the absolute cross sections, to be compared with the data. The risks of using relative cross section is that of overlooking limitations in the description of the reaction mechanism or in that of the structure description of the states involved in the reaction under study. Or in both.

In this connection, it is of notice that either one sets equal weight in correctly calculating the static and dynamic properties of the single-particles and of the collective modes respectively, and on their interweaving leading to dressed (renormalized), physical modes, than in working out the reaction mechanism, or the confrontation between theoretical predictions and experimental observation may not be fruitful⁷.

Echoing Heisenberg's requirement ⁸ that no concept enters the quantal description of a physical system which cannot, in principle, be observable, and Landau's findings that weakly excited states of a quantal many-body system may be regarded as a gas of weakly interacting elementary modes of excitation⁹, Bohr, Mottelson¹⁰ and coworkers developed a unified description of the nuclear structure. In particular, a Nuclear Field Theory (NFT)¹¹ in terms of Feynman diagrams describing the behavior of quasiparticles, vibrations and rotations, and of their couplings both in 3D¹² – as well as in gauge¹³ – and other “abstract” spaces, which had close connections with direct nuclear reactions¹⁴. Within this context one gives in Figs. 2.1.1–2.1.3 (insets) schematic representations of *unified NFT diagrams of structure and reactions (NFT_(s+r))*¹⁵, which microscopically describe the variety of structure and reaction processes in terms of Feynman diagrams in a basis of elementary modes of excitation. That is, in the present case in which the target is a closed shell sys-

dimensions, an angle θ is related to the length L of the circular arc it spans. $\theta = L/R$ rad.

⁷*Structure and Reactions*. Within this context one can ask how does one understands which the correct elements are to describe a reaction process, if one does not know in detail the structure of the initial and final states? In a nutshell: how can one understand reaction without knowing structure (eyes without object)? Vice versa, how can one understand what the elements needed for a correct description of the structure of nuclear states is, if one does not know how to observe them (specific probe), how to bring that information to the detector?. In other words, how can one understand structure without knowing reaction (object without eyes)? The answer to both questions is, likely, that one cannot.

⁸Heisenberg (1949).

⁹Landau (1941).

¹⁰Bohr (1964), Bohr and Mottelson (1969), Bohr (1976), Mottelson (1976), Bohr and Mottelson (1975), Bohr et al. (1958) and references therein.

¹¹Bès et al. (1974); Broglia et al. (1976); Bohr and Mottelson (1975); Bès and Kurchan (1990); Mottelson (1976) and refs. therein. See also Broglia (2019).

¹²Nilsson (1955); Bohr and Mottelson (1975) and refs. therein.

¹³Bohr et al. (1958); Belyaev (1959); Högaasen-Feldman (1961); Bès and Broglia (1966); Bjerregaard et al. (1966); Broglia and Riedel (1967); Bohr and Mottelson (1975).

¹⁴Alder et al. (1956), Alder and Winther (1975), Broglia and Winther (2004); concerning the general development of direct nuclear reactions see Austern (1970), Jackson (1970), Satchler (1980), Satchler (1983), Brink (1985), Glendenning (2004); Thompson and Nunes (2009), and refs. therein.

¹⁵Broglia (1975); Broglia and Winther (2004); Potel et al. (2013a); Broglia et al. (2016).

tem, a particle-hole (inelastic scattering), one-particle (single-particle stripping) and two-particle (Cooper pair transfer) modes.

In inset (I) of Fig. 2.1.1 a diagram describing elastic scattering is shown, while in inset (II) a NFT_(s+r) diagram describing the inelastic excitation of the low-lying octupole vibration of ^{208}Pb is displayed. A curved arrow on a line indicates propagation of the projectile in the continuum (asymptotic waves). The horizontal dashed line represents the action of the mean field or of the bare interaction (see Fig. 1.2.2).

In the insets of Fig. 2.1.2 a NFT_(s+r) diagrams describing the process $^{208}\text{Pb}(d, p)^{209}\text{Pb}$ are schematically shown. A standard pointed arrowed line indicates the neutron moving with the proton in the deuteron ((curved) arrowed double line), or around the (assumed, for simplicity, inert) ^{208}Pb core (bold face line). The jagged curve represents the recoil mode coupling the intrinsic and the relative motion, thus accounting for the different mass partition and the associated change in scaling between entrance and exit channel distorted waves. The corresponding momentum mismatch being taken care of by a Galilean transformation (recoil effect). The jagged curve can transfer the information of momentum mismatch to either the residual nucleus (inset (I)) or to the outgoing particle, i.e. the proton (inset (II)). Horizontal short arrowed lines stand for the proton-neutron (nucleon-nucleon) bare interaction inducing transfer. It is of notice that choosing in an appropriate way the (post or prior) representation to describe the reaction process (energy conservation), one can evidence the single-particle mean field or the proton-neutron interaction as inducing the transfer process.

In Fig. 2.1.3 a cartoon representation of the $^{208}\text{Pb}(t, p)^{210}\text{Pb}$ process is given, while in the inset the corresponding NFT diagram is displayed. The dineutron moving in the triton and around the ^{208}Pb core, (pair addition mode) is represented by a double arrowed line. Each individual transferred neutron is indicated with a single arrowed line. The curved arrows on the triton and on the proton indicate motion in the continuum with incoming and outgoing asymptotic waves, respectively. The pointed arrow encompassing the pair addition mode and the core ^{208}Pb , indicate intrinsic (structure) motion. In selecting this NFT diagram the assumption was made, following the results of detailed calculations, that the main contribution to the process arises from the successive transfer of the nucleons. Two jagged curves are shown. One connecting the first and the second transfer process at the particle-recoil coupling vertex (PRCV, dashed open square), in keeping with the fact that the channel $^{209}\text{Pb}+d$ has no asymptotic waves. The second jagged curve emerges from the second PRCV (inset (I)) or from a new one (inset (II)), and carries the mismatch information to the outgoing proton (see App. 2.B, in particular the paragraph after Eq. (2.B.36))¹⁶.

It is of notice that the successive transfer of neutrons described in insets (I) and (II) can be related to the successive tunneling of pairs of electrons involved

¹⁶Concerning the apparent non linearity of the recoil process displayed in inset (I) (two jagged lines associated with a single PRCV), we refer to Fig. 2.A.1 and associated discussion.

in the Josephson effect¹⁷, see Sect. 4.6, in particular the discussion following Eq. (4.6.1)).

As discussed in the following chapters, Cooper pairs are extended objects, the fermionic partners being correlated over distances much larger than nuclear dimensions (correlation length $\xi \approx 14$ fm $\gg R_0 \approx 6$ fm ($A = 120$)). Because the single particle potential acts on these pairs as a rather strong external field, this correlation length feature is not obvious in structure calculations, becoming apparent by connecting structure with Cooper pair transfer processes.

Elementary modes of excitation, that is single-particle and collective motion, are the way nuclei react to external probes, and thus closely connected with the physical observables. Namely, absolute transition probabilities and absolute differential cross sections. Within this context, bare elementary modes of excitation already contain an important fraction of nuclear many-body correlations, thus making the diagonalization of the nuclear Hamiltonian leading to dressed elementary modes of excitation, a low-dimension problem. For example, in terms of $10^2 \times 10^2$ matrices, each matrix element containing already much physical insight into nuclear structure at large¹⁸.

In keeping with the fact that all the nuclear degrees of freedom are exhausted by those of the nucleons, and that the different reactions, that is elastic, Coulomb and inelastic excitations, as well as one- and two-particle transfer reactions project particular, but somewhat overlapping components of the total wavefunction, the nuclear elementary modes of excitation give rise to an overcomplete, non orthogonal, Pauli principle violating basis, both concerning structure as well as reactions. Nuclear field theory¹⁹ provides the conserving sum rules protocol to diagonalize in this basis the three- and four-point vertex of the nuclear Hamiltonian (Eq. (1.2.10)) to any order of perturbation theory, also infinite if so required for specific processes (see Sect. 2.5). The dressed physical elementary modes resulting from the interweaving of the bare modes are orthogonal to each other and fulfill Pauli principle, providing a microscopic solution to the many-body nuclear problem.

NFT_(s+r) diagrams²⁰ (see e.g. Figs. 2.7.2, 2.7.3 (a)+(b) and 2.7.5) embody predictions in terms of absolute differential cross sections and transition probabilities. Quantities which can be directly compared with the observables obtained by studying the nuclear system with the variety of ever more precise arsenal of experimental probes (see e.g. Fig. 2.8.1).

At this point a proviso or two are in place. The original elementary modes of

¹⁷Josephson (1962); Anderson (1964b). See also Brink and Broglia (2005) App. L and refs. therein.

¹⁸Illuminating discussion with B. A. Brown during the 15th International Conference on Nuclear Reactions Mechanisms (Varenna, June 2018) regarding this, and related issues are gratefully acknowledged (RAB).

¹⁹Bès et al. (1974), Bès et al. (1976b), Bès et al. (1976c), Bès et al. (1976a), Bès and Broglia (1975), Broglia et al. (1976), Bès et al. (1975a), Mottelson (1976), Bès and Broglia (1977), Bortignon et al. (1977), Bortignon et al. (1978), Broglia and Winther (2004), Reinhardt (1975), Reinhardt (1978a), Reinhardt (1978b), Reinhardt (1980); see also Broglia (2019).

²⁰Broglia (1975); Broglia and Winther (2004); Broglia et al. (2016).

nuclear excitation melt together, due to their interweaving, into effective fields²¹. Each of them display properties which reflect that of the others with which it couples, their individuality resulting from the actual relative importance of each one of them. What one calls a physically (dressed) particle is only partially to be associated with that particle field alone. It is also partially to be associated with the vibrational fields (surface, density, spin²², pairing²³, etc. vibrational modes), because they are in interaction through the bare NN -interaction as well as the particle-vibration coupling vertices²⁴. And conversely, what one calls a nuclear vibration can couple to particle-hole (in the case of a surface vibration), two-particle (in the case of a pair addition) or a two-hole (in the case of a pair removal) configurations, and materialize in specific fermionic states.

Thus nucleons (fermions) couple to vibrational modes (quasi-bosons) and, eventually, can reabsorb them returning to the original state. The outcome of such processes, namely the dressed physical elementary modes of excitation, is closely connected with the renormalization program of quantum electrodynamics (QED)²⁵ implemented in NFT in terms of Feynman diagrams. Renormalized NFT, i.e (NFT)_{ren} implies that the intermediate, virtual states clothing the elementary modes of excitation, should be fully dressed²⁶. They can thus, in principle, be taken from experiment.

The specific experimental probes of the bare elementary modes of nuclear excitation reveal only one aspect, in most cases likely the most important one of the physical (dressed) elementary modes. Renormalized NFT reflects the physical unity of low-energy nuclear research requiring the melting not only of elementary modes of excitation but also of structure and reaction theory, let alone of the different experimental techniques developed to study the atomic nucleus. In other words the need for a “complete” set of experimental probes to reveal the multi faceted properties of dressed elementary modes of excitation.

As seen from the contents of the present monograph, the accent is set at relating theoretical predictions with experimental findings, through the unification of structure and reactions. In particular the unification of pairing and two-nucleon transfer processes, where the two subjects are blended together, which is what happens in nature. Once the NFT rules to work out the variety of elements (spectroscopic and, with the help of them, reaction amplitudes) have been laid out and/or the pertinent literature refer to, concrete embodiments are provided and eventual absolute cross sections and transition probabilities calculated and confronted with the experimental data.

As already stated in the previous chapter, an essential test theory has to pass

²¹Within this context see Dickhoff and Van Neck (2005) and references therein.

²²Bertsch and Broglia (2005) Chs. 6,7,8.

²³Brink and Broglia (2005) Ch. 5.

²⁴Bohr and Mottelson (1975).

²⁵Feynman (1975); Schwinger (2001).

²⁶Barranco et al. (2017), see also Broglia et al. (2016), see also Chs. 6 and 7.

is connected with some operator identities which are known as sum rules²⁷. An overcomplete (non-orthogonal) set of basis states, like that provided by elementary modes of excitation, will violate sum rules. For example, a hole and a pair addition is non orthogonal to a particle and a surface ($p - h$) vibration (see Sect. 2.5.4). As we shall see, taking into account the couplings between these modes following the rules of NFT, such violations are eliminated and sum rules fulfilled (Sect. 2.5).

A similar situation is found in the case of e.g. two-nucleon transfer, because the single-particle states in projectile and target are, as a rule, non-orthogonal. Taking into account the corresponding overlaps one can remove the overcounting (see Sects. 4.2, 6.2.9 and 6.5).

2.2 Sum rules revisited

A quantitative measure of the overcompleteness of the elementary modes of excitation basis is provided by the use of exact and of empirical²⁸ sum rules that the observables (cross sections) associated with the variety of probes to which the nucleus is subject, have to fulfill. Examples of the first type are given by the Thomas-Reiche-Kuhn (TRK) sum rule^{29,30}, and by a sum rule which relates one-with two-particle transfer processes³¹. Within the framework of simple, but still physically relevant models, connected with inelastic scattering processes, we refer to Sect. 2.5.2 in general and Eqs. (2.5.40)–(2.5.44) in particular. With regards to one-particle transfer sum rules we refer to Sect. 2.5.3 and Eqs. (2.5.75) and (2.5.78).

Concerning examples of the second type (empirical, physical sum rules), we refer to sum rules associated with two-nucleon transfer reactions (TNTR)³² discussed in Sect. 2.2.1 and³³ in Sect. 7.2.2.

Physically, they provide information concerning: 1) the maximum amount of energy which the quantal system can absorb from a beam of photons (γ -rays) shined on it; 2) the total two-nucleon transfer cross section (ring area fraction of the geometrical reaction cross section) exhausted by the final ($A \pm 2$) states populated in the transfer process.

²⁷See e.g. Bohr and Mottelson (1975); Ring and Schuck (1980), Bertsch and Broglia (2005) Ch. 4, Bortignon et al. (1998), Chs. 3 and 8 and refs. therein.

²⁸In this connection see paragraph after Eq. (2.B.44).

²⁹Bohr and Mottelson (1975), Sect. 6-4.

³⁰Bertsch and Broglia (2005), Chapter 3, in particular Sect. 3.3.

³¹Bayman and Clement (1972); Lanford (1977).

³²Broglia et al. (1972).

³³Within this context, the absolute two-nucleon transfer cross section populating the ground state of a superfluid nucleus is proportional to the number of Cooper pairs contributing to the nuclear condensate (modulus squared). This quantity is rather stable along a pairing rotational band (see however Fig. 3.1.3 and associated discussion in Sect. 3.1), in keeping with the fact that the “intrinsic” $|BCS\rangle$ -state of the deformed system in gauge space, is a coherent state and essentially the same for all members of the band. This fact is at the basis of a newly found physical sum rule (Potel et al. (2017)).

In other words, these sum rules provide a quantitative measure of the single-particle subspace the quantal system under study, in particular the nucleus, uses to correlate particle-hole excitations and thus induce the antenna-like motion of protons against neutrons or, to correlate pairs of nucleons moving in time reversal states around the Fermi energy, leading to a, static or dynamic, sigmoidal distribution of the associated pair ($\gamma, \bar{\nu}$) level occupancy around the Fermi energy.

As shown in Sect.. 1.7, the TRK sum rule can, in the nuclear case, be written as

$$S(E1) = \sum_{\alpha} |\langle \alpha | F | \tilde{0} \rangle|^2 (E_{\alpha} - E_0) = \frac{9}{4\pi} \frac{\hbar^2 e^2}{2m} \frac{NZ}{A}, \quad (2.2.1)$$

where $|\alpha\rangle$ labels a complete set of excited states populated by acting with the dipole operator F (Eq. (1.7.5)) on the correlated vacuum state $|\tilde{0}\rangle$. Each (quasi) bosonic elementary mode of excitation provides, within the harmonic approximation, a specific contribution to the total zero point fluctuations of the ground state, that is,

$$\langle \tilde{0} | F^2 | \tilde{0} \rangle = \sum_{\alpha} \frac{\hbar\omega}{2C_{\alpha}} = \sum_{\alpha} \frac{\hbar^2}{2D_{\alpha}} \frac{1}{\hbar\omega_{\alpha}}. \quad (2.2.2)$$

As mentioned in the previous chapter, ZPF perturb the static nucleon Fermi sea, that is the set of occupied levels of the mean field potential (see Fig. 2.2.1), inducing virtual particle-hole excitations (k, i , i.e. $\epsilon_i \leq \epsilon_F$ and $\epsilon_k > \epsilon_F$). Because $|\langle \alpha | F | \tilde{0} \rangle|^2$ measures the probability with which the state $|\alpha\rangle$ is populated, the α -sum in (2.2.1) gives a measure of the maximum energy that the nucleus can absorb from the γ -beam. It is customary to measure $|\langle \alpha | F | \tilde{0} \rangle|^2$ in single-particle (sp) units (Weisskopf (W) units)

$$\begin{aligned} B_{sp}(E1; j_1 \rightarrow j_2) &= \frac{3}{4\pi} e_{E1}^2 \langle j_1 \frac{1}{2} 1 0 | j_2 \frac{1}{2} \rangle^2 \times \langle j_2 | r | j_1 \rangle^2, \\ &\approx \frac{0.81}{4\pi} A^{2/3} e_{E1}^2 \text{ fm}^2 = B_W(E1), \end{aligned} \quad (2.2.3)$$

where $(e)_{E1} = (N/A)e$ for protons and $(e)_{E1} = -(Z/A)e$ for neutrons, in keeping with the fact that the motion of a nucleon is associated with a recoil of the rest of the nucleus, and that the center of mass of the system must remain, in an intrinsic excitation³⁴, at rest. It is of notice that,

$$\begin{aligned} S(E1) &= \sum_n |\langle n | F | \tilde{0} \rangle|^2 (E_n - E_0) \\ &= \sum_{k,i} |\langle k, i | F | \text{gs(MF)} \rangle|^2 (\epsilon_k - \epsilon_i), \end{aligned} \quad (2.2.4)$$

³⁴Bohr and Mottelson (1969).

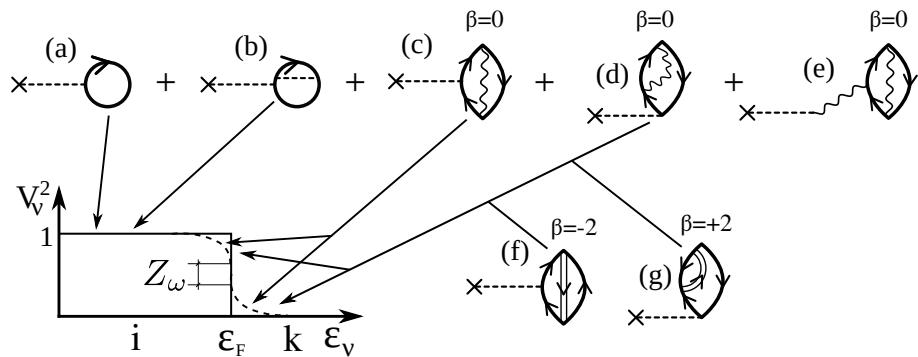


Figure 2.2.1: Schematic representation of the Fermi distribution. The sharp, (continuous line) step function schematically represents the Hartree-Fock occupation numbers. The associated nuclear density is measured with the help of an external field (cross attached to a dashed line) through processes of type (a) (Hartree: H) and (b) (Fock: F) is expected to display a diffusivity of the order of the strong force range. Zero Point Fluctuations (ZPF) associated with collective particle-hole vibrations, i.e. processes with transfer quantum number $\beta = 0$ and shown in (c), (d) and (e), and with pairing vibrations, i.e. pair removal (graph (f)) and pair addition modes (graph (g)), smooth out the occupation numbers around the Fermi energy (dashed curve; the discontinuity gives a measure of the single-particle content Z_ω) and lead to a nuclear density of larger (dynamical) diffusivity and radius than that associated with HF (see Sect. 1.8). One- and two-particle strengths which in this (mean field) approximation are found in a single A -mass system are, as a result of ZPF ($\beta = 0, \pm 1, \pm 2$) distributed over a number of nuclei ($A, A \pm 1, A \pm 2$) (see also Sect. 5.8, Fig. 5.8.1).

provided $|\tilde{0}\rangle$ contains the ground state correlations mentioned in connection with Eq. (2.2.2), and that $|\text{gs(MF)}\rangle$ those associated with HF (see Fig. 1.2.2). In other words, provided

$$|\text{HF}\rangle = |\text{gs(MF)}\rangle = \prod_{i \in \text{occup}} a_i^\dagger |0\rangle \quad (2.2.5)$$

where $|0\rangle$ is the particle vacuum state ($a_j|0\rangle = 0$), and $\Gamma_\alpha|\tilde{0}\rangle = 0$, Γ_α^\dagger being the creation operator of a dipole (RPA) correlated particle-hole like mode ($\Gamma_\alpha^\dagger = \sum_{ki} X_{ki}^\alpha a_k^\dagger a_i + Y_{ki}^\alpha (a_k^\dagger a_i)^\dagger$)³⁵.

Relation (2.2.4) is a consequence of the fact that $S(E1)$ is proportional to the average value of the double commutator $[[H, F], F]$ in the ground state of the system ($|\tilde{0}\rangle$ or $|\text{HF}\rangle$). Because F is a function of only the nucleon coordinates and, assuming $v(|\mathbf{r} - \mathbf{r}'|)$ to be velocity independent, the only contribution to the double commutator arises from the (universal) kinetic energy term of the Hamiltonian. Thus, the value (2.2.1) is model independent. In other words, this value does not depend on the correlations acting among the nucleons, but on the number of them participating in the motion and on their mass (inertia) as testified by the fact that $\sum_\alpha \hbar\omega_\alpha \left(\frac{\hbar\omega_\alpha}{2C_\alpha}\right) = \sum_\alpha \left(\frac{\hbar^2}{2D_\alpha}\right)$. *It is then not surprising that the TRK sum rule was used in the early stages of quantum mechanics, to determine the number of electrons in atoms.*

2.2.1 Empirical two-nucleon transfer sum rules

Let us now consider two-nucleon transfer processes. The absolute cross sections associated with the population of final states can be set essentially on equal footing with respect to each other concerning Q -value and recoil effects, with the help of empirically determined global functions³⁶ (see also Eqs. (2.B.38)–(2.B.40) and subsequent discussion). In this way, the theoretical absolute cross section associated with e.g. the $A(t, p)A + 2$ population (we assume N to be even) of the n th final state of spin J can be written as

$$\sigma^{(n)}(J, Q_0) = \left| \sum_{j_1 \geq j_2} B^{(n)}(j_1 j_2; J) S(j_1 j_2; J, Q_0) \right|^2, \quad (2.2.6)$$

where

$$S^2(j_1 j_2; J, Q_0) = \sigma(j_1, j_2; J, Q_0), \quad (2.2.7)$$

³⁵See footnote 12 Ch. 1. See also Bertsch and Broglia (2005), Ch. 4, and Brink and Broglia (2005) Ch. 8, Sect. 8.3, Bohr and Mottelson (1975) Sect. 6-5 h. As can be seen from Fig 1.2.3 (inset), in the RPA approximation no scattering vertices (Fig. 1.2.1) are present. Consequently, the coupling between one- and two- phonon states is, within this (harmonic) approximation, not possible.

³⁶see Broglia et al. (1972).

the quantity $\sigma(j_1, j_2; J, Q_0)$ being the absolute two-nucleon transfer cross section associated with the pure two-particle configuration $(j_1, j_2)_J$, while

$$B(j_1 j_2; J) = \left\langle \Phi_{J_f}(\xi_{A+2}) \left| \left[\Phi_{J_i}(\xi_A) \frac{[a_{j_1}^\dagger a_{j_2}^\dagger]_J}{[1 + \delta(j_1, j_2)]^{1/2}} \right]_{J_f} \right. \right\rangle, \quad (2.2.8)$$

is the two-nucleon spectroscopic amplitude, $\Phi_{J_i}(\xi_A)$ being the wavefunction describing the ground state of the initial nucleus, $\Phi_{J_f}(\xi_{A+2})$ that of the final state, ξ labeling the relative radial and spin intrinsic coordinates. Assuming A to be a closed shell system, and $J = 0$, one can write

$$|0_n^+\rangle = \sum_j c^{(n)}(j, j; J = 0) |j, j; J = 0\rangle, \quad (2.2.9)$$

where $n = 1, 2, 3, \dots$ labels the final nucleus states of spin and parity $J^\pi = 0^+$ in increasing energy order, and $B^{(n)} = c^{(n)}$. Making use of the completeness relation of the coefficients $c^{(n)}$ one obtains,

$$\sum_n \sigma^{(n)}(J = 0, Q_0) = \sum_j \sigma(j, j; J = 0, Q_0). \quad (2.2.10)$$

The above equation parallels (2.2.4), aside from the fact that the Q -value effect is, in connection with (2.2.4), analytically dealt with, while $\sigma(\{Q\})$ is a functional of Q . The complete separation of the relative and intrinsic motion coordinates taking place in e.g. (2.2.4) is in keeping with the fact that in elastic and inelastic processes the mass partition is equal in both entrance and exit channels. Thus, the intrinsic (structure) and the relative motion (reaction) coordinates can be treated separately. This is not the case for transfer processes, both intrinsic and reaction coordinates being interwoven through the recoil process (particle-recoil mode coupling (dashed open square), jagged curves (recoil mode), see e.g. Fig. 2.1.3 as well as Figs. 2.7.2 and 2.7.3; see also Sect. 2.B.1 of App. 2.B).

A parallel with the discussion carried out in connection with (2.2.3) regarding the TRK sum rule, can be drawn defining two-particle transfer units as,

$$\sigma_{2pu}^{max}(J, Q_0) = \max [\sigma(j_1, j_2; J, Q_0)], \quad (2.2.11)$$

where $\max[\]$ indicates that the largest two-particle absolute cross section in the single-particle subspace (*hot orbital*), is to be considered. In this way one can write the relation (2.2.10) in dimensionless units. Another quite useful, this time exact, two-particle transfer sum rule has been introduced in the literature³⁷, which relates two-nucleon stripping and pick-up reactions cross sections, with single-particle transfer processes³⁸.

³⁷Bayman and Clement (1972). In this connection, and within the context of a schematic model, see Eq. (2.5.75) and subsequent discussion.

³⁸Lanford (1977)

The above arguments carried out for nuclei around closed shells can, equally well, be applied to the case of open shell nuclei, making use of the corresponding two-nucleon spectroscopic amplitudes³⁹. In particular, in the case of phase-correlated, independent pair motion, i.e. the BCS mean field solution of the pair problem (pairing Hamiltonian). The summed spectroscopic pair transfer amplitudes, each term weighted with $(j + 1/2)^{1/2}$, is given by

$$\begin{aligned}\alpha'_0 &= \langle BCS(N+2) | P'^\dagger | BCS(N) \rangle \\ &= \sum_j \frac{2j+1}{2} U'_j(N) V'_j(N+2),\end{aligned}\quad (2.2.12)$$

where⁴⁰

$$P'^\dagger = \sum_{m>0} a'_{jm}^\dagger a'_{jm}^\dagger = \sum_j \sqrt{\frac{2j+1}{2}} T'^\dagger(j^2(0)), \quad (2.2.13)$$

creates two nucleons in time reversal states,

$$T'^\dagger(j^2(0)) = \frac{[a'_j^\dagger a'_j^\dagger]^0}{\sqrt{2}}, \quad (2.2.14)$$

being the two-nucleon transfer operator. The associated expectation value in the $|BCS\rangle$ ground state is the two-nucleon transfer spectroscopic amplitude

$$\begin{aligned}B(j^2(0), N \rightarrow N+2) &= \langle BCS(N+2) | T'^\dagger(j^2(0)) | BCS(N) \rangle \\ &= \sqrt{\frac{2j+1}{2}} U'_j(N) V'_j(N+2),\end{aligned}\quad (2.2.15)$$

implying that to create a pair of particles in a quantal state j , it has to be empty in the initial system (i.e. N , probability $U'^2_j(N)$) and has to result occupied in the final nucleus (i.e. $N+2$, probability $V'^2_j(N+2)$).

The ZPF associated with pairing vibrations smooth out the sharp HF Fermi surface (Fig. 2.2.1). The number of pairs in each level participating in this smoothing is $(2j+1)/2$, their occupancy being measured by the simultaneous, and apparently contradictory, property of being a particle (amplitude V'_j) and a hole (amplitude U'_j). In other words, α'_0 measures the number of pairs of nucleons (Cooper pairs) participating in the smoothing out of the Fermi surface⁴¹, being the spectroscopic amplitude associated with the population of pairing rotational bands in two-nucleon transfer processes. It is expected that α'_0 depends weakly on N and is about

³⁹See Yoshida (1962); see also App. 2 of Broglia et al. (1973).

⁴⁰The primed quantities are the particle creation operators and BCS occupation amplitudes referred to the intrinsic system of reference in gauge space (see Fig. 1.4.1 and Sect. 4.7). See also Potel et al. (2013b).

⁴¹Schrieffer (1964); Potel et al. (2017) and refs. therein.

conserved along a pairing rotational band. Because $d\sigma(gs(N) \rightarrow gs(N+2))/d\Omega \approx |\alpha'_0|^2$, conservation is also expected for these absolute cross section. But in this case, it is a conservation of a physical character, and not a mathematical one (see Sect. 7.2.1).

If one finds that at the angle where $L = 0$ two-nucleon transfer differential cross sections have the first maximum, as a rule close to 0° in (t, p) or (p, t) reactions, the two nucleon strength function is dominated by a single peak, that associated with the ground state, and this is so for a number of isotopes differing in mass number by two, it can be concluded that one is in presence of a pairing rotational band (see Figs. 3.1.3 and 3.1.4). This is one of the reasons why (2.2.12) can be viewed as the order parameter of the nuclear superfluid phase and, in keeping with (2.2.13) and (2.2.14), two-nucleon transfer reaction regarded as the specific tool to probe pairing in nuclei.

Because in a finite many-body system like the nucleus, quantal fluctuations in general and those of particle number in particular (pairing vibrations), are much larger than in bulk systems, a dynamic parallel to (the static deformation value) α_0 , namely α_{dyn} can be defined at profit (see Fig. 3.5.7)⁴².

2.3 Fluctuations and damping

In this section, we discuss some concepts and applications of the particle-vibration coupling mechanism⁴³ found at the basis of the development of NFT. The Hamiltonian describing a system of independent particles and of collective surface vibrations can be written as⁴⁴

$$H = H_{MF} + H_{coll} + H_c, \quad (2.3.1)$$

where

$$H_{MF} = T + U \quad (2.3.2)$$

is the mean field Hamiltonian, sum of the single-particle kinetic energy and of the self-consistent potential $U = f(\rho)$, functional of the density. That is,

$$U = U_H + U_x, \quad (2.3.3)$$

where the Hartree potential U_H was defined in Eq. (1.2.8) and,

$$U_x = - \sum_{i(\epsilon_i \leq \epsilon_F)} \varphi_i^*(\mathbf{r}') v(|\mathbf{r} - \mathbf{r}'|) \varphi_i^*(\mathbf{r}), \quad (2.3.4)$$

is the exchange (Fock) potential.

⁴²See also Potel et al. (2017).

⁴³See footnote 14 Ch. 1.

⁴⁴That is, essentially (1.2.10) without v , four-point vertex shown in Sect. 2.5 to be necessary for working out the (sum rule) conserving nuclear field theory rules. Within this framework, one can consider the present Section as a warm up to the introduction of NFT, aiming only at providing evidence of the importance of particle vibration coupling PVC renormalization processes.

To second order in the particle–vibration coupling strength one finds⁴⁵ that the nucleons move in an ω -dependent, non-local dielectric medium, that is

$$\begin{aligned} \left(-\frac{\hbar^2}{2m} \nabla_r^2 + U_H(r) \right) \varphi_j(\mathbf{r}) &+ \int d\mathbf{r}' U_x(\mathbf{r}, \mathbf{r}') \varphi_j(\mathbf{r}'), \\ &+ (\Delta E_j + iW_j) \varphi_j(\mathbf{r}) \\ &\approx \left(-\frac{\hbar^2}{2m_k} \nabla_r^2 + U''_H(r) + \Delta E''_j + iW''_j \right) \varphi_j(\mathbf{r}), \\ &= \varepsilon_j \varphi_j(\mathbf{r}), \quad \left(U''_H = \frac{m}{m_k} U \text{ and similarly for } \Delta E'' \text{ and } W'' \right), \end{aligned} \quad (2.3.5)$$

where $m_k = \left(1 + \frac{m}{\hbar^2 k} \partial U_x / \partial k \right)^{-1} \approx 0.7m$ is the k -mass⁴⁶, while

$$\Delta E_j(\omega) = \lim_{\Delta \rightarrow 0} \sum_{\alpha'} \frac{V_{v,\alpha'}^2(\omega - E_{\alpha'})}{(\omega - E_{\alpha'})^2 + (\frac{\Delta}{2})^2}, \quad (2.3.6)$$

and

$$W_j(\omega) = \lim_{\Delta \rightarrow 0} \sum_{\alpha'} \frac{V_{v,\alpha'}^2}{(\omega - E_{\alpha'})^2 + (\frac{\Delta}{2})^2}, \quad (2.3.7)$$

are the real and imaginary contributions to the self-energy calculated in second order of perturbation theory, ω standing for a generic value of the single-particle energy⁴⁷ (see Fig. 2.3.1; note that in this case $E_{\alpha'} = \epsilon_{\nu_2} + \hbar\omega_{\alpha}$).

⁴⁵see e.g. Brink and Broglia (2005), Mahaux et al. (1985) and references therein. See also Bernard and Giai (1981).

⁴⁶This is in keeping with the fact that the non-local component of the mean field can be parametrized at profit as $0.4E$, where $E = |(\hbar^2 k^2 / 2m) - \epsilon_F|$ (Perey and Buck (1962)) see also Sect. 2.7.1.

⁴⁷Given a Hamiltonian H_c , the contribution to the energy in second order perturbation theory can be written as

$$\Sigma_v(\omega) = \sum_{\alpha'} \frac{V_{v,\alpha'}^2}{\omega - E_{\alpha'}},$$

where $|\alpha'\rangle \equiv |n_\alpha = 1, v'\rangle$ are the intermediate states which can couple to the initial single-particle state v . Note that the expression above is not well defined, in that the energy denominator may vanish. Now, we are not confronted with an accidental degeneracy, but with a case where there can be many intermediate states with $E_{\alpha'} \approx \omega$. In other words, where the particle can decay into more complicated (doorway-) states $|\alpha'\rangle$ (Feshbach (1958)), starting in the single-particle level v of energy ω , without changing its energy (real, on-the-energy shell process). This is a typical dissipative process, and has to be solved by direct diagonalization (see Fig. 2.3.4). Another way around, is to extend the function $\Sigma_v(\omega)$ into the complex plane ($E_{\alpha'} \rightarrow E_{\alpha'} + i\frac{\Delta}{2}$) thus *regularizing the divergence* through a coarse grain approximation, determining the finite contributions and then taking the limit for $\Delta \rightarrow 0$ (Eqs.(2.3.6) and (2.3.7)). The resulting complex potential (*optical potential* from the *complex dielectric function* of optics), parametrizes in simple terms the shift of the centroid of the single-particle state and its finite lifetime.

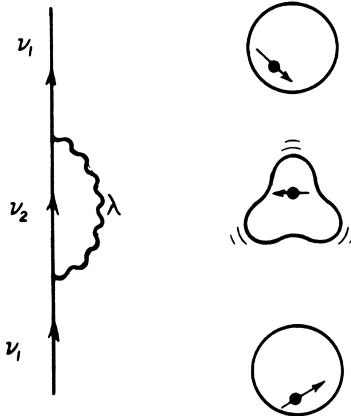


Figure 2.3.1: Self-energy (polarization, PO) graph renormalizing a single-particle. Time ordering leads to the corresponding correlation diagram (CO) (see e.g. Fig. 1.2.4; within this context see also Brink and Broglia (2005) Fig. 9.2 of this reference).

For many purposes ΔE can be treated in terms of an effective mass

$$m_\omega = m(1 + \lambda), \quad (2.3.8)$$

where

$$\lambda = -\left. \frac{\partial \Delta E}{\partial \omega} \right|_{\epsilon_F}, \quad (2.3.9)$$

is the *mass enhancement factor*, while

$$Z_\omega = m/m_\omega,$$

measures the single-particle content (discontinuity) at the Fermi energy.

Consequently, Eq.(2.3.5) can be rewritten as

$$\left(-\frac{\hbar^2}{2m^*} \nabla_r^2 + U'_H + iW'(\omega) \right) \varphi_j(\mathbf{r}) = \varepsilon_j \varphi_j(\mathbf{r}), \quad (2.3.10)$$

with

$$m^* = \frac{m_k m_\omega}{m}. \quad (2.3.11)$$

with $U'_H = (m/m^*)U$ and similarly for W' . Because $\lambda \approx 0.4$ (i.e. the dressed single-particle m_ω is heavier than the bare nucleon, as it has to carry a phonon along or, better, move through a cloud of virtual phonons) and $m_k = 0.7m$, $m^*/m \approx 1$ and $Z_\omega = m/m_\omega \approx 0.7$. Furthermore, due to the fact that $\hbar\omega_\alpha \approx 2 - 2.5 \text{ MeV}$, the range of single-particle energies $E = \varepsilon - \varepsilon_F$ over which the particle-vibration coupling process displayed in Fig.2.3.1 is effective is $\approx \pm 2\hbar\omega_\alpha \approx 4 - 5 \text{ MeV}$ around the Fermi energy (see Figs.2.3.2 and 2.3.3)

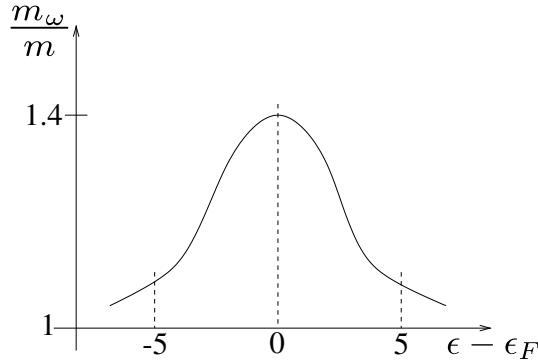


Figure 2.3.2: Schematic representation of the nucleon ω -mass as a function of the single-particle energy.

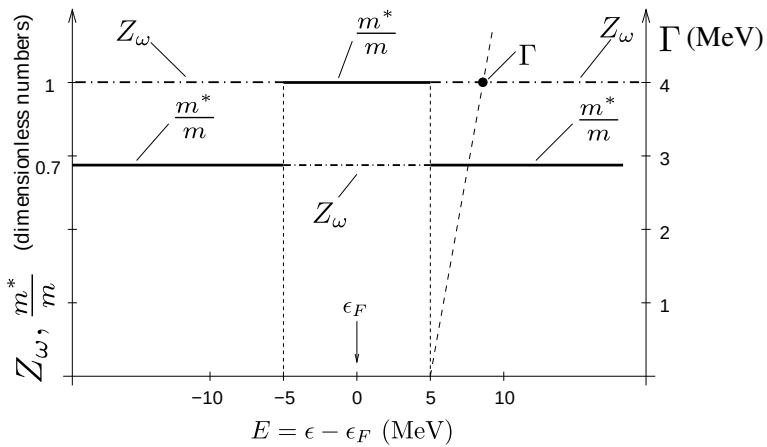


Figure 2.3.3: Schematic representation of the behaviour of m^*/m , $Z_\omega = (m_\omega/m)^{-1}$ and Γ (Fig. 2.3.7) as a function of $E = \epsilon - \epsilon_F$. Note the scales for Z_ω and m^*/m (left) and Γ (right).

It is of notice that ΔE_j indicates the shift of the centroid of the "dressed" single-particle state due to the coupling to the doorway states, while $\Gamma = 2W$ measures the energy range over which the single-particle state spreads due to this coupling (see Fig. 2.3.4). While a large number of states contribute to ΔE ("off-the-energy-shell process", i.e. intermediate, virtual processes in which energy is not conserved), only "on-the-energy-shell processes", that is processes which conserve energy, contribute to Γ . In fact

$$\lim_{\Delta \rightarrow 0} \frac{\Delta}{(\omega - E_{\alpha'})^2 + \left(\frac{\Delta}{2}\right)^2} = 2\pi\delta(\omega - E_{\alpha'}),$$

and

$$\Gamma(\omega) \approx 2\pi\bar{V}^2 n(\omega), \quad (2.3.12)$$

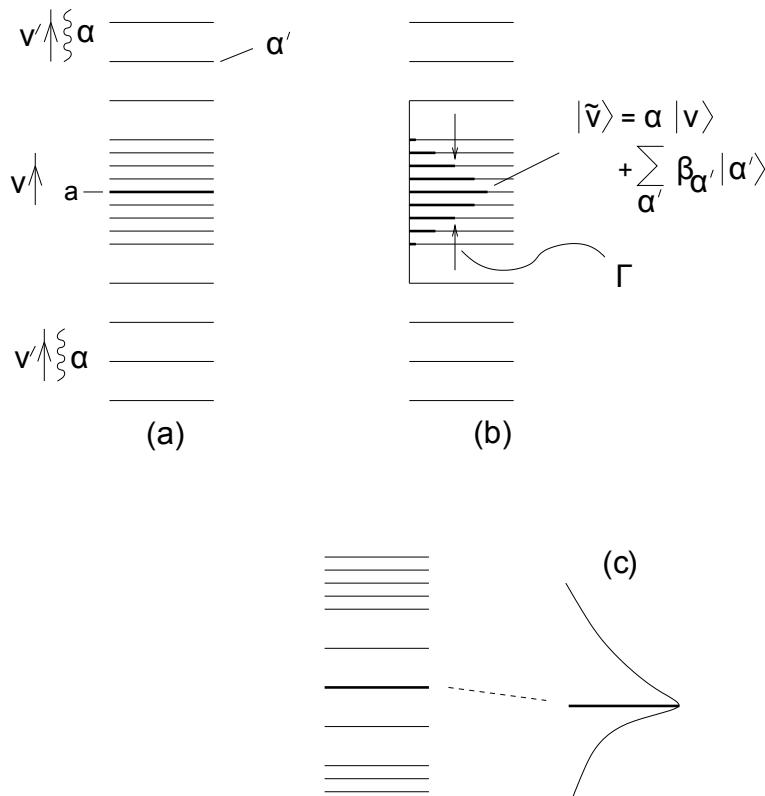


Figure 2.3.4: (a) Schematic representation of the result of the diagonalization of H_c in a basis consisting of the single-particle states $|v\rangle$ and the $|\alpha'\rangle = |\nu', \alpha\rangle$ doorway states. (b) Range of doorway mixing leading to the dressed single-particle state $|\tilde{\nu}\rangle$. In (c) we show a situation where there are more states $|\alpha'\rangle$ above $|a\rangle$ than below.

where \bar{V}^2 is the average value of $V_{\nu,\alpha'}^2$, while

$$n(\omega) = \sum_{\alpha'} \delta(\omega - E_{\alpha'}), \quad (2.3.13)$$

is the density of energy-conserving states α' . Eq.(2.3.12) is known as *Fermi Golden rule*.

Assuming the distribution of single-particle levels is symmetric with respect to the Fermi energy,

$$\Delta E(\omega) = \lim_{\Delta \rightarrow 0} \sum_{\alpha'} \frac{V_{\nu,\alpha'}^2(\omega - E_{\alpha'})}{(\omega - E_{\alpha'})^2 + \left(\frac{\Delta}{2}\right)^2} = 0 \quad (2.3.14)$$

as there are equally many states pushing the state down than up in energy (see Fig.2.3.4).

In the above discussion, the imaginary potential was introduced as an approximation to the breaking of a stationary state into many, more complicated stationary

states through the coupling to doorway states (Fig. 2.3.4(b)). This is the correct picture to describe the coupling of a nucleon moving in a single-particle state with more complicated configurations^{48,49}. However, such a description can become quite involved. On the other hand, to account for the change in the centroid energy and the spreading width of a single-particle state in terms of an *optical potential* ($\Delta E_\nu + iW_\nu$) is very economic. In any case Γ measures the range of energy over which the "pure" single-particle state $|a\rangle$ spreads due to the coupling to the more complicated doorway states $|\alpha'\rangle$. In other words, a stationary state

$$\varphi_\nu(\mathbf{r}, t) = e^{i\omega t} \varphi_\nu(\mathbf{r}), \quad (2.3.15)$$

has a probability density

$$\int d\mathbf{r} |\varphi_\nu(\mathbf{r}, t)|^2 = \int d\mathbf{r} |\varphi_\nu(\mathbf{r})|^2 = 1, \quad (2.3.16)$$

which does not depend on time. That is, if at $t = 0$, the probability that the particle is in a state ν is 1, it will have this probability also at $t = \infty$, implying an infinite lifetime. If however

$$\begin{aligned} \hbar\omega &= \varepsilon_\nu^{(0)} + \Delta E_\nu(\omega) + i\frac{\Gamma_\nu}{2}(\omega), & = \varepsilon_\nu + i\frac{\Gamma_\nu}{2}(\omega), & , (\varepsilon_\nu = \varepsilon_\nu^{(0)} + \Delta E_\nu; \Gamma_\nu = 2W_\nu) \\ \varphi_\nu(\mathbf{r}, t) &= e^{i\frac{\varepsilon_\nu t}{\hbar}} e^{-\frac{\Gamma_\nu t}{2\hbar}} \varphi_\nu(\mathbf{r}), \\ \int d\mathbf{r} |\varphi_\nu(\mathbf{r}, t)|^2 &= e^{-\frac{\Gamma_\nu t}{\hbar}}, \end{aligned} \quad (2.3.17)$$

implying a lifetime for the single-particle state given by

$$\tau = \Gamma_\nu/\hbar. \quad (2.3.18)$$

One may ask, how is it possible that the coupling to complicate (but still simple) states like $|\alpha'\rangle = |n_\alpha = 1, \nu'\rangle$ made out of a nucleon in the state ν' and a one phonon state of quantum numbers α , can explain the damping of a single-particle state lying 8-10 MeV from the Fermi energy ε_F , where the density of levels is expected to be conspicuous? This is because the Hamiltonian (2.3.1) contains the basic physics needed to describe the dressed single-particle motion as far as collective surface modes are concerned. Modes which, as a rule, play the central role in the dressing

⁴⁸See however Caldeira and Leggett (1981), Caldeira and Leggett (1983) and refs. therein.

⁴⁹To be noted that if we spread the strength of a stationary quantal state in a number of doorway stationary states over an energy range Γ (of the order of few MeV in the case of the GDR, see Sects. 1.5 and 5.5.1), and set all components in phase at $t = 0$, they will essentially be out of phase at time $t = \tau = \hbar/\Gamma$. In other words, each component will behave independent of each other and the original correlated state, created at $t = 0$ with probability 1 essentially ceases to exist at $t = \tau$. This does not imply that each of the incoherent members of the original coherent state cannot γ -decay at a much later stage ($\Gamma_{\gamma_0}/\Gamma \lesssim 10^{-2}$), see Bortignon et al. (1998). Discussions with B. Herskind on this issue are acknowledged.

of the single-particle motion in nuclei⁵⁰. Couplings to more complicated states go through a hierarchy of couplings. In other words, the variety of couplings should

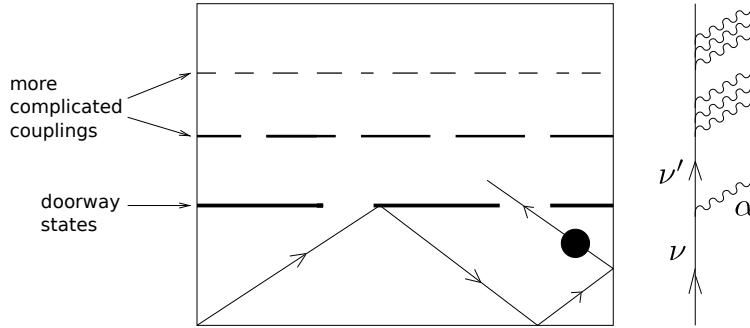


Figure 2.3.5: Schematic representation of the different levels of couplings leading to the damping of a single-particle state. It is essentially the first doorway coupling which controls the probability the ball (black dot) reflecting elastically on the walls of the box has to remain passed the first compartment. To the right, the coupling of the nucleon, moving originally in the state $|\nu\rangle$ (arrowed line), with an increasing number of phonons $|\alpha\rangle$ (wavy lines).

first go through the coupling to states of type $|\alpha'\rangle$ which act as proper doorway states (see Fig.2.3.5)⁵¹. Summing up, in the nuclear case, the *doorway coupling provides the basic mechanism to break the single-particle strength while, as a rule, higher-order couplings essentially fill in valleys* (see Fig.2.3.6).

In the case of the $1s_{1/2}$ orbital of ^{40}Ca ($\varepsilon - \varepsilon_F = -8$ MeV), simple estimates⁵² lead to $\bar{V}^2 \approx 0.3\text{MeV}$ for the coupling to an $L = 2$ phonon, and $n(\varepsilon_F) \approx 2 \text{ MeV}^{-1}$. Consequently

$$\Gamma \approx 4 \text{ MeV}, \quad (2.3.19)$$

in overall agreement with the experimental findings (see Fig. 2.3.7).

The estimate given in Eq.(2.3.19) is a particular example of the general (empirical) result (see Fig. 2.3.3)⁵³.

$$\Gamma_{sp}(E) = \begin{cases} 0.5|E| & |E| > 5 \text{ MeV}, \\ 0 & E \leq 5 \text{ MeV}, \end{cases} \quad (2.3.20)$$

where

$$E = |\varepsilon - \varepsilon_F|. \quad (2.3.21)$$

⁵⁰Something which was learned in the study of the nuclear spectrum in nuclei along the stability valley, and which have been found to remain correct also in connection with exotic nuclei. In particular light halo exotic nuclei regarding the phenomenon of parity inversion.

⁵¹Feshbach (1958).

⁵²Mahaux et al. (1985)

⁵³Bertsch et al. (1983)

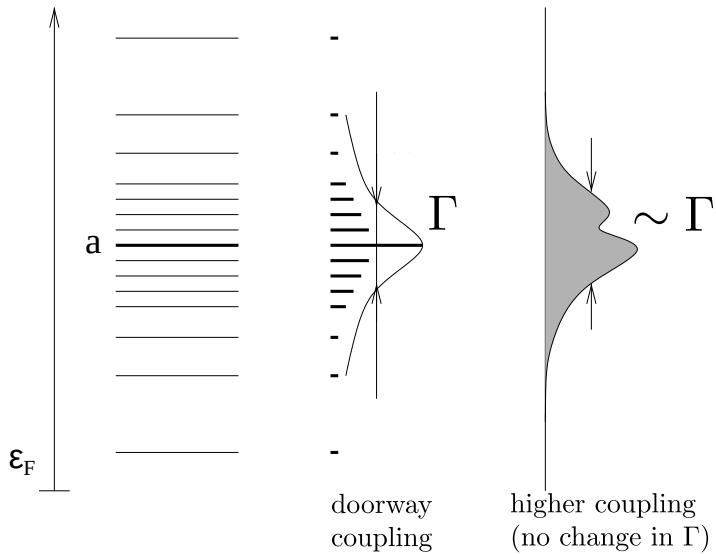


Figure 2.3.6: Schematic representation of the breaking of a single particle state $|a\rangle$ (heavy black horizontal line) through the coupling to doorway states ($|\alpha'\rangle = |n_\alpha = 1, \nu'\rangle$; thin horizontal lines) and eventually to increasingly more complicated (many-particle)-(many-hole) configurations.

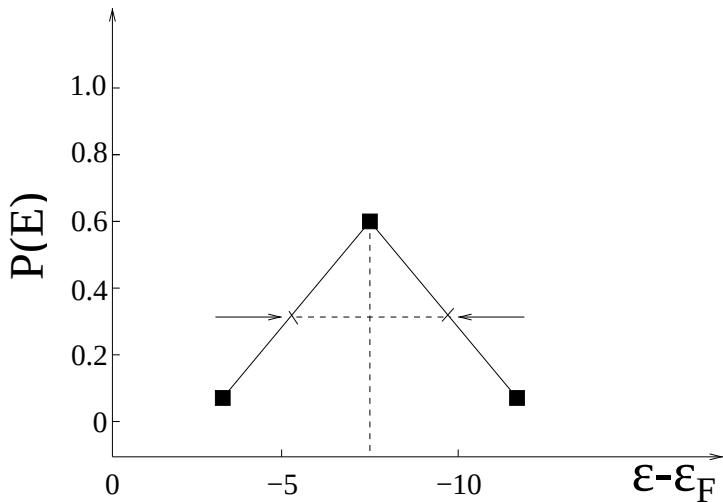


Figure 2.3.7: Schematic representation of the experimental strength function (solid squares) associated with the $2s$ state of ^{40}Ca . Also indicated is the full width at half maximum (FWHM).

2.3.1 Induced interaction

A nucleon close to the Fermi energy which, by bouncing inelastically off the nuclear surface excites a collective mode moving in the process to another, or remain-

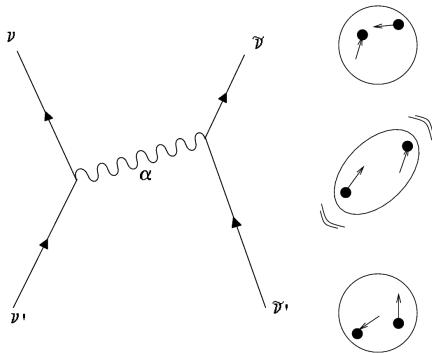


Figure 2.3.8: Schematic representation of the exchange of surface phonons between nucleons.

ing in the same state, has no other choices than to continue in such a state, or to reabsorb the vibration at a later instant of time (self-energy, Fig. 2.3.1). In the presence of another nucleon, the excited collective vibration by one nucleon may be absorbed by the second one (Fig. 2.3.8), the exchange of a vibration leading to an (induced) interaction.

Simple estimates of this induced interaction leads, in the case of ^{210}Pb , that is two neutrons above the $N = 126$ closed shell, to correlation energies for pairs of particles moving around it and coupled to angular momentum $J^\pi = 0^+$ of ≈ -1.5 MeV (see Fig. 2.3.9), when α is summed over the different multipolarities ($\lambda^\pi = 2^+, 3^-, 5^-$, label α in Fig. 2.3.8), and about half that value if one takes into account the fact that the single particle content for each of the interacting particles is about $Z_\omega \approx 0.7$. From this result one can conclude that the pairing interaction induced by the process depicted in Fig. 2.3.8, renormalizes in an important way the bare, $NN(^1\text{S}_0)$ -short range pairing interaction. This issue is taken up in more detail later on, in particular in connection with exotic halo nuclei (see e.g. Fig. 2.7.1 and Sect. 3.6). Concerning nuclei lying along the stability valley, see next section.

2.4 Well funneled nuclear structure landscape

Collective modes are composite. Being made out of fermions (nucleons) they contain in their microscopic structure, a consistent fraction of the nuclear correlations. One is then confronted to deal, in working out their interweaving through three-(PVC) and four- (ν) point vertices, with an overcomplete set of states. And, as a consequence, the need to correct for eventual Pauli-principle violations and non-orthogonality contributions.

The rules to do so have been cast into a graphical effective field theory, namely the nuclear field theory (NFT). In it, the free fields are to be calculated in the HF (HFB) approximation (particle (quasiparticle)) and in the RPA (QRPA) (vibrations). These elementary modes of excitation interact through the particle-vibration

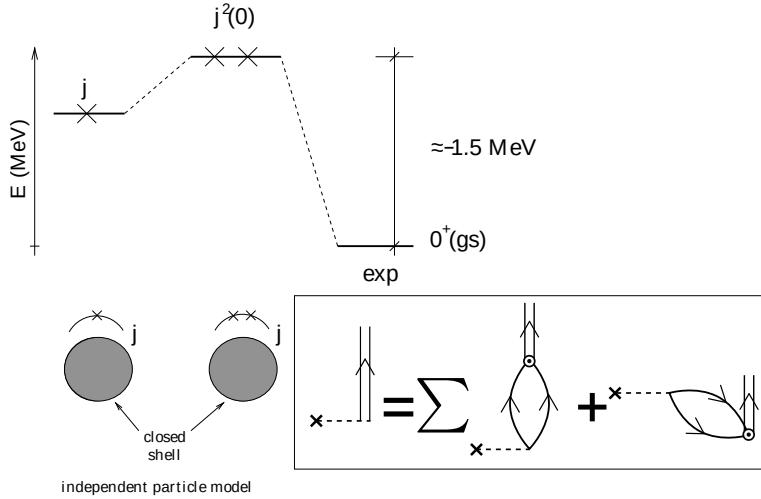


Figure 2.3.9: Schematic representation of the predictions of the independent particle model for one- and two-particles outside closed shell, in comparison with the experimental findings (e.g. for the case of ^{210}Pb , where $j = g_{9/2}$). In the inset, the graphical RPA dispersion relation of the monopole pair addition mode of ^{208}Pb , i.e. $|^{210}\text{Pb(gs)}$) is shown (see Sect. 3.5.2).

coupling vertices, while particles can also interact through four-point vertices (NN -bare interaction)⁵⁴.

A similar situation is found in the case of transfer processes in general, and of two-nucleon transfer in particular. One can work out the associated transfer amplitude by orthogonalizing, making use of second order perturbation theory, the single-particle wavefunctions of target and projectile. In other words, correct the corresponding expressions for non-orthogonality to the order of perturbation needed. This can be done both within the semiclassical approximation⁵⁵ or the DWBA⁵⁶.

The NFT rules for evaluating the effect of the couplings between fermions and (quasi) bosons involve a number of restrictions concerning initial and intermediate states as compared with the usual rules of perturbation theory. This is in keeping with the fact that the collective modes contain, from the start, the correlations arising from forwards and backwards going particle-hole ($\beta = 0$) as well as particle-particle ($\beta = +2$) and hole-hole ($\beta = -2$) bubbles, where β is the transfer quantum number.

The general validity of NFT rules have been demonstrated by proving the equivalence existing, to each order of perturbation theory, between the many-body

⁵⁴In connection with the reaction processes one finds again four-point (e.g. the proton-neutron interaction in the (p, d) reaction) and three-point vertices (e.g. particle-recoil coupling vertices).

⁵⁵See Sects. 6.1 and 6.5, and Sect. 2.B.1 of App. 2.B.

⁵⁶Sect. 4.2 and 6.1; see also the discussion connected with the origin of the tunneling Hamiltonian (4.6.1), in connection with weak coupled metallic superconductors and the Josephson effect.

finite nuclear system propagator calculated in terms of Feynman diagrams involving only the fermionic degrees of freedom i.e. explicitly respecting Pauli principle in a complete and not overcomplete basis, also known as Feynman-Goldstone propagator, and the propagator constructed in terms of Feynman diagrams involving fermion and phonon degrees of freedom (NFT Feynman diagrams) in the case of a general two-body interaction and an arbitrary distribution of single-particle levels⁵⁷.

Concerning the actual embodiment of NFT one can recognize the practical difficulties of respecting the corresponding rules. This is in keeping with the fact that at present there is not a bare, well behaved, low- k NN -force (eventually with $3N$ and higher order corrections) with which it is possible to generate a mean field, containing also the Fock component, to determine the single-particle states (HF) and, by introducing a periodic time-dependence calculate the collective modes (RPA and QRPA) associated with the variety of particle-hole ($\beta = 0$; density, spin, isospin, etc.) and pairing ($\beta = \pm 2$; monopole and multipole pair addition and pair subtraction) modes⁵⁸.

On the other hand, implementation of the NFT rules (renormalization)⁵⁹ have been carried out, making use of the bare Argonne v_{14} potential, and of the Skyrme force⁶⁰ SLy4 to determine the mean field and spin vibrational channels, and of multipole–multipole forces with self-consistent coupling constants for the variety of density vibrational channels.

The resulting structure predictions can, as a rule, together with the corresponding reaction computational codes, in particular COOPER and ONE (see App. 7.A), provide an overall account of “complete” sets of experimental data, obtained with the help of Coulomb, inelastic, γ -decay and one- and two-nucleon transfer data, which extensively maps out the nuclear structure and reaction landscape⁶¹, as discussed below.

Similar, but more accurate results are obtained making use of renormalized NFT, which parallels renormalized QED (see e.g. Sects. 5.2.2–5.2.4; see also Sect. 5.3.3 and Table 5.3.1). Namely: a) adjusting the bare potential parameters to be used together with a k -mass (similar to the renormalization of the bare mass in QED); b) making use of empirical values of the deformation parameters (PVC vertices) and phonon energies (parallel to renormalization of the bare charge)⁶².

2.4.1 The island of open-shell superfluid nuclei $^{118,119,120,121,122}\text{Sn}$

In what follows, we bring one- and two-particle transfer reactions together with an essentially “complete” set of other observables for the open-shell nucleus ^{120}Sn

⁵⁷Bès and Broglia (1975) and Bès et al. (1976a); see also Baranger (1969) and the lecture notes of McFarlane (1969).

⁵⁸See however footnote 10 Chapter 1.

⁵⁹Broglia et al. (2016).

⁶⁰Chabanat et al. (1997)

⁶¹Idini et al. (2015, 2014); Potel et al. (2013a).

⁶²Barranco et al. (2017, 2020). See also Sects. 5.2.2–5.2.4.

lying along the stability valley. From this example one can see that: 1) it is possible to predict, with few free parameters, most of them strongly constrained by empirical input, the experimental findings within a 10% level of accuracy; 2) the nuclear landscape, as it emerges from NFT based on elementary modes of excitation and of their interweaving, is well funneled (Fig. 2.4.1, see also table 2.4.1)⁶³, its minimum essentially coinciding with the global minimum resulting from the empirical renormalization choice of basic quantities (m_k , strength bare pairing, properties of few low-lying collective ($p - h$) modes and pairing vibrations). An important proviso concerning the above parlance is that one considers a group of homogeneous nuclei as e.g. open shell spherical superfluid nuclei (like the Sn isotopes)⁶⁴, or nuclei around closed shells (like $^{208,209,210}\text{Pb}$, ^{209}Bi , ^{210}Po , etc., or $^{10,11,12}\text{Be}$, $^{9,10,11}\text{Li}$, etc.)⁶⁵.

Let us now comment on Fig. 2.4.1. In it, the root mean square deviations $\sigma(x)$ (see also Table 2.4.1) between theoretical predictions and experimental values of the different structural properties which characterize ^{120}Sn are displayed. The calculations involve the island of open-shell superfluid nuclei $^{118,119,120,121,122}\text{Sn}$. The root mean square deviations are displayed as a function of the pairing coupling constant G (referred to the employed value $G_0 = 0.22$ MeV), the k -mass m_k ($(m_k)_0 = 0.7m$), the dynamical quadrupole deformation parameter β_2 ($(\beta_2)_0 = 0.13$) and in general as a function of x (x_0), measured with respect to the minimum value $\sigma_{\min} = \sigma(x_{\min})$, displayed in the interval $0.5 \leq x/x_0 \leq 1.5$ and *normalized* according to $0 \leq N(\sigma(x) - \sigma_{\min}) \leq 1$.

The curves represent: the deviation from experiment of the pairing gap associated with the $h_{11/2}$ orbital ($\Delta_{h_{11/2}}(G/G_0)$ (I); $\Delta_{h_{11/2}}(m_k/(m_k)_0)$ (II); $\Delta_{h_{11/2}}(\beta_2/(\beta_2)_0)$ (III); the deviation of the quasiparticle spectrum ($E_{qp}(G/G_0)$) (IV); $E_{qp}(\beta_2/(\beta_2)_0)$ (V); the deviation of the $h_{11/2} \otimes 2^+$ multiplet splitting $E_{h_{11/2} \otimes 2^+}(\beta_2/(\beta_2)_0)$ (VI); the deviation of the centroid position of the $d_{5/2}$ strength function $S_{d_{5/2}}(\beta_2/(\beta_2)_0)$ (VII); the deviation of the width of the $d_{5/2}$ strength function $S_{d_{5/2}}(\beta_2/(\beta_2)_0)$ (VIII); the deviation of the quadrupole transition strength $B(E2)(\beta_2/(\beta_2)_0)$ (IX); the deviation of the two-neutron transfer cross section $\sigma_{2n}(p, t)$ (G/G_0) (X). For an overview see Fig. 2.8.1. The basic feature displayed by Fig. 2.4.1 is the fact that, in spite of the fluctuations of the results typical of finite many-body systems, they clearly define a funnel in which all minima fall within a narrow window of x/x_0 values (1 ± 0.2).

This can be considered an emergent property of a description of structure and reactions carried out in a basis of elementary modes of excitation interacting, according to the NFT rules, through the NN -interaction as well as through PVC and PRC vertices.

The concept of a well funneled energy landscape is easy to understand in the case in which the number of particles $N \rightarrow \infty$ (thermodynamic limit). For example,

⁶³For details see Idini et al. (2015).

⁶⁴Within this context we refer to the conclusions of Idini et al. (2015).

⁶⁵In connection with light exotic halo nuclei we refer to Sect. 2.7 of the present Chapter and to Sects. 5.2.4 and 7.1.

a swing has a very simple and well funneled potential energy landscape. A similar concept which may still retain the (semi) classical viewpoint, but now referred to the free energy of large molecules, has been used in an attempt to describe protein folding⁶⁶. One has hypothesized that the results of an all atom, explicit solvent classical molecular dynamic simulations can be interpret in terms of a somewhat rugged, but still well funneled free energy landscape.

When we see such a behaviour in the nuclear case, even not so well defined, and somehow imperfect, we recognize that the nucleus is, after all, not macroscopic⁶⁷. Concepts strictly valid for $N \rightarrow \infty$ are strongly renormalized by quantal finite size effects, in particular zero point fluctuations (ZPF, Fig. 2.4.2)⁶⁸.

On the other hand, it is likely that the “imperfect” nature of the nuclear structure landscape funnel, an example of which is shown in Fig. 2.4.1, embodies more accurately the physics of quantal many-body systems than that of a smooth, more pedagogical construct essentially based on potential energy, even with the entropic contribution (free energy). This is in keeping with the fact that the interaction terms (potential energy) of the Hamiltonian contain the last vestiges of Newton’s conception of force or, causation, being thus too much anchored to classical mechanics⁶⁹.

Within this context we refer to Fig. 2.4.2, and to the fact that the ground state (nuclear vacuum) contains the physics of the system in terms of virtual processes. This is in keeping with the fact that acting on the system with the variety of probes available in the laboratory one obtains as on-shell final states, whose properties can eventually be observed with the help of the appropriate detectors, the variety of dressed (renormalized) elementary modes of excitation of the nucleus under study. In particular, diagrams (b), (c) of Fig 2.4.2 describe inelastic and two-nucleon transfer processes while diagrams (d)–(f) portray one-nucleon transfer reactions.

In all orders in the particle vibration coupling vertex ((PVC), solid dot), starting from second order (graph (f)), *NFT diagrams take care of the Pauli principle acting between the quasiparticles considered explicitly (continuous curves) and those*

⁶⁶Wolynes; Wolynes et al. (2012); see also cover figure of the issue of PNAS in which this reference was published.

⁶⁷Anderson (1972).

⁶⁸This nuclear result in turn, may be used at profit to shed light into the physics which is at the basis of protein folding. These molecules are after all quantal systems, and the associated quantal fluctuations are likely to play an important role in the associated fleeting transition states. The fact that average values and shapes of the transit-time distributions agree well with the simplest one-dimensional theory may only reflect the large uncertainties of the tunneling probabilities, small changes in the barrier’s parameters compensating for the lack of quantal effects, let alone the fact that being most high-dimensional models, as well as real processes, “sloppy”, their behavior depend on very few parameters (collective variables (CV), see Buchanan (2015), Transtrum et al. (2015), and references therein). It is unlikely that one or few of them are not related to quantal fluctuations. This is particularly true if one considers a NMR chemical shift biased molecular dynamic simulation, in keeping with the fact that NMR, based on the precession of nuclear spins is by its essence, a quantal phenomenon. Within this context one is reminded of α - and exotic-decay (see e.g. Ch. 7 of Brink and Broglia (2005) and references therein). While setting the formation probability of an alpha particle equal to one, could be compensated by modest changes in the associated tunneling probability (barrier), this was not possible in the case of e.g. exotic decay.

⁶⁹Born (1948) pp. 95,103; Pais (1986) p 258.

Observables	SLy4	$d_{5/2}$ shifted	Opt. levels	Fig. 2.4.1
Δ	10 (0.7%)	10 (0.7 %)	50 (3.5 %)	(I), (II), (III)
E_{qp}	190 (19%)	160 (16%)	45 (4.5 %)	(IV), (V)
Mult. splitt.	50 (7%)	70 (10%)	59 (8.4 %)	(VI)
$d_{5/2}$ strength (centr.)	200 (20%)	40 (4%)	40 (4%)	(VII)
$d_{5/2}$ strength (width)	160 (20%)	75 (9.3%)	8 (1%)	(VIII)
$B(E2)$	1.4 (14%)	1.34 (13%)	1.43 (14%)	(IX)
$\sigma_{2n}(p, t)$	40 (2%)	40 (2%)	40 (2%)	(X)

Table 2.4.1: Root mean square deviation σ between the experimental data and the theoretical values expressed in keV for the pairing gap, quasiparticle energies, multiplet splitting, centroid and width of the $5/2^+$ low-lying single-particle strength distribution (Fig. 5.2.3). In single-particle units B_{sp} for the γ -decay ($B(E2)$ transition probabilities) and in mb for $\sigma_{2n}(p, t)$ (Fig. 7.2.1). In brackets, the ratio $\sigma_{rel} = \sigma/L$ between σ and the experimental range L of the corresponding quantities, is given: 1.4 MeV (Δ), 1 MeV (E_{qp}), 700 keV (mult. splitting), 1 MeV ($d_{5/2}$ centroid), 809 keV (=1730–921) keV; $d_{5/2}$ width), 10 B_{sp} ($B(E2)$), 2250 μ b ($\sigma_{2n}(p, t)$)) is given. Columns 2,3, and 4 contain the results of NFT calculations making use of bare single-particle levels from Hartree-Fock with Sly4, same but for a 600keV shift towards the Fermi energy of the $\epsilon_{d_{5/2}}$ orbital, and optimal values of ϵ_j for all valence levels so that the dressed quasiparticle states provide the best fit to that data, respectively. In the last column reference to the corresponding diagrams shown in Fig. 2.4.1 are given. See also Fig. 2.8.1.

participating in the modes (double wavy lines), as well as between modes. As a consequence, self energy processes based on pure or little collective two quasiparticle excitations are correctly screened out.

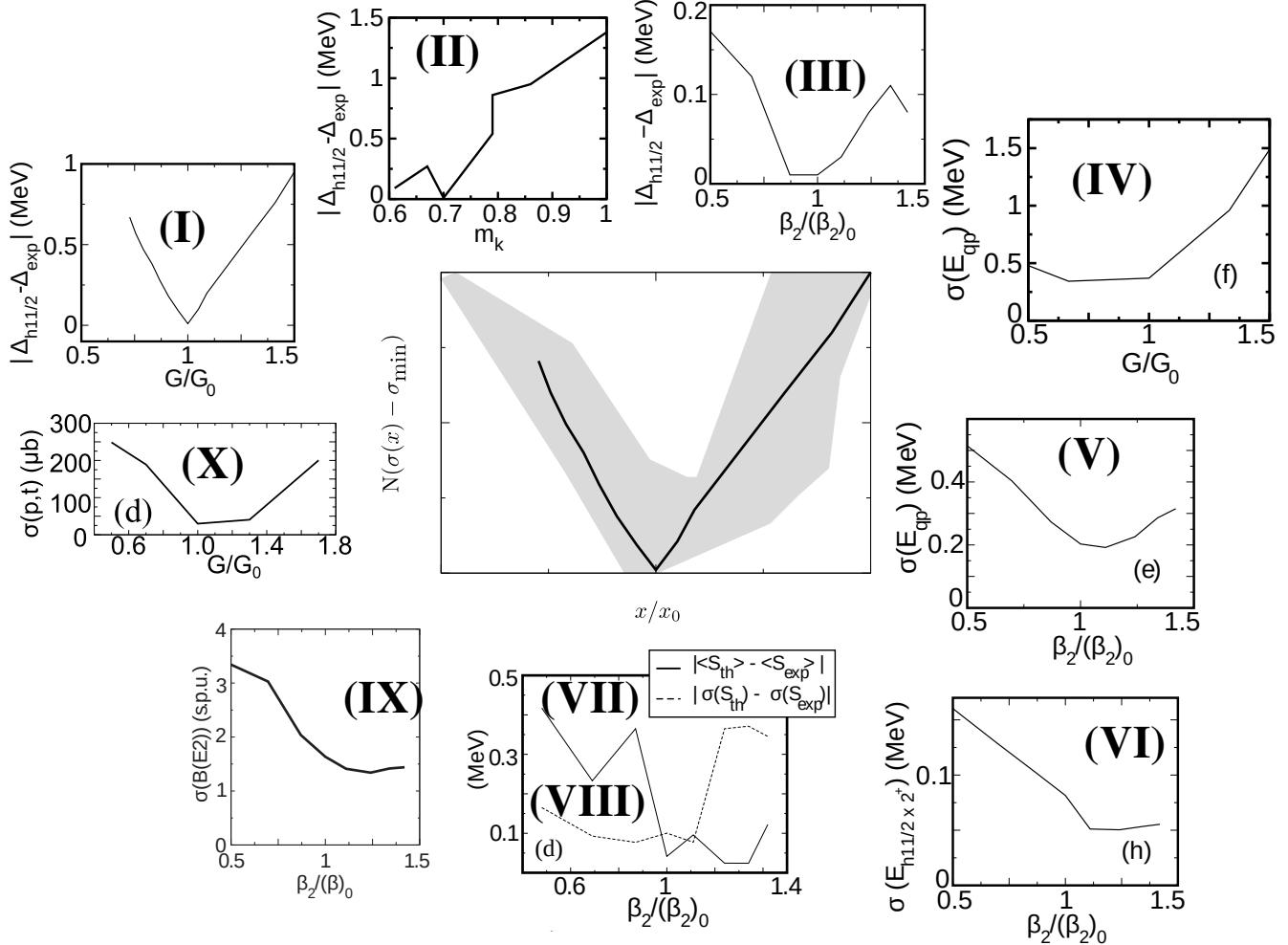


Figure 2.4.1: Root mean square deviations $\sigma(x)$ between theoretical predictions and experimental values of the different structural properties which characterize the open-shell nucleus ^{120}Sn (see text as well as Table 2.4.1 for the characterization of (I)–(X); see also Fig. 2.8.1); (central figure) the bold black line and grey area provide a schematic representation of $\langle \sigma \rangle$ and associated fluctuations.

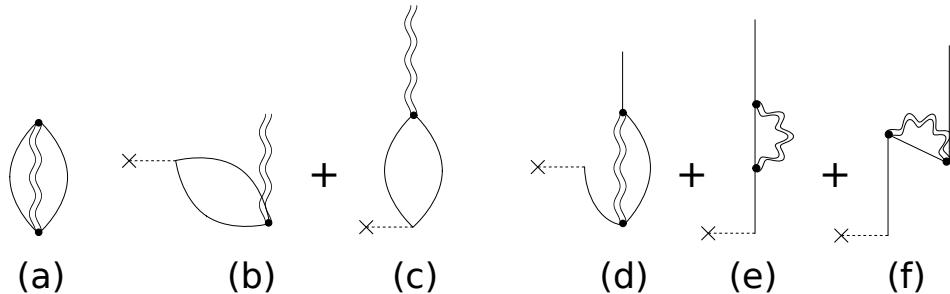


Figure 2.4.2: Schematic representation of the $\text{NFT}_{(s+r)}$ diagrams at the basis of the characterization of a superfluid nucleus like e.g. ^{120}Sn . **(a)** Nuclear structure ($\text{NFT}_{(s)}$). Zero point fluctuations (ZPF) characterizing the nucleus ground state. Continuous lines describe quasiparticle (qp) states, double wavy curves stand for correlated two quasiparticle ($2qp$) vibrational modes. Because $\alpha_\nu^\dagger = U_\nu a_\nu^\dagger - V_\nu a_\nu$ these modes encompass both particle-hole (ph) like vibrations, e.g. surface quadrupole vibrations, as well as correlated (pp) and (hh) monopole and multipole pairing vibrations. Intervening with an external field (cross followed by dashed line) one can excite **(b)** and **(c)** multipole (ph -like; inelastic scattering) and pairing ($2p$ -like, $2h$ -like) vibrations (two-particle transfer) as well as **(d)-(f)** single-quasiparticle states.

2.5 Nuclear Field Theory for pedestrians

The diagrams were intended to represent physical processes and the mathematical expressions to describe them... I would see electrons going along, being scattered at one point... emitting a photon and the photon goes over there... I thought that if they really turn out to be useful it would be fun to see them in the pages of Physical Review

R. P. Feynman

Nuclear Field Theory (NFT) was tailored after Feynman's graphical version of quantum electrodynamics (QED). It is then natural that in discussing NFT, analogies with QED are recurrent. Arguably, as a consequence of special relativity which put an end to the concept of ether, the field-free and matter-free vacuum was rightly considered as *bona fide* empty space. The advent of quantum mechanics changed this situation, the vacuum becoming populated.

In quantum mechanics an oscillator, for example, cannot be at rest. The os-

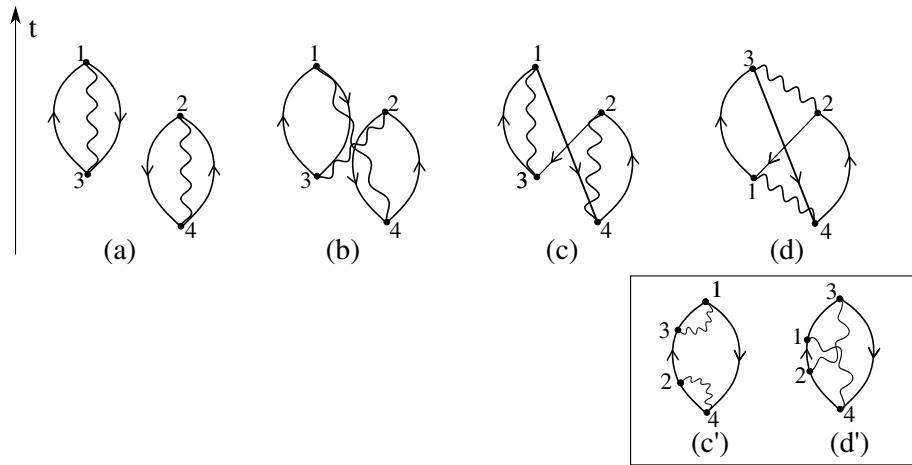


Figure 2.5.1: Oyster diagrams describing the correlation of the nuclear ground state associated with the ZPF of collective particle–hole–like excitations. In (a) we show two of such diagrams. In (b) and (c) we display a symmetrized (boson exchange), and antisymmetrized (fermion exchange) correction to (a), while (d) contains a simultaneous boson and fermion exchange. In all the diagrams shown, only ground state correlation vertices are present. They are connected with the $Y_{k_i}^\alpha$ –components ($\epsilon_k > \epsilon_F$, $\epsilon_i \leq \epsilon_F$) of the RPA wavefunction describing the collective mode (wavy line). While this is so for any time ordering, i.e., the sequence with which the particle-vibration coupling vertex (black dots) appear in the case of the processes shown in (a) and (b), this is not the case in connection with processes shown in (c) and (d) as can be seen from the corresponding diagrams (c') and (d') shown in the inset. Because of Pauli principle between particles (holes) present and those involved in the collective modes, the harmonic approximation has to be corrected. This is diagrammatically reflected by the presence of scattering vertices.

cillatory nature of the radiation field requires zero point fluctuations (ZPF) of the electromagnetic fields in the vacuum state. The occupation of the negative kinetic energy electron states and the subsequent calculation of the cross section for pair creation by photons, contributed another step in the understanding of the QED vacuum, let alone the Lamb shift⁷⁰.

When the fields are expressed in terms of creation and annihilation operators, the interaction between fermion and boson fields is proportional to the product of two fermion creation or destruction operator a^\dagger or a , and of one boson operator Γ^\dagger or Γ : e.g. $a_\nu^\dagger a_\nu \Gamma_\alpha^\dagger$, (see Fig. 2.5.1). That is, bilinear in the fermion fields and linear in the (quasi-) boson fields.

⁷⁰In high energy collisions and accelerator laboratories some of the original beam energy can be consumed by ripping electron-positron pairs out of the vacuum (Bruce et al. (2007)).

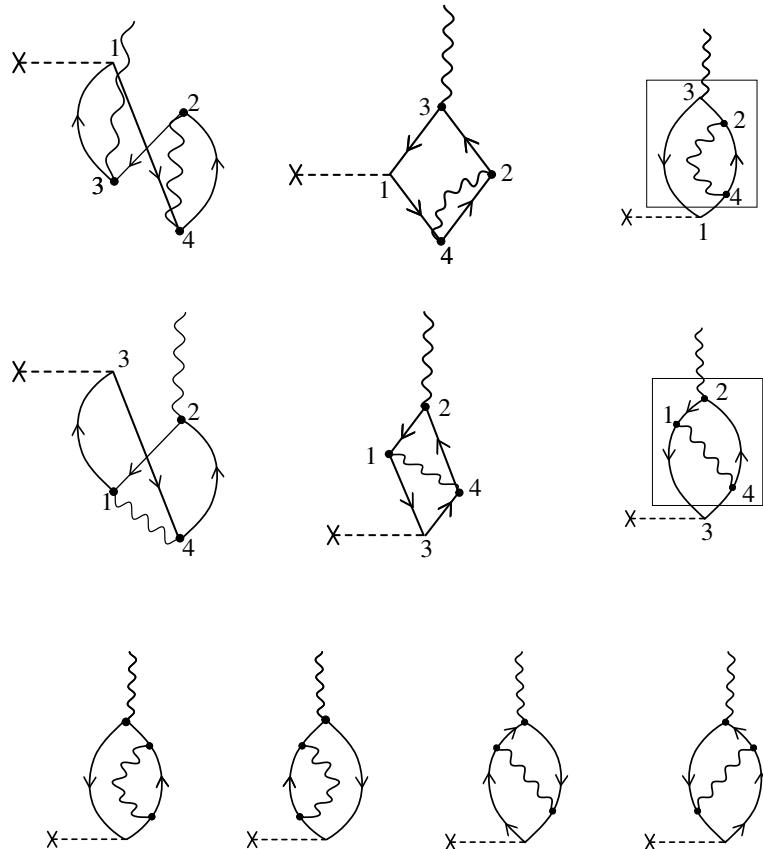


Figure 2.5.2: Some of the possible outcomes resulting from acting with an external single-particle field, i.e. that associated with inelastic processes (represented by a horizontal dashed line starting with a \times) on the ZPF of a nucleus ground state associated with particle-hole correlated vibrations. Within this context one returns to the question of renormalization mentioned in the text (see also Idini et al. (2015), Broglia et al. (2016), Barranco et al. (2017); see also Sect. 5.3) The diagrams of the first row result by intervening the virtual process shown in Fig. 2.5.1 (c) and eventual time orderings. Similar for those of the second row but in connection with diagram (d) of Fig. 2.5.1. The boxed processes correspond to particle self-energy (first row) and vertex correction (second row). Reversing the sense in which the fermions (arrowed lines) circle the loop from anticlockwise to clockwise, one obtains two new graphs. The set of processes obtained in this way are shown in the third and last row, and constitute a sum rule conserving set of diagrams (see first row Fig. 1.2.6).

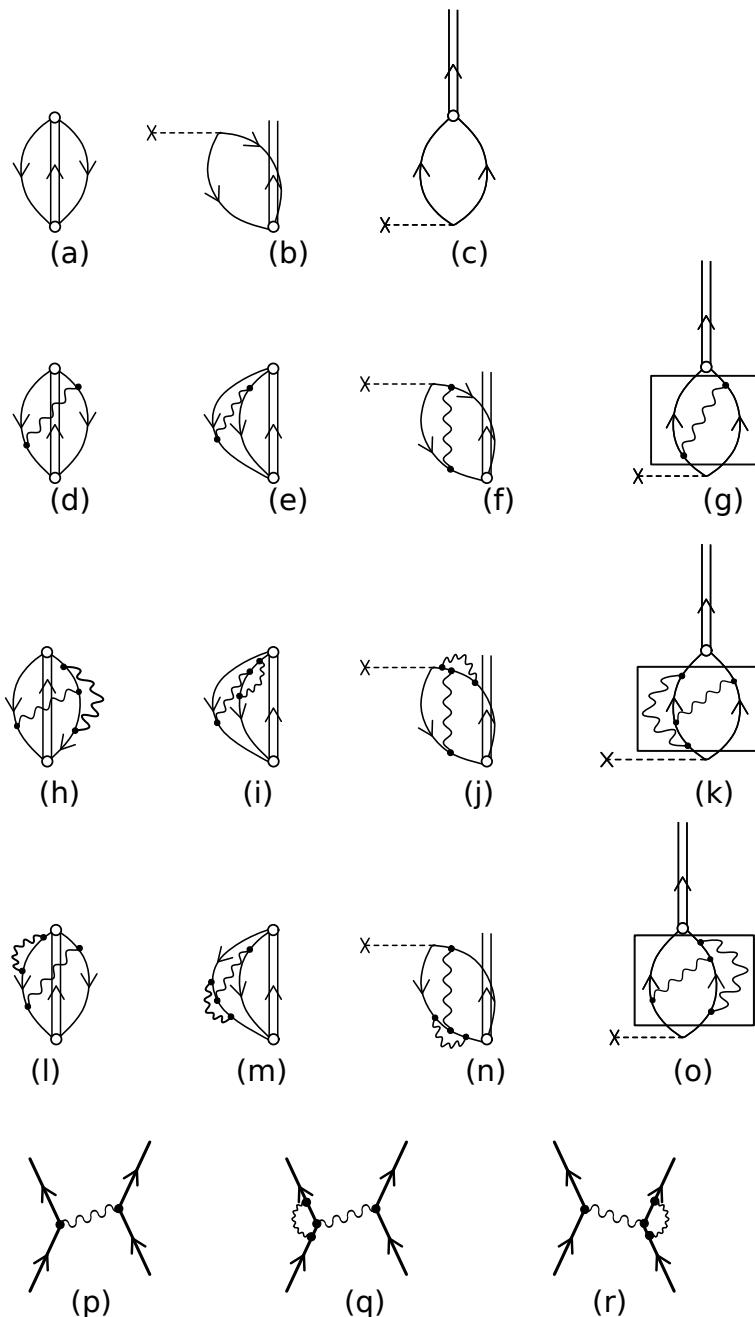


Figure 2.5.3: ZPF associated with the pair addition mode taking into account the interweaving of nucleons with density modes. The processes boxed in (g), (k) and (o), are associated with the induced pairing interaction (medium polarization effects; (p), (q), (r)) resulting from the exchange of density modes between nucleons moving in time reversal states, including also vertex corrections. The two-nucleon stripping and pickup external field is labeled by a dashed horizontal line which starts with a \times . The possibility of using pairing vibrational modes as intermediate bosons contributing to the induced pairing interaction, not only in 1S_0 channels but also in other channels (multipole modes) is discussed in App. 7.D. In particular, in connection with the possible presence of “vortex-like” pair addition modes, in exotic, halo nuclei, with $J^\pi = 1^-$ and $\beta = +2$ quantum numbers, connected also with the pygmy dipole mode in ^{11}Li (see e.g. Broglia et al. (2019a) and refs. therein).

A detailed graphical NFT treatment of the vacuum has an important consequence concerning the probing of nuclear structure with reactions. By intervening it with an external field one will excite the modes whose properties can be compared with experimental findings.

In other words, if one is in doubt of which are the properly dressed elementary, physical modes of excitation, one should find out how to specifically excite the mode in question, by acting with an external field on the ZPF of the vacuum (Hawking-like radiation⁷¹; see also Sect. 5.3 in particular Fig. 5.3.2). That is, carry out a NFT-like *gedankenexperiment*⁷² as in Fig. 2.5.2 for $p - h$ -vibrations and in Fig. 2.5.3 regarding pairing vibrations. Because the corresponding processes deal with physical states, they translate with ease into a laboratory setup. In keeping with the fact that the vacuum contains all the information (physical degrees of freedom) of the quantal system under study. Forcing virtual processes associated with vacuum ZPF to become real, one gets, in each instance, the real, dressed, physical mode of excitation.

The fact that one can treat fully quantum mechanically and on equal footing both structure and reaction processes is apparent from these figures and considerations⁷³. Thus unification of structure and reactions, and dressing of energies, vertices (interactions) and formfactors (single-particle radial wavefunctions and associated transition densities), results in a single vacuum correlated state (see e.g. Figs. 2.5.1 and 2.6.2). Vacuum states which through their fluctuations, reflect both single-particle, normal and abnormal (pairing) density vibrations and their interweaving. It also indicates the set of specific probes which make these virtual states to collapse into on-the-energy shell states, providing the corresponding physical information to the outgoing particles, also photons, which eventually interact with the corresponding detectors. Structure and reaction processes free of non-orthogonality, overcompleteness and Pauli principle violation contributions.

Summing up, the last line of Fig. 2.5.2 displays, together with the corresponding time orderings, the lowest order self-energy and vertex correction renormalising (particle-hole) modes. It thus gives rise to the physical collective vibrations whose properties can be directly compared with the experimental findings. In other words, the processes shown in the last line of Fig. 2.5.2 imply that the elementary modes participating in the virtual states have to display, exception made for energy (off-shell modes) the same properties of the physical, dressed (renormalized), on-shell modes whose properties can be directly compared with experiment⁷⁴ (renormalized NFT). This is because one can, through an experiment force such virtual states to become real (on-shell) on short call.

Let us now provide an introduction to NFT for pedestrians and see how the

⁷¹Barranco et al. (2019b).

⁷²H. C. Øersted, circa 1812.

⁷³Embodiments being found in e.g. Figs. 2.7.2, 2.7.3 and 5.3.2.

⁷⁴Examples of these processes in the case of giant resonances are found in Bortignon and Broglia (1981); Bertsch et al. (1983) and references therein. For low-lying states see Barranco et al. (2004).

above considerations become implemented⁷⁵.

2.5.1 The concept of elementary modes of excitation⁷⁶

The Hamiltonian of a many–body system of noninteracting particles, bosons or fermions, can be written as

$$H = \sum_i H_i, \quad (2.5.1)$$

where the summation is over all the particles of the system and where each H_i depends only on the variables of the i -th particle. The single-particle Schrödinger equation is

$$H_i \psi_k(\mathbf{r}_i) = \epsilon_k \psi_k(\mathbf{r}_i), \quad (2.5.2)$$

where ϵ_k is the single–particle energy eigenvalue and

$$\psi_k(\mathbf{r}_i) \equiv \langle \mathbf{r}_i | a_k^\dagger | 0 \rangle \quad (2.5.3)$$

is the corresponding wave function. The operator a_k^\dagger creates a particle in the state k when acting in the vacuum state $|0\rangle$. The energy levels of the system are given by the equation

$$E_n = \sum_k n_k \epsilon_k, \quad (2.5.4)$$

the corresponding eigenstates being

$$|n\rangle = \prod_k \frac{(a_k^\dagger)^{n_k}}{\sqrt{n_k!}} |0\rangle, \quad (2.5.5)$$

where $n_k = 0$ or 1 in the case of fermions and $n_k = 0, 1, 2, \dots$ in the case of bosons.

We now consider a system of interacting particles. The Hamiltonian will in this case be

$$H = \sum_i H_i + \frac{1}{2} \sum_{i,j} H_{ij}, \quad (2.5.6)$$

where i, j label the co–ordinates of the i -th and j -th particle.

In some cases it is possible to recast the two–body Hamiltonian in the form

$$H = \sum_\tau H'_\tau, \quad (2.5.7)$$

⁷⁵For details we refer to Bortignon et al. (1977) and refs. therein.

⁷⁶Bès and Broglia (1977)

with the associated Schrödinger equation

$$H'_\tau \psi_\tau(\zeta) = \epsilon_\tau \psi_\tau(\zeta), \quad (2.5.8)$$

ζ representing a general variable (e.g. the single-particle coordinate, the gap parameter, the shape of the nucleus, etc.). The wave function $\psi_\tau(\zeta)$ is the ζ -coordinate representation of the eigenstate $\Gamma_\tau^\dagger |\tilde{0}\rangle$. The operator Γ_τ^\dagger creates an excitation with the quantum number τ when acting in the state $|\tilde{0}\rangle$, the correlated vacuum of all the excitations τ .

The energy of the levels of the system, or at any rate of the most important ones to determine the physical response of it to external probes can be expressed in the form

$$E_m = \sum_\tau n_\tau \epsilon_\tau. \quad (2.5.9)$$

The corresponding eigenstate can be written in the same way as before, *i.e.*

$$|n\rangle = \prod_\tau \frac{(\Gamma_\tau^\dagger)^{n_\tau}}{\sqrt{n_\tau!}} |\tilde{0}\rangle. \quad (2.5.10)$$

Additivity features similar to (2.5.9) hold for other physical quantities, *i.e.*

$$\langle n | O | m \rangle = \sum_\tau A_\tau \sqrt{n_\tau} \delta(n_\tau, m_\tau + 1), \quad (2.5.11)$$

where

$$O = \sum_\tau A_\tau \Gamma_\tau^\dagger \quad (2.5.12)$$

is the operator which specifically excites the eigenstates described by $\psi_\tau(\xi)$. Because the excitation energies E_m and observables $|\langle m' | O | m \rangle|^2$ (e.g. absolute two-particle transfer cross-section, electromagnetic-transition probabilities, etc.) are linear combinations of ϵ_τ and A_τ , respectively, the eigenstates with energy ϵ_τ and associated observable A_τ are called the *elementary excitations of the system*.

The elementary modes of excitation of a many-body system represent a generalization of the idea of normal modes of vibration. They provide the building blocks of the excitation spectra, giving insight into the deep nature of the system one is studying, aside from allowing for an economic description of complicated spectra in terms of a gas of, as a rule, weakly interacting bosons and fermions. In the nuclear case they correspond to dressed particles and empirically renormalised vibrations (rotations).

There lie two ideas behind the concept of elementary modes of excitation⁷⁷. First, that one does not need to be able to calculate the total binding energy of

⁷⁷This concept was introduced by Landau (Landau (1941)) to describe the spectrum of He II. It was subsequently utilized by Bohr and Mottelson (Bohr and Mottelson (1975)) to obtain a unified description of the nuclear spectrum.

a nucleus to accurately describe the low energy excitation spectrum, in much the same way in which one can calculate the normal modes of a metal rod not knowing how to calculate its total cohesive energy. The second idea is that low-lying states ($\hbar\omega \ll \epsilon_F \ll BE$, i.e. binding energy) are of a particularly simple character, and amenable to a simple treatment, their interweaving being carried out at profit, in many cases, in perturbation theory⁷⁸. Within this context it is necessary to have a microscopic description of the ground state of the system which ensures that it acts as the vacuum state $|\tilde{0}\rangle$ of the elementary modes of excitation. In other words $a_\nu|\tilde{0}\rangle = 0, \Gamma_\alpha|\tilde{0}\rangle = 0$, where $a_\nu^+|\tilde{0}\rangle = |\nu\rangle$ and $\Gamma_\alpha^+|\tilde{0}\rangle = |\alpha\rangle$ represent a single-particle and a one-phonon state. This implies, in keeping with the indeterminacy relations⁷⁹ $\Delta x \Delta p \geq \hbar/2, \Delta I \Delta \Omega \geq 1, \Delta N \Delta \phi \geq 1$, etc., that $|\tilde{0}\rangle = |0\rangle_F|0\rangle_B$ displays the quantal zero point fluctuations (ZPF) of the many-body system under study.

Within the framework of nuclear field theory (NFT) used below, in which single-particle (fermionic, F) and vibrational (quasi-bosonic, B) elementary modes of excitation are to be calculated making use of HFB and QRPA respectively⁸⁰, $|\tilde{0}\rangle$ must display the associated ZPF (cf. App. 2.C). In particular for (harmonic) vibrational modes the indeterminacy relation achieves its lowest possible value $\Delta x \Delta p = \hbar/2$, the associated zero point energy amounting to $\hbar\omega/2$ for each degree of freedom. For example $5\hbar\omega/2$ for quadrupole vibrations, $\hbar\omega$ being the energy of the collective vibrational mode under consideration.

An illustrative example of the above arguments is provided by the low-lying quadrupole vibrational state of ^{120}Sn . Diagonalizing the effective SLy4 Skyrme interaction⁸¹ in QRPA leads to a value of $B(E2)(890 e^2 \text{ fm}^2)$ which is about a factor of 2 smaller than experimentally observed ($2030 e^2 \text{ fm}^2$). Taking into account renormalisation effects in NFT, namely in a conserving approximation (self-energy and vertex corrections), one obtains⁸² a value of $2150 e^2 \text{ fm}^2$.

If the collective phonons are not the main object of the study, but are to be used to dress the single-particle states and/or give rise to the induced pairing interaction, one can make use of phonons which account for the experimental findings (empirical renormalization⁸³)⁸⁴.

⁷⁸More precisely, and in keeping with the fact that (quasi) boson degrees of freedom have to decay through linear particle-vibration coupling vertices of strength Λ into their fermionic components to interact with another vibrational mode, the interweaving between the variety of many-body components dressing a single-particle state or a collective vibration will be described at profit in terms of an arrowed matrix which, assuming perturbation theory to be valid, can be transformed, neglecting contributions of the order of Λ^3 or higher, into a co-diagonal matrix, namely a matrix whose non-zero elements are $(i, i - 1)$ and $(i, i + 1)$, aside from the diagonal ones (i, i) .

⁷⁹The quantities I and N are the angular momentum and particle number, conjugate variables to the Euler and gauge angles Ω, ϕ respectively.

⁸⁰That is, Hartree-Fock (HF) and Random Phase (RPA) approximations extended to open shell superfluid nuclei. Namely Hartree-Fock-Bogoliubov and Quasi Random Phase Approximation.

⁸¹Ring and Schuck (1980).

⁸²Barranco et al. (2004).

⁸³Idini et al. (2015); Broglia et al. (2016); Barranco et al. (2017).

⁸⁴As already mentioned, with the help of experimental probes which couple weakly to the nucleus, i.e. in such a way that the system can be expressed in terms of the properties of the excitation

2.5.2 NFT rules and applications

A field theory can be formulated in which the nuclear elementary modes of excitation play the role of the free fields and in which their mutual interweaving takes place through four-point (v) as well as three-point, particle-vibration coupling vertices⁸⁵. This theory provides a graphical perturbative approach to obtain the exact solution of the many-body nuclear-structure problem in the product basis $\psi_\tau(\zeta)\psi_\eta(\Delta)\dots\psi_\gamma(\Gamma)$

In what follows we state and apply the nuclear-field-theory rules, to calculate the interactions between the nuclear free fields and the reaction processes between the resulting physical states making use of a simple model.

Schematic model

The model considered consists of two single-particle levels, each with degeneracy⁸⁶ Ω and with a schematic monopole particle-hole interaction coupling the particles in the two levels.

The total Hamiltonian is equal to

$$H = H_{sp} + H_{TB} \quad (2.5.13)$$

where

$$H_{sp} = \frac{\epsilon}{2}N_0, \quad N_0 = \sum_{\sigma=\pm 1,m} \sigma a_{m,\sigma}^\dagger a_{m,\sigma}, \quad (2.5.14)$$

and

$$H_{TB} = -\frac{V}{2} (A^\dagger A + AA^\dagger), \quad A^\dagger = \sum_m a_{m,1}^\dagger a_{m,-1}. \quad (2.5.15)$$

in the absence of probes (see however Sect. 7.4), it has been possible to identify, among others, the following elementary modes of excitation in systems around closed shells:

- a) single-particles and -holes,
- b) shape vibrations,
- c) spin and isospin vibrations and charge exchange modes,
- d) pairing vibrations.

Away from closed shells one has to add to the above modes:

- e) rotations and $2qp$ -like vibrations in 3D-space (e.g. quadrupole rotations and β - and γ -vibrations)
- f) rotations and $2qp$ -like vibrations in gauge space (pairing rotations and pairing vibrations).

Different probes have been utilized in the process of the identification of the different modes. In particular two-neutron transfer reactions induced by tritons and protons have played a central role in unraveling the basic features of the pairing modes.

⁸⁵Bès et al. (1974); Broglia et al. (1976); Bohr and Mottelson (1975); Mottelson (1976).

⁸⁶It is of notice the difference of a factor of 2 in the degeneracy of each level as compared to Sect. 2 of Bortignon et al. (1977) in which case it is 2Ω . This is in keeping with the fact that, as a rule, $\Omega = (2j + 1)/2$. See also Eq. (2.5.85) and related discussion.

The index σ ($=\pm 1$) labels the two levels, while m labels the degenerate states within each level. The strength of the monopole coupling is denoted V and the energy difference between the two levels ϵ . The matrix element of (2.5.15) is given by

$$\langle m, 1; m', -1 | H_{TB} | m'', 1; m''', -1 \rangle = -V\delta(m, m')\delta(m'', m'''). \quad (2.5.16)$$

Field-theoretical solutions

The (quasi-) boson fields are defined through the random-phase approximation, in terms of particle-hole excitations. The basis utilized to describe the nuclear systems is a product of the different free fields. The closed-shell system of the model under consideration corresponds to the lowest ($\sigma = -1$) level filled with Ω particles, while the upper ($\sigma = 1$) level remains empty. The basis particle and hole states are obtained by adding or removing a single particle to/from this closed-shell configuration. The corresponding wave functions and energies, which should include the Hartree-Fock corrections (see Fig. 1.2.2 (b), (c)) generated by the residual interaction⁸⁷, are

$$\begin{cases} |m, 1\rangle = a_{m,1}^\dagger |0\rangle, & E(m, 1) = \frac{1}{2}(\epsilon + V), \\ |m, -1\rangle = a_{m,-1}|0\rangle, & E(m, -1) = \frac{1}{2}(\epsilon + V). \end{cases} \quad (2.5.17)$$

Thus the unperturbed energy for producing a particle-hole excitation with respect to the ground state is

$$\epsilon' = E(m, 1) + E(m, -1) = \epsilon + V. \quad (2.5.18)$$

The term V in (2.5.18) is the Hartree-Fock contribution to the particle-hole excitation.

If we define the creation operator of the normal modes as

$$\Gamma_\nu^\dagger = \sum_m \lambda_m^\nu a_{m,1}^\dagger a_{m,-1}, \quad (2.5.19)$$

the linearization equation (see Eq. (1.1.15)),

$$[H, \Gamma_\nu^\dagger] = \omega_\nu \Gamma_\nu^\dagger, \quad (2.5.20)$$

yields

$$\begin{cases} \omega_1 = \epsilon' - V\Omega, \\ \omega_\nu = \epsilon' \quad (\nu = 2, 3, \dots, \Omega). \end{cases} \quad (2.5.21)$$

⁸⁷The Hartree-Fock energy associated with the Hamiltonian (2.5.13) can be obtained from the linearization relation $[H, a_{\sigma,m}^\dagger] = E(m, \sigma)a_{\sigma,m}^\dagger$ acting on the Hartree-Fock vacuum (see Eq. (1.1.16)), which in this case coincides with the single-particle vacuum defined by $a_{m,-1}^\dagger |0\rangle = a_{m,1}^\dagger |0\rangle = 0$.



Figure 2.5.4: Graphical representation of the amplitude of the collective phonon (wavy line) on a given particle-hole excitation $((m, 1), (m, -1))$. This amplitude can be written in terms of the interaction vertex denoted Λ_i , and the energy denominator $\omega_i - \epsilon'$. The particles (holes) are depicted by upward- (downward-) going arrowed lines. Time is assumed to run upwards.

Utilizing (2.5.20) and the normalization condition

$$[\Gamma_\nu, \Gamma_{\nu'}^\dagger] = \delta(\nu, \nu'), \quad (2.5.22)$$

we obtain for the amplitudes associated with the lowest mode

$$\lambda_m^1 = \frac{1}{\sqrt{\Omega}}. \quad (2.5.23)$$

One can also write this amplitude as the ratio between a coupling matrix element and an energy denominator, i.e.

$$\lambda_m^1 = \frac{\Lambda_1}{\omega_1 - \epsilon'}. \quad (2.5.24)$$

From (2.5.21), (2.5.23) and (2.5.24) we obtain

$$\Lambda_1 = -V \sqrt{\Omega}, \quad (2.5.25)$$

which is the strength with which a particle hole excitation $(m, 1; m, -1)$ couples to the collective phonon (see Fig. 2.5.4). This can also be seen by calculating the matrix element of the interaction Hamiltonian (2.5.15) between the normal modes and the single particle-hole state

$$\Lambda_\nu = \langle n_\nu = 1 | H_{TB} | m, 1; m', -1 \rangle = -V \sqrt{\Omega} \delta(m, m') \delta(\nu, 1). \quad (2.5.26)$$

Note that the particle-vibration coupling strengths associated with the other normal modes lying at an energy ϵ' (see (2.5.21)) are equal to zero. The exact solution of (2.5.13) is reproduced by utilizing as the basic degrees of freedom both the vibrations (see (2.5.21)) and the particles (see (2.5.17)) coupled through the interaction

(2.5.16) (four-point vertex) and (2.5.26) (particle-vibration coupling). A significant part of the original interaction has already been included in generating the collective mode (2.5.21). This implies that the rules for evaluating the effect of the couplings (2.5.16) and (2.5.26) between fermions and bosons, involve a number of restrictions as compared with the usual rules of perturbation theory that are to be utilized in evaluating the effect of the original interaction (2.5.15) acting in a fermion space. They read as follows:

- I) In initial and final states, proper diagrams involve collective modes and particle modes, but not any particle configuration that can be replaced by a combination of collective modes. This restriction permits an initial state comprising the configuration $(n_v = 1; m)$, but excludes $(m', 1; m', -1; m, 1)$.
- II) The couplings (2.5.16) and (2.5.26) are allowed to act in all orders to generate the different diagrams of perturbation theory; the restriction I) does not apply to internal lines of these diagrams.
- III) The internal lines of diagrams are, however, restricted by the exclusion of diagrams in which a particle-hole pair is created and subsequently annihilated without having participated in subsequent interactions.
- IV) The energies of the uncoupled particle and phonon fields are to be calculated by utilizing the Hartree-Fock approximation (see eq. (2.5.17)) and the RPA (see eq. (2.5.21)), respectively. The contributions of all allowed diagrams are evaluated by the usual rules of perturbation theory.

We note that the external fields acting on the system are allowed to create any state which may generate the different diagrams of perturbation theory. The corresponding matrix elements should be weighted with the amplitude of the component through which the final state is excited.

The above rules are also valid for those situations which cannot be treated in perturbation theory and where a full diagonalization is called for. Thus, *e.g.*, when the system displays a spurious state (see Sect. 2.5.3).

In what follows we discuss the energy of the $2p - 1h$ -like excitations, simplest modes which can display spuriousity. We distinguish between two types of states, namely

$$|n_i = 1; m, 1\rangle, \quad \begin{cases} \omega_1 = \epsilon' - V\Omega, & \Lambda_1 = -\sqrt{\Omega}V \\ \omega_i = \epsilon', & \Lambda_i = 0 \end{cases} \quad (i = 1; m = 1, 2, \dots, \Omega), \quad (2.5.27)$$

and

$$|m', 1; m', -1; m, 1\rangle, \quad \epsilon' \quad (m, m' = 1, 2, \dots, \Omega). \quad (2.5.28)$$

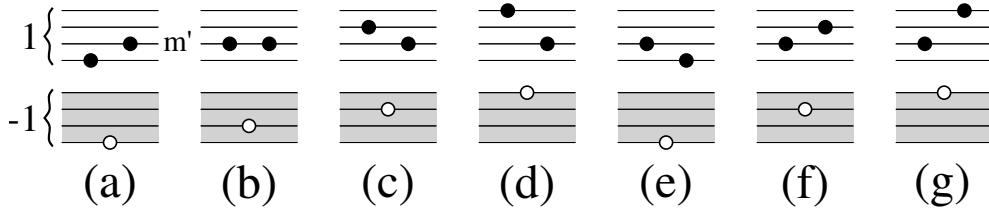


Figure 2.5.5: Schematic two-level model. Count of the states $|m, 1; m - 1; m', 1\rangle$ in the case of $j = 3/2$ and thus $\Omega = 2j + 1 = 4$. State (b) is not allowed because of Pauli principle. The states ((a),(e)), ((c),(f)) and ((d),(g)) are pairwise identical, in keeping with the indistinguishability of the particles. Thus, the states (a), (c) and (d) (equivalent (e), (f) (g)) exhaust the degrees of freedom of states of type (2.5.28). In other words, there are only $\Omega - 1 = 3$ two-particle one-hole states in which the odd particle is in the state $(m', 1)$. It is of notice that all substates m of levels $\sigma = \pm 1$ are degenerate. One displays them equally spaced only for didactical purposes.

The physical states are to be written as

$$|qm\rangle = \sum_i \xi_{iqm} |n_i = 1; m, 1\rangle, \quad (2.5.29)$$

in keeping with the fact that (2.5.28) cannot be basis states according to rule I), but only intermediate states. The quantities ξ_{iqm} are the amplitudes of the physical state in the different components of the product basis of elementary excitations. Rule I) eliminates most of the double counting of two-particle, one-hole states. The model space contains Ω “proper” states of the form $|n_i; m, 1\rangle$, in which case the odd particle is in the state $(m, 1)$. That is $|n_1; m, 1\rangle$ ($\omega_1 = \epsilon' - V\Omega$) and $|n_i; m, 1\rangle$ ($\omega_i = \epsilon'$, $i = 2, \dots, \Omega$). However, there are only $\Omega - 1$ two-particle, one-hole states in which the odd particle is in the state $(m, 1)$ (Fig. 2.5.5). Therefore, a spurious state remains in the spectrum based on elementary modes of excitation. In other words, allowing the quantum number m to run over all possible Ω -states, the model space contains Ω^2 states (one for each value of m), while the correct number is $\Omega(\Omega - 1)$. Thus the basis $|n_1 = 1; m, 1\rangle$ contains Ω spurious states. Its origin can be traced back to the violation of the Pauli principle.

To obtain the energy of $|qm\rangle$ we have to allow the states $|n_1 = 1; m, 1\rangle$ to interact through the vertices (2.5.16) and (2.5.26) and generate all the different perturbation theory diagrams (see rule II) except those containing bubbles (see rule III).

The different graphical contributions to the energy, calculated within the framework of the Brillouin-Wigner perturbation theory are displayed in Fig. 2.5.6. There is only one (diagonal) matrix element given by a single summation, which can be

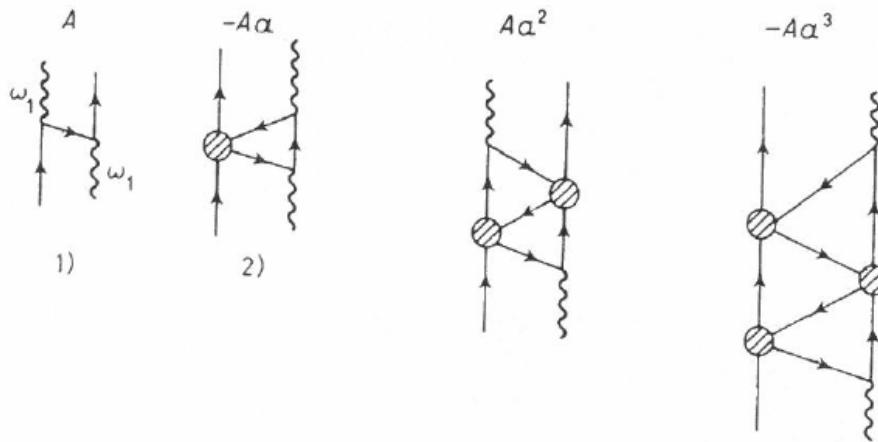


Figure 2.5.6: Contributions to the interaction of a fermion and a collective boson ω_i to order $1/\Omega^4$. The secular equation $E - E^{(0)} = A \sum_n a^n (-1)^n$ is given in terms of the quantities $A = 4\Omega V^2/(3\epsilon' - 2E)$ and $a = 2V/(3\epsilon' - 2E)$. The hatched circle stands for the four-point vertex (2.5.16) (see also Fig. 2.5.8 (I)).

carried to all orders in the interaction vertices⁸⁸, and can be written as

$$\begin{aligned} X_{ii'} &= A \sum_n (-1)^n a^n \delta(i, i') = \\ &= \frac{A}{1+a} \delta(i, i') \delta(n, 1) = -K(E) (\sqrt{\Omega} V)^2 \delta(i, i') \delta(i, 1), \end{aligned} \quad (2.5.30)$$

where a and A are defined in the caption to the figure and

$$K(E) = \left(\frac{3}{2}\epsilon' - E + V \right)^{-1}, \quad (2.5.31)$$

is the effective coupling strength. The associated secular equation

$$|(\omega_i - E)\delta(i, i') + X_{ii'}| = 0 \quad (2.5.32)$$

is equivalent to the dispersion relations ($i = 1$ and $i \neq 1$)

$$\frac{1}{K(E)} = \frac{(\sqrt{\Omega} V)^2}{\frac{1}{2}\epsilon' + \omega_i - E}. \quad (2.5.33)$$

Thus the energies of the system are determined by the equation

$$E = \frac{1}{2}\epsilon' + \omega_1 + \frac{\Omega V^2}{\frac{3}{2}\epsilon' - E + V}. \quad (2.5.34)$$

⁸⁸Concerning the proper expansion parameter see Eq. (2.5.85). See also Broglia (2020) App. A.

It admits the two solutions

$$E_{qm} = \begin{cases} \frac{3}{2}\epsilon', \\ \frac{1}{2}\epsilon' + \omega_1 + V = \frac{3}{2}\epsilon' - \Omega V + V, \end{cases} \quad (2.5.35)$$

which agree with the exact values⁸⁹.

Because $\Lambda_i = 0$ for $i \neq 1$, there is no summation in (2.5.29) and

$$|qm\rangle = N_{qm}|n_1 = 1; m, 1\rangle, \quad (2.5.36)$$

where

$$1 = N_{qm}^2 \left(1 - \frac{\partial X_{11}}{\partial E} \right) = N_{qm}^2 \left(1 - \frac{\Omega V^2}{\left(\frac{3}{2}\epsilon' - E + V \right)^2} \right). \quad (2.5.37)$$

For $E_{qm} = \frac{1}{2}\epsilon' + \omega_1 + V$ we obtain

$$N_{qm}^2 = \frac{\Omega}{\Omega - 1}, \quad (2.5.38)$$

while for $E_{qm} = \frac{3}{2}\epsilon'$ the state is non-normalizable as the quantity in parentheses in (2.5.37) is either negative ($\Omega > 1$) or zero ($\Omega = 1$). The state defined by

$$|q, m\rangle = \sqrt{\frac{\Omega}{\Omega - 1}} |n_1 = 1; m, 1\rangle, \quad (2.5.39)$$

and

$$E_{qm} = \frac{1}{2}\epsilon' + \omega_1 + V = \frac{3}{2}\epsilon' - V(\Omega - 1), \quad (2.5.40)$$

exhausts the inelastic sum rule in agreement with the exact results. Note that (2.5.39) is specifically excited in inelastic processes, as can be seen by direct inspection. The external inelastic field can act in two ways, exciting either a particle-hole pair or a phonon, with amplitudes

$$\langle m, 1; m', -1 | A_1^\dagger | 0 \rangle = \delta(m, m'), \quad (2.5.41)$$

and

$$\langle n_1 = 1 | A_1^\dagger | 0 \rangle = \sqrt{\Omega} \delta(i, 1), \quad (2.5.42)$$

respectively. The different graphical contributions to the inelastic-scattering process are displayed in Fig. 2.5.7, and can again be summed to all orders in the interaction vertices giving

$$\langle n_1 = 1; m, 1 | A_1^\dagger | m, 1 \rangle = \sqrt{\Omega} + \frac{\Lambda_1}{\frac{3}{2}\epsilon' - E_{qm} + V}. \quad (2.5.43)$$

⁸⁹The exact solutions can be obtained by noting that the operators A^\dagger, A and $\frac{1}{2}N_0$ are generators of the SU_2 group (see Bortignon et al. (1977)).

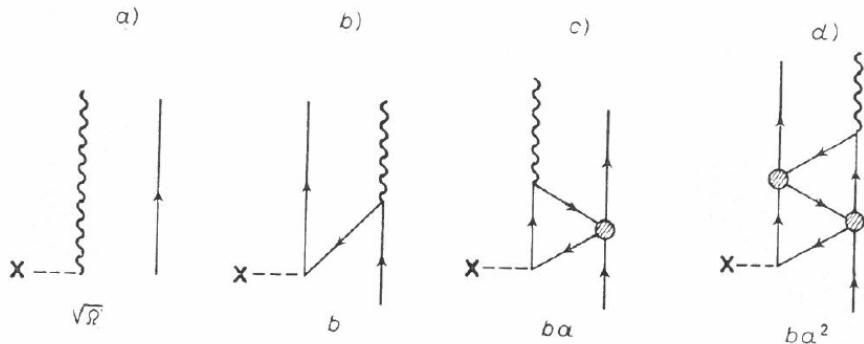


Figure 2.5.7: Graphical representation of the different terms contributing to the matrix element of the inelastic operator $\sqrt{\Omega} A^\dagger$ up to order $1/\Omega^3$. Note that the different contributions b), c), etc. have a one-to-one correspondence with the different contributions to E (see Fig. 2.5.6); $a = -2V/(3\epsilon' - 2E)$, $b = 2\Lambda_1/(3\epsilon' - 2E)$.

For $E_{qm} = \frac{3}{2}\epsilon'$ this quantity is equal to zero. Thus, the corresponding states do not carry any inelastic strength, a feature which is closely related to the fact that they cannot be normalized and that they do not display any correlation energy. On the other hand, the matrix element associated with (2.5.39) is

$$\langle qm|A^\dagger|m, 1\rangle = \sqrt{\frac{\Omega}{\Omega-1}} \frac{\Omega-1}{\sqrt{\Omega}} = \sqrt{\Omega-1}, \quad (2.5.44)$$

value which agrees with the exact answer.

The results (2.5.40) and (2.5.44) can be traced down to Pauli-principle corrections. In fact, the state $|n_i = 1; m, 1\rangle$ has a nonvanishing matrix element, implying a single particle-vibration coupling vertex, with the state $|m, 1; m, -1; m, 1\rangle$. This component, which is spurious, is removed by the different graphs displayed in Figs. 2.5.6 and 2.5.7. *The presence of the odd particle $(m, 1)$ blocks the particle-hole excitation $(m, 1; m, -1)$ which was present in the uncoupled system. Thus the system increases its energy by a quantity V . The reduction of the inelastic amplitude from $\sqrt{\Omega}$ to $\sqrt{\Omega-1}$ also indicates that there is one less particle-hole excitation responding to the external probe.*

2.5.3 Spurious states

While the model space product of elementary modes of excitation discussed in the last section contains Ω^2 states, only $\Omega(\Omega-1)$ are physically possible, the number of spurious states being Ω . On the other hand, the agreement between the exact and the nuclear-field-theoretical results shows that the effects of those spurious states are eliminated from all the matrix elements associated with physical observables.

In what follows we show that, in fact, the spurious states are isolated in an ex-

plicit way in nuclear field theory⁹⁰. Their energy coincides with the initial unperturbed energy, while all physical operators have zero off-diagonal matrix elements between any physical state and a spurious state, in particular the unit operator, which measures the overlap of the two types of states.

For this purpose we use again a schematic model consisting in a number, Ω , of single-particle levels in which particles interact by means of a “monopole” force,

$$H = H_{sp} + H_{int}, \quad (2.5.45)$$

where

$$H_{sp} = \frac{1}{2} \sum_{m=1}^{\Omega} \epsilon_m (a_{m,1}^\dagger a_{m,1} - a_{m,-1}^\dagger a_{m,-1}), \quad (2.5.46)$$

and

$$H_{int} = -VA^\dagger A, \quad (2.5.47)$$

with

$$A^\dagger = \sum_{m=1}^{\Omega} a_{m,1}^\dagger a_{m,-1}. \quad (2.5.48)$$

The energy of the i -th phonon is determined by the RPA dispersion relation (see rule IV))

$$\sum_{m=1}^{\Omega} \frac{1}{\epsilon_m - \omega_i} = \frac{1}{V}. \quad (2.5.49)$$

The eigenfunction corresponding to the different modes is

$$|n_i = 1\rangle = \sum_m \frac{\Lambda_i}{\epsilon_m - \omega_i} a_{m,1}^\dagger a_{m,-1} |0\rangle. \quad (2.5.50)$$

The particle-vibration coupling constant is given by⁹¹

$$\Lambda_i = -\langle n_i = 1 | H_{int} | m, 1; m', -1 \rangle = \left[\sum_m \frac{1}{(\epsilon_m - \omega_i)^2} \right]^{-\frac{1}{2}} \delta(m, m'), \quad (2.5.51)$$

where $|n_i = 1\rangle$ denotes a state containing one phonon, while $|m, 1; m, -1\rangle$ is the eigenstate associated with a particle-hole excitation.

⁹⁰Broglia et al. (1976)

⁹¹The normalization of $|n_i = 1\rangle$ implies $\langle n'_i = 1 | n_i = 1 \rangle = 1 = \sum_m \Lambda_i^2 / (\epsilon_m - \omega_i)^2$ and thus $\Lambda_i = [1/(\epsilon_m - \omega_i)^2]^{-1/2}$. In other words, the normalization is determined by the sum squared of the amplitudes, as the coupling matrix element does not depend on the energy. The HF solution can be obtained from the relation $[H, a_{m,1}^\dagger] |HF\rangle = \epsilon_{m,1}^{HF} |HF\rangle = \left(\frac{1}{2} \epsilon_m a_{m,1}^\dagger - VA^\dagger a_{m,-1}^\dagger \right) |HF\rangle = \frac{1}{2} \epsilon_m a_{m,1}^\dagger |HF\rangle$, as the state $(m, -1)$ is occupied in $|HF\rangle$. Thus $\epsilon_{m,1}^{HF} = \frac{1}{2} \epsilon_m$.

The other interaction to be included (rule II) is the four-point vertex which has the value

$$\langle m, 1; m', -1 | H_{int} | m'', 1; m''', -1 \rangle = -V \delta(m, m') \delta(m'', m'''). \quad (2.5.52)$$

The single-particle energies to be used in calculating the different graphs are $\frac{1}{2}\epsilon_m$, as the Hartree-Fock contribution (see rule IV)) of H_{int} is zero.

Similarly to H_{int} the “inelastic operator” has two different matrix elements, namely

$$\langle n_i = 1 | a_{m',1}^\dagger a_{m',-1} | 0 \rangle = \frac{\Lambda_i}{\epsilon_{m'} - \omega_i} \quad (2.5.53)$$

and

$$\langle m', 1; m'', -1 | a_{m,1}^\dagger a_{m,-1} | 0 \rangle = \delta(m, m') \delta(m', m''). \quad (2.5.54)$$

In what follows we discuss again the system comprising an odd particle, in the orbit $(m, 1)$, in addition to a single phonon excitation of the vacuum. According to rule I) initial and final states may involve both collective modes and particle modes, but not any particle configuration that can be replaced by a combination of collective modes. The exclusion of the states $|m, 1; m', 1; m', -1\rangle$ eliminates most of the double counting of two-particle, one-hole states. The Ω “proper” states of the form $|n_i = 1; m, 1\rangle$ are allowed. However, there are only $\Omega - 1$ (two-particle, one-hole) states in which the odd particle is in the state $(m, 1)$ (see Fig. 2.5.5). Therefore, a spurious state remains in the spectrum of the elementary modes of excitation. If we displace the zero-point energy of the odd system to $(1/2)\epsilon_m$, the unperturbed energy of the basis states $|n_i = 1; m, 1\rangle$ is ω_i .

The lower-order corrections to this energy which do not contain bubbles are drawn in Fig. 2.5.8 (I). Iterating these processes to infinite order we obtain the secular equation

$$|(\omega_i - E)\delta(i, i') + X_{ii'}(E)| = 0, \quad (2.5.55)$$

where

$$X_{ii'} = -\frac{\Lambda_i \Lambda_{i'}}{E - \epsilon_m - V}. \quad (2.5.56)$$

The different contributions calculated within the framework of the Brillouin–Wigner perturbation theory are energy dependent, and take into account renormalization effects of the states not explicitly included in the calculations. The dispersion relation fixing the energies E_{qm} of the physical states is

$$E - \epsilon_m - V = \sum_{i=1}^{\Omega} \frac{\Lambda_i^2}{\omega_i - E} = F(E). \quad (2.5.57)$$

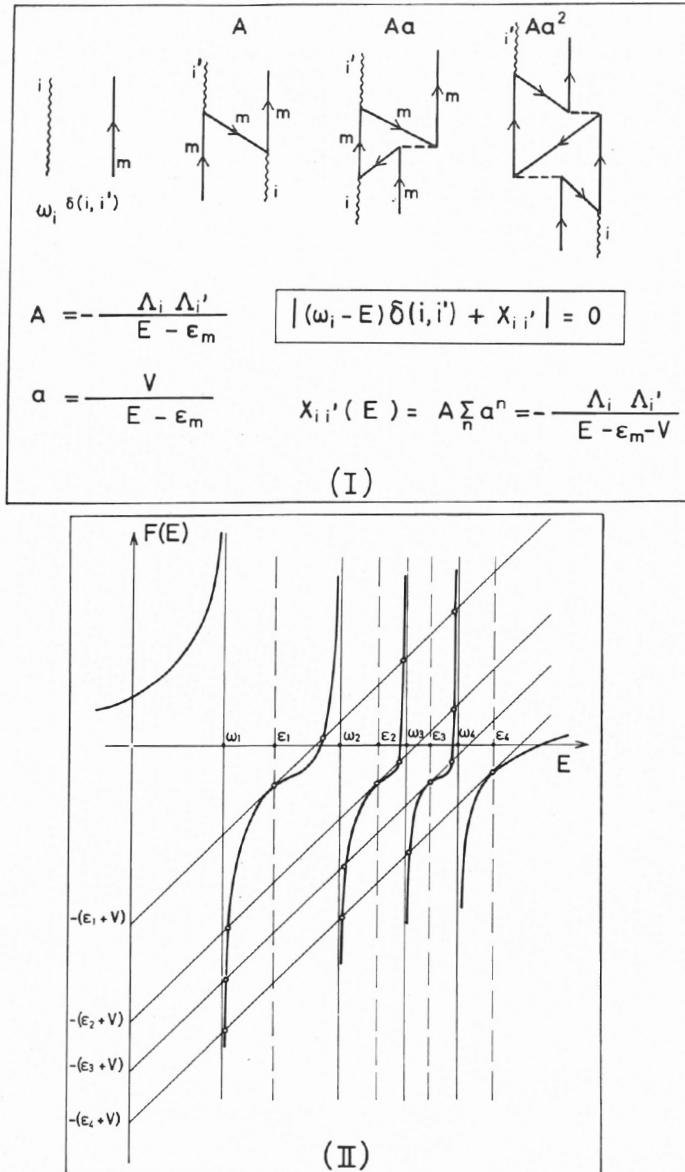


Figure 2.5.8: (I) Lower order contributions to the energy matrix element between the basis states $|n_i = 1; m, 1\rangle$. The dashed line stands for the model bare interaction (see Eq. (2.5.52)). The quantity $X_{ii'}(E) = A \sum_n a^n = -\Lambda_i \Lambda_{i'}/(E - \epsilon_m - V)$, where $A = -\Lambda_i \Lambda_{i'}/(E - \epsilon_m)$ and $a = V/(E - \epsilon_m)$, is the matrix element iterated to all orders in $1/\Omega$. The secular equation of the problem is $|(\omega_i - E)\delta(i, i') + X_{ii'}| = 0$, and is equivalent to the dispersion relation (2.5.57). (II) Graphical solution of the dispersion relation (2.5.57), for the case $\Omega = 4$. The function $F(E) = \sum_i \Lambda_i^2/(\omega_i - E)$ is displayed as a continuous thick line, while the parallel lines $E - \epsilon_m - V$ have been drawn as thin continuous lines intersecting the ordinates axis at $-(\epsilon_m + V)$. The intersections between the two functions give the eigenvalues of the secular equation. For each value of ϵ_m there are $\Omega + 1$ roots, the root at $E = \epsilon_m$ being double.

There is one equation for each single-particle level because the monopole force cannot change the m-state of the odd particle. The relation (2.5.57) can be solved graphically as shown in Fig. 2.5.8 (II). The energy $E = \epsilon_m$ is always a root of (2.5.57), in fact a double root since

$$\left[\frac{dF(E)}{dE} \right]_{E=\epsilon_m} = \sum_i \frac{\Lambda_i^2}{(\omega_i - \epsilon_m)^2} = 1, \quad (2.5.58)$$

and the line $E - \epsilon_m - V$ is at 45° . The remaining intersections of this line and the function $F(E)$ give rise to $\Omega - 1$ additional roots denoted by (qm) , whose energy E_{qm} agrees with the physical eigenvalues obtained from the exact solution of the model. The eigenvectors associated with the physical states (qm) are

$$|qm\rangle_F = \sum_i \xi_{iqm} |i; m, 1\rangle, \quad (2.5.59)$$

where

$$\xi_{iqm} = -N_{qm} \frac{\Lambda_i}{\omega_i - E_{qm}} = \langle i; m, 1 | qm \rangle_F. \quad (2.5.60)$$

The normalization condition which determines N_{qm} is⁹²

$$\begin{aligned} {}_F\langle qm | qm \rangle_F &= 1 = \sum_{i,i'} \left(\delta(i, i') - \frac{\partial X_{ii'}}{\partial E} \right) \xi_{iqm}^* \xi_{i'qm} = \\ &= N_{qm}^2 \left[\sum_i \frac{\Lambda_i^2}{(\omega_i - E_{qm})^2} - \frac{1}{(E_{qm} - \epsilon_m - V)^2} \sum_{i,i'} \frac{\Lambda_i^2 \Lambda_{i'}^2}{(\omega_i - E_{qm})(\omega_{i'} - E_{qm})} \right] = \\ &= N_{qm}^2 \left[\sum_i \frac{\Lambda_i^2}{(\omega_i - E_{qm})^2} - 1 \right], \end{aligned} \quad (2.5.61)$$

where the dispersion relation (2.5.57) has been utilized, and where $X_{ii'}$ is the matrix element appearing in (2.5.55) and defined in (2.5.56). For $E_{qm} = \epsilon_m$ the factor multiplying N_{qm}^2 is zero (see eq. (2.5.58)). Thus, there are only $\Omega - 1$ states which can be normalized when solving the Hamiltonian (2.5.45) within the framework of nuclear field theory. The full spuriosity of the elementary-mode product basis is concentrated in a single state⁹³.

⁹²This is in keeping with the fact that one is using Brillouin-Wigner perturbation theory (see e.g. App. A of Bès et al. (1976b)), where the differences between the exact energy E and the unperturbed energy replace the differences between unperturbed energies in the denominators of the Rayleigh-Schrödinger expansion. Furthermore, in the Brillouin-Wigner treatment the initial state never appears as an intermediate state, situations which are explicitly considered in the Rayleigh-Schrödinger treatment, and which in the Brillouin-Wigner one is taken into account through the use of the exact energy in the energy denominators. In the diagrams corresponding to transition matrix elements they are taken into account, in addition, by the normalization coefficients ξ .

⁹³Note that the mathematical relation $N^2 f(E) = 1, N^2$, being the norm of the state with energy E , implies that such state is spurious if $f(E) = 0$ or $f(E) < 0$ (see Eq. (2.5.37) and subsequent discussion).

The subscript F has been utilized above to indicate that we are dealing with the nuclear-field solution of the Hamiltonian (2.5.45) (for simplicity it will not be used in the following). Note that these eigenvectors are expressed in terms of only allowed initial or final states (see rule I)

$$|i; m, 1\rangle \equiv a_{m,1}^\dagger |i\rangle, \quad (2.5.62)$$

which are assumed to form an orthonormal basis, in particular in deriving the relation (2.5.61). *This is equivalent to the basic assumption of nuclear field theory of the independence of the different modes of excitation, i.e., in the present case,*

$$[\Gamma_i, a_{m,1}^\dagger] = 0. \quad (2.5.63)$$

Rules I–IV) discussed in the last section give the proper mathematical framework to this ansatz, which has played a basic role in developing a unified theory of nuclear structure. The above discussion can be illuminated by utilizing a conventional treatment of the residual interaction. Expanding the states $|n_i = 1; m, 1\rangle$ in terms of particle and hole states, we can write, with the help of (2.5.50),

$$a_{m,1}^\dagger |n_i = 1\rangle = a_{m,1}^\dagger \sum_{m' \neq m} \frac{\Lambda_i}{\epsilon' - \omega_i} a_{m',1}^\dagger a_{m',-1} |0\rangle. \quad (2.5.64)$$

The overlap between the states $|n_i = 1; m, 1\rangle$ is thus given by,

$$\begin{aligned} Z(i, i') &= \langle i' | a_{m,1}^\dagger a_{m,1}^\dagger | i \rangle \\ &= \sum_{m' \neq m} \frac{\Lambda_i \Lambda_{i'}}{(\epsilon_{m'} - \omega_i)(\epsilon_{m'} - \omega_{i'})} = \delta(i, i') - \frac{\Lambda_i \Lambda_{i'}}{(\epsilon_m - \omega_i)(\epsilon_m - \omega_{i'})}, \end{aligned} \quad (2.5.65)$$

where the orthogonality relation,

$$\sum_{m'} \frac{\Lambda_i \Lambda_{i'}}{(\epsilon_{m'} - \omega_i)(\epsilon_{m'} - \omega_{i'})} = \delta(i, i'), \quad (2.5.66)$$

of the RPA solutions in the even system has been utilized. Because of the non-orthogonality of the basis, the eigenvalues of the system are determined by the relation

$$|Z(E)(H - E)| = 0. \quad (2.5.67)$$

This is fulfilled for

$$|H - E| = 0, \quad (2.5.68)$$

which yields the $\Omega - 1$ physical roots, as well as for

$$|Z(E)| = 0. \quad (2.5.69)$$

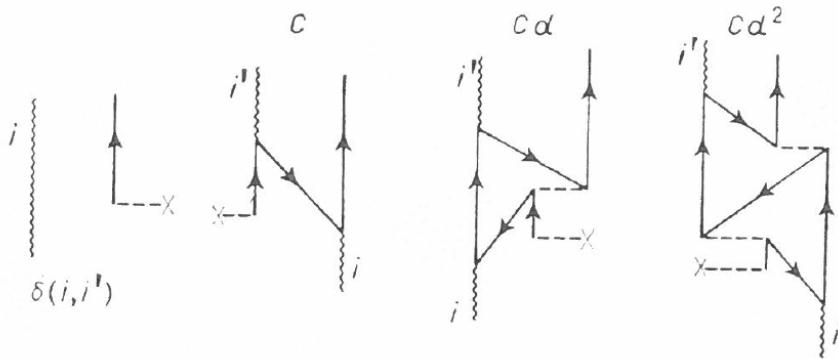


Figure 2.5.9: Lower order contributions to the one-particle transfer reaction induced by $a_{m,1}^\dagger$. The result of iterating the different contributions to all orders in $1/\Omega$ is equal to $T_{qm}(i, i') = C \sum_n d^n = -\Lambda_i \Lambda_{i'} / ((\omega_i - \epsilon_m)(E_{qm} - \epsilon_m - V))$, $C = -\Lambda_i \Lambda_{i'} / ((\omega_i - \epsilon_m)(E_{qm} - \epsilon_m))$, $d = V/(E_{qm} - \epsilon_m)$.

This solution corresponds to the spurious root $E_{qm} = \epsilon_m$. In fact⁹⁴,

$$\begin{aligned} \lim_{\delta \rightarrow 0} \sum_i \xi_{iqm}(E_{qm} = \epsilon_m + \delta) Z_{ii'} &= \lim_{\delta \rightarrow 0} N_{qm}(E_{qm} = \epsilon_m + \delta) \\ &\times \sum_i \frac{\Lambda_i}{\omega_i - (\epsilon_m + \delta)} \sum_{m \neq m'} \frac{\Lambda_i \Lambda_{i'}}{(\epsilon_{m'} - \omega_i)(\epsilon_{m'} - \omega_{i'})} = 0, \end{aligned} \quad (2.5.70)$$

since

$$\sum_i \frac{\Lambda_i^2}{(\epsilon_m - \omega_i)(\epsilon_{m'} - \omega_i)} = \delta(m, m'). \quad (2.5.71)$$

Note that this solution in terms of the overlap Z gives the exact answer in the present case, because of the simplicity of the model. In a general case which includes ground-state correlations this may not be true any longer.

We now calculate the one-particle stripping process leading to the odd system. This calculation illustrates the explicit concentration of the whole spuriousness into a single state which has zero correlation energy⁹⁵ and zero amplitude for the different physical processes exciting the $\Omega - 1$ physical states.

One has first to calculate the amplitude for the transition to a basis component ($n_i = 1; m, 1$) including only those graphs in which all intermediate states are excluded from appearing as initial or final states. This exclusion reflects the fact that

⁹⁴Within the context of renormalization, one first calculates the expressions for a finite value of δ and then takes the limit.

⁹⁵This is because the spurious state has zero phase space to correlate.

the diagonalization procedure has included all interaction effects that link these allowed states. The final amplitude for the transition to the state $|qm\rangle$ is obtained by summing the amplitudes ($n_i = 1; m, 1$) each weighted by ξ_{iqm} given in Eq. (2.5.60).

The lower-order contributions to the one-particle transfer amplitude between the state $|n_i = 1\rangle$ and the state $|qm\rangle$ are displayed in Fig. 2.5.9. They can be summed up to all orders of $1/\Omega$, the result being equal to (for the definition of T_{qm} see the corresponding caption)

$$\begin{aligned}
 & \langle qm | a_{m,1}^\dagger | n_i = 1 \rangle \\
 &= \sum_{i'} \xi_{i'qm} \left\{ \delta(i, i') - \frac{\Lambda_i \Lambda_{i'}}{(\omega_i - \epsilon_m)(E_{qm} - \epsilon_m)} \left[\frac{1}{1 - V/(E_{qm} - \epsilon_m)} \right] \right\} \\
 &= \sum_{i'} \xi_{i'qm} \{ \delta(i, i') - T_{qm}(i, i') \} \\
 &= -N_{qm} \left[\frac{\Lambda_i}{\omega_i - E_{qm}} - \frac{\Lambda_i}{(\omega_i - \epsilon_m)(E_{qm} - \epsilon_m - V)} \sum_{i'} \frac{\Lambda_{i'}^2}{\omega_{i'} - E_{qm}} \right] \\
 &= \frac{N_{qm}(E_{qm} - \epsilon_m)\Lambda_i}{(E_{qm} - \omega_i)(\omega_i - \epsilon_m)}. \tag{2.5.72}
 \end{aligned}$$

This quantity is zero for the spurious roots⁹⁶ (*i.e.* $E_{qm} = \epsilon_m$) and agrees with the exact result for the $\Omega - 1$ remaining physical roots.

Utilizing the relations

$$\frac{1}{V} = \sum_{m'} \frac{1}{\epsilon_{m'} - \omega_i}, \tag{2.5.73}$$

and

$$\frac{1}{V} = \sum_{m' \neq m} \frac{1}{\epsilon_{m'} - E_{qm}}, \tag{2.5.74}$$

we obtain

$$\sum_{m' \neq m} \frac{1}{(\epsilon_{m'} - E_{qm})(\epsilon_{m'} - \omega_i)} = \frac{1}{(E_{qm} - \omega_i)(\epsilon_m - \omega_i)}. \tag{2.5.75}$$

With the help of this relation we can derive the *one-particle transfer sum rule*. Note that (2.5.73) is the dispersion relation for the free phonon field. The second relation is, however, alien to the field theory results. Nevertheless, one can show that the solutions E_{qm} of (2.5.74) and of the nuclear-field-theory dispersion relation (2.5.57) are identical, except for the root $E_{qm} = \epsilon_m$. One can, therefore, utilize

⁹⁶Making use of (2.5.61) one obtains $\lim_{\delta \rightarrow 0} [(E_{qm} - \epsilon_m)N_{qm}]_{E_{qm}=\epsilon_m+\delta} = \lim_{\delta \rightarrow 0} \left\{ \sqrt{2} \delta^{3/2} / [\sum_i \frac{\Lambda_i}{\omega_i - \epsilon_m}]^{1/2} \right\} = 0$.

(2.5.74) as a mathematical relation without further justifications in the present context. One obtains

$$\begin{aligned} \sum_{qm} \left| \langle qm | a_{m,1}^\dagger | n_i = 1 \rangle \right|^2 &= \sum_{qm} \Lambda_{qm}^2 \Lambda_i^2 \sum_{m' \neq m} \frac{1}{(\epsilon_{m'} - E_{qm})(\epsilon_{m'} - \omega_i)} \\ &\quad \times \sum_{m'' \neq m} \frac{1}{(\epsilon_{m''} - E_{qm})(\epsilon_{m''} - \omega_i)}, \end{aligned} \quad (2.5.76)$$

where

$$\Lambda_{qm} = -N_{qm}(E_{qm} - \epsilon_m) = \left[\sum_{m' \neq m} \frac{1}{(\epsilon_{m'} - E_{qm})^2} \right]^{-\frac{1}{2}}. \quad (2.5.77)$$

Thus

$$\sum_{qm} \left| \langle qm | a_{m,1}^\dagger | n_i = 1 \rangle \right|^2 = \Lambda_i^2 \sum_{m' \neq m} \frac{1}{(\epsilon_{m'} - \omega_i)^2} = 1 - \frac{\Lambda_i^2}{(\epsilon_m - \omega_i)^2}, \quad (2.5.78)$$

where use has been made of the orthogonality relation

$$\sum_{qm} \frac{\Lambda_{qm}^2}{(\epsilon_{m'} - E_{qm})(\epsilon_{m''} - E_{qm})} = \delta(m', m''), \quad (m', m'' \neq m). \quad (2.5.79)$$

The result (2.5.78) coincides with the exact result. Physically it means that the single-particle orbital $(m, 1)$ is blocked by the amount $\Lambda_i^2/(\epsilon_m - \omega_i)^2$, which is the probability that the phonon $(n_i = 1)$ is in the particle-hole configuration $(m, 1; m, -1)$, *i.e.* with its particle in the orbital $(m, 1)$.

2.5.4 Applications

The nucleus $^{209}_{83}\text{Bi}_{126}$ can be viewed as a proton outside the double closed shell nucleus $^{208}_{82}\text{Pb}_{126}$. ^{209}Bi displays 12 states below 3 MeV. Within the framework of NFT, these states can be interpreted in terms of the elementary excitations of the ground state of ^{208}Pb , (gs) which plays the role of the “vacuum” state.

In what follows we discuss some aspects of this low-lying spectrum in terms of fermions (protons; a_j, a_j^\dagger), surface ($\Gamma^\dagger(\beta = 0\lambda)$) and pairing ($\Gamma^\dagger(\beta = 2\lambda)$) proton pair addition vibrational modes.

The unperturbed states of the closed-shell-plus-one-particle system can be written in terms of the free fields as

$$|n2\lambda, j; IM\rangle = [\Gamma_n^\dagger(2\lambda)a_j]_{IM}|0\rangle, \quad (2.5.80)$$

and

$$|n0\lambda, j; IM\rangle = [\Gamma_n^\dagger(0\lambda)a_j^\dagger]_{IM}|0\rangle. \quad (2.5.81)$$

This constitutes the basis set of states $\{\alpha_i\}$. All other states give rise to the complementary Hilbert space $\{a_i\}$.

The elementary modes of excitation interact through the particle–vibration and four-point vertices displayed in Fig. 2.5.10 giving rise to the matrix elements

$$M_1(nj, n'j') \equiv \langle [\Gamma_{n'}^\dagger(0\lambda)a_{j'}^\dagger]_{IM}|h_{eff}(E)|[\Gamma_n^\dagger(0\lambda)a_j^\dagger]_{IM} \rangle, \quad (2.5.82)$$

$$M_2(nj, n'j') \equiv \langle [\Gamma_{n'}^\dagger(2\lambda)a_{j'}]_{IM}|h_{eff}(E)|[\Gamma_n^\dagger(2\lambda)a_j]_{IM} \rangle, \quad (2.5.83)$$

and

$$M_3(nj, n'j') \equiv \langle [\Gamma_{n'}^\dagger(2\lambda)a_{j'}]_{IM}|h_{eff}(E)|[\Gamma_n^\dagger(0\lambda)a_j^\dagger]_{IM} \rangle. \quad (2.5.84)$$

They are to be calculated by utilizing the graphical techniques of perturbation theory and the rules discussed in Sect. 2.5.2. *There are two parameters on which to expand upon in carrying out a perturbative calculation. The first one is the strength of the interaction vertices measured in terms of the average distance between single-particle levels. The second is $1/\Omega$, where $\Omega = \sum_j(j + \frac{1}{2})$ is the effective degeneracy of the valence shells (in connection to this “standard” definition of Ω we refer to in the footnote⁸⁶ of this Chapter). These two parameters are in general connected through involved expressions. In the simple model discussed in Sect. 2.5.2, however, their relation is explicit and can be expressed as*

$$\epsilon = O(1), \quad \Lambda = O\left(\frac{1}{\sqrt{\Omega}}\right) \quad \text{and} \quad V = O\left(\frac{1}{\Omega}\right). \quad (2.5.85)$$

Another feature which determines the family of diagrams to select to a given order of perturbation is the number of internal lines which can be freely summed up. Each of these summations introduces a multiplicative factor Ω . Based on a wealth of detailed calculations for realistic distributions of levels one can conclude that relations (2.5.85) are also valid in such cases⁹⁷.

In what follows, we provide an example of a NFT calculation in a realistic situation, namely ²⁰⁹Bi. This nucleus has been investigated by means of high-resolution anelastic process⁹⁸. Through these experiments a septuplet of states around 2.6 MeV of excitation was identified, with spins ranging from $\frac{3}{2}^+$ to $\frac{15}{2}^+$.

⁹⁷ An alternative way to argue concerning the expansion parameter, in this case in connection with pairing vibrations, is to use the dimensionless quantity $x = 2G\Omega/D$ appropriate of a model made of two j -shells separated by an energy $D = 2\epsilon$ in which pairs of nucleons moving in time-reversal states interact through a pairing interaction of strength G . Phase transition takes place at $x \geq 1$, while situations away from phase transitions but still displaying consistent fluctuations typical of nuclei with low-lying collective vibrations, correspond to $x \approx 0.5$. Assuming $D(O(\epsilon)) = O(1)$, one obtains $G(O(V)) = O(1/\Omega)$ and naturally $\Lambda = O(1/\sqrt{\Omega})$, in keeping with the fact that the induced interaction can be written as $\Lambda^2/(\epsilon - \omega)$.

⁹⁸Ungrin et al. (1971); see also Bohr and Mottelson (1975).

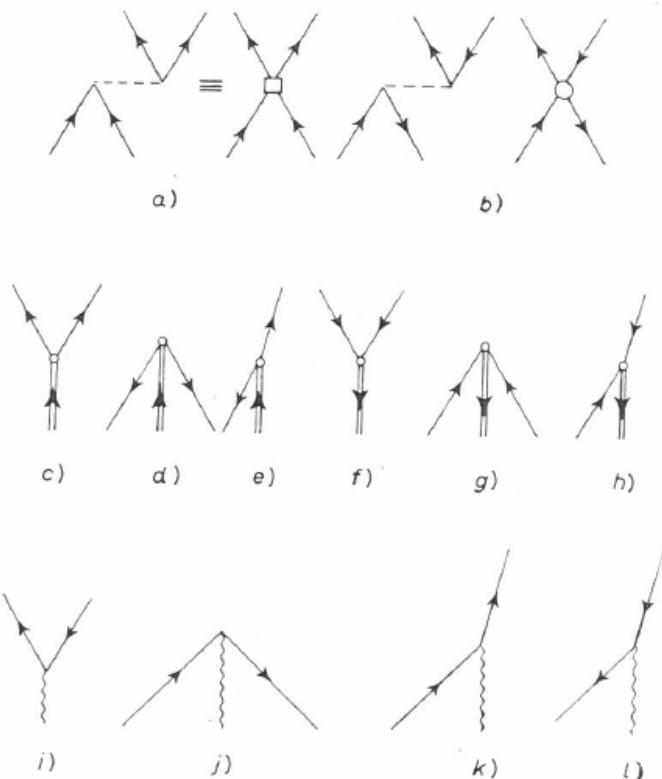


Figure 2.5.10: Interactions coupling the fermion fields with the pairing and surface vibrations. The different fermion and boson free fields are particles, holes, pairing vibrations ($\beta = \pm 2$) and surface vibrations ($\beta = 0$), β being the transfer quantum number. The two possible four-point vertices are given in a) and b). They correspond to the pairing and particle-hole model bare interactions. In graphs c)–h) all possible couplings between the fermion fields (arrowed lines) and the pairing vibrational fields (double lines arrowed) are displayed. Graphs i)–l) are all the possible coupling vertices between the surface vibrations (wavy line) and the fermion fields. Note that there is no direct coupling between the two boson fields, as the field theory we are dealing with is linear in the different (quasi) boson field coordinates.

In zeroth order these states can be interpreted in terms of a proton moving in the $h_{9/2}$ orbital coupled to the lowest octupole vibration of ^{208}Pb . The $\frac{3}{2}^+$ of this multiplet displays also a large parentage based on the proton pair addition mode of ^{208}Pb , and a proton hole moving in the $d_{3/2}$ orbital, as revealed by the (t, α) reaction⁹⁹ on ^{210}Po . The above results indicate that the low-lying (two-particle, one-hole) type of states of ^{209}Bi , are amenable to a simple description in term of the basis states¹⁰⁰

$$|j_1^{-1}, (\beta = 2), \lambda^\pi; IM\rangle \equiv |j_1^{-1} \otimes \lambda^\pi(^{210}\text{Po}); IM\rangle \quad (\lambda^\pi = 0^+, 2^+, 4^+) \quad (2.5.86)$$

and

$$|j_2, (\beta = 0), \lambda^\pi; IM\rangle \equiv |j_2 \otimes \lambda^\pi(^{208}\text{Pb}); IM\rangle \quad (\lambda^\pi = 3^-) \quad (2.5.87)$$

Only the lowest states of each spin and parity λ^π are included in the basis states, while all the RPA solutions are included in the intermediate states. The quadrupole surface vibrational modes were allowed only as intermediate states. The single hole and particle (proton) states j_1^{-1} and j_2 respectively, correspond to experimentally known levels around the $Z = 82$ ($N = 126$) shell closures. In what follows, the two $\frac{3}{2}^+$ states will be discussed.

The two basis states

$$|\alpha\rangle \equiv |d_{3/2}^{-1} \otimes \text{gs}(^{210}\text{Po}); 3/2^+\rangle, \quad (2.5.88)$$

and¹⁰¹

$$|\beta\rangle \equiv |h_{9/2} \otimes 3^-(^{208}\text{Pb}); 3/2^+\rangle, \quad (2.5.89)$$

are 118 keV apart. They mix strongly through the couplings depicted by the graphs a) and b) of Fig. 2.5.11.

Because of the energy dependence of h_{eff} (Eqs. (2.5.82)–(2.5.84)), there is a different matrix element for each final state. The diagonalization of the matrices was carried out self-consistently, *i.e.* the energy denominators of the different graphs are to be calculated by utilizing the exact energies¹⁰². The corresponding graphical contributions to the single-particle spectroscopic amplitudes, and thus eventually absolute (t, α) differential cross sections¹⁰³, and absolute inelastic scattering (d, d') cross-sections are also collected in Fig. 2.5.11. To be noted is the very different ratio of the (d, d') and (t, α) cross sections associated with the two

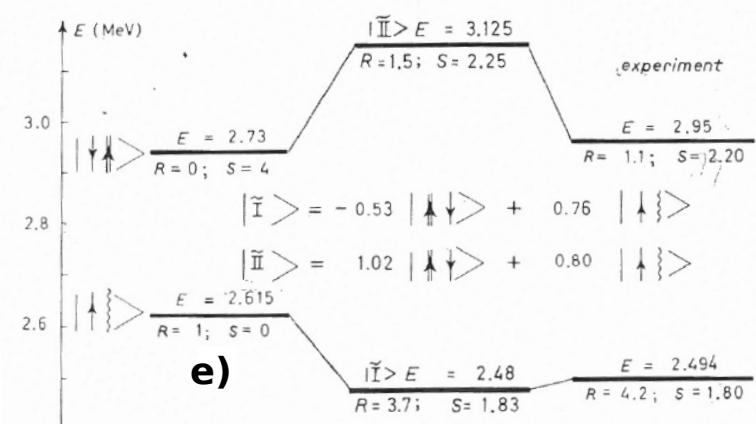
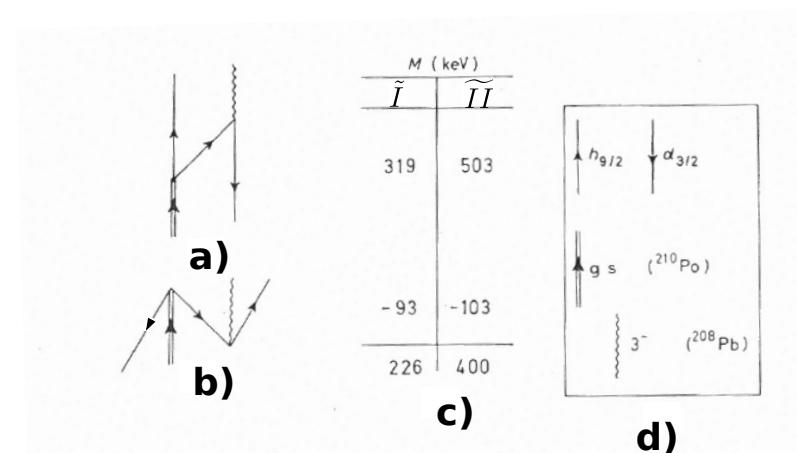
⁹⁹Regarding this pickup reaction, the connection between theory and experiment will be carried out in terms of spectroscopic factors.

¹⁰⁰Barnes et al. (1972).

¹⁰¹Although not likely, the reader is advised not to confuse the label of the state $|\beta\rangle$ with the transfer quantum number β used before.

¹⁰²for more details, see ref. Bortignon et al. (1977); see also Bortignon et al. (1976).

¹⁰³It is of notice that in the present case, and at variance with the rest of the monograph, use will be made, for the sake of didactics, of the concept of spectroscopic factors (see end of Sect. 3.1 as well as Sect. 5.8).



$$\left\{ \begin{array}{c} -0.53 \\ 1.02 \\ -0.103 \\ -0.103 \end{array} \right\} + \left\{ \begin{array}{c} 0.76 \\ 0.80 \\ 0.135 \\ 0.135 \end{array} \right\}^2 = \frac{2 \times 10^{-2}}{1 \times 10^{-5}}$$

f)

$$4 \times \left\{ \begin{array}{c} -0.53 \\ 1.02 \\ -0.103 \\ -0.103 \end{array} \right\}^2 = \frac{1.12}{4.16}$$

g)

$$4 \times \left\{ \begin{array}{c} -0.53 \\ 1.02 \\ -0.010 \\ -0.011 \end{array} \right\} + \left\{ \begin{array}{c} 0.76 \\ 0.80 \\ -0.211 \\ -0.333 \end{array} \right\}^2 = \frac{1.82}{2.27}$$

h)

$$\frac{1}{10} \left\{ \begin{array}{c} 0.76 \\ 0.80 \\ -0.577 \end{array} \right\}^2 = \frac{1.92 \times 10^{-2}}{2.13 \times 10^{-2}} e^2 b^3 \quad (3.3\%)$$

i)

$$\begin{aligned}
 & \frac{1}{10} \left\{ -0.53 \left[\begin{array}{c} \text{Diagram 1} \\ \text{Diagram 2} \end{array} \right] + 1.02 \left[\begin{array}{c} \text{Diagram 3} \\ \text{Diagram 4} \end{array} \right] \right\} \\
 & + 0.76 \left[\begin{array}{c} \text{Diagram 5} \\ \text{Diagram 6} \\ \text{Diagram 7} \\ \text{Diagram 8} \\ \text{Diagram 9} \\ \text{Diagram 10} \end{array} \right] + 0.80 \left[\begin{array}{c} \text{Diagram 11} \\ \text{Diagram 12} \\ \text{Diagram 13} \\ \text{Diagram 14} \\ \text{Diagram 15} \\ \text{Diagram 16} \end{array} \right] \\
 & = \left\{ \begin{array}{ll} 0.0216 & (3.7\%) \\ 0.0087 & (1.5\%) \end{array} \right. \quad \mathbf{j)
 \end{aligned}$$

Figure 2.5.11: In **a**, **b** and **c**) we give the two graphical contributions and the corresponding numerical values to the matrix element $M(E) = \langle d_{3/2}^{-1} \otimes gs(^{210}\text{Po}) | h_{eff}(E) | h_{9/2} \otimes 3^-(^{208}\text{Pb}); 3/2 \rangle$ in lowest order in $1/\Omega$. The resulting wave functions $|\tilde{I}\rangle$ and $|\tilde{II}\rangle$ are displayed in **e**) normalized according to (2.5.61). In **e**) we also give the unperturbed, and the renormalized theoretical energies of the levels. The (t, α) spectroscopic factor corresponding to the reaction $^{210}\text{Po}(t, \alpha)^{209}\text{Bi}$ is denoted by S , while

$$R = \frac{d\sigma(h_{9/2} \rightarrow J)}{d\sigma(gs(^{208}\text{Pb}) \rightarrow 3^-(^{208}\text{Pb}))}.$$

is the ratio of inelastic cross sections. In **d**) we display the free fields, while in **e**) we provide a summary of the results of the calculations in comparison with the data. The zeroth and order $1/\Omega$ contributions to the electromagnetic excitations are collected in **i**) and **j**). The value $0.577e^2b^3$ is the $B(E3; 0 \rightarrow 3)$ value associated with the 2.615 MeV state in ^{208}Pb . In **g**) and **h**) we give the zeroth and order $1/\Omega$ contributions to the spectroscopic factor associated with the $^{210}\text{Po}(t, \alpha)^{209}\text{Bi}$ reaction. Finally in **f**) we display the lowest contribution to the spectroscopic factor associated with the $^{208}\text{Pb}(^3\text{He}, d)$ reaction, which gives a measure of the ground state correlations of ^{208}Pb associated with the existence in its low-lying spectrum, of an octupole and a proton pair addition mode (see also Tables 2.5.1–2.5.3).

	$E_x(\text{MeV})$	$S(t, \alpha)(2j + 1)$	$S(^3\text{He}, d)$
$3/2^+$	2.49	$1.8 \pm 0.3 (4)$	< 0.01
$3/2^+$	2.95	$2.2 \pm 0.3 (4)$	< 0.01
$1/2^+$	2.43	$1.8 (2)$	< 0.02
$11/2^-$	3.69	$10 (12)$	< 0.05

Table 2.5.1: Observed single-particle strength (and corresponding errors) associated with the transfer reactions $^{210}\text{Po}(t, \alpha)^{209}\text{Bi}$ and $^{208}\text{Pb}(^3\text{He}, d)^{209}\text{Bi}$ (see Bortignon et al. (1977)).

states. While $R_1 = B(E3; (\frac{3}{2})_1)/B(E3; (\frac{3}{2})_2)$ is approximately equal to 2.5, the ratio $R_2 = \sigma((t, \alpha); (\frac{3}{2})_2)/\sigma((t, \alpha); (\frac{3}{2})_1)$ is close to one (see Tables 2.5.1–2.5.3).

Because the component $|\beta\rangle$ carries the inelastic-scattering strength, while the (t, α) reaction proceeds mainly through the component of type $|\alpha\rangle$, *the difference between R_1 and R_2 can be traced back to the corrections associated with the overcompleteness of the unperturbed basis states which give rise to rather different normalizations (see Sect. 2.5.3) of the two physical states $|\tilde{I}\rangle$ and $|\tilde{II}\rangle$ (see Fig. 2.5.11 (e)).*

Said it differently, in a conventional two-state shell model calculation implying a single matrix one would obtain

$$|\tilde{I}\rangle = A|\alpha\rangle + B|\beta\rangle \quad (2.5.90)$$

and

$$|\tilde{II}\rangle = -B|\alpha\rangle + A|\beta\rangle \quad (2.5.91)$$

with $A^2 + B^2 = 1$. The model would predict the value $(B/A)^2$ for both the ratio $R_1 = \sigma((d, d'); (3/2)_{\tilde{I}})/\sigma((d, d'); (3/2)_{\tilde{II}})$ and $R_2 = \sigma((t, \alpha); (3/2)_{\tilde{II}})/\sigma((t, \alpha); (3/2)_{\tilde{I}})$. The fact that $(R_1)_{th} = (0.0376/0.0156) = 2.41$ (against $(R_1)_{exp} = (0.042/0.011) = 3.8$) and $(R_2)_{th} = (2.25/1.83) = 1.22$ (against $(R_2)_{exp} = (2.2/1.8) = 1.22$), is a direct consequence of the overcompleteness of the basis which is taken care by NFT. This field theory provides a systematic mathematical procedure to deal with the spurious states. In the present case, associated with the overcompleteness of the basis ((2.5.88), (2.5.89)).

Within the framework of conventional shell model calculation ((2.5.90), (2.5.91)) the asymmetry between R_1 and R_2 can be related to the finite overlap¹⁰⁴ between the basis states $|\alpha\rangle$ and $|\beta\rangle$.

2.6 Competition between ($\beta = 0$) and ($\beta = \pm 2$) ZPF

Particle-hole like vibrations, as e.g. collective surface quadrupole vibrations, induce dynamical distortions of the mean field which virtually break the magnetic

¹⁰⁴Making use of the experimental values displayed in Tables 2.5.1 and 2.5.2, this overlap is estimated to be ≈ 0.25 .

	$E_x(\text{MeV})$	$\frac{\sigma(^{209}\text{Bi}(9/2^-;gs) \rightarrow ^{209}\text{Bi}(3/2^+;E))}{\sigma(^{208}\text{Pb}(gs) \rightarrow (3^-;2.615 \text{ MeV}))}$
3/2	2.49	0.042 ± 0.003
3/2	2.95	0.011 ± 0.002

Table 2.5.2: Observed total inelastic cross section σ^{oct} (and corresponding errors) associated with the lowest octupole vibrational state of ^{208}Pb can be written in terms of that associated with a single magnetic substate σ' as $\sigma_{3^-}^{oct} = 7\sigma'$. That associated with the multiplet $(h_{9/2} \otimes 3^-)_{J^+}(J = 3/2 - 15/2)$ as $\sigma_{3^-}^{oct} = 70\sigma'$, in keeping with the fact that the $h_{9/2}$ state has 10 magnetic substates. Thus, the strength associated with the 3/2 channel is $4/70 = 0.057$ to be compared with the observed summed (percentage) strength 0.053 ± 0.005 ($= (0.042 \pm 0.003) + (0.011 \pm 0.002)$) associated with the 2.45 MeV and the 2.95 MeV 3/2 $^+$ states.

	$E_n(\text{MeV})$		$\frac{\sigma(h_{9/2} \rightarrow 3/2^+)}{\sigma(0^+ \rightarrow 3^-)} (\%)$		$S(t, \alpha)$		$S(^3\text{He}, d)$	
	Theory	Exp	Theory	Exp	Theory	Exp	Theory	Exp
3/2	2.480	2.49	3.76	4.2 ± 0.3	1.83	1.8 ± 0.3	0.02	< 0.01
3/2	3.125	2.95	1.56	1.1 ± 0.2	2.25	2.2 ± 0.3	10^{-5}	< 0.01

Table 2.5.3: Summary of NFT predictions concerning the structure of the two lowest 3/2 $^+$ states of ^{209}Bi , in comparison with the experimental data.

degeneracy of levels into two-fold (Kramer's) degenerate (Nilsson-like) levels and to a reduction of the size of the discontinuity at the Fermi surface typical of non-interacting Fermi systems (see Fig. 2.2.1 (a)–(e)). Pairing vibrations does similarly, through dynamical ($U_j V_j$) weighting factors (see Fig. 4.4.2), which are operative in an energy region¹⁰⁵ $\epsilon_F \pm E_{corr}(\beta = \pm 2)$ (see Figs. 2.2.1 (f), (g) and 2.6.1 (g)). At the same time the dressing of nucleons by particle–hole and pairing vibrations, leads to an effective ω -mass making nucleons heavier, thus approaching the centroid of the valence orbitals lying above and in the Fermi sea towards the Fermi surface. In Fig. 2.6.1 a schematic representation of the effects the interweaving of single-particle motion and collective vibrations has on pairing correlations, is displayed (see also Fig. 5.8.1).

Within this context, one can mention that there is a remarkable similarity between the variation of V_ν^2 around the Fermi surface for the BCS ground state at $T = 0$ for strong coupled superconductors, and the normal-metal Fermi function at $T \gtrsim T_c$ (critical temperature) as a result of the renormalization of the single-electron levels due to the coupling to lattice phonons¹⁰⁶, something which is not unknown within the nuclear scenario for both closed and open shell nuclei (e.g. ^{208}Pb (see Fig. 1.2.5), Sn-isotopes (see Fig. 5.2.3, as well as Idini (2013) and Idini et al. (2015)). Thus, the change in the Fermi system, metal or nucleus, in

¹⁰⁵ E_{corr} is the correlation energy associated with the pair addition ($\beta = +2$) or with the pair removal modes ($\beta = -2$).

¹⁰⁶ Tinkham (1996).

going from the normal to the superconducting phase, cannot be solely described in terms of changes in the occupation numbers of single-particle states, as testified by the fact that through this mechanism no gap opens in the single-particle spectrum. Rather, the main mechanism is that of Off-Diagonal-Long-Range-Order (ODLRO)¹⁰⁷, that is the fact that $|BCS(\phi)\rangle_{\mathcal{K}'} = \prod_v (U'_v + e^{-2i\phi} V'_v P_v^\dagger |0\rangle)$ describes a system in which there is a gap for both single-pair translation and pair dissociation¹⁰⁸, as a result of the gauge phase coherence of all the Cooper pairs participating in the BCS condensate. Furthermore, and similarly important, are the consequences of such phasing in connection with the Cooper pair tunneling probability (Sects. 4.3, 4.7.1 and 4.12) between two weakly coupled superconductors (Josephson junction, Sect. 4.6), as well as between two superfluid nuclei in a heavy ion collision, eventually at energies below the Coulomb barrier (Sect. 7.3). Namely, the fact that successive tunneling is the dominant mechanism.

Zero point fluctuations induced by particle-hole like vibrations and by pairing modes compete with each other for phase space, through Pauli principle (see Fig. 2.6.2), thus eventually leading to a single ground state containing all of the dressed renormalized ZPF. The Pauli principle NFT diagrams shown in Figs. 2.6.2 (b) and (d) are at the basis of the stabilization of the ground state in general and of the competition between (as a rule quadrupole) deformations in 3D-space which breaks single-particle degeneracy (Nilsson potential), and in gauge space which thrives on large degeneracies¹⁰⁹. It is also the reason why single open shell nuclei are usually spherical. When tidal-like polarization effects in doubly open shell nuclei become overwhelming, the nucleus makes use of a Jahn-Teller mechanism to profit at best and simultaneously, of the quadrupole-quadrupole (alignment) and of the pairing (independent pair motion in Kramers degenerate levels) interactions. In other words, of potential energy (quadrupole deformation, localization) and of pairs of nucleons entangled with each other, over distances $\xi (\gg R_0)$ resulting in strongly overlapping entities and thus little sensitive to the orientation of the quadrupole deformed field (small moment of inertia, smaller than the rigid one), effect weakened in turn by confinement (mean field), and spatial quantization ($\mathcal{J} \approx 5\mathcal{J}_{irrot}$).

¹⁰⁷ See footnote 42 Ch. 1.

¹⁰⁸ See Eqs. (1.4.53) and (1.4.54) and following discussion.

¹⁰⁹ Within this context see Bayman (1961); Bès and Sorensen (1969); Mottelson (1962); Bohr and Mottelson (1975), and references therein.

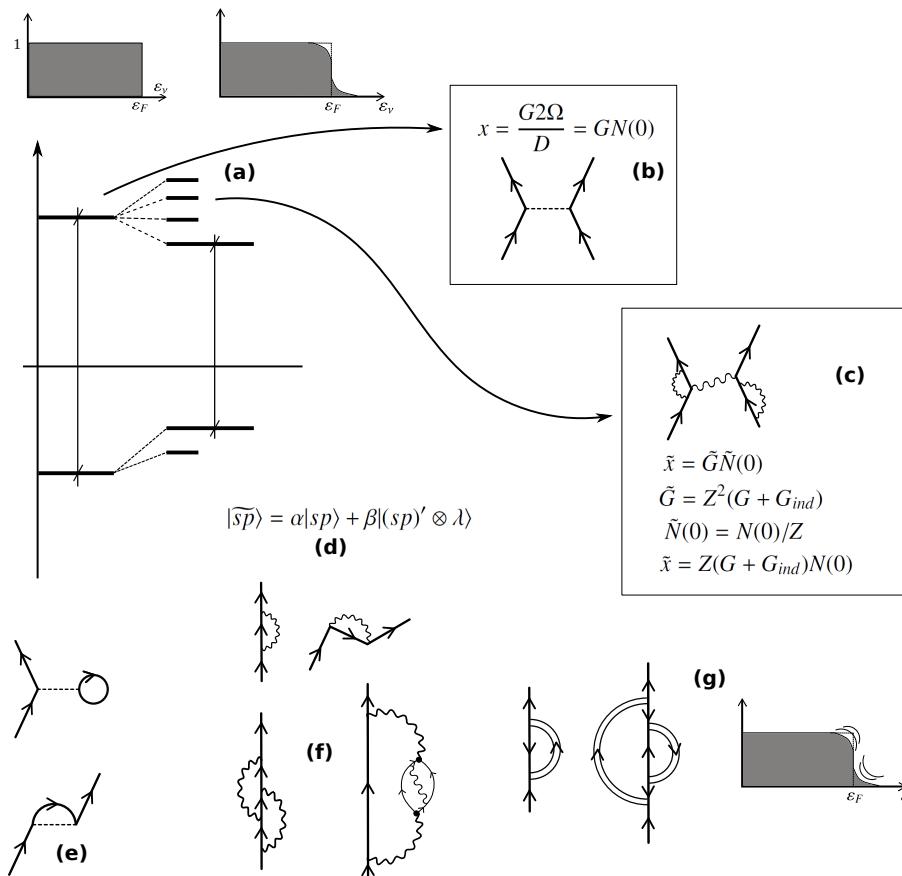


Figure 2.6.1: Schematic representation of some of the consequences the interweaving of the elementary modes of excitation with varied transfer quantum number ($\beta = 0, \pm 1, \pm 2$; (f), (g)) have in the (mainly single-particle) nuclear spectrum (a)–(e). In particular pair correlations (b, c), as measured by the (two-level) dimensionless parameter $x = G2\Omega/D = GN(0)$, product of the bare coupling constant G and the density of states (DOS) at the Fermi energy (ratio of the single-particle degeneracy $2\Omega = (2j+1)$, and the single-particle energy separation; see Högaasen-Feldman (1961); Broglia et al. (1968)). Coupling with surface modes (f) reduce the effective value of D leading to an increase of $N(0)$ as measured by Z ; see e.g. Barranco et al. (2005) and refs. therein). The eventual increase of x , as reflected by \tilde{x} , results from a delicate balance of the two effects eventually overwhelmed by the induced pairing interaction resulting from the exchange of collective ($\beta = 0$) vibrations between pairs of nucleons moving in time reversal states close to the Fermi energy, as compared with the increase in confinement kinetic energy reflected by the dynamical smoothing of the Fermi energy through the coupling of single-particle states to $\beta = 0$ and $\beta = \pm 2$ vibrations.

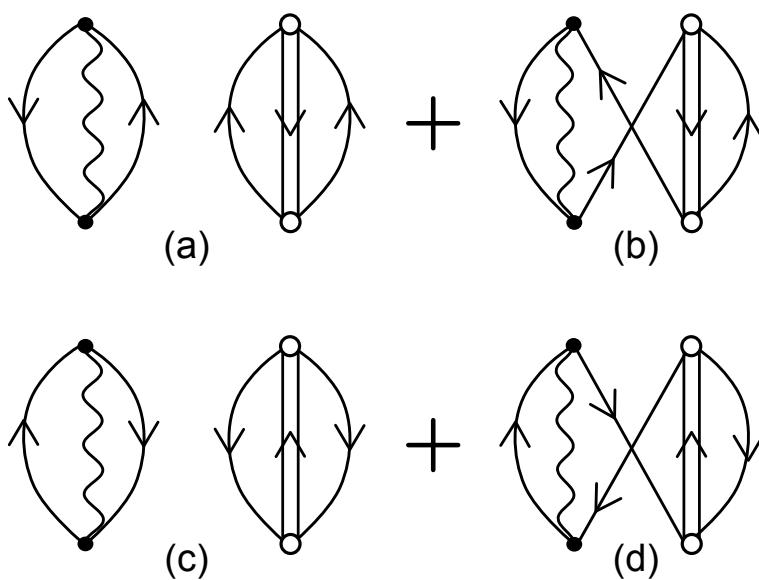


Figure 2.6.2: (a) (c) ZPF associated with $p-h$ and pairing vibrations (pair subtraction and pair addition modes) make use of the same nucleon degrees of freedom to simultaneously, and independently, correlate $p-h$, $p-p$ and $h-h$ excitations, thus violating Pauli principle. The NFT processes (b) and (d), which contribute to the correlation energy of the nucleus with opposite sign to that contributed by (a) and (c) (each unavoidable crossing of fermion lines contributes a minus sign), remove Pauli violating contributions to the corresponding order of perturbation in $1/\Omega$ (see Eq. (2.5.85) and related text).

2.7 Coupling between structure and reaction channels

In this Section the unification of NFT of structure and reactions is further developed using as example the light exotic two-neutron halo nucleus ^{11}Li . In particular, we dwell upon the variety of renormalization processes and associated form factors needed to calculate one- and two- neutron transfer reactions. The use of the same elements in the eventual calculation of the corresponding polarization contribution to the optical potential¹¹⁰, a subject not treated in the present monograph, is briefly mentioned.

2.7.1 Bare particles and Hartree-Fock field

Nucleon elastic scattering experiments at energies of tens of MeV can be accurately described in terms of an optical potential in which the real component is parametrized according to the (Woods–Saxon) potential¹¹¹,

$$U(r) = Uf(r), \quad (2.7.1)$$

$f(r)$ being a Fermi (sigmoidal) function, of radius $R_0 = r_0 A^{1/3}$ ($r_0 = 1.2$ fm), diffusivity $a = 0.65$ fm, and strength

$$U = U_0 + 0.4E \quad (2.7.2)$$

where

$$U_0 = V_0 + 30 \frac{N - Z}{A} \text{ MeV}, \quad V_0 \approx -51 \text{ MeV}, \quad (2.7.3)$$

while E is the energy of the scattered particle $\epsilon_k = \hbar^2 k^2 / 2m$, measured from the Fermi energy, m being the nucleon mass. One can replace the k -dependence in (2.7.2) by the so-called k -mass¹¹²

$$m_k = m \left(1 + \frac{m}{\hbar^2 k} \frac{dU}{dk} \right)^{-1}, \quad (2.7.4)$$

where the energy independent Woods-Saxon potential has a depth given by $\left(\frac{m}{m_k}\right) U_0 = U'_0$ ¹¹³. For the nucleons of the ^{11}Li core, namely ^9Li , $m_k = m(1 +$

¹¹⁰See footnote 23 Ch 1.

¹¹¹cf. e.g. Bohr and Mottelson (1969) and refs. therein.

¹¹²What in nuclear matter is called the k -mass and is a well defined quantity, in finite systems like the atomic nucleus, in which linear momentum is not a conserved quantity, is introduced to provide a measure of the spatial non-locality of the mean field, and is defined for each state as the expectation value of the quantity inside the parenthesis in Eq. (2.7.4), calculated making use of the corresponding single-particle wavefunction (see e.g. ref. Bernard and Giai (1981), in which case m_k is referred to as the non-locality effective mass)

¹¹³See e.g. Fig. 2.14 Mahaux et al. (1985).

$0.4)^{-1} \approx 0.7m$. For the halo neutrons¹¹⁴, $m_k/m = (1 + O \times 0.4)^{-1}$, where $O (= (R_0/R)^3)$ is the overlap between the core and the halo neutrons. Making use of the values $R_0 = 2.66$ fm and $R = 4.58$ fm (see Eq. (4.10.4)) one obtains $O \approx 0.2$ (see (3.6.8)) and thus, $m_k \approx 0.93 m$.

2.7.2 Physical particles

Within the framework of the above scenario, the full complexity of the many-body nuclear Hamiltonian

$$H = T + v, \quad (2.7.5)$$

has been reduced to

$$H_{HF} = T + U(r) + U_x(\mathbf{r}, \mathbf{r}'). \quad (2.7.6)$$

In other words, and making use of the expression,

$$H = T + v(|\mathbf{r} - \mathbf{r}'|) = H_{HF} + (v(|\mathbf{r} - \mathbf{r}'|) - (U(r) + U_x(|\mathbf{r} - \mathbf{r}'|))) \quad (2.7.7)$$

the full many-body nuclear Hamiltonian has been approximated by the Hartree-Fock Hamiltonian by neglecting the term in parenthesis. That is, by assuming that the sum of the direct and exchange potential gives a sensible approximation to the full two-body interaction v . This approximation (adding an appropriate spin-orbit potential), although providing a number of important insights into the nuclear structure and reactions as the sequence of single-particle levels¹¹⁵ (and associated magic numbers), and sensible nucleon-nucleus elastic phase shifts, disagrees with experiment on a number of points. In particular, leading to a too low level density of states around the Fermi energy and to infinite mean free paths, also for nucleons moving in states far removed from the Fermi energy, e.g. deep hole states.

To move further, one has to go beyond independent particle as well as potential scattering motion. That is, one has to allow the particles moving in both bound and continuum states to interact with each other and among themselves through four point vertices (see e.g. diagrams a) and b) of Fig. 2.5.10), and by coupling to $\beta = \pm 2$ pairing, and $\beta = 0$ surface mode, (diagrams c)–h) and i)–l) respectively). Let us exemplify the consequences of the interaction between elementary modes of excitation in the case of $^{11}\text{Li} + p$.

¹¹⁴ Assuming a velocity independent interaction v , the k -dependence of the mean field stems from the exchange (Fock) potential $U_x(\mathbf{r}, \mathbf{r}') = -\sum_i \varphi_i^*(\mathbf{r}') v(|\mathbf{r} - \mathbf{r}'|) \varphi_i(\mathbf{r})$ (linear in O), while the central potential is written as $U(r) = \sum_i \int d\mathbf{r}' |\varphi_i(\mathbf{r}')|^2 v(|\mathbf{r} - \mathbf{r}'|)$, (independent of O). It is of notice that the coupling between e.g. the low-lying quadrupole vibration of the core (^8He) and a halo neutron is also linear in O , i.e. $\langle H_c \rangle_{2^+((\text{core}), n(\text{halo}))} = \beta_2 \left\langle \frac{R_0}{\sqrt{5}} \frac{\partial U}{\partial r} \right\rangle O \langle j || Y^2 || 1/2 \rangle$, where $\langle j || Y^2 || 1/2 \rangle \approx 0.7$ ($j = 5/2, 3/2$), and $\left\langle R_0 \frac{\partial U}{\partial r} \right\rangle \approx 1.44 U_0 \approx -60$ MeV, see Brink and Broglia (2005) Eqs. (D20) p. 303 and (D26) p. 304. See also Fig. 2.7.1.

¹¹⁵ Mayer and Jensen (1955). It is of notice however the exceptions found in the case of light, halo nuclei, displaying parity inversion.

2.7.3 $^{11}_3\text{Li}_8$ structure in a nutshell

The sequence of single-particle levels for $^{10}_3\text{Li}_7$ associated with the mean field potential (2.7.3) implies that the distance between the last occupied neutron state $0p_{1/2}(\epsilon_{1/2^-} = -1.2 \text{ MeV})$ and the first empty one, $1s_{1/2}(\epsilon_{1/2^+} = 1.5 \text{ MeV})$ is 2.7 MeV (see Fig. 2.7.1, see also Fig. 1.1.2). In other words, within the framework of the Mayer and Jensen sequence of levels, in ^{10}Li the $0s_{1/2}, 0p_{3/2}$ neutron orbitals are fully occupied, while $0p_{1/2}$ carries one neutron, making ^{10}Li a neutron closed shell system ($N = 8$) with a hole in it¹¹⁶.

There is experimental evidence¹¹⁷ which testifies to the fact that the first unoccupied states of ^{10}Li are a virtual $1/2^+$ ($\epsilon_{1/2^+} \approx 0.2 \text{ MeV}$) and a resonant $1/2^-$ ($\epsilon_{1/2^-} \approx 0.5 \text{ MeV}$) state. According to NFT, this is a consequence of the self-energy renormalization of the bare $1/2^+$ state and of the $1/2^-$ state through a mainly PO (polarization) and CO (correlation) process respectively¹¹⁸ (Fig. 2.7.1). Thus parity inversion and the melting of the $N = 8$ closed shell in favor of the new magic number $N = 6$.

While ^{10}Li is not bound, ^{11}Li displays a two-neutron separation energy $S_{2n} \approx 400 \text{ keV}$. This value is the result of a subtle bootstrap mechanism. Being at threshold and basically not feeling a centrifugal barrier, the $1/2^+$ and $1/2^-$ states, are essentially not available for the short range bare NN -pairing interaction¹¹⁹, requiring a Cooper pair binding mechanism mediated by the exchange of long wavelength collective modes. This is the natural scenario of a very low-lying collective dipole mode for two main reasons. First, the presence of an eventual dipole particle-hole transition $1/2^+ \rightarrow 1/2^-$, with energy $\lesssim 1 \text{ MeV}$. The second one is related to the fact that the neutron halo in ^{11}Li can hardly sustain multipole surface vibrations, e.g. quadrupole vibrations, aside from displaying a very large radius as compared to that of the closed shell core $^9_3\text{Li}_6$, and thus a small overlap with it. This (halo) controlled phenomenon¹²⁰ has a two-fold hit: i) to screen the bare $NN^{-1}S_0$ short

¹¹⁶It is of notice that in the characterization of nuclear single-particle levels in terms of number of nodes n , orbital angular momentum l and total angular momentum $j(l, 1/2)$ (i.e. (nlj)) while univoque for l and j , allows for the flexibility whether to include or less the node at the origin. In the first case $n = 0$ is the first level of a given (l, j) , in the second $n = 1$.

¹¹⁷See footnote 30 of Ch. 5. See also Sect. 5.2.4.

¹¹⁸Barranco et al. (2001, 2020), CO (involving ground state correlation vertices), PO (particle moving around closed shells and polarizing the core).

¹¹⁹Bennaceur et al. (2000); Hamamoto and Mottelson (2003, 2004).

¹²⁰Within this context one is essentially forced to make a subtle extension of the statement according to which the single-particle motion is the most collective of all nuclear motions (Mottelson (1962)), emerging from the same properties of the nuclear interaction (both bare and induced) as collective motion does, and in turn at the basis of the detailed properties of each collective mode, acting as scaffolds and filters of the variety of embodiments. In fact, one has to add the characterisation of “physical” to “single-particle motion” (i.e. dressed) to englobe in the above statement also the present situation (parity inversion). In other words, while the bare $s_{1/2}$ and $p_{1/2}$ orbitals could never lead to a low-lying dipole strength, the corresponding clothed, physical states do so in a straightforward manner, the dressing effects being in this case of the same order of magnitude of the mean field effects. Consequently: “physical, dressed single-particle motion, is one of the most collective of nuclear motions”, seems to be the right statement. Within this context, one can point

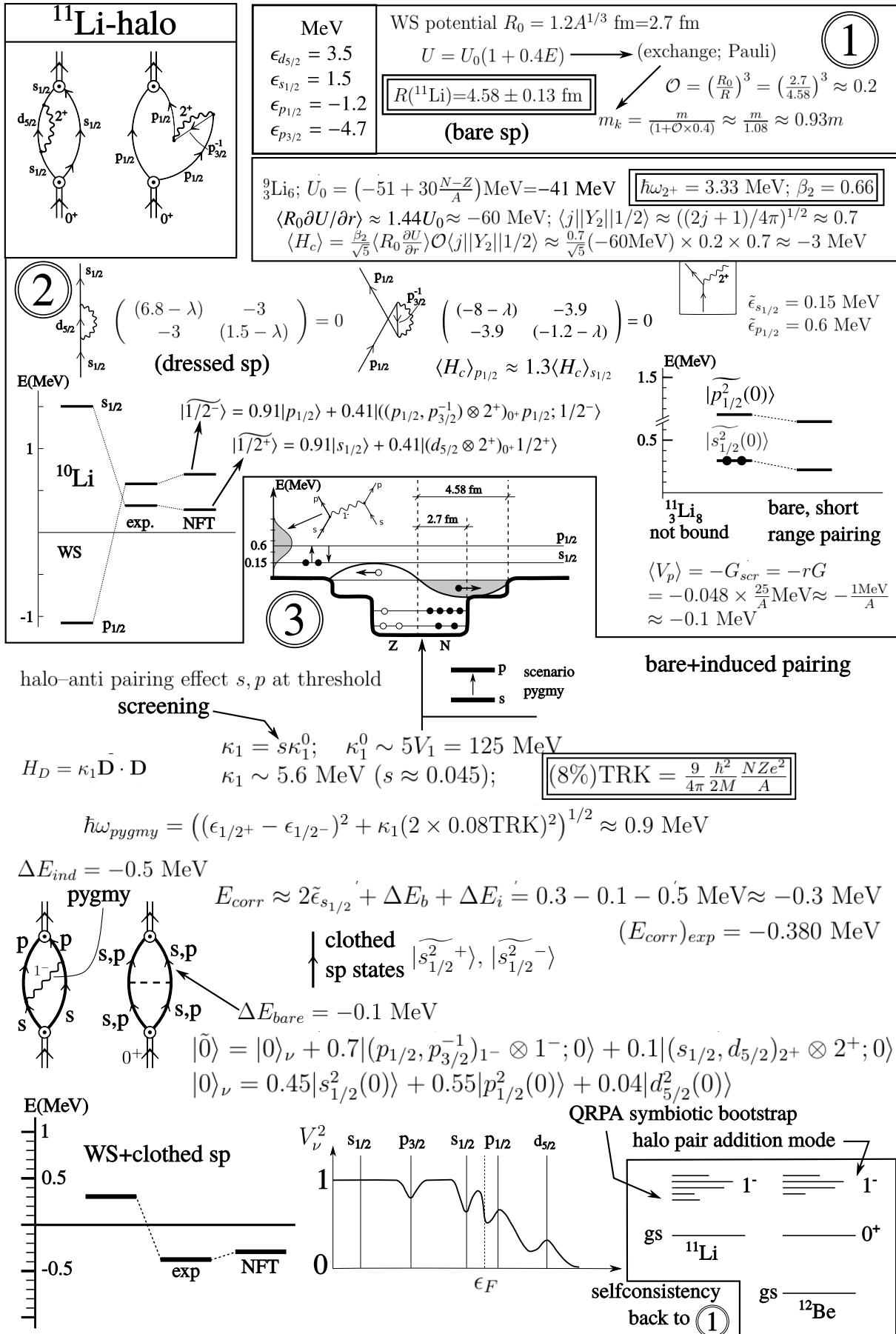


Figure 2.7.1

range pairing interaction making it subcritical, ii) to screen the (repulsive) symmetry interaction, and (consequently), to allow for the presence, at low energies, of a consistent fraction of the TRK-sum rule.

In fact, a 1^- resonance carrying $\approx 8\%$ of the TRK sum rule and with energy centroid $\lesssim 1$ MeV have been observed¹²¹. Thus, the connotation of pygmy dipole resonance (PDR). Exchanged between the heavily dressed neutrons moving in parity inverted states it provides, together with the contribution of the strongly screened, bare NN -pairing interaction, the glue needed to bind the neutron halo Cooper pair to the ${}^9\text{Li}$ core. With some amount of experimental input¹²², NFT allows to propagate dressing effects which not only renormalizes mean field, but overwhelms it, resulting in an overall account of the experimental findings. The discussion of the many-body effects at the basis of these phenomena is carried out in three steps, as schematically displayed in Fig. 2.7.1 and described below¹²³ (for details see Sects. 3.6 and 5.2.4):

1) Starting with well defined elements: Woods-Saxon (WS) potential, and the parameters characterizing the low-lying quadrupole vibration of the core ${}^9\text{Li}$ (**input**, double boxed quantities), calculate the single-particle levels and collective vibration (separable interaction) and determine the corresponding particle-vibration coupling vertices (strength and form factors). From the ratio of the WS radius (R_0) and of the observed one ($R({}^{11}\text{Li})$ **input**) determine the overlap O . Because $O \ll 1$, the contribution of the exchange (Fock) potential to the empirical WS potential is small concerning the halo neutrons¹²⁴. Consequently the neutron halo k -mass m_k has a value close to the observed neutron mass m ;

to a tumbling-like chain of incipient phase transitions: parity-inversion \rightarrow dipole (dynamic) distortion \rightarrow pygmy dipole resonance \rightarrow (single) BCS Cooper pair formation, similar to that observed in the phase transition between normal-superconducting phases in metals, spontaneously breaking of translational invariance (crystal) \rightarrow phonons \rightarrow Cooper pairs \rightarrow BCS condensation (Nambu (1991), see also Broglia et al. (2019a)).

¹²¹Zinser et al. (1997), Nakamura et al. (2006), Shimoura et al. (1995), Ieki et al. (1993), Sackett et al. (1993), Kobayashi et al. (1989); Kanungo et al. (2015); Aumann (2019).

¹²²See e.g. the corresponding discussion in Barranco et al. (2017).

¹²³The information contained in this figure (see also Broglia et al. (2019b)), finds its origin going back to predictions made in 2001 (Barranco et al. (2001); see Fig. 3.6.3). Results from a *tour de force*-like experiment carried out at TRIUMF (Vancouver) by Tanihata et al. (2008) (see also Tanihata et al. (2013) and Fig. 7.1.2) provided evidence concerning the predicted phonon mediated pairing in nuclei (see Potel et al. (2010)). Figure 2.7.1 displays an analytic synthesis of a nuclear field theory solution –(NFT) presented and discussed at the pedestrian level in Sect. 2.5– of the structure of one- and two- neutron halo nuclei lying at the dripline, namely ${}^{10}\text{Li}$ (Barranco et al. (2020)) and ${}^{11}\text{Li}$ (Barranco et al. (2001)). Thus, of the unified interpretation of the results of almost two decades of experimental research carried out and being carried out at major laboratories around the world (TRIUMF (Vancouver; Tanihata et al. (2008)), ISOLDE (CERN; Jeppesen et al. (2006)), TRIUMF (Cavallaro et al. (2017)), etc.). Figs. 2.7.1, 2.7.2 and 2.7.3 condense also a number of facets about pygmy resonances in nuclei and of quantal dipole fluctuations mediated pairing, an intermediate boson also found at the basis of the van der Waals interaction (see Broglia et al. (2019a)). Within this scenario, see Section 3.6 *Halo pair addition mode and pygmy*, as well as the parallel carried out in connection with the role the van der Waals interaction plays in molecular systems in general, and proteins in particular (see Sect. 3.7 and App. 3.B).

¹²⁴See footnote 114 of the present Chapter.

2) Making use of the above elements one can dress the bare single-particle states, in particular the $s_{1/2}$ and $p_{1/2}$ states. Parity inversion ensues, with $1/2^+$ and $1/2^-$ at threshold. As a consequence the $N = 8$ shell closure melts away, $N = 6$ becoming a new magic number, testifying to the fact that large amplitude fluctuations can, in nuclei, be as important as static mean field effects. As a result ${}^3_3\text{Li}_7$ is not bound. Adding one more neutron and switching on the bare pairing interaction (e.g. a contact force $V(r_{12}) = -4\pi V_0 \delta(\mathbf{r}_1 - \mathbf{r}_2)$ with constant matrix element¹²⁵ $G = 1.2 \text{ fm}^{-3} V_0/A \approx (25/A) \text{ MeV}$), the screening $r = \frac{(M_j)\text{halo}}{(M_j)\text{core}} \approx \frac{2}{2j+1} \left(\frac{R_0}{R}\right)^3 \approx 0.048$ (see Eq. (3.6.3)) resulting from the poor overlap between halo and core neutrons leads to a value of the strength of the pairing interaction ($G_{scr} = r \times G$) which is subcritical, and thus to an unbound system. In fact, $G_{scr} = 0.048 \times 25/A \text{ MeV} \approx 0.1 \text{ MeV}$ and $\Delta E = 2\tilde{\epsilon}_{s_{1/2}} - G_{scr} \approx 0.3 \text{ MeV} - 0.1 \text{ MeV} \approx 0.2 \text{ MeV}$. Summing up, $G_{scr} < G_c$, G_c being the minimum value of the pairing coupling constant leading to a bound state¹²⁶;

3) The sloshing back and forth of the halo neutrons against the core nucleons moving in phase, leads to a pygmy dipole mode which feels a strongly screened repulsive symmetry potential. In keeping with the fact that the strength of the bare dipole-dipole interaction $\kappa_1^0 \sim 1/R^2({}^{11}\text{Li})$, the screening factor can be estimated as $((R({}^{11}\text{Li})/\xi)^2(R = 4.58 \text{ fm}, \xi = 20 \text{ fm} \text{ (correlation length of the halo Cooper pair), } s \approx 0.052)$, the value obtained in Sect. 3.6 being $s = 0.043$. In other words, while it takes a quantity proportional to $5V_1 = 125 \text{ MeV}$ to separate protons from neutrons in the core, this value is reduced to $s \times 5V_1 = 5.4 \text{ MeV}$ ($(V_1)_{scr} = sV_1 \approx 1 \text{ MeV}$) for halo neutrons. This, as well as the small energy difference between the parity inverted levels, is at the basis of the fact that $\approx 8\%$ of the Thomas-Reiche-Kuhn

¹²⁵Brink and Broglia (2005), pp 40-42.

¹²⁶At the basis of superconductivity one finds the result obtained by Cooper (1956). He worked out the problem of two electrons interacting through an attractive interaction above a quiescent Fermi sea. Thus, all but two of the electrons are assumed to be noninteracting. The background of electrons enter the total problem only through the Pauli principle by blocking states below the Fermi surface from participating in the remaining two-particle problem. If one measures the kinetic energy ϵ_k relative to its value at the Fermi surface only states with $\epsilon_k > 0$ are available to the interacting pair of electrons. Cooper found that a bound state of the pair always exists for arbitrarily weak coupling as long as the potential is attractive near the Fermi surface, a mechanism which implies the instability of the normal phase. Cooper pair binding is a rather remarkable result for the usual two-body problem. If one has only two particles coupled by an attractive interaction, they would not form bound states unless the attractive interaction exceeds a certain critical value. In fact, for a two-body system with an attractive interaction, a bound state exists for any value of the strength, provided the density of levels $N(E)$ ($E = |\epsilon - \epsilon_F|$) is different from zero around the Fermi energy ($N(0) \neq 0$). This is so for one- and two-dimensional systems, in which case $N(E) \sim E^{-1/2}$ and E^0 respectively, while for 3D-systems $N(E) \sim E^{1/2}$ (Fermi gas model). Cooper model reduces the 3-dimensional to a lower-dimensional (essentially 2D) quantal system (Gor'kov (2012), see also Cohen-Tannoudji and Guéry-Odelin (2011)). In ${}^{11}\text{Li}$ the last two neutrons are very weakly bound. Consequently they move away from the nucleus core ${}^9\text{Li}$, lowering in the process their relative momentum and forming a misty cloud or halo (Austin and Bertsch (1995)). One can thus view this system as the nuclear embodiment of a Cooper pair. The question then arises, why is there a critical value for the pairing interaction? The answer is likely related to spatial quantization and a not quiescent Fermi sea, features associated with the fact that the nucleus is a finite quantal many-body system.

sum rule (**input**) is found at low energy ($\lesssim 1$ MeV). Another way to say the same thing is that $(V_1)_{scr} = sV_1$ is at the basis of the fact that the $s_{1/2} - p_{1/2}$ energy difference ($\Delta\tilde{\epsilon} \approx 0.45$ MeV) is only increased by a modest value ($\hbar\omega_{pygmy} \lesssim 1$ MeV $\approx 10^{21}$ Hz), and that at the same time, the $E1$ -strength remains essentially unchanged, not being shifted to higher energies by the GDR. Typical values of the dipole strength in the case of nuclei lying along the stability valley being $\approx 10^{-4}B_W(E1)$ for low-energy single-particle transitions, while in the present case one finds a value close to $1B_W(E1)$.

The two halo neutrons dressed by the vibrations of the core (heavy arrowed lines in lowest left corner of Fig. 2.7.1) and interacting through the bare NN -pairing force are not bound. Consequently, the pygmy resonance will fade away almost as soon as it is generated (essentially lasting the neutron transversal time $\approx 10^{-21}$ s in e.g. a ${}^9\text{Li}(t, p)$ reaction). This, unless it is exchanged between the two neutrons forcing them to jump from the configuration $s_{1/2}^2(0)$ at threshold ($2 \times \tilde{\epsilon}_{s_{1/2}} \approx 0.3$ MeV) into $p_{1/2}^2(0)$, also close to threshold ($2 \times \tilde{\epsilon}_{p_{1/2}} \approx 1.2$ MeV). In other words, one finds the dipole pygmy resonance acting as an intermediate boson which couples to the halo neutrons with a strength $\Lambda \approx 0.5$ MeV (QRPA calculation). Thus, the corresponding $|gs({}^{11}\text{Li})\rangle$ correlation energy $E_{corr} \approx -0.3$ MeV is mainly due to the dipole pygmy exchange process.

The resulting symbiotic halo pair addition mode of ${}^{11}\text{Li}$ can, in principle, be used as a building block of the nuclear spectrum, amenable to be moved around. A possible candidate for such a role is the first excited 0^+ state of ${}^{12}\text{Be}$, together with the associated dipole state built on it and eventually other fragments of the associated $E1$ -low energy-strength.

To calculate the pygmy dipole resonance (PDR) of ${}^{11}\text{Li}$ one needs to know the ground state of this nucleus (halo-pair addition mode) so as to be able to determine microscopically the occupation factors of the $1s_{1/2}, 1p_{3/2}, \epsilon s_{1/2}, \epsilon p_{1/2}, \epsilon d_{5/2}, \dots$, etc. states. This is the basic input needed to work out the corresponding QRPA equations, whose diagonalization provide the energy, transition density and $E1$ -transition probability of the mode. But to do so one is required to know the PDR. Arrived to this point, the protocol schematically presented in Fig. 2.7.1 requires to go back to 1 and repeat the whole procedure until convergence is achieved.

2.7.4 ${}^{11}\text{Li}(p, p){}^{11}\text{Li}$ and transfer reaction channels

Because NFT rules have, a priori, no limitations concerning whether the excitations studied lie or not in the continuum, or whether the single-particle motion displays asymptotic waves, it allows for a unified description of structure (s) and reactions (r) ($NFT_{(s+r)}$). An example of the above statement is provided by Fig. 2.7.2. Graph (a) describes one of the processes contributing to the elastic reaction ${}^{11}\text{Li}(p, p){}^{11}\text{Li}$ as the system propagates in time. Namely, a polarization contribution to the global (mean field) optical potential describing proton elastic scattering off ${}^{11}\text{Li}$.

In what follows we describe the processes taking place in the interval of time

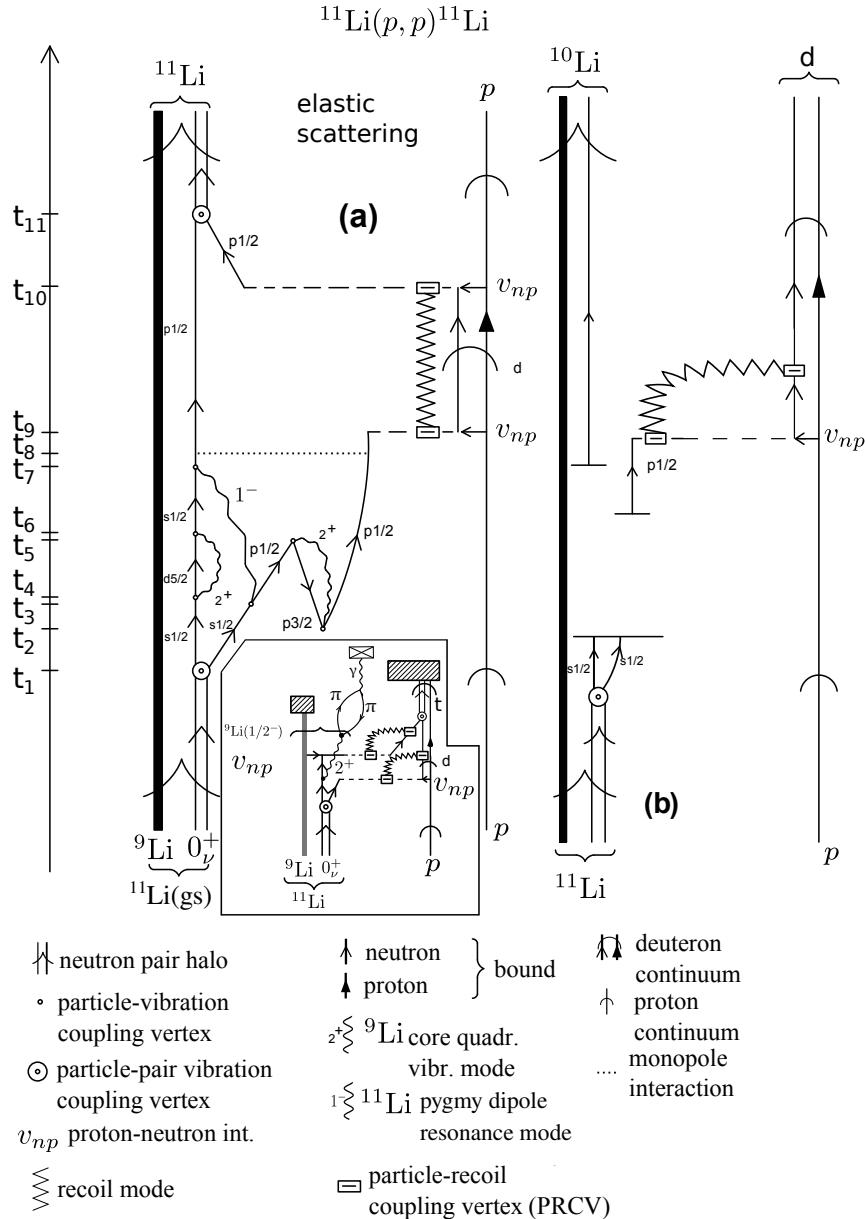


Figure 2.7.2: (a) NFT-diagram describing one of the processes contributing to the elastic reaction $^{11}\text{Li}(p, p)^{11}\text{Li}$ as the system propagates in time (polarization contribution to the global (mean field) optical potential). In the inset, a schematic NFT diagram describing the process $^{11}\text{Li}(p, t)^9\text{Li}(1/2^-)$ is displayed. A dashed open rectangle indicates the particle-recoil coupling vertex. A crossed box represents a γ -detector, while hatched rectangles stand for particle detectors. (b) Schematic NFT diagram describing the reaction $^{11}\text{Li}(p, d)^{10}\text{Li}$, i.e. same as in (a) up to time t_8 (reason for which no details are repeated between t_2 and t_8). From there on the deuteron continues to propagate to the detector bringing to it, aside from nuclear structure information, the information resulting from its interaction with the recoil mode.

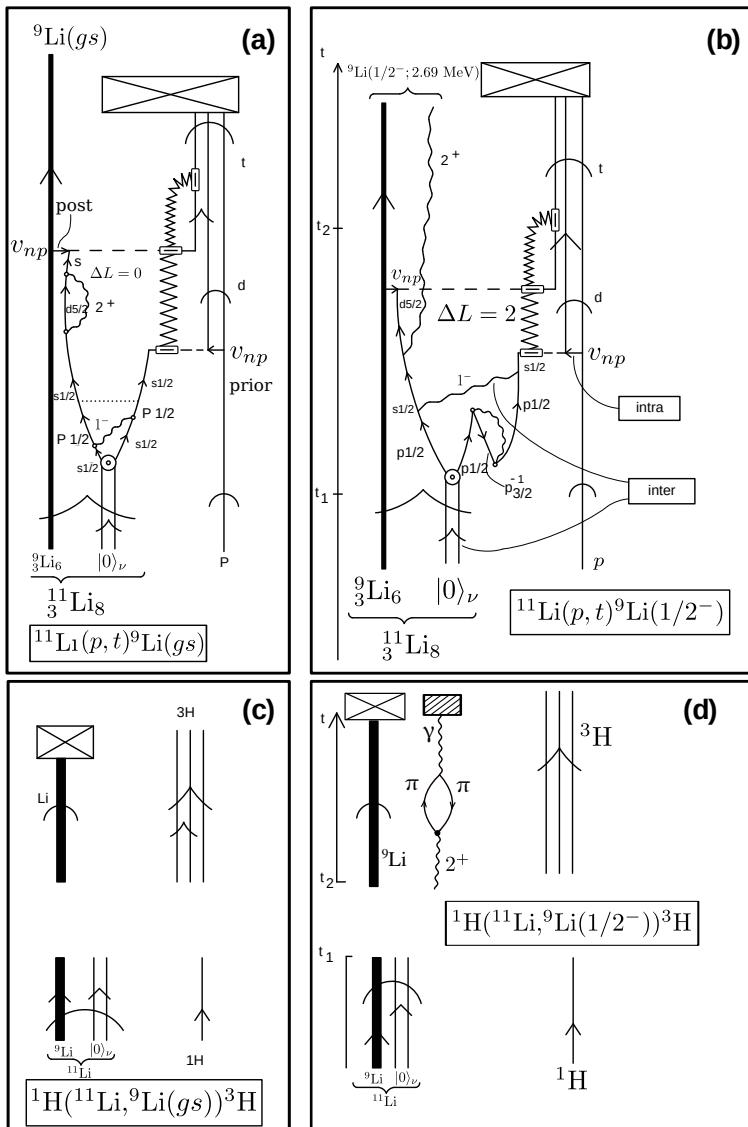


Figure 2.7.3: NFT-diagram describing a contribution to the reaction amplitude **(a)** $^{11}\text{Li}(p, t)^9\text{Li(gs)}$, **(b)** $^{11}\text{Li}(p, t)^9\text{Li}(1/2^-)$, **(c)** $^1\text{H}(^{11}\text{Li}, ^9\text{Li(gs)})^3\text{H}$ and **(d)** $^1\text{H}(^{11}\text{Li}, ^9\text{Li}(1/2^-))^3\text{H}$. These last two processes making it concrete, through inverse kinematics, the gedankenexperiments (a) and (b). Time is assumed to run upwards. A single arrowed line represents a fermion (proton) (π) or neutron (ν). A double arrowed line two correlated nucleons. In the present case two correlated (halo) neutrons (halo-neutron pair addition mode $|0\rangle_\nu$). A heavy arrowed line represents the core system $|^9\text{Li(gs)}\rangle$. A standard pointed arrow refers to structure, while "round" arrows refer to reaction. A wavy line represents (particle-hole) collective vibrations, like the low-lying quadrupole mode of ^9Li , or the dipole pygmy resonant state which, together with the bare pairing interaction (horizontal dotted line) binds the neutron halo Cooper pair to the core. A short horizontal arrow labels the proton-neutron interaction v_{np} responsible for the single-particle transfer processes, represented by an horizontal dashed line. A dashed open square indicates the particle-recoil coupling vertex. The jagged line represents the recoil mode resulting from the mismatch between the relative centre of mass coordinates associated with the mass partitions $^{11}\text{Li}+p$, $^{10}\text{Li}+d$ (virtual) and $^9\text{Li}+t$. The γ -detector is represented by a hatched box, the particle detector by a crossed rectangle.

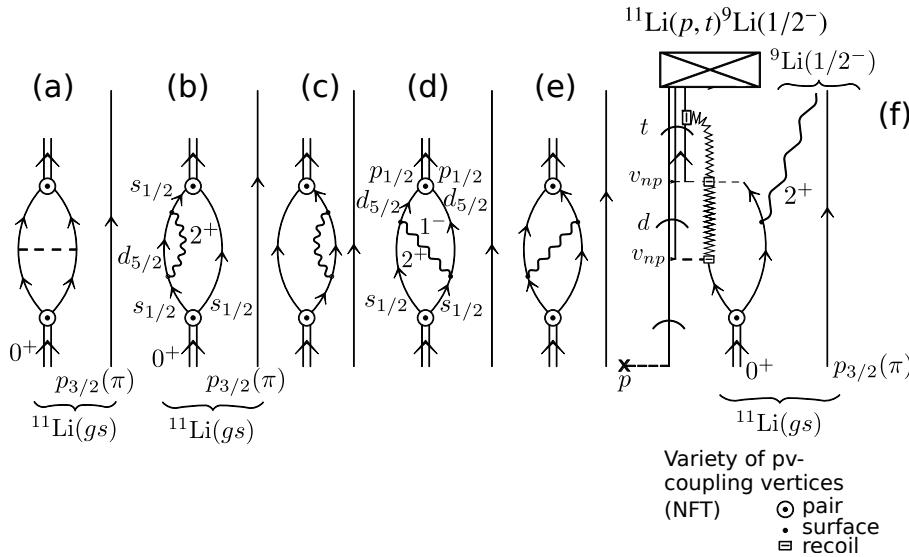


Figure 2.7.4: (a–e) Lowest order, NFT diagrams associated with the processes contributing to the binding of the neutron halo Cooper pair (double arrowed line) of ^{11}Li to the core ^9Li through the bare pairing interaction (dashed line) as well as the exchange of the core quadrupole phonon and of the soft dipole mode of ^{11}Li (wavy line). Single arrowed lines describe the nucleon independent-particle motion of neutrons ($s_{1/2}, d_{5/2}$, etc.) as well as of the $p_{3/2}(\pi)$ proton considered as a spectator; (a) Bare pairing interaction, four-point vertex (horizontal dotted line); (b, c) self energy, effective mass polarization (PO) process dressing the $s_{1/2}(v)$ single-particle state (a similar diagram, but corresponding to correlation (CO) processes dressing the $p_{1/2}$ state is not shown, see Fig. 2.7.2); (d, e) vertex correction (induced pairing interaction) renormalizing the vertex with which the pair addition mode couples to the fermions (dotted open circle); (f) NFT diagram describing the reaction $^{11}\text{Li}(p, t)^{9}\text{Li}(1/2^-)$ populating the first excited state of ^9Li , the dashed horizontal line starting with a cross stands for the (p, t) probe. The successive transfer of the two halo neutrons ($^{11}\text{Li}(gs)+p \rightarrow ^{10}\text{Li}+d \rightarrow ^9\text{Li}(1/2^+) + t$) is shown, in keeping with the fact that this process provides the largest contribution to the absolute differential cross section. The jagged line represents the recoil mode carrying to the outgoing particle the effect of the momentum mismatch associated with the transfer process (recoil).

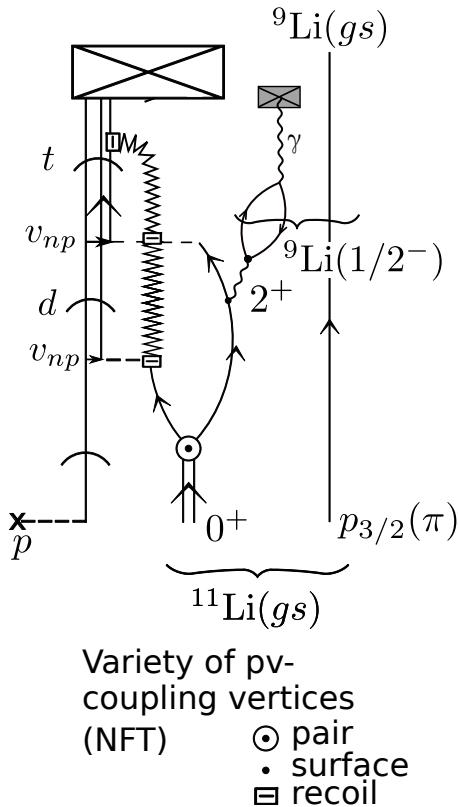


Figure 2.7.5: Gedankenexperiment: γ -ray coincidence between $^1\text{H}(^{11}\text{Li}, ^9\text{Li})^3\text{H}$ and $^9\text{Li}(\text{gs}) + \gamma$ ($E2; 2.69 \text{ MeV}$) decay process. The virtual quadrupole phonon associated with self-energy and vertex correction processes becomes real through the action of the $^1\text{H}, ^3\text{H}$ (p, t) external field. Thus, it is not only that recoil modes are “measured” by particle detectors in connection with outgoing particles which have asymptotic wavefunctions, but also the quadrupole vibration, whose eventual γ -decay (see Fig. 5.3.2 (II)) can be measured by the γ -detector. The apparent bubble process made out of a proton particle-hole component of the quadrupole mode does not contradict NFT rule III (Sect. 2.5). In fact, the initial vertex (solid dot) corresponds to a nuclear PVC, while the second vertex the proton excitation to the electromagnetic field, leading to γ -decay.

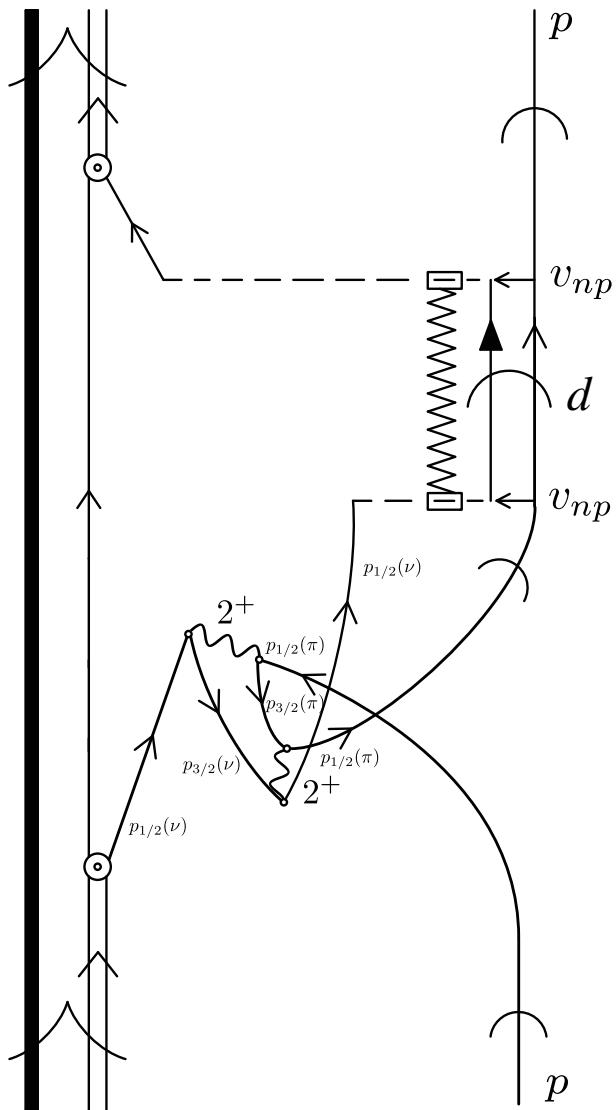


Figure 2.7.6: In keeping with standard direct reaction praxis, neither in Fig. 2.7.2 nor in 2.7.3 antisymmetrization is carried out between the impinging proton and the protons of ^{11}Li . Within the present discussion ($(^{11}\text{Li}(p, p))^{11}\text{Li}$), an example of such processes corresponds to the exchange of a proton participating in the quadrupole vibration of the core, with the projectile, as shown in the figure. Such a process will not only be two orders higher in perturbation in the particle-vibration coupling vertex than the original one shown in graph (a) of Fig. 2.7.2. It will also be strongly suppressed by the square of the overlap between a proton moving in the continuum, and a $p_{1/2}$ proton of the ^9Li core.

$t_1 - t_{11}$, starting from the $t = 0$ situation in which a proton impinges¹²⁷ on ^{11}Li .

¹²⁷To be carried out in inverse kinematics, i.e. a beam of ^{11}Li on a ^1H target.

At time t_1 , the halo pair addition mode $|0_\nu\rangle$ couples to a pure, bare configuration $s_{1/2}^2(0)$. At time t_2 , due to the zero point fluctuations associated with the quadrupole vibration of the ^{11}Li core, namely ${}^9\text{Li}$, the virtual state $((p_{1/2}, p_{3/2}^{-1})_{2^+} \otimes 2^+)_{0^+}$ is created. At time t_4 , one of the $s_{1/2}$ neutrons excites the quadrupole vibration of the core reabsorbing it at time t_6 . As a result of this self energy process, its energy is lowered to threshold, becoming a virtual state. The other $s_{1/2}$ neutron state excites at t_3 the PDR and moves into the $p_{1/2}$ orbital after which, due to Pauli principle, becomes exchanged¹²⁸ with the homologous $p_{1/2}$ of the $(p_{1/2}, p_{3/2}^{-1})_{2^+}$ configuration, exchange process which is completed by time t_5 . As a result the $p_{1/2}$ state undergoes a conspicuous repulsion, becoming a resonant state¹²⁹. At time t_7 the $s_{1/2}$ neutron absorbs the PDR and moves into a $p_{1/2}$ state. At time t_8 it interacts, through the bare 1S_0 -pairing interaction, with the other $p_{1/2}$ neutron. Before the two $p_{1/2}$ neutrons couple to the $|0_\nu\rangle$ state one of them is picked up at time t_9 under the action of the proton-neutron interaction v_{np} , by the projectile (proton), to form a virtual deuteron. The recoil effect associated with the new mass partition being taken care of by the particle-recoil coupling vertex and associated recoil mode (jaggy line). At time t_{10} , under the effect of v_{np} , the neutron of the virtual deuteron is transferred back to the virtual ${}^{10}\text{Li}$ nucleus to form again ${}^{11}\text{Li}$ and thus the original mass partition, as testified by the fact that the recoil phonon (jaggy line initiated at time t_9) is reabsorbed at a second particle-recoil coupling vertex. The resulting $(p_{1/2}^2(0))$ configuration couples, at time t_{11} , to the $|0_\nu\rangle$ state leading to the $|{}^{11}\text{Li(gs)}\rangle$ state. The system has thus returned to the entrance channel configuration, namely that of ${}^{11}\text{Li} + p$, which propagates to $t \rightarrow +\infty$, the proton eventually bringing the nuclear structure information to the detector.

In other words, the real and imaginary non-local and ω -dependent processes shown in Fig 2.7.2 (a) contribute to the polarization component of the optical potential¹³⁰ and to its A -dependence. It is of notice that Fig. 2.7.2 exemplifies the elements needed to extend and formalize NFT rules of structure so as to be able to deal, on equal footing or, more properly, in an unified fashion, with reactions.

As schematically shown in Fig. 2.7.1 (see also Sect. 3.6.1), to each nuclear structure process displayed in diagram (a) of Fig. 2.7.2, it corresponds a specific

¹²⁸The (inevitable) crossing between the two $p_{1/2}$ arrowed lines: see Fig. 2.7.1, diagram at the center of the panel indicated by a (double circled) number 2, and the corresponding one in the upper left corner (${}^{11}\text{Li-halo}$); see also diagrams (a) and (b) of Fig. 1.2.4.

¹²⁹It is of notice that the fact that to follow the dressing of the $p_{1/2}$ neutron by the quadrupole vibration of the core ${}^9\text{Li}$, one has to go backwards in time, in a N-like trajectory testifies to the fact that the hole states play, in nuclei, the role anti-particles do in field theory, in particular in QED.

¹³⁰Concerning the optical potential we refer to footnote 23 of Ch. 1. In relation with the program of NFT_(s+r) one can mention that Landau felt that Feynman diagrams have an independent basic importance, because of the possibility of relating them directly to physical observables. Feynman diagrams allow to describe processes where one set of particles with given energies, momenta, angular momenta, go in and another set (or the same) comes out. At the basis of this approach one finds vertex processes and dispersion relations. Now, vertex processes can simply mean the variety of processes connecting the incoming particles with the outgoing ones. In other words, within the present framework the processes taking place between times $t_2 - t_{11}$ (Figs. 2.7.2 (a) and (b)) and $t_1 - t_2$ (Fig. 2.7.3 (b)) (Landau (1959); ter Haar (1969)). See also Sect. 5.3.5.

equation, amplitude, etc., and thus a number with appropriate units. This is also so regarding the reaction aspects of the diagram, and thus reaction amplitudes, formfactors and absolute differential cross sections. Within this context we refer to Fig. 4.1.2 in connection with the inset to Fig. 2.7.2 (a), as well as to graph (a) and (b) of Fig. 2.7.3.

In Fig. 2.7.2 (b) one assumes the same processes to take place as in (a) up to time t_8 (reason for which no details are repeated between t_2 and t_8). From there on the deuteron continues to propagate to the detector, and the effect of the particle-recoil coupling vertex is to be worked out and the corresponding outgoing distorted waves modified accordingly. Likely, the halo neutron in ^{10}Li will break up before the system can be recorded by the particle detector. Summing up, in the center of mass reference frame both p and ^{11}Li display asymptotic states in entrance as well as in exit channels in case (a), and only in the entrance channel in case (b), while in the exit channel only $^{10}\text{Li}({}^9\text{Li}+n)$ and the deuteron do so.

Another examples of the NFT diagrams of structure and reactions are given in Fig. 2.7.3. In (a) a contribution associated with the reaction $^{11}\text{Li}(p,t){}^9\text{Li}(\text{gs})$ is shown, while in (b) one associated with the population of the first excited $1/2^-$ (2.69 MeV) state of ${}^9\text{Li}$ is displayed. We refer to Figs. 2.7.4 and 2.7.5 for a compact graphical representation of this last process. The importance of such process is that it provided direct evidence of phonon mediated pairing interaction in nuclei as theoretically predicted¹³¹.

Concerning the question of Pauli principle in reaction processes, e.g. between the incoming proton and the collective modes of the core (${}^9\text{Li}$) we refer to Fig. 2.7.6.

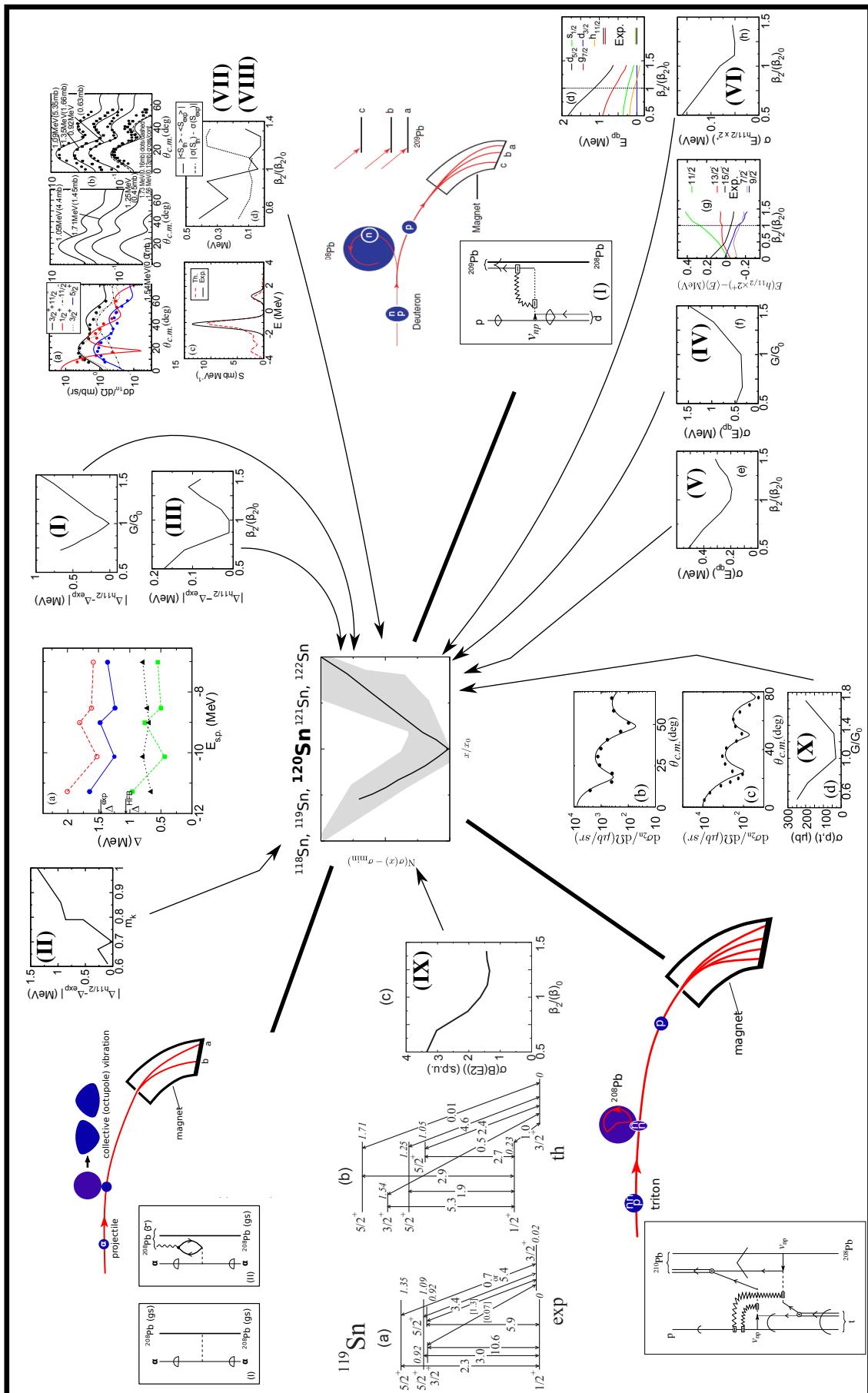
2.8 Characterization of an open-shell nucleus: ^{120}Sn

In this section we aim at giving an overall view of the versatility of $(\text{NFT})_{\text{ren(s+r)}}$ to describe both structure as well as the results of experimental probes which provide an essentially complete characterization of an atomic nucleus or, better, of a small island of closely related nuclei (isotopes). Details of the structure and reaction calculations, techniques and computer codes employed to achieve such characterization are provided in the following chapters.

In keeping with the fact that $^{120}_{50}\text{Sn}$ is a typical example of single open shell superfluid nuclei, it has been studied extensively with a variety of probes. Elastic, anelastic (Coulomb excitation and subsequent γ -decay as well as inelastic scattering). Also one- and two-particle transfer reactions. The corresponding absolute differential cross sections and transition probabilities involve as targets and residual systems the island of superfluid nuclei $^{118,119,120,121,122}\text{Sn}$.

A theoretical description of the variety of observables have been carried out solving the Nambu-Gorkov equation to propagate the different $(\text{NFT})_{\text{ren(s)}}$ pro-

¹³¹Barranco et al. (2001), Tanihata et al. (2008); Potel et al. (2010); Tanihata et al. (2013); Beceiro-Novoa et al. (2015).

Figure 2.8.1: Characterization of ^{120}Sn .

cesses dressing the single-particle states and renormalizing the bare pairing interaction, as well as spectroscopic and transition amplitudes¹³². With the help of the computer codes SINGLE and COOPER, tailored to propagate the $(\text{NFT})_{\text{ren(s)}}$ spectroscopic amplitude content to the detector, one- and two- nucleon transfer absolute differential cross sections were worked out. In Fig. 2.8.1 theory is compared to experiment in terms of renormalized energies, absolute differential cross sections and electromagnetic transition probabilities.

In the **upper left** part of this figure, the cartoon representation displayed in Fig. 2.1.1 is here used to schematically illustrate¹³³ anelastic processes, Coulomb excitation and the quadrupole γ -decay of ^{119}Sn (**middle left** (a) experiment, (b) theory, (c) standard deviation between theory and experiment calculated as a function of β_2 and displayed normalized with respect to the experimental value $(\beta_2)_0$ (**IX**)). This value together with those also labeled by roman numerals, referred to below, were used to draw the center boxed nuclear structure landscape (see Fig. 2.4.1 and Table 2.4.1).

In the **bottom left** part of the figure, a schematic representation of an experimental setup to measure two-nucleon transfer processes is given. Also a $(\text{NFT})_{\text{ren(s+r)}}$ diagram describing successive transfer is displayed (Fig. 2.1.3). In the **bottom middle** part of the figure, the theoretical absolute differential cross sections $d\sigma_{2n}/d\Omega$ (continuous curves) associated with the reactions $^{120}\text{Sn}(p, t)^{118}\text{Sn}(\text{gs})$ and $^{122}\text{Sn}(p, t)^{120}\text{Sn}(\text{gs})$ are displayed ((b) and (c)), in comparison with the experimental data (solid dots). The theoretical cross sections were recalculated as a function of the pairing strength G and the corresponding standard deviation with respect to the experimental value determined. The results are shown in (d) (**X**) as a function of G , normalized with respect to the (equivalent) value G_0 of v_{14} (Argonne 1S_0 bare NN -interaction; only the values corresponding to $^{120}\text{Sn}(p, t)^{118}\text{Sn}$ are displayed). It is of notice that the corresponding relative standard deviation σ/L for $G = G_0$ ($L \equiv \sigma_{2n}$ being the experimental cross section), is quite small ($70\mu\text{b}/2250\mu\text{b} \approx 0.03$)¹³⁴.

In the **upper middle** part of the figure and under the label (a), the state dependent pairing gap of the valence orbitals of ^{120}Sn are displayed, in comparison with the experimental findings ($\Delta^{\text{exp}} \approx 1.45$ MeV, arrow left). The gap associated with the lowest quasiparticle state $h_{11/2}$ calculated as a function of m_k (different Skyrme interaction) as well as of G/G_0 and $\beta_2/(\beta_2)_0$ have been used to work out the corresponding standard deviations with respect to the experimental findings and are displayed to the right (**I**) and (**III**) and the left (**II**) of (a).

In the **upper right** part of the figure, absolute differential cross sections and

¹³²See Idini et al. (2015) and refs. therein, see also Broglia et al. (2016).

¹³³While in Fig. 2.1.1 one refers to the excitation of an octupole mode, in the present case we deal with a quadrupole vibration.

¹³⁴This result is related with the fact that $\sigma \sim |\alpha_0|^2$, α_0 being the BCS order parameter. That is, the number of Cooper pairs participating in the BCS condensate which measures the deformation in gauge space. Because the state $|BCS\rangle$ describing phase-correlated, independent pair motion (Cooper pair condensate) is a coherent state displaying off-diagonal-long-range-order (ODLRO), it is not surprising that $|\alpha|^2 = \left| \sum_{\nu>0} U_\nu V_\nu \right|^2 = \left| \langle BCS | P^\dagger | BCS \rangle \right|^2 = \left| \langle BCS | P | BCS \rangle \right|^2$ plays the role of a physical, non-energy weighted sum rule (Potel et al. (2017)). See Sect. 7.2.1.

strength functions associated with the one-particle transfer processes $^{120}\text{Sn}(d, p)^{121}\text{Sn}$, are displayed. In (a), the absolute differential cross sections associated with the low-lying states $h_{11/2}, d_{3/2}, s_{1/2}$ and $d_{5/2}$ are shown (theory: continuous curve; data: solid dots). In (b) (left) the calculated $^{121}\text{Sn}(5/2^+)$ absolute differential cross sections (continuous curves) are shown and compared with the experimental data (right, solid dots; also given here are DWBA fits used in the analysis of the experimental data). In (c) the calculated strength function associated with the $5/2^+$ state (dashed curve) is compared to the data (continuous curve), while in (d) the difference between the centroid and width of the experimental and calculated $d_{5/2}$ strength function are shown as a function of the ratio $\beta_2/(\beta_2)_0$ in terms of solid and dashed curves ((VII) and (VIII)). In the **middle right** part of the figure, a cartoon representation of a setup to measure one-nucleon transfer reactions is displayed. Also shown is a $(\text{NFT})_{\text{ren (s+r)}}$ diagram describing the process (see Fig. 2.1.2).

In (d) (**lower right** part of the figure), the lowest quasiparticle energy values are displayed as a function of $\beta_2/(\beta_2)_0$ in comparison with the data. The root mean-square deviation between the experimental and theoretical levels as a function of $\beta_2/(\beta_2)_0$ and of G/G_0 are shown in (e) and (f) respectively ((V), (IV)). In (g) the experimental energies of the members of the $h_{11/2} \otimes 2^+$ multiplet, are compared with the theoretical values calculated as a function of the ratio $\beta_2/(\beta_2)_0$. Finally in (h) the root mean-square deviation between the experimental and theoretical energies of the members of the $h_{11/2} \otimes 2^+$ multiplet shown in (g) are given as a function of $\beta_2/(\beta_2)_0$ (VI).

Summing up, the nuclear structure landscape is well funneled and theory provides an overall account of the data, when the physical values of β_2, G and m_k are used¹³⁵.

2.9 Summary

In Fig. 2.8.1 we report on the $(\text{NFT})_{\text{ren (s+r)}}$ analysis of a rather complete set of data aimed at characterizing the superfluid nucleus ^{120}Sn , namely: **1**) $^{120}\text{Sn}(p, t)^{118}\text{Sn}(\text{gs})$, $^{122}\text{Sn}(p, t)^{120}\text{Sn}(\text{gs})$; **2**) $^{120}\text{Sn}(p, d)^{119}\text{Sn}$, $^{121}\text{Sn}(p, d)^{120}\text{Sn}$; **3**) $^{119}\text{Sn}(\alpha, \alpha')^{119}\text{Sn}$ (γ -decay); **4**) energies; **5**) strength functions; **6**) pairing gaps and **7**) $B(E2)$ transition probabilities.

Arbitrarily forcing the particle-vibration coupling (PVC) strength, the strength of the bare pairing force and the value of the k -mass to depart from their “physical” values, one can test the robustness of the $\text{NFT}_{\text{ren (s+r)}}$ picture of ^{120}Sn given, and of the funneled character of the associated nuclear structure and reaction landscape.

In a very real sense this, namely the results collected in Fig. 2.8.1, is a nucleus. That is, the summed experimental and theoretical structural information accessed through asymptotic states, outcome of probing the system with a “complete” array of experiments interpreted with the help of a unified structure and reaction theoretical scheme as, for example, that provided by $(\text{NFT})_{\text{ren (s+r)}}$.

¹³⁵For details, see Idini et al. (2015), and Broglia et al. (2016), see also Fig. 2.4.1.

Appendix 2.A NFT_(s+r): linear theory

NFT is linear in the particle-vibration coupling vertices associated with the variety of modes (see e.g. Fig. 2.5.10). This is also true concerning its extension to describe reaction processes, as can be seen from graph (a) of Fig. 2.A.1. For simplicity, this diagram and similar ones are drawn as displayed in Fig. 2.A.1 (b) (see e.g. Fig. 2.1.3). However, in all cases the effects of recoil are properly taken into account following (a), as shown in e.g. Fig. 4.1.2 as well as in Sect. 6.1 within the framework of second order DWBA, and in Sect. 6.5 (Eq. (6.5.7)) in the semiclassical approximation.

Appendix 2.B NFT and reactions

Nuclear Field Theory was systematically developed to describe nuclear structure processes. This fact did not prevent the translation into this graphical language, expressions which embodied the transition amplitude of a variety of reaction processes, in particular second order (in v_{np}) transition amplitudes associated with two nucleon transfer reactions¹³⁶.

The new feature to be considered regarding transfer processes and not encountered neither in structure, nor in elastic or anelastic processes, is the graphical representation of recoil effects. That is, a physical phenomenon associated with the change in the coordinate of relative motion reflecting the difference in mass partition between entrance (also intermediate, if present) and exit channels. Phenomenon which mixes in a non-analytically separable fashion, structure and reaction degrees of freedom.

Nuclear structure processes do not affect the center of mass of the nucleus, with a proviso. In fact, the shell model potential violates translational invariance of the total nuclear Hamiltonian and, thus, single-particle excitations can be produced by a field proportional to the total center-of-mass coordinate. The translational invariance can be restored by including the effects of the collective field generated by a small displacement α of the nucleus. Such a displacement, e.g. in the x -direction, gives rise to a coupling which can be written as,

$$H_c = \kappa\alpha F, \quad (2.B.1)$$

where

$$F = -\frac{1}{\kappa} \frac{\partial U}{\partial x}, \quad (2.B.2)$$

and

$$\kappa = \int \frac{\partial U}{\partial x} \frac{\partial \rho_0}{\partial x} d\tau = -A \left\langle \frac{\partial^2 U}{\partial x^2} \right\rangle, \quad (2.B.3)$$

¹³⁶Broglia (1975); Broglia and Winther (2004).

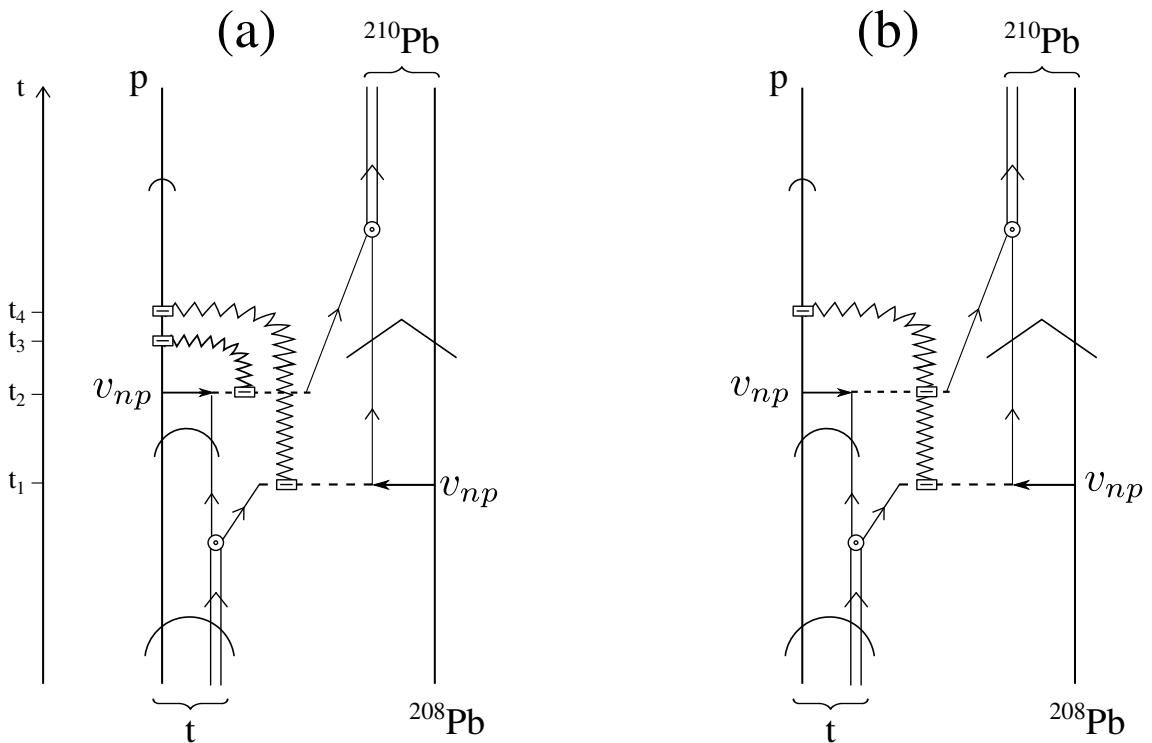


Figure 2.A.1: NFT_(s+r) diagram describing the successive transfer contribution to the reaction $^{208}\text{Pb}(t, p)^{210}\text{Pb}(\text{gs})$. That is the population of the lowest energy, monopole pair addition mode of ^{208}Pb . Concerning the different symbols used, we refer to Figs. 2.1.3 and 2.7.2. In particular concerning the recoil mode (jagged line) and the associated particle-recoil coupling vertex (dashed open rectangle). (a) of the variety of contributions associated with the different sequence of events taking place at times t_1, t_2, t_3 and t_4 , one has chosen that in which the information of a mass $2m$ transfer process to be carried out to the detector is conveyed to the outgoing particle at $t_3(t_4)$ by the recoil process associated with the nucleon transfer at $t_2(t_1)$; diagram (b) is used only for simplicity when one is interested in discussing aspects of the (s+r)-diagram other than recoil.

corresponding to a normalization of α such that $\langle F \rangle = \alpha$. It is of notice that both κ and U are negative for attractive fields (p. 444 of¹³⁷)

The spectrum of normal modes generated by the field coupling (2.B.1), namely by a Galilean transformation of amplitude α , contains an excitation mode with zero energy for which zero point fluctuations diverge in just the right way to restore translational invariance to leading order in α . In fact, while the zero point fluctuations (ZPF)

$$\lim_{\omega_\alpha \rightarrow 0} \left(\frac{\hbar\omega_\alpha}{2C_\alpha} \right)^{1/2} = \lim_{\omega_\alpha \rightarrow 0} \left(\frac{\hbar^2}{2D_\alpha \hbar\omega_\alpha} \right)^{1/2}, \quad (2.B.4)$$

diverge the inertia remains finite and equal to $D_\alpha = AM$, as expected, C_α being the restoring force constant. The additional dipole roots include, in particular, the isoscalar dipole modes associated with $\hat{D} = \sum_{i=1}^A r_i^3 Y_{1\mu}(\hat{r}_i)$, which can be viewed as a non-isotropic compression mode¹³⁸.

The operators leading to transformations associated with the change in the coordinates of relative motion (recoil effects) in nuclear reaction processes, are Galilean operators ($\sim \exp(\mathbf{k}_{\alpha\beta} \cdot (\mathbf{r}_\beta - \mathbf{r}_\alpha))$). Their action (on e.g. the entrance channel distorted waves), as that of (2.B.1) on the shell model ground state, can be graphically represented in terms of NFT diagrams (or eventual extensions of them). In Figs 2.1.2 and 2.1.3 as well as 2.7.2–2.7.6 they are drawn in terms of jagged lines. How do we calculate such couplings? Let us elaborate on this point.

When one states that the small displacement α of the nucleus leads to a coupling (2.B.1) one means a coupling between the single-particle and the collective displacement of the system as a whole. When one talks about the spectrum of normal modes associated with such a coupling, one refers to the harmonic approximation (RPA). Thus, to the solutions of the dispersion relation¹³⁹,

$$-\frac{2\kappa}{\hbar} \sum_i \frac{|F_i|^2 \omega_i}{\omega_i^2 - \omega_a^2} = 1, \quad (2.B.5)$$

where the sum is over dipole particle-hole excitations. This dispersion relation can be represented graphically through the diagrams shown in Fig. 1.2.3 (inset). In particular, α acting on the vacuum creates the collective mode. This can also be seen by expressing α in second quantization, namely

$$\alpha = \sqrt{\frac{\hbar\omega_\alpha}{2C_\alpha}} (\Gamma_\alpha^\dagger + \Gamma_\alpha), \quad (2.B.6)$$

where $\sqrt{\hbar\omega_\alpha/2C_\alpha} = \sqrt{\frac{\hbar^2}{2D_\alpha} \frac{1}{\hbar\omega_\alpha}}$ is the zero-point amplitude of the collective (displacement) mode. Now, none of the above arguments lose their meaning in the

¹³⁷Bohr and Mottelson (1975).

¹³⁸See e.g. Colò et al. (2000).

¹³⁹Bohr and Mottelson (1975); Eq. (6-244), Brink and Broglia (2005), Sect. 8.3.1.

case in which there is a root with $\omega_\alpha = 0$, also in keeping with the fact the inertia remains finite. In Figs. 2.7.2–2.7.6 we do something similar to what is done in Fig. 1.2.3 (inset). The dot, which in this figure represents the particle-vibration coupling, is replaced by a small dashed open square, which we label “particle-recoil coupling vertex” (see labels Fig. 2.7.2). It constitutes a graphical mnemonic for counting the degrees of freedom that are at play. In this case the coordinates of relative motion. Also the fact that in connection with the appearance of such vertices one has to calculate matrix elements of precise form factors which involve the recoil phases.

Concerning the actual calculation of a particle-mode vertex in which $\omega_a \rightarrow 0$, an empirical way out is that of a coarse-grained-like symmetry restoration. In this case κ is adjusted in such a way that the lowest solution of Eq. (2.B.5), although being smaller than the rest of them, remains finite¹⁴⁰.

2.B.1 Transfer

We now consider, within the framework of the semiclassical¹⁴¹ approximation a general reaction



in which the nucleus a impinges on the nucleus A in the entrance channel $\alpha(a, A)$ and where the two nuclei in the exit channel β , namely b and B may differ from those in α , by the transfer of one or more nucleons.

In the center-of-mass system, the total Hamiltonian may be written as

$$\begin{aligned} H &= T_{aA} + H_a + H_A + V_{aA}, \\ &= T_{bB} + H_b + H_B + V_{bB}, \end{aligned} \quad (2.B.8)$$

where T_{aA} is the kinetic energy of the relative motion in channel α

$$T_{aA} = -\frac{\hbar^2}{2m_{aA}} \nabla_{aA}^2, \quad (m_{aA} = \frac{m_a m_A}{m_a + m_A}), \quad (2.B.9)$$

and similarly for T_{bB} . Assuming the nuclei in (2.B.7) to be heavy ions, one can solve the time-dependent Schrödinger equation

$$i\hbar \frac{\partial \Psi}{\partial t} = H\Psi \quad (2.B.10)$$

¹⁴⁰Within this context we refer to Bohr and Mottelson (1975), p. 446. With no coupling H_c (Eq. (2.B.1)), the ZPF $\alpha_0^{(0)}$ of the nuclear CM are small ($\sim A^{-2/3}$). Thus, it is possible to tune κ so as to make the ZPF associated with the lowest root large as compared to $\alpha_0^{(0)}$, but still compatible with the ansatz at the basis of RPA (small amplitude harmonic vibrations).

¹⁴¹Assuming Eq. (2.B.7) to describe a transfer process induced by a heavy ion reaction with bombarding energy lying somewhat above the Coulomb barrier, the associated de Broglie reduced wavelength $\lambda = \sqrt{2mE}$ is much smaller than the nuclear dimension, and the relative motion can be described in terms of classical trajectories. On the other hand, the particle transfer process is described fully quantum mechanically. The coupling between the intrinsic (structure) nucleon motion and the relative motion associated with the recoil effect, is dealt with in terms of narrow wave packets. See e.g. Broglia and Winther (2004) and references therein.

with the initial condition that a and A are in their ground states, and that the relative motion is described as a narrow wave-packet of rather well-defined impact parameter and velocity. Because of the quantal nature of the process under consideration, we study the quantal description in the limit of small wavelength of relative motion (semiclassical approximation). One expands Ψ on the channel wavefunctions

$$\Psi_\beta(t) = \Psi_m^b(\xi_b)\Psi_n^B(\xi_B)e^{i\delta_\beta} \quad (2.B.11)$$

where Ψ^b and Ψ^B describe the structure of the two nuclei and satisfy the equations

$$H_b\Psi_m^b(\xi_b) = E_m^b\Psi_m^b(\xi_b) \quad (2.B.12)$$

and

$$H_B\Psi_n^B(\xi_B) = E_n^B\Psi_n^B(\xi_B), \quad (2.B.13)$$

while ξ_b and ξ_B denote the intrinsic coordinates. The phase δ_β is defined by

$$\delta_\beta = \frac{1}{\hbar} \left\{ m_\beta \mathbf{v}_\beta(t)(\mathbf{r}_\beta - \mathbf{R}_\beta(t)) - \int_0^t \left(U_\beta(R_\beta(t')) - \frac{1}{2}m_\beta v_\beta^2(t') \right) dt' \right\}. \quad (2.B.14)$$

The index β labels both the partition of nucleons into b and B , as well as the quantal states of the two nuclei. The quantity U_β is the ion-ion potential in this channel. It is equal to the expectation value of $V_\beta = V_{bB}$ in the channel β . The distance between the centers of mass of the two systems is denoted by

$$\mathbf{r}_\beta = \mathbf{r}_{bB} = \mathbf{r}_b - \mathbf{r}_B. \quad (2.B.15)$$

The quantity \mathbf{R}_β and its derivative $\mathbf{v}_\beta = \dot{\mathbf{R}}_\beta$ describe the motion of the centers of mass of the wavepackets, and satisfy the corresponding classical equation of motion,

$$m_\beta \dot{\mathbf{v}}_\beta = -\nabla U_\beta(\mathbf{R}_\beta). \quad (2.B.16)$$

The phase factor $e^{i\delta_\beta}$ in the channel wavefunction, is essentially a Galilean transformation where an additional phase (related to the Q -value) has been added to eliminate, as far as possible, the diagonal matrix elements of the coupled equations. Using the notation $E_\beta = E_m^b + E_n^B$ and inserting the ansatz

$$\Psi = \sum_\beta c_\beta((r_\beta - R_\beta), t)\Psi_\beta(t)e^{-iE_\beta t/\hbar} \quad (2.B.17)$$

in eq. (2.B.10) one obtains, assuming narrow wavepackets, product of an amplitude $a_\beta(t)$ and a shape¹⁴² $\chi_\beta(\mathbf{r} - \mathbf{R}_\beta(t), t)$, ($c_\beta = a_\beta \chi_\beta$),

$$\begin{aligned} i\hbar \sum_\beta \dot{a}_\beta(t) \langle \Psi_\xi | \Psi_\beta \rangle_{\mathbf{R}_\xi} e^{-iE_\beta t/\hbar} \\ = \sum_\gamma \langle \Psi_\xi | V_\gamma - U_\gamma(r_\gamma) | \Psi_\gamma \rangle_{\mathbf{R}_\xi} a_\gamma(t) e^{-iE_\gamma t/\hbar}, \end{aligned} \quad (2.B.18)$$

¹⁴²See Broglia and Winther (2004) p. 295.

where the sub-index on the matrix elements indicate that the integration over the degree of freedom of the two nuclei, the average center-of-mass coordinate $\mathbf{r}_{\beta\gamma} = (\mathbf{r}_\beta + \mathbf{r}_\gamma)/2$ should be identified with the average classical coordinate, i.e.

$$\mathbf{r}_{\beta\gamma} \rightarrow \mathbf{R}_{\beta\gamma} = \frac{1}{2}(\mathbf{R}_\beta + \mathbf{R}_\gamma), \quad (2.B.19)$$

and the functions $\langle \Psi_\xi | V_\gamma - U_\gamma(r_\gamma) | \Psi_\gamma \rangle_{\mathbf{R}_{\xi\gamma}}$ are the form factors. The coupled equations (2.B.18) can be written in a more compact way by an orthogonalization procedure, which makes use of the *adjoint channel wavefunctions*

$$\omega_\xi = \sum_\gamma g_{\xi\gamma}^{-1} \Psi_\gamma, \quad (2.B.20)$$

where g^{-1} is the inverse of the overlap matrix

$$g_{\xi\gamma} = \langle \Psi_\xi | \Psi_\gamma \rangle, \quad (2.B.21)$$

that is

$$\sum_\xi g_{\gamma\xi} g_{\xi\beta}^{-1} = \sum_\xi g_{\gamma\xi}^{-1} g_{\xi\beta} = \delta(\gamma, \beta). \quad (2.B.22)$$

With this definition,

$$(\omega_\xi, \Psi_\beta) = \delta(\xi, \beta), \quad (2.B.23)$$

which takes care of non-orthogonality. Making use of the above relations one can rewrite (2.B.18) in the form

$$i\hbar \dot{a}_\beta(t) = \sum_\gamma \langle \omega_\beta | V_\gamma - U_\gamma | \Psi_\gamma \rangle_{\mathbf{R}_{\beta\gamma}} e^{i(E_\beta - E_\gamma)t/\hbar} a_\gamma(t). \quad (2.B.24)$$

That is, the proper transfer (tunneling) equations are obtained from (2.B.18) by a basis orthogonalization process¹⁴³. Solving these coupled equations with the condition that at $t = -\infty$ the system is in the ground state of a and A (entrance channel α), that is $a_\gamma(-\infty) = \delta_{\gamma,\alpha}$, one can calculate the differential cross section

$$\frac{d\sigma_{\alpha \rightarrow \beta}}{d\Omega} = P_{\alpha \rightarrow \beta} \sqrt{\left(\frac{d\sigma_\alpha}{d\Omega} \right)_{el} \left(\frac{d\sigma_\beta}{d\Omega} \right)_{el}}, \quad (2.B.25)$$

where $P_{\alpha \rightarrow \beta}$ is the absolute value squared of the transition amplitude $|a_\beta(t = +\infty)|^2$. It gives the probability that the system at $t = +\infty$ is in the final channel. The quantities $(d\sigma/d\Omega)_{el}$ are the (semiclassical) elastic cross sections.

¹⁴³Cf. with Eq. (4.6.1) and following discussion, in connection with the derivation of the Cooper pair tunneling Hamiltonian (Cohen et al. (1962)) between two weakly coupled superconductors (Josephson junction).

We now solve the coupled equations in first order perturbation theory. For this purpose we insert $\delta(\gamma, \alpha)$ at the place of $a_\gamma(t)$ obtaining

$$\begin{aligned} a_\beta(t) &= \frac{1}{i\hbar} \int_{-\infty}^t \langle \omega_\beta | V_\alpha - U_\alpha | \Psi_\alpha \rangle_{\mathbf{R}_{\beta\alpha}(t)} \exp^{i(E_\beta - E_\alpha)t'/\hbar} dt' \\ &= \frac{1}{i\hbar} \int_{-\infty}^t dt' \langle \Psi_\beta | V_\alpha - U_\alpha | \Psi_\alpha \rangle_{\mathbf{R}_{\beta\alpha}(t)} \exp^{i(E_\beta - E_\alpha)t'/\hbar} \end{aligned} \quad (2.B.26)$$

where the expansion,

$$\omega_\beta = \Psi_\beta - \langle \Psi_\alpha | \Psi_\beta \rangle_{\mathbf{R}_{\beta\alpha}(t)}, \quad (2.B.27)$$

has been used, and the ansatz made, that the global optical potentials (U : real part), and standard nucleon–nucleon interactions V fulfill the relation

$$\langle \Psi_\alpha | V_\alpha - U_\alpha | \Psi_\alpha \rangle = 0. \quad (2.B.28)$$

Let us consider for simplicity the one-particle transfer reaction¹⁴⁴

$$a (= b + 1) + A \rightarrow b + B (= A + 1). \quad (2.B.29)$$

Making use of (2.B.11), that is,

$$\Psi_\alpha = \Psi^a \Psi^A e^{i\delta_\alpha}, \quad (2.B.30)$$

and

$$\Psi_\beta = \Psi^b \Psi^B e^{i\delta_\beta}, \quad (2.B.31)$$

one can write

$$\begin{aligned} \langle \Psi_\beta | V_\alpha - U_\alpha | \Psi_\alpha \rangle_{\mathbf{R}_{\alpha\beta}} &= \langle \Psi^b \Psi^B | (V_\alpha - U_\alpha) e^{i\sigma_{\alpha\beta}} | \Psi^a \Psi^A \rangle_{\mathbf{R}_{\alpha\beta}} e^{i\gamma_{\alpha\beta}} \\ &= \langle \phi^{B(A)}(S^B(n), r_{1A}), U(r_{1b}) e^{i\sigma_{\alpha\beta}} \phi^{a(b)}(S^a(n), r_{1b}) \rangle_{\mathbf{R}_{\alpha\beta}} e^{i\gamma_{\alpha\beta}}. \end{aligned} \quad (2.B.32)$$

To obtain the above relation, we have separated the difference $\delta_{\alpha\beta} = \delta_\alpha - \delta_\beta$ between the phases δ_α and δ_β into a part $\gamma_{\alpha\beta}$ which only depends on time and is related to the effective Q -value of the reaction process, and a phase $\sigma_{\alpha\beta}$ which also depends on the center-of-mass coordinate of the transferred particles. That is

$$\begin{aligned} \gamma_{\alpha\beta}(t) &= \int_0^t \left\{ U_\alpha(R_\alpha(t)) - \frac{1}{2} m_\alpha v_\alpha^2(t') - U_\beta(R_\beta(t')) + \frac{1}{2} m_\beta v_\beta^2(t') \right\} \\ &\quad + \mathbf{k}_{\alpha\beta}(t)(\mathbf{R}_\alpha - \mathbf{R}_\beta), \end{aligned} \quad (2.B.33)$$

¹⁴⁴Concerning the question of non-orthogonality in one-particle transfer processes, the first contribution arises in second order of perturbation theory, within the framework of DWBA (Thompson and Nunes (2009)).

where $\mathbf{k}_{\alpha\beta}$ is the average wave vector

$$\mathbf{k}_{\alpha\beta} = \frac{1}{2\hbar} (\mathbf{m}_\alpha \mathbf{v}_\alpha(t) + \mathbf{m}_\beta \mathbf{v}_\beta(t)). \quad (2.B.34)$$

Similarly

$$\sigma_{\alpha\beta} = \mathbf{k}_{\alpha\beta}(t) \cdot (\mathbf{r}_\beta - \mathbf{r}_\alpha). \quad (2.B.35)$$

The phase σ is characteristic for transfer processes since the dynamical variables \mathbf{r}_α and \mathbf{r}_β are identical for elastic and inelastic scattering. It arises from the change in the center-of-mass coordinate taking place when mass is transferred from one system to other. It gives rise to the recoil effect. Summing up, the one-particle transfer amplitude reads

$$\begin{aligned} (a_\beta(t = +\infty))^{(1)} &= \int_{-\infty}^{\infty} \langle \phi^{B(A)}(S^B(n), \mathbf{r}_{1A}), U_{1b}(r_{1b}) e^{\sigma_{\alpha\beta}} \phi^{a(b)}(S^a(n), \mathbf{r}_{1b}) \rangle_{\mathbf{R}_{\alpha\beta}} \\ &\times \exp \left\{ i \left[(E_\beta - E_\alpha)t'/\hbar + \gamma_{\alpha\beta} \right] \right\} \end{aligned} \quad (2.B.36)$$

The phases $\delta_{\alpha\beta}$, $[\sigma_{\alpha\beta} + \gamma_{\alpha\beta}]$ play a similar role in the determination of transfer processes reaction amplitudes as δ_l does in connection with elastic scattering cross sections. In fact, $\delta_{\alpha\beta}$ determines the shift between incoming and outgoing waves and thus the interference process which is at the basis of the absolute value of the transfer differential cross section. In other words, the reaction part of the elastic and one-nucleon-transfer reaction cross section are embodied in δ_l and $\delta_{\alpha\beta}$ respectively. The nuclear structure part is contained in the reduced mass μ and potential U in the case of elastic scattering, and in the single-particle content and radial wave-functions (formfactors), mean field potential U_{1b} and Q -value phase in the transfer case. Within the diagrammatic representation of particle-transfer reaction theory, the recoil phase is represented by a jagged line. Similar to δ_l , $\delta_{\alpha\beta}$ and $\sigma_{\alpha\beta}$ cannot be measured directly but can, in principle, be inferred from the absolute differential cross section¹⁴⁵. In other words, the jagged line does not display asymptotic behavior, representing in all cases a virtual process.

Let us conclude this section by making some comments concerning two nucleon transfer processes

$$\alpha \equiv a (= b + 2) + A \rightarrow b + B (= A + 2) \equiv \beta, \quad (2.B.37)$$

and associated sum rules. For such reactions, the overlap appearing in Eq. (2.B.27) contributes with an amplitude $\langle \Psi_\beta | \mathbb{1} | \Psi_\gamma \rangle \langle \Psi_\gamma | V_\alpha - U_\alpha | \Psi_\alpha \rangle$, where $\mathbb{1}$ is the unit operator, while $\gamma \equiv f (= b + 1) + F (= A + 1)$, denotes the mass partition of the intermediate channel. The above expression indicates that a consistent description of two-nucleon transfer reactions in a non-orthogonal basis involves three strongly

¹⁴⁵For more details see Broglia and Winther (2004), in particular Section V.4 p. 308 and subsequent ones, and references therein.

interweaved reaction channels¹⁴⁶, namely α, γ and β . Consequently it has to be calculated, at least, up to second order of perturbation theory. Thus the need to calculate also $a^{(2)}(t)$, that is to add to simultaneous, also successive transfer and non-orthogonality corrections.

Regarding the sum-rules we consider, for simplicity, the simultaneous transfer amplitude (2.B.26) (see also Sect. 6.5, Eq. (6.5.4)). That is,

$$\begin{aligned} a^{(1)}(t = +\infty) &= \frac{1}{i\hbar} \int_{-\infty}^{\infty} dt \exp\left[\frac{i}{\hbar}(E^{bB} - E^{aA})t + \gamma_{\beta\alpha}(t)\right] \\ &\times \langle \phi^{B(A)}(\mathbf{r}_{1A}, \mathbf{r}_{2A}) | U(r_{1b}) | e^{i\sigma_{\beta\alpha}} \phi^{a(b)}(\mathbf{r}_{1b}, \mathbf{r}_{2b}) \rangle_{\mathbf{R}_{\alpha\beta}(t)}, \end{aligned} \quad (2.B.38)$$

where¹⁴⁷

$$\sigma_{\beta\alpha} = \frac{1}{\hbar} \frac{2m_n}{m_A} (m_{aA}\mathbf{v}_{aA}(t) + m_{bB}\mathbf{v}_{bB}(t)) \cdot (\mathbf{r}_\alpha - \mathbf{r}_\beta), \quad (2.B.39)$$

takes care of recoil effects, the phase factor $e^{i\sigma_{\beta\alpha}}$ being a generalized Galilean transformation associated with the mismatch between entrance and exit channels, while the phase

$$\begin{aligned} \gamma_{\beta\alpha}(t) &= \int_0^t dt' \left\{ U_\beta(\mathbf{R}_\beta(t')) - \frac{1}{2}m_\beta v_\beta^2(t') - U_\alpha(\mathbf{R}_\alpha(t')) + \frac{1}{2}m_\alpha v_\alpha^2(t') \right\} \\ &+ \frac{1}{2\hbar} (m_\alpha \mathbf{v}_\alpha(t) + m_\beta \mathbf{v}_\beta(t)) \cdot (\mathbf{R}_\beta(t) - \mathbf{R}_\alpha(t)), \end{aligned} \quad (2.B.40)$$

is related to the effective Q -value of the reaction.

The rate of change of the formfactor $\langle \phi^{B(A)}, U(r_{1b})e^{i\sigma_{\alpha\beta}}\phi^{a(b)} \rangle$ with time is slow, being completely overshadowed by the rapidly varying phase factor $\exp\left[\frac{i}{\hbar}(E^{bB} - E^{aA})t + \gamma_{\beta\alpha}(t)\right]$. Similar relations concerning recoil and Q -value effects can be obtained from the amplitudes associated to successive and to non-orthogonality terms i.e. $a^{(2)}$ and $a^{(NO)}$ (see Ch. 6, Sect. 6.5).

Summing up, to compare two-nucleon transfer cross sections on equal structural footing, one has to eliminate the kinematical oscillating phase which can completely distort the “intrinsic” (reduced matrix element) value of the two-nucleon cross section. And for that, one has to work on each of the three amplitudes to extract at best the phases which couple relative and intrinsic motion (see also Sect. 2.2.1).

Let us make a parallel with the sum rule associated with electromagnetic decay (Coulomb excitation, γ -decay). The absolute transition probability for absorption (emission) of a photon is measured in sec^{-1}

$$T(E1; I_1 \rightarrow I_2) = (1.59 \times 10^{15}) \times (E)^3 \times B(E1; I_1 \rightarrow I_2), \quad (2.B.41)$$

¹⁴⁶This is also testified by the fact that their formal expression can be shifted around by changing representation, among post-post, prior-prior, and prior-post, see e.g. Eq. (2.B.8) as well as Fig. 6.5.2.

¹⁴⁷See Eqs. (13)–(15) p. 310 Broglia and Winther (2004).

where E is the energy of the transition and $B(E1) = \langle I_2 || \mathcal{M}(E1) || I_1 \rangle / \sqrt{3}$ is¹⁴⁸ the reduced transition probability¹⁴⁹. The TRK-sum rule is written as

$$S(E1) = \langle 0 | [[H, \mathcal{M}(E1)], \mathcal{M}(E1)] | 0 \rangle / 2, \quad (2.B.42)$$

In this case the Q -value dependence of the observed absolute transition probabilities can be eliminated analytically (E^3 dependence), in keeping with the fact that the mass partition ($a + A \rightarrow a + A^*$) does not change between entrance and exit channels or, equivalently, the coordinate of relative motion $\mathbf{R}_{aa'}(t)$ is always that describing the relative center of mass position of the system in the entrance channel.

Expressed it differently, and returning to the expression of the first order (simultaneous) two-nucleon transfer amplitude $a^{(1)}(t = +\infty)$ one can only devise empirical protocols to try to extract the γ - and σ -phase dependence from it and set differential absolute two-nucleon transfer cross sections $d\sigma/d\Omega \sim |a|^2$ on equal footing regarding kinematics, so as to be able to compare intrinsic, reduced transition probabilities (structure). That is, extract the structure information contained in, e.g.,¹⁵⁰

$$\phi^{B(A)}(\mathbf{r}_{1A}, \mathbf{r}_{2A}) = \langle \mathbf{r}_{1A}, \mathbf{r}_{2A} | \Gamma_1^\dagger(\beta = +2) | \tilde{0} \rangle, \quad (2.B.43)$$

as well as in

$$\phi^{B(A)}(\mathbf{r}_{1A}, \mathbf{r}_{2A}) = \langle \mathbf{r}_{1A}, \mathbf{r}_{2A} | [a_k^\dagger a_k^\dagger]_0 | 0 \rangle. \quad (2.B.44)$$

Namely, in the RPA pair addition mode representation and in the pure two-particle configuration $|j_k^2(0)\rangle$ describing two nucleons moving in time reversal states around the close shell system $|0\rangle$. If it was possible to disentangle the γ and σ dependence of $a^{(1)}$ (as well as that of $a^{(2)}$ and $a^{(NO)}$) from its formfactor dependence, the comparison between the quantities¹⁵¹ $\sum_{n,k} |c_k^{(n)}|^2$, and

$\sum_{n,j} \left| \sum_k X_k^{(n)} \delta(j, k) - \sum_i Y_i^{(n)} \delta(j, i) \right|^2$ could eventually be phrased in terms of exact sum rules. This not being the case, one has to deal with approximate TNTR sum rules. With this proviso in mind, such sum rules are quite useful (see Sect. 7.2.2; also Sect. 2.2.1).

Appendix 2.C NFT vacuum polarization

The role zero point fluctuations play in the nuclear ground state, i.e. in the NFT vacuum can be clarified by relating it to the polarization of the QED vacuum. Let

¹⁴⁸ $\mathcal{M}(E1) = e \sum_k \left(\left(\frac{N-Z}{A} - t_Z \right) r Y_{1\mu} \right)_k$.

¹⁴⁹ Bohr and Mottelson (1969) p. 382.

¹⁵⁰ Where $\Gamma_n^\dagger(\beta = +2) = \sum_k X_k^n [a_k^\dagger a_k^\dagger]_0 - \sum_i Y_i^n [a_i^\dagger a_i^\dagger]_0$, is the RPA pair creation addition mode acting on a correlated closed shell system $|\tilde{0}\rangle$.

¹⁵¹ See e.g. Broglia and Riedel (1967).

us briefly dwell on the "reality" of such a phenomenon by recalling the fact that to the question of Rabi of whether the polarization of the QED vacuum could be measured¹⁵², in particular the change in charge density felt by the electrons of an atom, e.g. the electron of a hydrogen atom, due to virtual creation and annihilation of electron-positron pairs, Lamb gave a quantitative answer, both experimentally and theoretically¹⁵³. The corresponding correction (Lamb shift) implies that the $2s_{1/2}$ level lies higher than the $2p_{1/2}$ level by about 1000 MHz.

In connection with the discussion of Feynman of vacuum polarisation, where a field produces a pair, the subsequent pair annihilation producing a new field, namely a close loop, he implemented in his space-time trajectories, Wheeler's idea of electrons going backwards in time (positrons). Such trajectories would be like an **N** in time, that is electrons which would back up for a while, and go forward again. Being connected with a minus sign, these processes are associated with Pauli principle in the self-energy of electrons (see Fig. 2.4.2 (f)). The divergences affecting such calculations could be renormalised by first computing the self-energy diagram in second order and finding the answer which is finite, but contains a cut-off to avoid a logarithmic divergence¹⁵⁴. Expressing the result in terms of the experimental mass, one can take the limit (cut-off $\rightarrow \infty$) which now exists (see also Sect. 5.3.3).

In the nuclear case, for example Skyrme effective interactions give rise to particle-vibration coupling vertices which, because of the contact character of these interactions may lead to divergent zero point energies¹⁵⁵, unless a cut-off is introduced¹⁵⁶. The Gogny force being finite range does not display such problems. The question still remains concerning the stability of the results to be obtained by carrying out a complete summation over both collective and non collective contributions. Likely, one ensures it by going to higher orders in the oyster diagrams (see Figs. 1.8.1 (a) and (b), 1.8.3 (a), 2.4.2 (a), 2.5.1 and 2.6.2, as well as Sect. 6.3). The fermion exchange between two of these diagrams (Pauli principle) eliminates non-collective contributions, leading eventually to convergent results.

An economic and reliable method to achieve a similar result, is that of using renormalization. That is, to calculate the lowest order diagrams but introducing, in the intermediate states, the dressed physical (empirical) states.

¹⁵²Pais (1986) pp. 450, 451; Pais (2000) pp. 255-267.

¹⁵³Lamb Jr and Rutherford (1947); Kroll and Lamb (1949)

¹⁵⁴See Bethe (1947); Feynman (1961); Weinberg (1996a) p. 578; see also Mehra (1996) p. 295, and Bjorken and Drell (1998).

¹⁵⁵See Hellemans et al. (2013); Pastore et al. (2015) and refs. therein.

¹⁵⁶The question emerges of which are the provisos to be taken in the use of effective forces to higher orders in the PVC. Within this context see Mahaux et al. (1985), also Broglia et al. (2016); Barranco et al. (2017) concerning the implementation of renormalization in both configuration and 3D-spaces within the framework of NFT. In a nutshell, the bare mean field exists but its properties cannot be measured (not any more than the bare electron mass in renormalized quantum electrodynamics), and corresponds to a set of parameters of a Fermi-like function which ensure that the dressed states reproduce all of the experimental findings, for both structure and reactions.

Chapter 3

Collective pairing modes

I wrote down the trial ground state as a product of operators –one for each pair state– acting on the vacuum... $\Psi_0 = \prod_{\mathbf{k}} (U_{\mathbf{k}} + V_{\mathbf{k}} b_{\mathbf{k}}^\dagger) |0\rangle$... Since the pair creation operators $b_{\mathbf{k}}^\dagger$ commute for different \mathbf{k} 's... Ψ_0 represents uncorrelated occupancy of the various pair states.... The operators $b_{\mathbf{k}}^\dagger = c_{\mathbf{k}\uparrow}^\dagger c_{\mathbf{k}\downarrow}^\dagger$, being a product of two fermions (quasiparticle) creation operators, do not satisfy Bose statistics, since $b_{\mathbf{k}}^{\dagger 2} = 0$.

J. R. Schrieffer

3.1 Nuclear Structure in a nutshell

The low-energy properties of the finite, quantal, many-body nuclear system, in which nucleons interact through the strong force of strength $v_0 (\approx -100 \text{ MeV})$ and range $a (\approx 0.9 \text{ fm})$ are controlled, in first approximation, by independent particle motion. This is a consequence of the fact that nucleons display a sizable value of the quantal zero point (kinetic) energy of localization ($\hbar^2/Ma^2 \approx 50 \text{ MeV}$) as compared to the absolute value of the strength of the NN -potential¹ $|v_0| = 100$

¹ The corresponding ratio $q = \left(\frac{\hbar^2}{Ma^2}\right) \frac{1}{|v_0|}$ is known as the quantality parameter and was first used in connection with the study of condensed matter (de Boer (1948, 1957); de Boer and Lundbeck (1948); Nosanow (1976)). It was introduced in nuclear physics in Mottelson (1998) where its value $q = 0.5$ testifies to the validity of independent particle motion. It is of notice that questions like the one posed in connection with localization and long mean free path were already discussed by Lindemann (1910) in connection with the study of the stability or less of crystals. The generalization to aperiodic crystals, like e.g. proteins (Schrödinger (1944)) was carried out in Stillinger and Stillinger (1990).

MeV.

The corresponding ground state $|HF\rangle = \Pi_i a_i^\dagger |0\rangle$ describes a step function in the probability of the occupied ($\epsilon_i \leq \epsilon_F$) and empty ($\epsilon_k > \epsilon_F$) states, displaying a sharp discontinuity at the Fermi surface. Pushing the system it reacts with an inertia AM , sum of the nucleon masses (App. 2.B). Setting it into rotation, assuming the density $\rho(r) = \sum_i |\langle \mathbf{r}|i\rangle|^2$ ($|i\rangle = a_i^\dagger |0\rangle$) to be spatially deformed, it responds with the rigid moment of inertia. This is because the single-particle orbitals are solidly anchored to the mean field (Fig. 3.3.2).

Pairing acting on nucleons moving in time reversal states v, \bar{v} ($v \equiv (nlj)$), in configurations of the type $((l)_{L=0}^2, (s)_{S=0}^2)$, and lying close to the Fermi energy, alter the above picture in a conspicuous way². Within an energy range of the order of the absolute value of the pair correlation energy³ $|E_{corr}| (\approx 3\text{ MeV})$ centered around ϵ_F ($|E_{corr}|/\epsilon_F \ll 1$), the role of independent particles is taken over by pairs of nucleons, correlated over distances $\xi \approx \hbar v_F / (\pi \Delta) (\approx 14 \text{ fm}; \Delta = 1.2 \text{ MeV})$, and being equally phased (see Sect. 3.4). In other words, and bringing the argument to a previous step, one can posit that since the creation operators a_i^\dagger commute for different i 's, $|HF\rangle$ represents uncorrelated occupancy of the different single-particle states. Similarly, because the pair operators P_v^\dagger commute for different v 's, $|BCS\rangle$ represents uncorrelated occupancy of the various pair states. The (single-particle)-Cooper pair⁴ parallel is summarized in Fig. 3.4.1.

For intrinsic⁵ nuclear excitation energies and rotational frequencies⁶ sensibly smaller than $|E_{corr}/2|$ and than $\hbar\omega_{rot} \approx 0.5 \text{ MeV}$ respectively, the system can be described in terms of gauge phase correlated, pair motion. This is a consequence of the fact that the kinetic energy of (Cooper) pair confinement ($\hbar^2/(2M\xi^2) \approx 10^{-2} \text{ MeV}$), is much smaller than the absolute value of the pair binding energy $|E_{corr}|$, implying that each pair behaves as an entity⁷ of mass $2M$ and intrinsic structure

Its possible application to the atomic nucleus is discussed in App. 3.A

²Bohr et al. (1958); see also Broglia and Zelevinsky (2013) and references therein.

³In BCS, $E_{corr} \approx -\frac{N(0)}{2} \Delta^2$, where $N(0) = \frac{g}{2}$ is the density of nuclear states at the Fermi energy and for one spin orientation, $g_i = i/16 \text{ MeV}^{-1}$ ($i = N, Z$) being the result of an empirical estimate which takes surface effects into account (Bohr and Mottelson (1975) Eq. (2-47); Bortignon et al. (1998) Eq. (7.15)), while Δ is the pairing gap, 2Δ being the binding energy of the Cooper pair. For a typical superfluid, quadrupole deformed nucleus like ^{170}Yb , $N(0) = 5.3 \text{ MeV}^{-1}$, $\Delta \approx 1.1 \text{ MeV}$ and $E_{corr} = -3.2 \text{ MeV}$ (Shimizu et al. (1989)).

⁴Cooper (1956).

⁵As opposed to collective excitations which do not alter the temperature of the system, as all the energy is concentrated in a single mode.

⁶Coriolis force acts oppositely on each member of a Cooper pair. When the difference in rotational energy between superfluid and normal rotation becomes about equal to the correlation energy, the nucleon moving opposite to the collective rotation becomes so much retarded in its revolution period with respect to the partner nucleon, that eventually it cannot correlate any more with it and “align”, its motion (and spin) with the rotational motion, becoming again a pair of fermions and not participating in the BCS condensate. This happens for a (critical) angular momentum $I_c \approx (120 \times |E_{corr}|)^{1/2} \approx 20\hbar$, corresponding to a rotational frequency $\hbar\omega_c \approx 0.5 \text{ MeV}$ (see Bohr and Mottelson (1975), Broglia and Zelevinsky (2013) and references therein).

⁷The ratio $q_\xi = \frac{\hbar^2}{2M\xi^2} \frac{1}{|E_{corr}|} \approx 0.02$ provides a generalized quantality parameter. It testifies to the stability of nuclear Cooper pairs in superfluid nuclei.

$L = S = 0$. Cooper pairs behave like (quasi) bosons, the single-particle orbits on which they correlate become dynamically detached from the mean field, leading to a BCS condensate. It is however very different from a standard condensate of real bosons. Within this context see Sect. 3.4 (Figs. 3.3.3 and 3.3.4) and App. 4.A.

Cooper pairs exist also in situations in which the environmental conditions are above critical. For example, in metals at room temperature, in closed shell nuclei as well as in deformed open shell ones at high values of the angular momentum⁸. However, in such circumstances, they break essentially as soon as they are generated (pairing vibrations). While these pair addition and subtraction fluctuations have little effect in condensed matter systems with the exception than at⁹ $T \approx T_c$ (critical temperature), they play an important role in normal (non-superfluid) nuclei. In particular around closed shells¹⁰ (Fig. 3.1.1), and specially in the case of light, highly polarizable, exotic halo nuclei¹¹. From this vantage point one can posit that it is not so much, or, at least not only, the superfluid phase which is abnormal in the nuclear case, but the normal state around the closed shell systems, displaying a consistent component of dynamical abnormal density. The importance of pairing vibrations is also found in connection with the self-energy of nucleons moving around closed shells¹² as well as concerning the binding energy of the closed shell systems¹³.

Furthermore, it is of notice the role pairing vibrations play in the transition between superfluid and normal nuclear phases (cf. Fig. 3.1.2) as a function of the rotational frequency (angular momentum) as emerged from the experimental studies of high spin states carried out by, among others, Garrett and collaborators¹⁴. While the (dynamic) pairing gap associated with pairing vibrations leads to $\approx 20\%$ increase of the static pairing gap for low rotational frequencies, it becomes the overwhelming contribution above the critical frequency¹⁵.

The central role played by pairing vibrations within the present circumstances is that to restore particle-number conservation, another example after that provided by the quantity parameter and by its generalization to pair motion, of the fact that potential functionals are, as a rule, best profited by special arrangements of particles (spontaneous symmetry breaking), while fluctuations favor symmetry¹⁶.

⁸See e.g. Shimizu and Broglia (1990) and references therein.

⁹See Schmidt (1968), Schmid (1969) Abrahams and Woo (1968); concerning superfluid ^3He see Wölfe (1978).

¹⁰See Sects. 3.5 and 3.6; Bohr (1964), Bès and Broglia (1966), Högaasen-Feldman (1961), Schmidt (1972), Schmidt (1968), Barranco et al. (2001), Potel et al. (2013a), Potel et al. (2014), Schmidt (1964).

¹¹See Potel et al. (2013a) and refs. therein. Also Potel et al. (2013b) in connection with the closed shell system ^{132}Sn .

¹²See e.g. Bès and Broglia (1971a); Flynn et al. (1971).

¹³Baroni et al. (2004).

¹⁴Garrett (1985); Garrett et al. (1986), see also Shimizu et al. (1989), Barranco et al. (1987) and Ch. 6 of Brink and Broglia (2005).

¹⁵Shimizu et al. (1989), Shimizu and Broglia (1990), Shimizu (2013), Dönau et al. (1999), Shimizu et al. (2000).

¹⁶Anderson and Stein (1984); Anderson (1976).

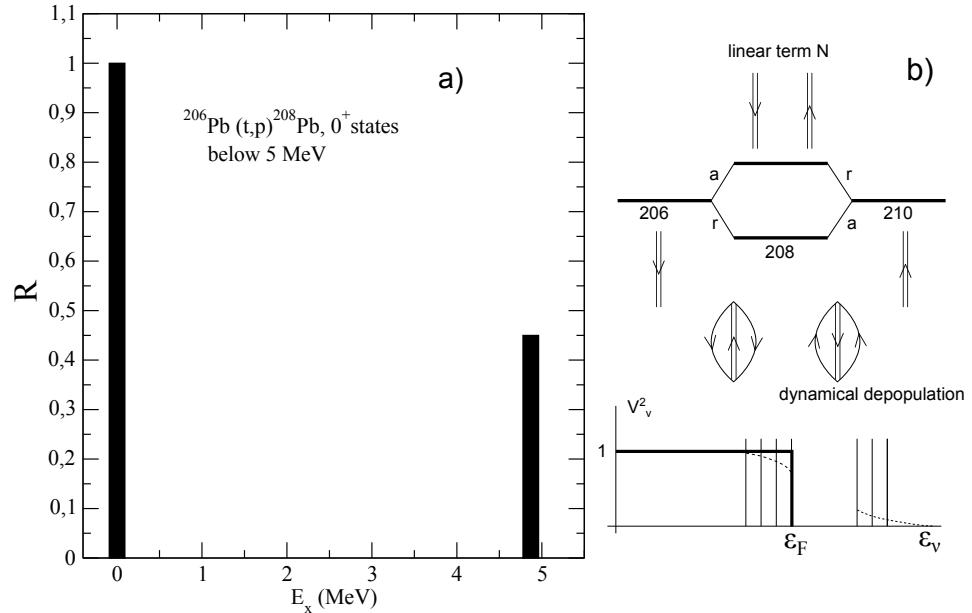


Figure 3.1.1: (a) Ratio of the absolute $L = 0$ differential cross sections $d\sigma(E_x, \theta = 59^\circ)/d\sigma(gs, \theta = 59^\circ)$ ($= (0.05 \text{ mb/sr})/(0.12 \text{ mb/sr})$) below 5 MeV for the reaction $^{206}\text{Pb}(t, p)^{208}\text{Pb}$ at the second maximum ($\theta = 59^\circ$; Bjerregaard et al. (1966)). It is of notice the large experimental errors of the corresponding angular distributions associated with the poor statistics of the cross section at the first maximum $\theta = 5^\circ$. This is the reason why the maximum at 59° was preferred to report the ratio of the cross sections. (b) Schematic representation of the pairing vibrational spectrum around ^{208}Pb . Also shown is a cartoon representation of the softening of the sharp mean field Fermi surface due to the ZPF of the pairing vibrational modes. The label a and r indicate the pair addition and pair removal modes. It is to be noted that a linear term in N has been added to the binding energy to make the binding energy values associated with ^{206}Pb ($N = 124$) and ^{210}Pb ($N = 128$), equal, in an attempt to emphasize a harmonic picture for the two-phonon state. Concerning the anharmonicities of the modes see last paragraph Sect. 3.5.

Within this context see Eqs. (1.4.55) and (1.4.59).

There are a number of methods which allows one to go beyond BCS mean-field approximation, or of its generalization known as the Hartree-Fock-Bogolyubov approximation (HFB). Generally referred to as number projection methods¹⁷(NP), they make use of a variety of techniques like Generator Coordinate Method, etc. as well as protocols like variation after projection, etc. The advantages of NP methods over the RPA is to lead to smooth functions for both the correlation energy and the pairing gap at the pairing phase transition between normal and superfluid phases.

In Fig. 3.1.3 we display the excitation function associated with the reaction $^{122}\text{Sn}(p, t)^{120}\text{Sn}(J^\pi)$, populating the low-energy states of the single open shell superfluid nucleus ^{120}Sn . The angle selected to report the value of the absolute differential cross sections, that is 5° , corresponds to the first, and largest, peak of the absolute $L = 0$ differential two-nucleon transfer cross section. Essentially all the strength in the $L = 0$ channel is concentrated in the ground state, the strongest 0^+ -excited state carrying a cross section of the order of 3% of that of the ground state. Within this context, the difference with the results displayed in Fig. 3.1.1 is apparent¹⁸.

In the inset to Fig. 3.1.3 a quantity closely related to the Sn-isotopes binding energy is displayed (bold face levels). Namely $B(^{50+N}\text{Sn}_N)-8.124N$ MeV+46.33 MeV, resulting from the subtraction of the contribution of the single nucleon addition to the nuclear binding energy (linear function in N) obtained by a linear fitting of the binding energies of all the Sn-isotopes. Also displayed is the parabolic fit to these energies, a quantity to be compared with $E_N = (\hbar^2/2I)(N - N_0)^2$, namely the energy associated with the members of the pairing rotational band. The difference with the spectrum of pair addition and subtraction modes displayed in Fig. 3.1.1 b) is again evident. Concerning the parallel one can draw between the 3D- and pairing-vibrations and -rotations, see Fig. 3.1.5.

3.1.1 Pairing rotational and vibrational bands probed with one- and two-nucleon transfer

A simple estimate of the pairing rotational band moment of inertia is provided by the single j -shell model¹⁹, namely $(\hbar^2/2I)=G/4 \approx 25/(4N_0)$ MeV. This estimate turns out to be rather accurate. Certainly better than one can expect. On the other hand, one is reminded of the fact that we are discussing properties which specifically characterize a coherent state²⁰, namely $|BCS\rangle$ (see Eq. (1.4.39)).

¹⁷cf. Ring and Schuck (1980), Egido (2013), Robledo and Bertsch (2013); cf. also Frauendorf (2013), Ring (2013), Heenen et al. (2013), and references therein.

¹⁸ While this “distortion” of the (t, p) excitation function is useful to emphasize the parallels between vibrational and rotational bands in 3D-and in gauge-spaces, it has to be used with care concerning the parallel with Cooper pair tunneling between weakly coupled superconductors (Sect. 4.6 and in particular, Sect. 7.3). This is also in keeping with the fact that at the basis of nuclear BCS one finds an interdisciplinary connection (see Bohr et al. (1958)).

¹⁹See e.g. Brink and Broglia (2005) App. H and refs. therein.

²⁰See Sect. 7.2; see also Potel et al. (2017).

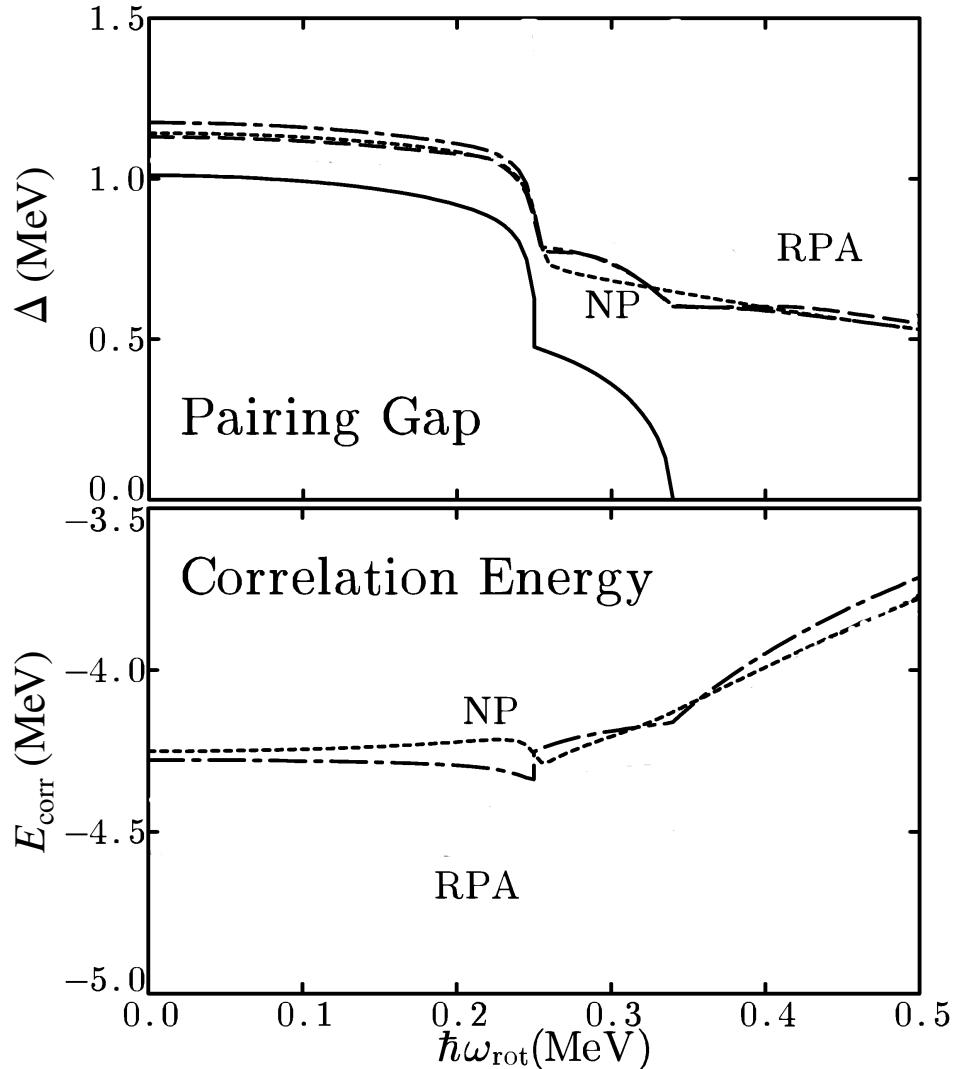


Figure 3.1.2: Pairing gap calculated taking into account the correlation associated with pair vibrations in the RPA approximation ($\Delta = (\Delta_{BCS}^2 + \frac{1}{2}G^2S_0(RPA))^{1/2}$) (upper panel) and RPA correlation energy (lower panel) for neutrons in ^{164}Er as a function of the rotational frequency (Shimizu et al. (1989); Shimizu (2013)). Both quantities are in MeV (*dashed-dotted curves*). The value of the static (mean-field) pairing gap Δ , which vanishes at $\hbar\omega_{\text{rot}} = 0.34$ MeV, is also displayed in the upper panel (continuous curve). The results of the number-projection (NP) calculations are shown as dotted curves; $S_0(RPA) = \sum_{n \neq AGN} [< n|P|0 > + < n|P^\dagger|0 >]_{RPA}^2$, $P^\dagger = \sum_{\nu>0} a_\nu^\dagger a_{\bar{\nu}}^\dagger$; $\Delta_{BCS} = G < BCS | P^\dagger | BCS >$ (*continuous line*) is the standard, static BCS pairing gap, while G is the pairing force strength. The non-energy weighted sum rule $S_0(RPA)$ describes the contribution of pairing fluctuations to the effective (RPA) gap, and is intimately associated with projection in particle number. It is of notice that $\sum_{n \neq AGN}$ means that the divergent contribution from the zero energy mode (Anderson, Goldstone, Nambu mode, see e.g. Broglia et al. (2000) and references therein), associated with the lowest ($\hbar\omega_0$) solution of the $H = H_p + H_p''$ is to be excluded (see Sect 4.7.1, discussion before Eq. (4.7.26) as well as Brink and Broglia (2005) App. J).

Also reported in Fig. 3.1.3 (inset), are the integrated values of the measured absolute two-neutron transfer cross sections, quantities which are reproduced by the theoretical predictions within experimental errors (Fig. 3.1.7). In principle, one could have expected a sensible constancy of these cross sections (transitions) as the pairing rotational model implies a common intrinsic deformed state in gauge space, namely $|BCS\rangle$ (see Eq. (1.4.65)). On the other hand, the number of Cooper pairs α'_0 which defines deformation in gauge space is rather small (≈ 6) and thus subject to conspicuous fluctuations ($\Delta\alpha'_0/\alpha'_0 \approx \sqrt{6}/6 \approx 0.4$). Because $\sigma \sim \alpha'^2_0$, fluctuations in σ of the order of 100%, can be expected.

In keeping with the analogies discussed in connection with Figs. 3.1.5 and 3.3.2 between pairing and quadrupole rotational bands, we note that in the electromagnetic decay of these last bands one expects, in the case of heavy nuclei, fluctuations of the order of $(\sqrt{200}/200)$, in keeping with the magnitude of the associated $B(E2)$ values, when measured in terms of single-particle units. Within this context, the average value of the absolute experimental cross sections displayed in the inset of Fig. 3.1.3 is $1762 \mu\text{b}$, while the average difference between experimental and predicted values²¹ is $94 \mu\text{b}$. Thus, the discrepancies between theory and experiment are bound in the interval $0 \leq |1 - \sigma_{th}(i \rightarrow f)/\sigma_{exp}(i \rightarrow f)| \leq 0.09$, the average discrepancy being 5%.

In Fig. 3.1.4 the excited pairing rotational bands based on 0^+ pairing vibrational modes are displayed as a single band, resulting from the average value of the 0^+ excited states with energy ≤ 3 MeV. The best parabolic fit is shown. Also given are the relative (p, t) absolute integrated cross sections normalized to the corresponding values of the ground state rotational band. The cross talk between bands is in all cases $\leq 8\%$, the single j -shell value estimate being²² 6%.

The above results underscore the fact that, at the basis of an operative coarse grained approximation to the nuclear many-body problem, one finds a judicious choice of the collective coordinates²³ which in turn is closely connected with the probe used to study the system. In other words, pairing vibrations are elementary modes of excitation containing the right physics to restore gauge invariance through their interweaving with quasiparticle states (Eqs. (1.4.59)–(1.4.62)). They project from the deformed state ($\alpha'_0 \neq 0$), the members (Eq. (1.4.65)) of pairing rotational bands which, similarly to pairing vibrational bands, are specifically excited in two-nucleon transfer reactions. Examples are provided by Cooper pair tunneling in heavy ion collisions between superfluid nuclei, as well as (t, p) , (p, t) , etc reactions²⁴. From other fields of physics, the Josephson²⁵ effect between weakly coupled metallic superconductors is rightly considered the paradigmatic example.

²¹Potel et al. (2013b).

²²Brink and Broglia (2005) App. H.

²³In this connection, we quote allegedly from S. Weinberg: “In solving a problem you may choose to use the degrees of freedom you like. But if you choose the wrong ones you will be sorry”.

²⁴See e.g. Yoshida (1962), Glendenning (1965), Bohr (1964), Bayman (1971), Broglia et al. (1973), Hansen (2013) and Potel et al. (2013a) and references therein.

²⁵Josephson (1962).

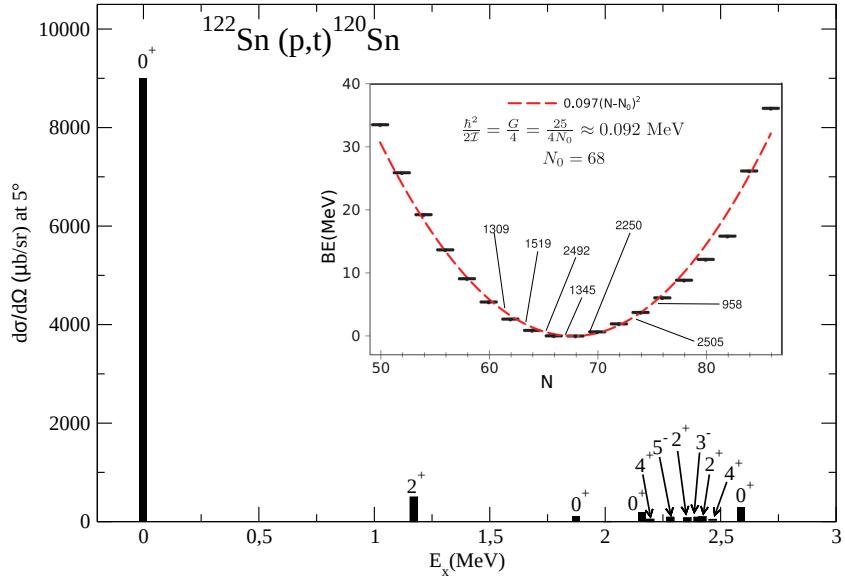


Figure 3.1.3: Excitation function associated with the reaction $^{122}\text{Sn}(p, t)^{120}\text{Sn}(J^\pi)$. The absolute experimental values of $d\sigma(J^\pi)/d\Omega|_{5^\circ}$ are given as a function of the excitation energy E_x (after Guazzoni et al. (2011)). In the inset the neutron pairing rotational band between magic numbers $N = 50$ and $N = 82$ is displayed (see also caption to Fig. 1.4.5), the absolute $^{A+2}\text{Sn}(p, t)^A\text{Sn}$ experimental cross sections taken from Guazzoni et al. (1999), Guazzoni et al. (2004), Guazzoni et al. (2006), Guazzoni et al. (2008), Guazzoni et al. (2011), Guazzoni et al. (2012), are also given in μb (number connected by lines to the corresponding levels of the pairing rotational band).

Within this context one can now take the basic consequence of pairing condensation in nuclei regarding reaction mechanisms. For this purpose let us consider a *gedankenexperiment* in which the superfluid target and the projectile can at best come in weak contact. Because $(\hbar^2/2M\xi^2)/|E_{corr}| \approx 10^{-2}$, Cooper pairs in superfluid nuclei behave as (potentially) very extended entities of mean square radius ($\xi \gg R_0$) and mass $2M$, even in the case in which the 1S_0 , NN -potential vanishes in the zone between the weakly overlapping densities of the two interacting nuclei. One then expects Cooper pair transfer to be observed. Not only. One also expects the associated absolute differential cross section to be of the same order of magnitude than one-nucleon transfer ones, and to be dominated by successive

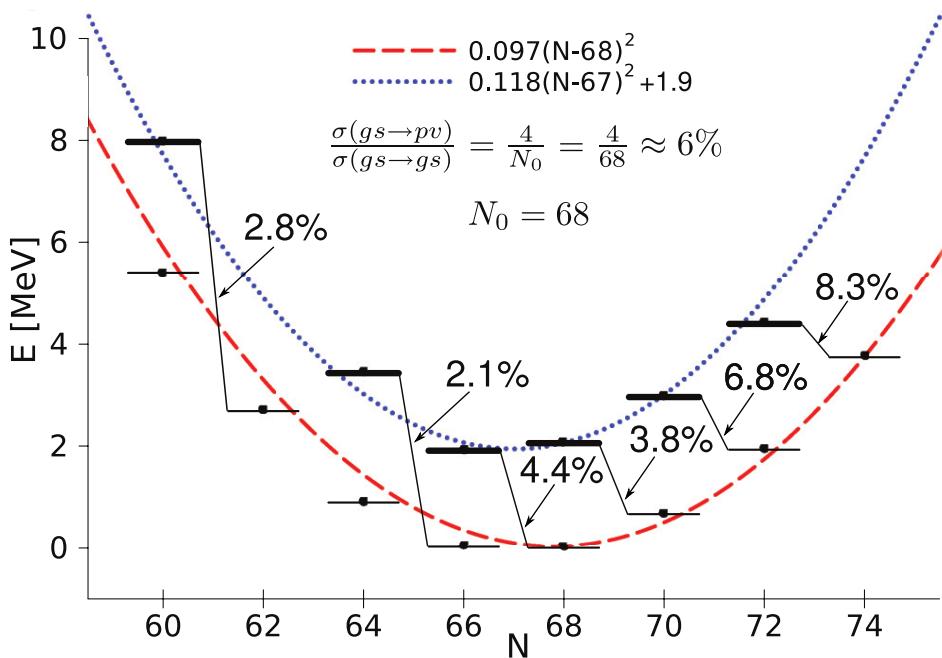


Figure 3.1.4: The weighted average energies ($E_{exc} = \sum_i E_i \sigma_i / \sum_i \sigma_i$) of the excited 0^+ states below 3 MeV in the Sn isotopic chain are shown on top of the ground state pairing rotational band, already displayed in Fig. 3.1.3 in the case of ^{120}Sn . Also indicated is the percentage of cross section for two-neutron transfer to excited states, normalized to the cross sections populating the ground states. The estimate of the ratio of cross sections displayed on top of the figure was obtained making use of the single j -shell model (Brink and Broglia (2005) App. H).

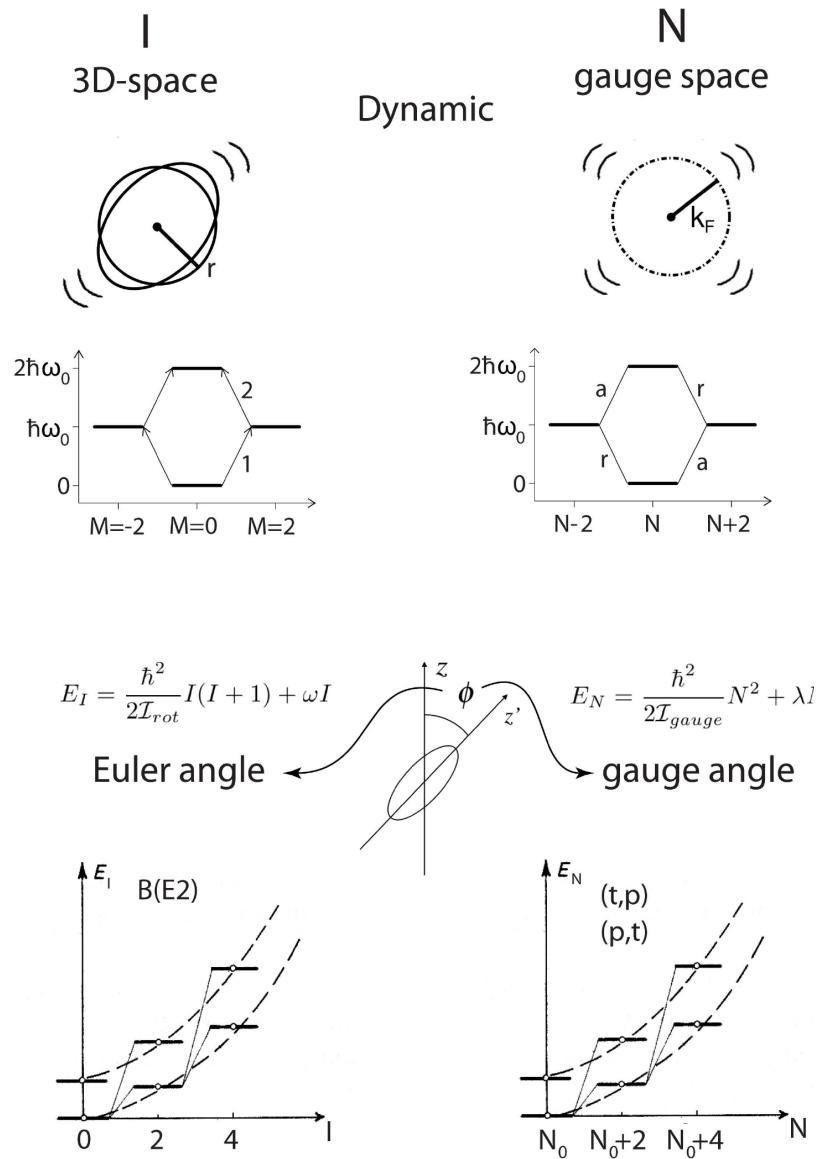


Figure 3.1.5: Parallel between dynamic and static deformations in 3D- and in gauge-space for the nuclear finite many-body system. In the first case, the angular momentum \mathbf{I} and the Euler angles are conjugate variables. In the second, particle number N and gauge angle. While the fingerprint of static (quadrupole and gauge) deformations are quadrupole and pairing rotational bands, vibrational bands are the expression of such phenomena in non deformed systems.

transfer. These expectations have been confirmed experimentally²⁶. The above parlance, being at the basis of the Josephson effect, reflects both one of the most solidly established results in the study of BCS pairing, and explains the workings of a paradigmatic probe of spontaneous symmetry breaking phenomena.

Due to the fact that, away from the Fermi energy pair motion becomes independent particle motion, one-particle transfer reactions like e.g. (d, p) and (p, d) can be used together with (t, p) and (p, t) processes, as valid tools to cross check pair correlation predictions. In particular, to shed light on the origin of pairing in nuclei: in a nutshell, the relative importance of the bare NN -interaction and the induced pairing interaction.

While the calculation of two-nucleon transfer spectroscopic amplitudes and differential cross sections are, a priori, more involved to be worked out than those associated with one-nucleon transfer reactions, the former are, as a rule, more “intrinsically” accurate than the latter ones. This is because, in the case of two nucleon transfer reactions, the quantity (order parameter α'_0) which expresses the collectivity of the members of a pairing rotational band, reflects the properties of a coherent state ($|BCS\rangle$). In other words, it results from the sum over many two-nucleon transfer spectroscopic amplitudes ($\sqrt{j_\nu + 1/2} U'_\nu V'_\nu$), all phased in the same way (phase correlated).

There is a further reason which confers $\alpha'_0 = \sum_j (j + 1/2) U'_j V'_j$ a privileged position with respect to the single contributions $(j + 1/2) U'_j V'_j$. It is the fact that $\alpha'_0 = e^{2i\phi} \sum_j (j + 1/2) U_j V_j = e^{2i\phi} \alpha_0$ defines a privileged orientation in gauge space, α_0 being the order parameter referred to the laboratory system which makes an angle ϕ in gauge space with respect to the intrinsic system to which α'_0 is referred²⁷. In other words, the quantities α'_0 measures the deformation of the superfluid nuclear system in gauge space.

Similar arguments can be used regarding the excitation of pairing vibrations in terms of Cooper pair transfer around closed shells as compared to one-particle transfer. As seen from Fig. 3.1.6 (c)–(b), the random phase approximation (RPA) amplitudes X_ν^a and Y_ν^a are coherently summed over pairs of time reversal states to give rise to the spectroscopic amplitudes associated with the direct excitation of the pair addition mode displayed in (d). Because of the (dispersion) relation

²⁶ Both in heavy ion reaction between superfluid nuclei (Montanari et al. (2014), see also Fig. 4.6.1 and Sect. 7.3) and light ions (d, p) and (t, p) reaction. See e.g. Cavallaro et al. (2017) $d\sigma(^9\text{Li}(d, p)^{10}\text{Li}(1/2^-))/d\Omega|_{\theta_{max}} \approx 0.8$ mb/sr, as compared to Tanihata et al. (2008) $d\sigma(^{11}\text{Li}(p, t)^9\text{Li}(1/2^-))/d\Omega|_{\theta_{max}} \approx 1$ mb/sr, Fortune et al. (1994) $^{10}\text{Be}(t, p)^{12}\text{Be}(\text{gs})$ ($\sigma = 1.9 \pm 0.5$ mb, $4.4^\circ \leq \theta_{cm} \leq 54.4^\circ$) as compared to Schmitt et al. (2013) $^{10}\text{Be}(d, p)^{11}\text{Be}(1/2^+)$ ($\sigma = 2.4 \pm 0.013$ mb, $5^\circ \leq \theta_{cm} \leq 39^\circ$) in the case of light nuclei around closed ($N = 6$) shell, and Bassani et al. (1965) $^{120}\text{Sn}(p, t)^{118}\text{Sn}(\text{gs})$ ($\sigma = 3.024 \pm 0.907$ mb, $5^\circ \leq \theta_{cm} \leq 40^\circ$) as compared to Bechara and Dietzsch (1975) $^{120}\text{Sn}(d, p)^{121}\text{Sn}(7/2^+)$ ($\sigma = 5.2 \pm 0.6$ mb, $2^\circ \leq \theta_{cm} \leq 58^\circ$).

²⁷ It is of notice that the square root of $\alpha_0 = e^{-2i\phi} \alpha'_0$, normalized to a proper volume element V , that is, $\Psi = e^{-i\phi} \sqrt{n_S}$, where $n_S = \alpha'_0/V$ is the density of superconducting electrons, constitutes the order parameter at the basis of the Ginzburg-Landau theory of superconductivity (Ginzburg and Landau (1950)), called by Ginzburg the Ψ -theory of superconductivity (Ginzburg (2004)), published seven years before than BCS.

(b)+(c) \equiv (d), the X_ν - and Y_ν -amplitudes are, in the Cooper pair transfer process, correlated, the situation being not so in the case of one-particle transfer (see e.g. Fig. 2.4.2 (e), (f)). In other words, while the (renormalized) spectroscopic amplitude and (renormalized) radial form factor associated with a one-nucleon transfer process depend on a single renormalized energy and associated wavefunction, the corresponding quantities for Cooper pair transfer depend on a distribution of (renormalized) states and wavefunctions of levels around the Fermi energy, leading to a coherent state²⁸.

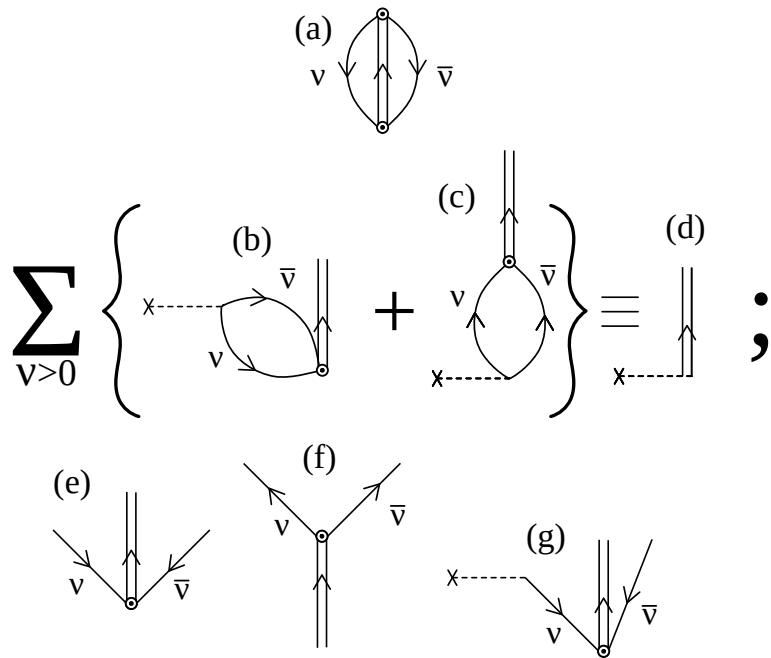


Figure 3.1.6: NFT diagrams associated with one- and two-particle transfer from closed shell associated with ZPF of pair addition modes. (a) ZPF associated with the virtual excitation of a pair addition mode and two uncorrelated holes. (b) two-particle transfer filling a two hole state associated with the ZPF, (c) two-particle transfer to time-reversal states lying above the Fermi energy. These processes receive contribution from all (v, \bar{v}) pairs (sum over $v > 0$), leading to (d), the direct excitation of the collective pair addition mode. The relation (b)+(c) \equiv (d) is the NFT graphical representation of the random phase approximation (RPA) dispersion relation used to calculate the properties of the pair addition mode in the harmonic approximation (Section 3.5). The backwards and forwards going RPA amplitudes are displayed in (e) and (f) respectively. (g) one-particle transfer proceeding through the filling of a hole associated with the ZPF.

²⁸We note that, in spite of the fact that one is dealing with the connection between structure and direct transfer reactions, no mention has been made of spectroscopic factors in relation with one-particle transfer processes, let alone when discussing two-particle transfer processes.

We write now, in somewhat more detail, the spectroscopic amplitudes (B -coefficients) for the two-nucleon transfer processes mentioned above. Namely, one for the case in which A and $B (= A + 2)$ are members of a pairing rotational band. A second one, for the case in which they are members of a pairing vibrational band. That is,

$$\begin{aligned} \text{1), } B((nlj)^2(0)) &= \langle BCS(N+2) | \frac{[a_{nlj}^\dagger a_{nlj}^\dagger]_0^0}{\sqrt{2}} | BCS(N) \rangle \\ &= \sqrt{j+1/2} U_{nlj}(N) V_{nlj}(N+2), \end{aligned} \quad (3.1.1)$$

and

$$\begin{aligned} \text{2), } B((nlj)^2(0)) &= \langle (N_0 + 2)(gs) | \frac{[a_{nlj}^\dagger a_{nlj}^\dagger]_0^0}{\sqrt{2}} | N_0(gs) \rangle \\ &= \begin{cases} \sqrt{j_k + 1/2} X^a(n_k l_k j_k) & (\epsilon_{j_k} > \epsilon_F) \\ \sqrt{j_k + 1/2} Y^a(n_i l_i j_i) & (\epsilon_{j_i} \leq \epsilon_F). \end{cases} \end{aligned} \quad (3.1.2)$$

Where the X and Y coefficients are the forwardgoing and backwardsgoing RPA amplitudes of the pair addition mode. For actual numerical values, see Tables 3.1.1 and 3.5.2–3.5.5. Making use of these spectroscopic amplitudes and treating explicitly the intermediate deuteron channel in terms of successive transfer, correcting both this and the simultaneous transfer channels for non-orthogonality contributions, makes the above picture the quantitative probe of Cooper pair correlations in nuclei²⁹ (Fig. 3.1.7).

Summing up, one will use throughout the present monograph, exception made when explicitly mentioned, absolute cross sections as the link between spectroscopic amplitudes and experimental observations³⁰.

3.2 Renormalization and spectroscopic amplitudes

As a result of the interweaving of single-particle and collective motion, the nucleons acquire a state dependent self energy $\Delta E_j(\omega)$. Consequently, the single-particle potential which was already non-local in space (exchange potential, related to the Pauli principle) becomes also non-local in time. A possible technique to make it local is that of the effective mass approximation. In it one describes the single-particle motion in terms of a local potential given by $U'(r) = (m/m^*)U(r)$. The quantity $m^* = m_k m_\omega / m$ is the effective nucleon mass, m_k being the so-called k -mass associated with the Fock potential, and $m_\omega = m(1 + \lambda)$ being the ω -mass (non-locality in time, as implied by the relation $\Delta\omega\Delta t \geq 1$). The quantity

²⁹Bayman and Chen (1982) and Potel et al. (2013a), see also Götz et al. (1975) and Broglia et al. (1977).

³⁰It is of notice that concerning nuclear structure information, it is also contained in the optical potential used in the calculation of the corresponding cross sections. See footnote 23 Ch. 1.

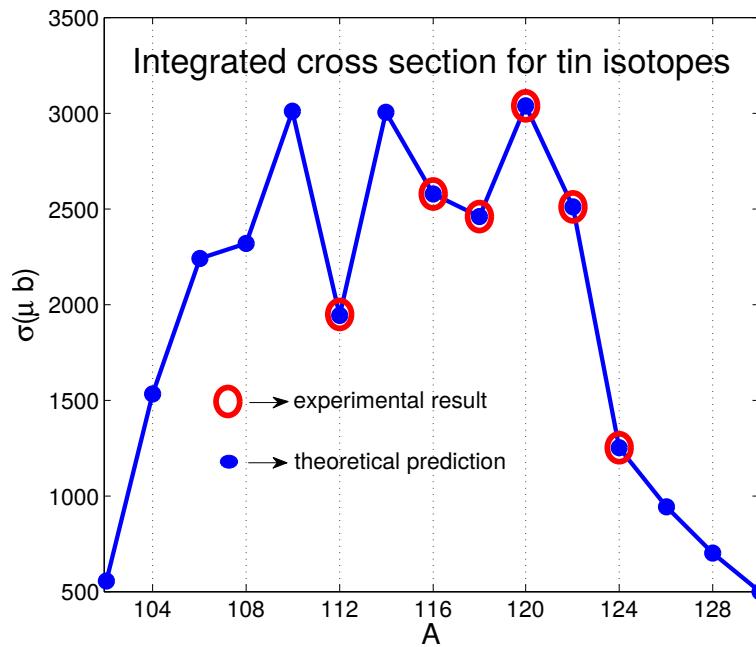


Figure 3.1.7: Absolute value of the two-nucleon transfer cross section $^{A+2}\text{Sn}(p, t)^A\text{Sn}(\text{gs})$ ($A = 112, 116, 118, 120, 122, 124$ cf. Potel et al. (2013a) Potel et al. (2013b); see Sect. 7.2) calculated taking into account successive and simultaneous transfer in second order DWBA, properly corrected for non-orthogonality contributions in comparison with the experimental data (Guazzoni et al. (1999), Guazzoni et al. (2004), Guazzoni et al. (2006), Guazzoni et al. (2008), Guazzoni et al. (2011), Guazzoni et al. (2012)).

$nlj^{\text{a})}$	$^{112}\text{Sn}(p,t)^{110}\text{Sn}(\text{gs})$	$^{124}\text{Sn}(p,t)^{122}\text{Sn}(\text{gs})$	$^{124}\text{Sn}(p,t)^{122}\text{Sn}(\text{gs})$	
$1g_{7/2}$	0.96	-1.1073	0.44	0.63
$2d_{5/2}$	0.66	-0.7556	0.35	0.60
$2d_{3/2}$	0.54	-0.4825	0.58	0.72
$3s_{1/2}$	0.45	-0.3663	0.36	0.52
$1h_{11/2}$	0.69	-0.6647	1.22	-1.24

Table 3.1.1: Two-nucleon transfer spectroscopic amplitudes associated with the reactions $^{112}\text{Sn}(p,t)^{110}\text{Sn}(\text{gs})$ and $^{124}\text{Sn}(p,t)^{122}\text{Sn}(\text{gs})$. **a)** quantum numbers of the two-particle configurations $(nlj)_{J=0}^2$ coupled to angular momentum $J = 0$. **b)** and **d)** $\langle \text{BCS} | T_\nu | \text{BCS} \rangle = \sqrt{(2j_\nu + 1)/2} U_\nu(A) V_\nu(A + 2)$ ($A + 2 = 112$ and 124 respectively), where $T_\nu = [a_\nu^\dagger a_\nu^\dagger]_0 / \sqrt{2}$ ($\nu \equiv nlj$) (cf. Potel et al. (2011, 2013a,b)) **c)** two-nucleon transfer spectroscopic amplitudes calculated making use of initial and final state wavefunctions obtained by diagonalizing a v_{low-k} , that is a renormalized, low-momentum interaction derived from the CD-Bonn nucleon-nucleon potential (see Guazzoni et al. (2006) and references therein). **e)** Two-neutron overlap functions obtained making use of the shell-model wavefunctions for the ground state of ^{122}Sn and ^{124}Sn calculated with the code NuShell (Brown and Rae, 2007). The wavefunctions were obtained starting with a G -matrix derived from the CD-Bonn nucleon-nucleon interaction Machleidt et al. (1996). These amplitudes were used in the calculation of $^{124}\text{Sn}(p,t)^{122}\text{Sn}$ absolute cross sections carried out by I.J. Thompson (Thompson, 2013).

$\lambda = -\partial\Delta E(\omega)/\partial\hbar\omega$ is the so-called mass enhancement factor. It reflects the ability with which vibrations dress single-particles. In other words, it measures the probability with which a nucleon moving at $t = -\infty$ in a “pure” orbital j can be found at a later time in a $2p - 1h$ like (doorway state) $|j'L; j\rangle$, L being the multipolarity of a vibrational state. Within this context, the discontinuity taking place at the Fermi energy in the dressed particle picture ($Z_\omega = (m/m_\omega)$) reflects the associated single-particle occupancy probability³¹.

It is of notice that dressed particles imply also an induced pairing interaction v_p^{ind} which renormalizes the bare $NN^{-1}S_0$ interaction v_p^{bare} , and results from the exchange of the dressing vibrations between pairs of nucleons moving in time reversal states close to the Fermi energy³². *In other words, fluctuations in the normal density $\delta\rho$ and the associated δU (Time-Dependent HF) and associated particle-vibration coupling vertices lead to abnormal (pairing) density.* Whether this results in a dynamic or static phenomenon, depends on whether the parameter (see

³¹See e.g. Brink and Broglia (2005) Ch. 9.

³²See e.g. Fig. 4.8.3, 7.E.1 (II) (b),(d),(e).

Fig. 3.5.7³³⁾

$$x' = G' N'(0), \quad (3.2.1)$$

product of the effective pairing strength,

$$G' = Z_\omega^2 (v_p^{bare} + v_p^{ind}), \quad (3.2.2)$$

and of the renormalized density of levels $N'(0)$ is considerably smaller (larger) than $\approx 1/2$. The quantity G' is the sum of the bare and induced pairing interaction, renormalized by the degree of single-particle content of the levels in which nucleons correlate. The quantity

$$N'(0) = Z_\omega^{-1} N(0) = (1 + \lambda) N(0) \quad (3.2.3)$$

is the similarly renormalized density of levels at the Fermi energy. From the above relations one obtains

$$x' = Z_\omega (v_p^{bare} + v_p^{ind}) N(0). \quad (3.2.4)$$

It is of notice that typical values of $Z_\omega \approx 0.7$, while for nuclei along the stability valley bare and induced pairing contributions are about equal which, according to Eq. (3.2.2) implies $G' \approx G = v_p^{bare}$, as in the case of a non-renormalized situation. On the other hand the physics is radically different, particles being a consistent fraction of the time in excited states coupled to collective vibrations, pairing acquiring a state dependence.

All of the above many-body, ω -dependent effects which imply in many cases a coherent sum of amplitudes, together with the corresponding renormalizations of the single-particle radial wavefunctions (formfactors) not discussed within the present framework, are not simple to capture in a spectroscopic factor³⁴ neither in connection with one-particle transfer, nor regarding two-nucleon transfer processes³⁵.

constituents	M/M_n	$a(\text{cm})$	$v_0(\text{eV})$	q	phase($T = 0$)
${}^3\text{He}$	3	2.9(-8)	8.6(-4)	0.19	liquid ^{a)}
${}^4\text{He}$	4	2.9(-8)	8.6(-4)	0.14	liquid ^{a)}
H_2	2	3.3(-8)	32(-4)	0.06	solid ^{b)}
${}^{20}\text{Ne}$	20	3.1(-8)	31(-4)	0.007	solid ^{b)}
nucleons	1	9(-14)	100(+6)	0.5	liquid ^{a),c),d)}

Table 3.3.1: Zero temperature phase for a number of systems of mass M (M_n : nucleon mass), the first four depending on atomic interactions (range Å, strength eV), the last one referring to the atomic nucleus. a) delocalized (condensed), b) localized, c) non-Newtonian solid (cf. e.g. Bertsch (1988), de Gennes (1994)), that is, systems which react elastically to sudden solicitations and plastically under prolonged strain, d) paradigm of quantal, strongly fluctuating, finite many-body systems. Delocalization or less does not seem to depend much on whether one is dealing with fermions or bosons (Mottelson (1998) and refs. therein; cf also Ebran et al. (2014b), Ebran et al. (2014a), Ebran et al. (2013), Ebran et al. (2012)).

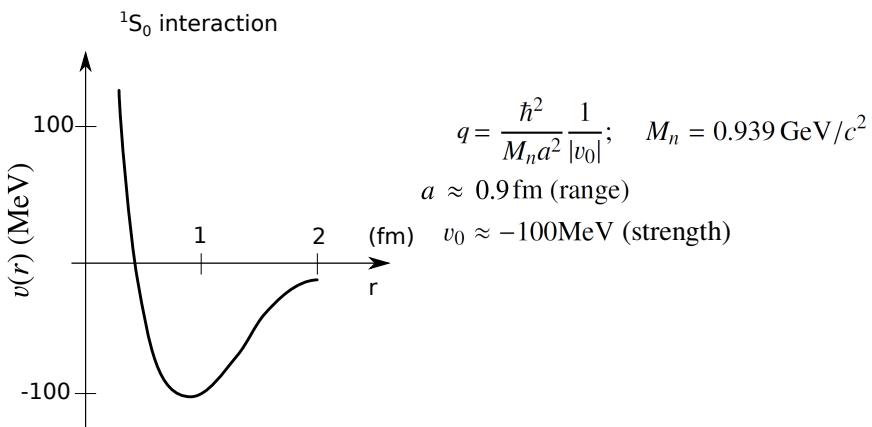


Figure 3.3.1: Schematic representation of the bare NN -interaction acting among nucleons displayed as a function of the relative coordinate $r = |\mathbf{r}_1 - \mathbf{r}_2|$, used to estimate the quantality parameter q , ratio of the zero point fluctuations (ZPF) of confinement and the potential energy.

3.3 Quantality Parameter

The quantality parameter³⁶ is defined as the ratio of the quantal kinetic energy of localization (confinement) and the potential energy, (cf. Fig. 3.3.1 and Table 3.3.1). Fluctuations, quantal or classical, favor symmetry: gases and liquids are homogeneous. Potential energy on the other hand prefers special arrangements: atoms like to be at specific distances and orientations from each other (spontaneous breaking of translational and of rotational symmetry reflecting the homogeneity and isotropy of empty space³⁷).

When q is small, quantal effects are small and the lower state for $T < T_c$ will be a solid (crystalline structure), T_c denoting the critical temperature. For sufficiently large values of $q (> 0.15)$ the system will display particle delocalization and, likely, be amenable, within some approximation, to a mean field description (Figs. 1.2.2 and 3.3.2). In fact, the step *delocalization* \rightarrow *mean field* is not automatic, neither guaranteed. In any case, not for all properties neither for all levels of the system. Let us elaborate on these points.

As already stated, independent particle motion can be viewed as the most collective of all nuclear properties, reflecting the effect of all nucleons on a given one resulting in a macroscopic effect. Consequently, it should be possible to calculate the mean field in an accurate manner. Arguably, as accurately as one can calculate collective vibrations, e.g. quadrupole vibrations. But this does not mean that one knows how to correctly calculate the energy and associated dynamical deforma-

³³Brink and Broglia (2005) App. H. Sect. H4 and refs. therein; Barranco et al. (2005).

³⁴In keeping with the fact that $m_k \approx 0.6 - 0.7m$ and that $m^* \approx m$, as testified by the satisfactory fitting standard Saxon-Woods potentials provides for the valence orbitals of nucleons of mass m around closed shells, one obtains $m_\omega \approx 1.4 - 1.7m$. Thus $Z_\omega \approx 0.6 - 0.7$. It is still an open question how much of the observed single-particle depopulation can be due to hard core effects, which shifts the associated strength to high momentum levels (see Dickhoff and Van Neck (2005), Jenning (2011), Kramer et al. (2001), Barbieri (2009), Schiffer et al. (2012), Duguet and Hagen (2012), Furnstahl and Schwenk (2010)). An estimate of such an effect of about 20% will not quantitative change the long wavelength estimate of Z_ω given above. Arguably, a much larger depopulation through hard core effects remains an open problem within the overall picture of elementary modes of nuclear excitation and of medium polarization effects. Within this context it remains an open question the role the renormalization of the radial dependence of the single-particle wavefunctions (form factors), due to many-body effects, can play in the determination of spectroscopic factors. Discussions with Augusto Macchiavelli on this subject are gratefully acknowledged.

³⁵See Barranco et al. (2005, 1999).

³⁶Nosanow (1976), de Boer (1957), de Boer (1948), de Boer and Lundbeck (1948), Mottelson (1998).

³⁷As already stated in footnote 1 of this Chapter, within such a general context, the physics embodied in the quantality parameter is closely related to that which is at the basis of the classical Lindemann criterion (Lindemann (1910)) to measure whether a system is ordered (e.g. a crystal) or disordered (e.g. a melted system) (Bilgram (1987), Löwen (1994), Stillinger and Stillinger (1990); Stillinger (1995)). The above statement concerning the competition between potential energy and fluctuations, is also valid for the generalized Lindemann parameter (Stillinger and Stillinger (1990), Zhou et al. (1999)), used to provide similar insight into inhomogeneous finite systems like e.g. proteins (aperiodic crystals Schrödinger (1944), see also Ehrenfest's theorem (Basdevant and Dalibard (2005) pag. 138 see also App. 3.A).

tion parameter of each single state of the quadrupole response function. Within this context one may find, through mean field approximation a good description for the energy of the valence orbitals of a nucleus but for specific levels (e.g. the $d_{5/2}$ level of the isotopes $^{119-120}\text{Sn}$, see Fig 5.2.3). It is not said that including particle-vibration coupling corrections, a process which in average makes theory come closer to experiment³⁸, single specific quasiparticle energies will satisfactorily agree with the data³⁹. Cases like this one constitute a sobering experience concerning the intricacies of the many-body problem in general, and of the nuclear one (finite many-body system, FMBS) in particular, where spatial quantization plays a central role. In other words, one is dealing with a self-bound, strongly interacting, finite many-body system generated from collisions originally associated with a variety of astrophysical events and thus with the coupling and interweaving of different scattering channels and resonances, a little bit as e.g. the Hoyle monopole resonance ($\alpha + \alpha + \alpha \rightarrow ^{12}\text{C}$). Within the anthropomorphic scenario such phenomena are found in the evolution of the Universe to eventually allow for the presence of organic matter and, arguably, life on earth⁴⁰ more likely than to make mean field approximation, also the renormalized one, an “exact” description of nuclear structure and reactions, for every single nuclear level.

3.4 Generalized quantity parameter

Remaining within the framework of independent particle motion, but allowing for the dressing of particles one obtains a sigmoidal distribution around the Fermi energy reflecting the ω -dependence of $Z_\nu(\omega)$. In the case of the Sn-isotopes where the density of levels is $N(0) \approx 4\text{MeV}^{-1}$ and the energy of the lowest $\beta = 0$ collective vibration ($\lambda^\pi = 2^+$) is $\hbar\omega_2 \approx 1.3$ MeV, one expects the width of the distribution to be ≈ 3 MeV around the Fermi energy. Allowing nucleons not only to emit and reabsorb $\beta = 0$ vibrational modes, but also to exchange them with the other nucleons, as well as to interact through v_p^{bare} , one is confronted with a many-body problem like the one discussed in Sect. 3.2 for $x \gg 1/2$, which can be solved at profit within the BCS framework. It leads, again for Sn-isotopes, to a sigmoidal occupation probability extending over an energy range also of ≈ 3 MeV ($\approx 2\Delta$) around the Fermi energy⁴¹.

In spite of the similitudes in occupancy between the two physical situations discussed above, the associated $L = 0$ (gs) \rightarrow (gs) two-nucleon transfer cross sections are quite different. In the first case, it is proportional to $(Z_{\nu_1} \sqrt{\sigma_{\nu_1}})^2 \approx (1/2) \times \sigma_{\nu_1}$

³⁸See e.g. Bohr and Mottelson (1975), Bohr and Broglia (1977), Hamamoto (1977), Bès et al. (1977), Barnes et al. (1972), Reich et al. (1974), Hamamoto and Siemens (1976), Bortignon et al. (1977), Bernard and Giai (1981), Mahaux et al. (1985), Bès and Broglia (1971a); Flynn et al. (1971), Bortignon et al. (1976), Bès et al. (1988), Barranco et al. (1987), Vinh Mau (1995), Barranco et al. (2001); Orrigo and Lenske (2009) and references therein.

³⁹See Tarpanov et al. (2014).

⁴⁰See e.g. Rees (2000), Meißner (2015) and references therein.

⁴¹Within this context see discussion following Eqs. (1.4.54) and (1.4.53) and Tinkham (1996).

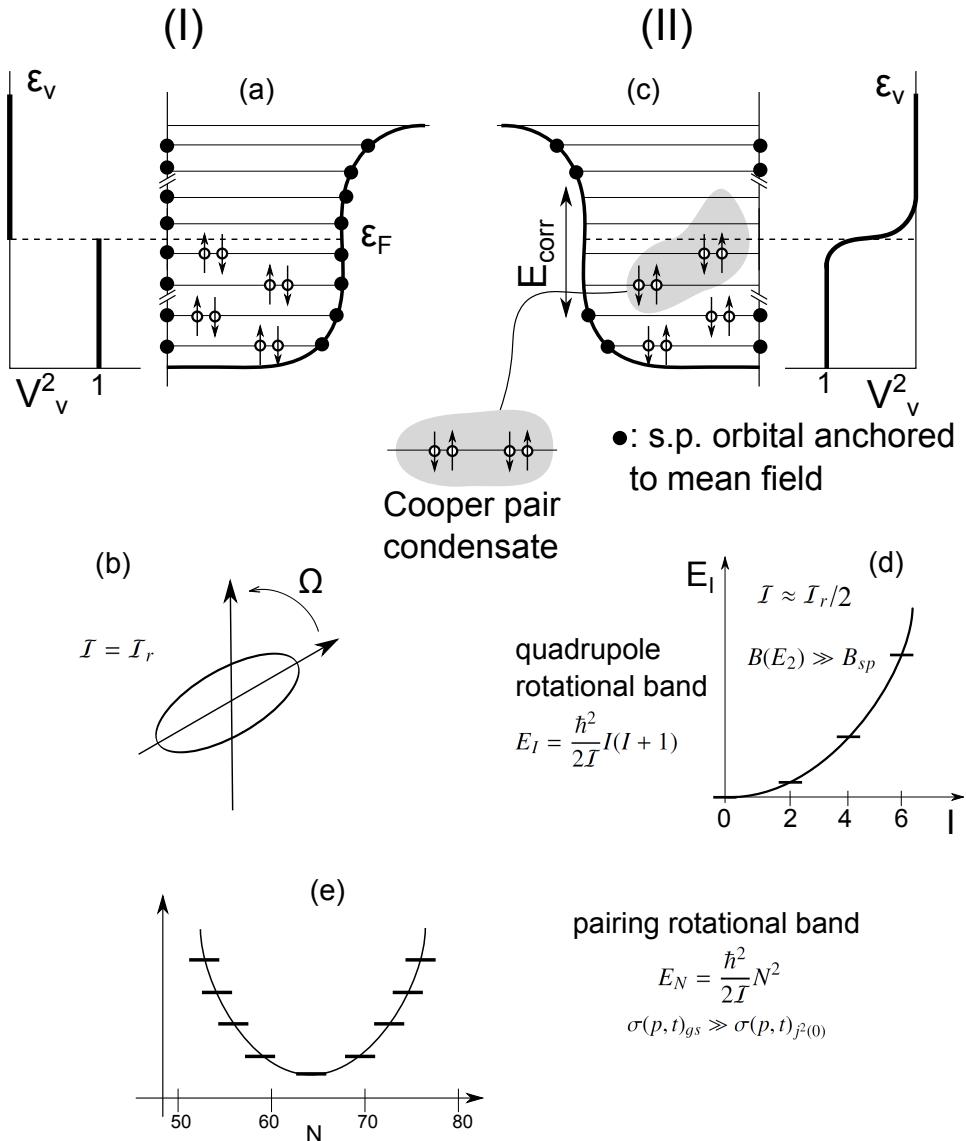


Figure 3.3.2: (I) (a) Schematic representation of “normal” (independent-particle) motion of nucleons in two-fold degenerate (Kramers, time-reversal degeneracy) orbits solidly anchored to the mean field (solid dots at the ends of the single-particle levels) and displaying a sharp, step-function-like, discontinuity in the occupancy at the Fermi energy can lead to a deformed (Nilsson (1955)) rotating nucleus with a rigid moment of inertia I_r . (b)

(II) (c) Schematic representation of phase correlated, independent nucleon Cooper pair motion in which few (of the order of 5-8) pairs lead to (d) a sigmoidal occupation function at the Fermi energy and, having mainly uncoupled themselves from the fermionic mean field (no solid dot at the end of the Kramers invariant single-particle levels), contribute in a reduced fashion to (d) the moment of inertia of quadrupole rotational bands leading to $I \approx I_r/2$ (cf. Belyaev (2013), Belyaev (1959), Bohr and Mottelson (1975) and references therein), (e) pairing rotational bands in gauge space, an example of which is provided by the ground states of the superfluid Sn-isotopes.

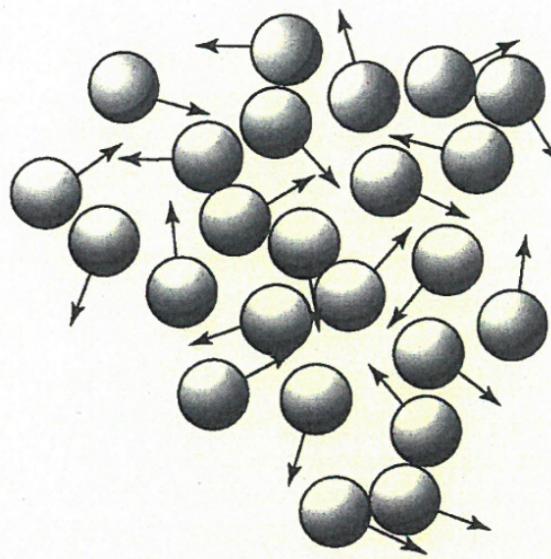


Figure 3.3.3: A system of localized pairs (Schafroth pairs). Such that the size of the pair bound state is small compared to the average spacing between pairs.

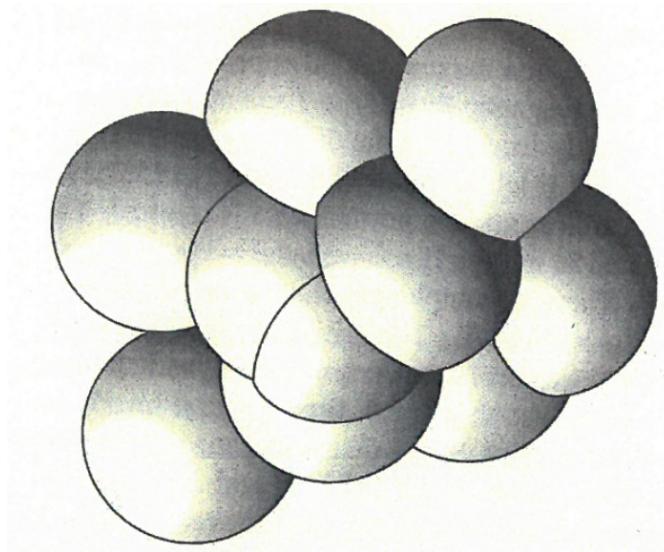


Figure 3.3.4: At the basis of BCS superconductivity one finds condensation of very extended ($\approx 10^4 \text{ \AA}$), strongly overlapping (within the volume of a given pair the centers of $\approx 10^6$ other pairs is found), weakly bound ($\approx 2\Delta \approx \text{meV}$) Cooper pairs corresponding to ordering in occupying momentum space, and not space-like condensation of strongly bound pairs which undergo Bose condensation (cf. Schrieffer (1964)).

and to $(\sum_{v>0} U'_v V'_v \sqrt{\sigma_v})^2 \approx (\alpha'_0)^2 \bar{\sigma} \approx 50 \times \bar{\sigma}$ in the second. The quantity σ_{v_1} is the cross section associated with the lowest $j_{v_1}^2(0)$ two-particle configuration. The summation $\sum_{v>0}$ extends over the orbitals around ϵ_F within a range $\approx 2\Delta$, and $\bar{\sigma}$ is the average of the cross sections σ_v associated with the valence configurations $j_v^2(0)$.

In connection with the Cooper pair (BCS solution) one should, more correctly, use the wavefunction (1.4.65) to calculate the two-nucleon transfer cross section. In keeping with the definition $c'_v = V'_v/U'_v$ ad the phasing (1.4.41), it is found that

$$\sum_{v>0} c'_v P_v^\dagger |0\rangle_F = e^{2i\phi} \sum_{v>0} c_v P_v^\dagger |0\rangle_F, \quad (3.4.1)$$

leading to,

$$|\Psi_{N_1}\rangle \sim \left(e^{2i\phi} \sum_{v>0} c_v P_v^\dagger \right)^{N_1/2} |0\rangle_F. \quad (3.4.2)$$

That is, independent pair motion, each Cooper pair contributing a phase ϕ to $|N\rangle$.

Note that in a metallic superconductor as described by BCS, there are of the order of 10^6 Cooper pairs around the Fermi surface. Because the phase ϕ is common to so many pairs, its effects can become felt on the macroscopic scale. By the same token, the quantal behaviour of the BCS condensate must also be the same as the quantal behaviour of a single Cooper pair. A feature which reflects itself in superfluid nuclei, although in this case one is talking of only 6–8 Cooper pairs.

If one finds, as one does⁴², that the tunneling energy is a function of the relative phase, it will adjust so as to minimize the energy through tunneling of electrons. The generalized quantality parameter, ratio of the kinetic energy of confinement ($T_\xi = \hbar^2/(2m_e\xi^2) \approx 4 \times 10^{-5}$ meV; $\xi \approx 10^4$) of the electrons forming the Cooper pairs undergoing tunneling, and the Cooper pair correlation (binding) energy $E_{corr}(\approx 3$ meV, $\Delta \approx 1.4$ meV, Pb), is much smaller than 1 ($q_\xi \approx 10^{-5}$). A result which implies that for $T < T_c$ ($T_c \approx 7.1$ K, critical temperature of Pb), the electron Cooper pair partners are solidly anchored to each other. Said it differently, the intrinsic Cooper pair motion is frozen because of the pairing gap. Consequently, if under such conditions ($T < T_c$) one observes a current across the unbiased junction between the two superconductors, it must be a current of carriers of mass $2m_e$ and charge $2e$. If one electron partner of a Cooper pair tunnels, the other entangled electron does it also. The parallel between independent-particle and independent-Cooper pair motion, all with the same phase, is summarized in⁴³ Fig. 3.4.1. *It is of notice that in the present case this result emerges not because of the large role played by potential effects, but because of the very large dimensions, low relative intrinsic momentum of the Cooper pair.*

⁴²Josephson (1962); Anderson and Rowell (1963).

⁴³In connection with Fig. 3.4.1, the estimate $R = 8/k_F$ was carried out with the help of the Fermi gas model (see e.g. Bohr and Mottelson (1969), for $A \approx 120$). The Fermi momentum is written as $k_F \approx (3\pi^2 A/2V)^{1/3} \approx (\frac{3\pi^2}{2} \rho(0))^{1/3}$. It is of notice that this expression leads for $\rho(0) \approx 0.17 \text{ fm}^{-3}$ to the Fermi momentum $k_F \approx 1.36 \text{ fm}^{-1}$.

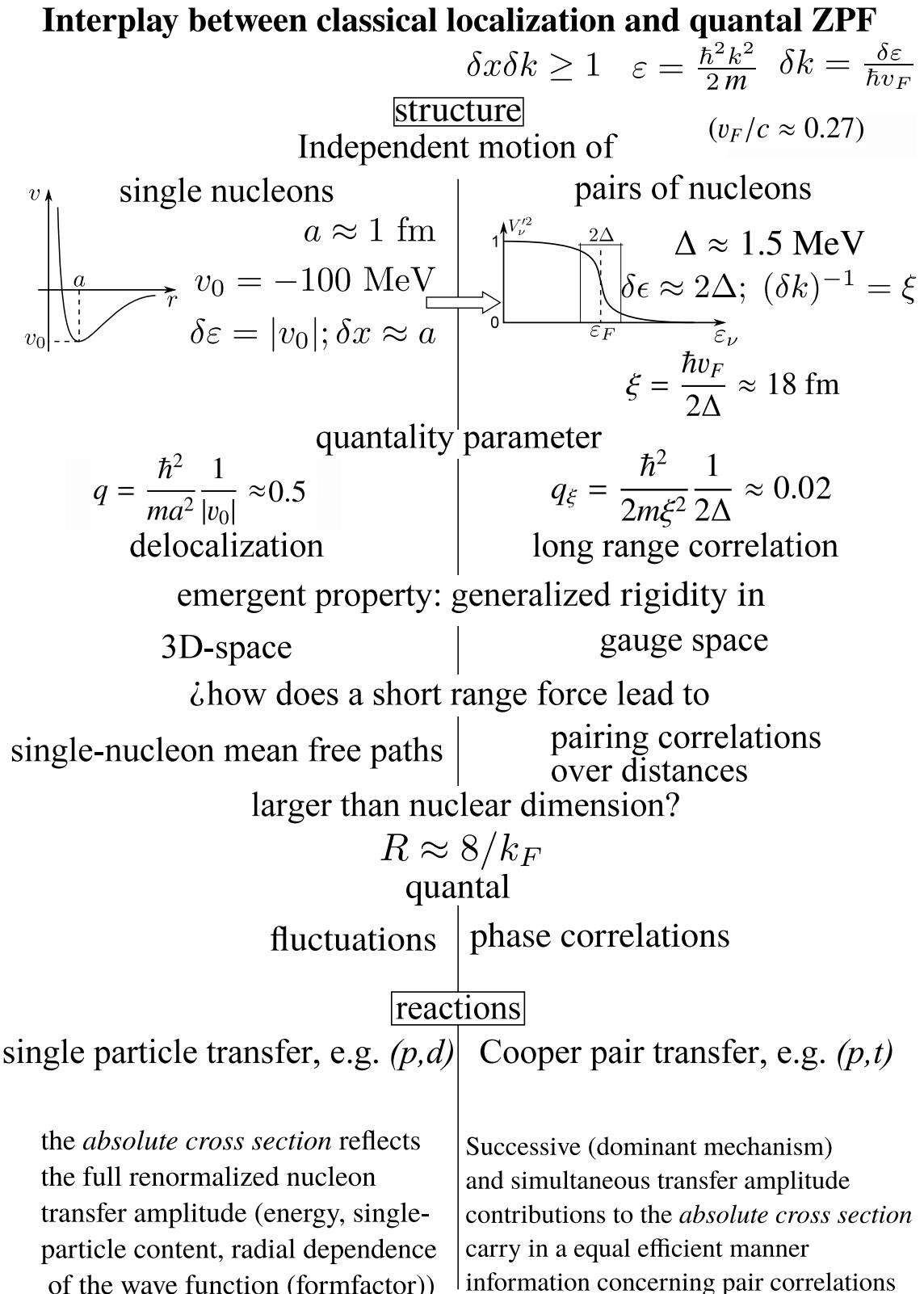


Figure 3.4.1: Classical localization and zero point fluctuations, associated with independent-particle (normal density) and independent-pair (abnormal density) motion.

We conclude this section by remarking once more that BCS condensation, that is condensation of strongly overlapping, very extended, weakly bound Cooper pairs corresponds to ordering in occupying momentum space, and not space-like condensation of strongly bound clusters which undergo Bose condensation. In BCS condensation, the inner, intrinsic structure of the pair, that is, the fact that it is made out of fermions entangled in time reversal states, is the characterizing feature, at the basis of ODLRO. Cooper pairs (see Eqs. (1.4.48) and (3.4.1)) are not bosons, as $(P_\nu^\dagger)^2 = 0$ testifies (see also App. 4.A). However, it is true that, under certain circumstances, namely $T < T_c$, $q \ll 1/\xi$ and $d \ll \xi$ ($D_0 \lesssim \xi$ in the case of heavy ion reactions between superfluid nuclei), where ξ is the correlation length and d is the width of the Josephson junction (D_0 being the distance of closest approach), Cooper pairs show properties which strongly resemble those of bosons⁴⁴.

3.5 Two-nucleon spectroscopic amplitudes associated with pairing vibrational modes in closed shell systems: the ^{208}Pb case.

The solution of the pairing Hamiltonian leads, in the case of closed shell systems and within the harmonic approximation (RPA), to pair addition (*a*) pair removal (*r*) two-particle, two-hole correlated modes. The associated creation operator have been defined in Eqs. (1.3.7)–(1.3.9) for the pair addition mode. Similarly, in the case of the pair removal mode,

$$\Gamma_r^\dagger(n) = \sum_i X_n^r(i)\Gamma_i^\dagger + \sum_k Y_n^r(k)\Gamma_k,$$

X, Y fulfilling

$$\sum X^2 - Y^2 = 1,$$

and

$$\Gamma_k^\dagger = a_k^\dagger a_{\bar{k}}^\dagger, \quad (\epsilon_k > \epsilon_F).$$

Similarly,

$$\Gamma_i^\dagger = a_{\bar{i}}^\dagger a_i, \quad (\epsilon_i \leq \epsilon_F).$$

The relations

$$[H, \Gamma_a^\dagger(n)] = \hbar W_n(\beta = +2),$$

and

$$[H, \Gamma_r^\dagger(n)] = \hbar W_n(\beta = -2),$$

where β is the transfer quantum number, lead to the dispersion relations,

$$\frac{1}{G(\pm 2)} = \sum_k \frac{(\Omega_k/2)}{2\epsilon_k \mp W_n(\pm 2)} + \sum_i \frac{(\Omega_i/2)}{2\epsilon_i \pm W_n(\pm 2)}, \quad (3.5.1)$$

⁴⁴ter Haar (1995b), ter Haar (1995a) Ch. 4.9 and Ehrenfest and Oppenheimer (1931).

orbit	ϵ_j	$\epsilon_{p_{1/2}} - \epsilon_k \equiv \epsilon_k - \epsilon_{p_{1/2}} $
$0h_{9/2}$	-10.62	3.47
$1f_{7/2}$	-9.50	2.35
$0i_{13/2}$	-8.79	1.64
$2p_{3/2}$	-8.05	0.90
$1f_{5/2}$	-7.72	0.57
$2p_{1/2}$	-7.15	0
$\epsilon_F = -5.825$ MeV		$\epsilon_k - \epsilon_{g_{9/2}} \equiv \epsilon_{g_{9/2}} - \epsilon_k $
$1g_{9/2}$	-3.74	0.
$0i_{11/2}$	-2.97	0.77
$0j_{15/2}$	-2.33	1.41
$2d_{5/2}$	-2.18	1.56
$3s_{1/2}$	-1.71	2.03
$1g_{7/2}$	-1.27	2.47
$2d_{3/2}$	-1.23	2.51

Table 3.5.1: Valence single-particle levels of ^{208}Pb . In the upper part the occupied levels ($\epsilon_i \leq \epsilon_F$) are displayed while in the lower part the empty levels are shown ($\epsilon_k \geq \epsilon_F$). Of notice that $\epsilon_{p_{1/2}} - \epsilon_{g_{9/2}} = 3.41$ MeV, is the single-particle gap associated with $N = 126$ shell closure.

n labeling the corresponding solutions in increasing order of energy. In the above equation, $\Omega_j = j + 1/2$ is the pair degeneracy of the orbital with total angular momentum j .

For the case of the (neutron) pair addition and pair subtraction modes of ^{208}Pb the above equations are graphically solved in Fig 3.5.1 (see also Table 3.5.1). The minimum of the dispersion relation defines the Fermi energy of the system under study. This is in keeping with the fact that *in the case in which* $W_1(\beta = +2) = W_1(\beta = -2) = 0$, *situation corresponding to the transition between normal and superfluid phases*⁴⁵, the energy value at which the dispersion relation touches for the first time the energy axis, coincides with the BCS λ variational parameter. It is of notice that, as a rule, the Fermi energy of closed shell nuclei is empirically defined as half the energy difference between the last occupied and the first empty single particle state⁴⁶. Making use of the values

$$\begin{cases} E_{corr}(+2) = BE(208) + BE(210) - 2BE(209) = 1.248 \text{ MeV}, \\ E_{corr}(-2) = BE(208) + BE(206) - 2BE(207) = 0.640 \text{ MeV}, \end{cases}$$

one obtains $W_1(-2) + W_1(+2) = (BE(208) - BE(206)) - (BE(210) - BE(208))$

⁴⁵See footnote 27 Ch. 1.

⁴⁶cf. e.g. Mahaux et al. (1985).

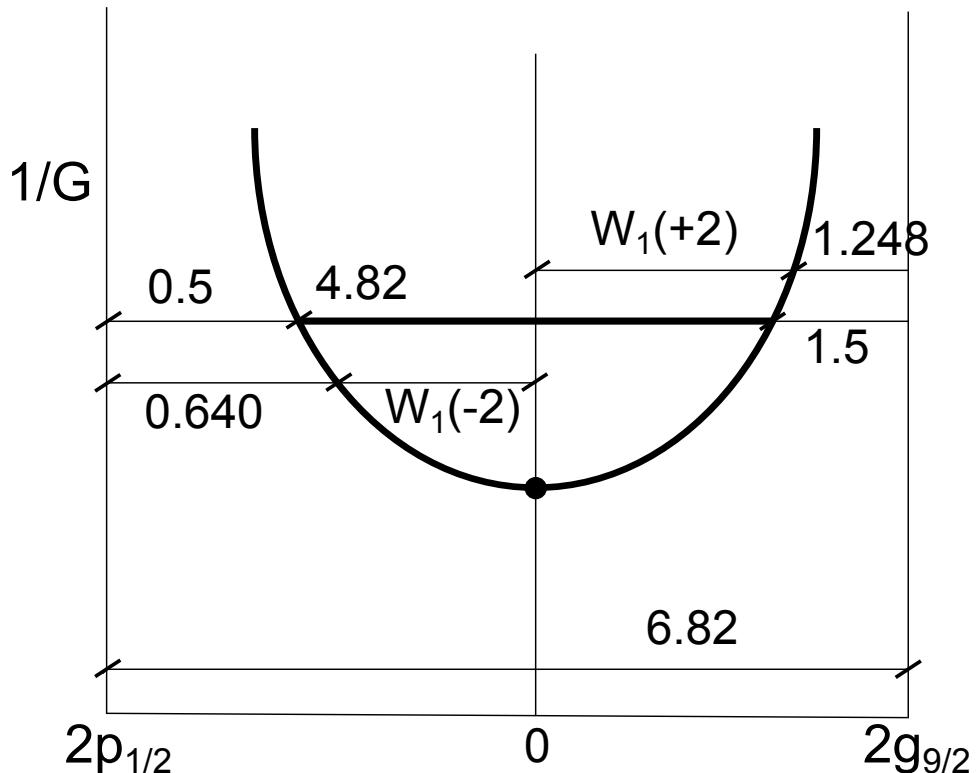


Figure 3.5.1: The right hand side of the RPA pairing vibrational dispersion relation for neutrons in the case of the closed shell system ^{208}Pb in the region between the two neighboring shells ($p_{1/2}$ and $g_{9/2}$). All quantities are in MeV. For each value of G there is a straight horizontal line, which is divided by the the curve in three sections. The first one from the left corresponds to the pairing correlation energy of the nucleus ^{206}Pb (two correlated neutron hole states) while the last segment to the right measures the pairing correlation energy of ^{210}Pb (two correlated neutrons above closed shell) the intermediate segment measures the energy of the two phonon (correlated $(2p - 2h)$) pairing vibrational state of ^{208}Pb .

$= 14.11 - 9.115 = 4.995$ MeV. Notice that in the above calculations all energies differences are positive. In particular (see Table 3.5.1)

$$\epsilon_i < \epsilon_F \Rightarrow \epsilon_F - \epsilon_i = -|\epsilon_F| + |\epsilon_i| = |\epsilon_i| - |\epsilon_F| > 0,$$

and

$$\epsilon_k > \epsilon_F \Rightarrow \epsilon_k - \epsilon_F = -|\epsilon_k| + |\epsilon_F| = |\epsilon_F| - |\epsilon_k| > 0.$$

Thus,

$$\begin{cases} 2(\epsilon_F - \epsilon_{p_{1/2}}) = W_1(-2) + E_{corr}(-2) > 0, \\ 2(\epsilon_{g_{9/2}} - \epsilon_F) = W_1(+2) + E_{corr}(+2) > 0. \end{cases}$$

From Fig. 3.5.1 and Table 3.5.1 one can then write,

$$2 \times (-5.825 - (-7.15)) \text{ MeV} = 2.650 \text{ MeV} = W_1(-2) + 0.640 \text{ MeV}$$

and

$$2 \times (-3.74 \text{ MeV} - (-5.825) \text{ MeV}) = 4.17 \text{ MeV} = W_1(+2) + 1.248 \text{ MeV}.$$

Consequently,

$$W_1(-2) = 2.01 \text{ MeV} \quad \text{and} \quad W_1(+2) = 2.92 \text{ MeV},$$

leading to,

$$W_1(+2) + W_1(-2) = 4.93 \text{ MeV}. \quad (3.5.2)$$

3.5.1 Pair removal mode

In Fig. 3.5.2, a graphical representation of the forwards going RPA amplitude of the pair removal mode is shown. Its expression for $n = 1$, is

$$X_1^r(i) = \frac{\frac{1}{2}\Omega_i^{1/2}\Lambda_1(-2)}{2(\epsilon_F - \epsilon_i) - W_1(-2)},$$

where

$$\begin{aligned} 2 \times (\epsilon_F - \epsilon_i) - W_1(-2) &= 2 \times (\epsilon_F - \epsilon_i) - 2 \times (\epsilon_F - \epsilon_{p_{1/2}}) + E_{corr}(-2) \\ &= 2 \times (\epsilon_{p_{1/2}} - \epsilon_i) + E_{corr}(-2) = 2 \times (|\epsilon_i| - |\epsilon_{p_{1/2}}|) + E_{corr}(-2). \end{aligned}$$

Thus,

$$X_1^r(i) = \frac{\frac{1}{2}\Omega_i^{1/2}\Lambda_1(-2)}{2(|\epsilon_i| - |\epsilon_{p_{1/2}}|) + E_{corr}(-2)}.$$

Making use of the empirical value of $E_{corr}(-2)$ worked out above one obtains,

$$X_1^r(i) = \frac{\frac{1}{2}\Omega_i^{1/2}\Lambda_1(-2)}{2(|\epsilon_i| - |\epsilon_{p_{1/2}}|) + 0.640 \text{ MeV}}.$$

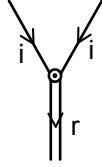


Figure 3.5.2: NFT representation of the forwards going RPA amplitude of the pair removal mode (double downward going arrowed line) describing a two correlated hole state (single downward going arrowed line for each hole with quantum numbers collectively labeled i).

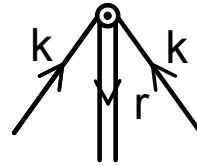


Figure 3.5.3: Same as Fig. 3.5.2 but for the backwards going amplitudes.

In Fig. 3.5.3 we display the graphical process associated with the backwards going ($n = 1$) RPA amplitude,

$$Y_1^r(k) = \frac{\frac{1}{2}\Omega_k^{1/2}\Lambda_1(-2)}{2(\epsilon_k - \epsilon_F) + W_1(-2)}.$$

Making use of

$$2 \times (\epsilon_F - \epsilon_{p_{1/2}}) - E_{corr}(-2) = W_1(-2),$$

one can write

$$2 \times (\epsilon_F - \epsilon_{p_{1/2}}) + 2 \times (\epsilon_k - \epsilon_F) - E_{corr}(-2) = 2 \times (\epsilon_k - \epsilon_F) + W_1(-2),$$

leading to

$$2 \times (|\epsilon_{p_{1/2}}| - |\epsilon_k|) - E_{corr}(-2) = 2 \times (|\epsilon_{p_{1/2}}| - |\epsilon_{g_{9/2}}|) + 2 \times (|\epsilon_{g_{9/2}}| - |\epsilon_k|) - E_{corr}(-2).$$

Thus,

$$Y_1^r(k) = \frac{\frac{1}{2}\Omega_k^{1/2}\Lambda_1(-2)}{2(|\epsilon_{g_{9/2}}| - |\epsilon_k|) + 2(|\epsilon_{p_{1/2}}| - |\epsilon_{g_{9/2}}|) - E_{corr}(-2)}.$$

With the help of $2 \times (|\epsilon_{p_{1/2}}| - |\epsilon_{g_{9/2}}|) - E_{corr}(-2) = 6.82 \text{ MeV} - 0.640 \text{ MeV} = 6.18 \text{ MeV}$, one obtains,

$$Y_1^r(k) = \frac{\frac{1}{2}\Omega_k^{1/2}\Lambda_1(-2)}{2(|\epsilon_{g_{9/2}}| - |\epsilon_k|) + 6.18 \text{ MeV}}.$$

The above expressions of $X_1^r(i)$ and $Y_1^r(k)$ contain the experimental values of the 2-hole correlation energies (0.640 MeV). Because (see Fig. 3.5.1) the associated values of G does not lead to the observed correlation energy of the pair addition mode (1.248 MeV), we prefer to choose a single value of G and use the resulting $E_{corr}(-2)$ (=0.5 MeV) and $E_{corr}(+2)$ (=1.5 MeV), correlation energies, to calculate the corresponding X, Y amplitudes for both the lowest removal and lowest addition pairing modes. Making use of,

$$2 \times (|\epsilon_{p_{1/2}}| - |\epsilon_{g_{9/2}}|) = 6.82 \text{ MeV} \quad \text{and} \quad 2 \times (|\epsilon_{p_{1/2}}| - |\epsilon_{g_{9/2}}|) - E_{corr}(-2) \\ = (6.82 - 0.5) \text{ MeV} = 6.32 \text{ MeV},$$

one can write

$$X_1^r(i) = \frac{\frac{1}{2}\Omega_i^{1/2}\Lambda_1(-2)}{2(|\epsilon_i| - |\epsilon_{p_{1/2}}|) + 0.5 \text{ MeV}},$$

$$Y_1^r(k) = \frac{\frac{1}{2}\Omega_k^{1/2}\Lambda_1(-2)}{2(|\epsilon_{g_{9/2}}| - |\epsilon_k|) + 6.32 \text{ MeV}}.$$

Tables 3.5.2 and 3.5.3 contain the amplitudes of the pair removal mode of ^{208}Pb ($\Gamma_r^\dagger(1) = \sum X_1^r(i)\Gamma_i^\dagger + \sum Y_1^r(k)\Gamma_k$), that is of the two neutron correlated hole state describing $|^{206}\text{Pb}(\text{gs})\rangle = \Gamma_r^\dagger(1)|0\rangle$.

It is of notice that the coupling strength $\Lambda_1(-2)$ with which the pair removal mode couples to the two single-particles (-holes) states is calculated by normalizing the amplitudes: **1)** Tamm Dancoff, TD implies no ground state correlations⁴⁷, i.e. to assume that $Y \equiv 0$; $\sum_i A^2(i) = 1.5549 \text{ MeV}^{-2}$ and thus $\Lambda_1(-2) = 0.802 \text{ MeV}$, $(\sum X(i)_T^2 = 1)$; **2)** RPA, $\Lambda_1^2(-2) \times (\sum_i A^2(i) - \sum_k B^2(k)) = \Lambda_1^2(-2) \times 1.45073 = 1$. Thus $\Lambda_1(-2) = 0.830 \text{ MeV}$.

The above results show that there is a few percentage difference between the two values of Λ (TD and RPA), as well as for the corresponding X amplitudes (Table 3.5.2). Nonetheless, ground state correlations as expressed by the Y amplitudes (Table 3.5.3), gives rise to a 53% increase in the $^{206}\text{Pb}(t, p)^{208}\text{Pb}(\text{gs})$ absolute cross section, from 0.34 mb to 0.52 mb to be compared with experimental data $\sigma = 0.68 \pm 0.24 \text{ mb}$ (see Fig. 4.5.4).

3.5.2 Pair addition mode

In Fig. 3.5.4 the X -amplitude of the pair addition mode is shown (NFT diagram). The associated expression ($n = 1$),

$$X_1^a(k) = \frac{\frac{1}{2}\Omega_k^{1/2}\Lambda_1(+2)}{2(\epsilon_k - \epsilon_F) - W_1(+2)},$$

can, making use of the relation

⁴⁷See footnote 27 Ch. 1.

units		MeV	MeV ⁻¹	RPA	TD
nlj	Ω_i	$ \epsilon_i - \epsilon_{p_{1/2}} $	$A(i) = \frac{\frac{1}{2}\Omega_i^{1/2}}{2(\epsilon_i - \epsilon_{p_{1/2}}) + 0.5 \text{ MeV}}$	$X_1^r(i)$	$X_1^r(i)$
$2p_{1/2}$	1	0	1	0.83	0.80
$1f_{5/2}$	3	0.57	0.528	0.44	0.42
$2p_{3/2}$	2	0.90	0.307	0.25	0.25
$0i_{13/2}$	7	1.64	0.350	0.29	0.28
$1f_{7/2}$	4	2.35	0.192	0.16	0.15
$0h_{9/2}$	5	3.47	0.150	0.12	0.12

Table 3.5.2: Forwards going RPA amplitudes $X_1^r(i)$ of the lowest pair removal mode of ^{208}Pb (i.e. $|^{206}\text{Pb}\rangle$ state), cf. Table XVI Broglia et al. (1973).

units		MeV	MeV ⁻¹	RPA
nlj	Ω_k	$ \epsilon_{g_{9/2}} - \epsilon_k $	$B(k) = \frac{\frac{1}{2}\Omega_k^{1/2}}{2(\epsilon_{g_{9/2}} - \epsilon_k) + 6.23 \text{ MeV}}$	$Y_1^r(i)$
$1g_{9/2}$	5	0	0.179	0.15
$0i_{11/2}$	6	0.77	0.158	0.13
$0j_{15/2}$	8	1.41	0.156	0.13
$2d_{5/2}$	3	1.56	0.093	0.08
$3s_{1/2}$	1	2.03	0.046	0.04
$1g_{7/2}$	4	2.47	0.090	0.07
$2d_{3/2}$	2	2.51	0.063	0.05

Table 3.5.3: Same as Table 3.5.2 but for the backwards amplitudes $Y_1^r(k)$ of the lowest energy pair removal mode.

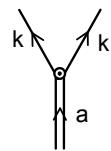


Figure 3.5.4: Same as Fig. 3.5.2 but for the pair addition mode

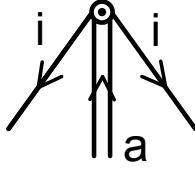


Figure 3.5.5: Same as Fig. 3.5.3 but for the pair addition mode

$$\begin{aligned} 2 \times (\epsilon_k - \epsilon_F) - W_1(+2) &= 2 \times (\epsilon_k - \epsilon_F) - 2 \times (\epsilon_{g_{9/2}} - \epsilon_F) + E_{corr}(+2) \\ &= 2 \times (\epsilon_k - \epsilon_{g_{9/2}}) + E_{corr}(+2) = 2 \times (|\epsilon_{g_{9/2}}| - |\epsilon_k|) + E_{corr}(+2), \end{aligned}$$

be expressed as

$$X_1^a(k) = \frac{\frac{1}{2}\Omega_k^{1/2}\Lambda_1(+2)}{2(|\epsilon_{g_{9/2}}| - |\epsilon_k|) + E_{corr}(+2)}.$$

Similarly (Fig. 3.5.5),

$$Y_1^a(i) = \frac{\frac{1}{2}\Omega_i^{1/2}\Lambda_1(+2)}{2(\epsilon_F - \epsilon_i) + W_1(+2)},$$

can be written, with the help of the relation

$$\begin{aligned} 2 \times (\epsilon_F - \epsilon_i) + W_1(+2) &= 2 \times (\epsilon_F - \epsilon_i) - 2 \times (\epsilon_{g_{9/2}} - \epsilon_F) - E_{corr}(+2) \\ &= 2 \times (\epsilon_{p_{1/2}} - \epsilon_i) + 2 \times (\epsilon_{g_{9/2}} - \epsilon_{p_{1/2}}) - E_{corr}(+2) \\ &= 2 \times (|\epsilon_i| - |\epsilon_{p_{1/2}}|) + 2 \times (|\epsilon_{p_{1/2}}| - |\epsilon_{g_{9/2}}|) - E_{corr}(+2), \end{aligned}$$

as

$$Y_1^a(i) = \frac{\frac{1}{2}\Omega_i^{1/2}\Lambda_1(+2)}{2(|\epsilon_i| - |\epsilon_{p_{1/2}}|) + 2\Delta\epsilon_{sp} - E_{corr}(+2)}.$$

Making use of $E_{corr}(+2) = 1.5$ MeV (cf. Fig. 3.5.1) and

$$2 \times \Delta\epsilon_{sp} = 2 \times (|\epsilon_{p_{1/2}}| - |\epsilon_{g_{9/2}}|) = 6.82 \text{ MeV},$$

one can write $2\Delta\epsilon_{sp} - E_{corr}(+2) = (6.82 - 1.5)$ MeV = 5.32 MeV, leading to

$$\begin{cases} X_1^a(k) = \frac{\frac{1}{2}\Omega_k^{1/2}\Lambda(-2)}{2(|\epsilon_{g_{9/2}}| - |\epsilon_k|) + 1.5 \text{ MeV}}, \\ Y_1^a(i) = -\frac{\frac{1}{2}\Omega_i^{1/2}\Lambda(+2)}{2(|\epsilon_i| - |\epsilon_{p_{1/2}}|) + 5.32 \text{ MeV}}. \end{cases}$$

The corresponding numerical values are displayed in Tables 3.5.4 and 3.5.5, while in Fig. 3.5.6 we display a schematic summary of the graphical solution of the dispersion relations.

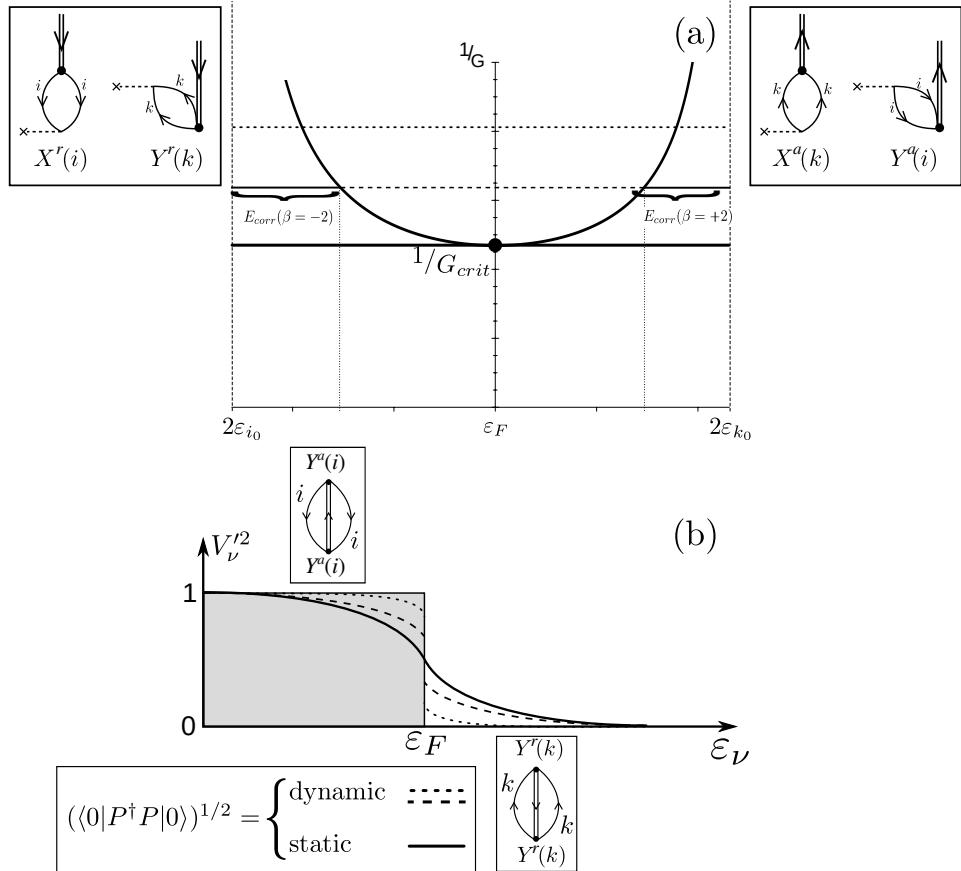


Figure 3.5.6: Schematic representation of the quantal phase transition taking place as a function of the pairing coupling constant in a closed shell nucleus. (a) dispersion relation associated with the RPA diagonalization of the Hamiltonian $H = H_{sp} + H_p$ for the pair addition and pair removal modes. In the insets are shown the two-particle transfer processes exciting these modes, which testify to the fact that the associated zero point fluctuations (ZPF) which diverge at $G = G_{crit}$, blur the distinction between occupied and empty states typical of closed shell nuclei. (b) occupation number associated with the single-particle levels. For $G < G_{crit}$ there is a dynamical depopulation (population) of levels $i(k)$ below (above) the Fermi energy. For $G > G_{crit}$, the deformation of the Fermi surface becomes stable, although with a non-vanishing dynamic component (see Fig. 3.1.2).

units		MeV	MeV ⁻¹	
<i>nlj</i>	Ω_k	$ \epsilon_{g_{9/2}} - \epsilon_k $	$C(k) = \frac{\frac{1}{2}\Omega_k^{1/2}}{2(\epsilon_{g_{9/2}} - \epsilon_k) + 1.5 \text{ MeV}}^a$	$X_1^a(k)$
$1g_{9/2}$	5	0	0.745	0.82
$0i_{11/2}$	6	0.77	0.403	0.44
$0j_{15/2}$	8	1.41	0.327	0.36
$2d_{5/2}$	3	1.56	0.187	0.21
$3s_{1/2}$	1	2.03	0.090	0.10
$1g_{7/2}$	4	2.47	0.155	0.17
$2d_{3/2}$	2	2.51	0.108	0.12

Table 3.5.4: Forwards going RPA amplitudes associated with the pair addition mode of ^{208}Pb (cf. Table XVI Broglia et al. (1973)). a), in the TD approximation (amplitudes not shown) the particle-pair addition coupling constant is calculated from the normalization of the amplitudes $C(k)$. That is, $\sum_k C^2(k) = 0.903 \text{ MeV}^{-2}$ and $\Lambda_1(+2) = (0.903)^{-1/2} \text{ MeV} = 0.950 \text{ MeV}$.

units		MeV	MeV ⁻¹	
<i>nlj</i>	Ω_i	$ \epsilon_i - \epsilon_{p_{1/2}} $	$D(i) = \frac{\frac{1}{2}\Omega_i^{1/2}}{2(\epsilon_i - \epsilon_{p_{1/2}}) + 5.32 \text{ MeV}}^a$	$Y_1^a(i)$
$2p_{1/2}$	1	0	0.094	-0.1
$1f_{5/2}$	3	0.57	0.134	-0.15
$2p_{3/2}$	2	0.90	0.099	-0.11
$0i_{13/2}$	7	1.64	0.154	-0.17
$1f_{7/2}$	4	2.35	0.100	-0.11
$0h_{9/2}$	5	3.47	0.091	-0.10

Table 3.5.5: Same as Table 3.5.4 but for the backwards going amplitude. a) $\sum_i D^2(i) = 0.079$ and $\Lambda_1^2(+2)(\sum_k C^2(k) - D^2(i)) = \Lambda_1^2(+2)(0.903 - 0.079) \text{ MeV}^{-2} = 0.824 \text{ MeV}^{-2}$; $\Lambda_1(+2) = (0.824)^{-1/2} \text{ MeV}$, thus $\Lambda_1(+2) = 1.102 \text{ MeV}$.

Let us conclude this Section by noting that while the harmonic (RPA) description of the pair vibrational mode of ^{208}Pb provides a fair picture of the two neutron transfer spectroscopic amplitudes, in keeping with the collective character of these (coherent) states, conspicuous anharmonicities in the multi-phonon spectrum have been observed and calculated⁴⁸. Within the framework of Fig. 3.1.5, we schematically emphasize in Fig. 3.5.7 the relative importance of dynamic and static pairing distortions, in comparison with the corresponding quantities in the case of quadrupole surface distortions in 3D-space⁴⁹. These results underscore the major role pairing vibrations play in nuclei around closed shells, while those shown in Fig. 3.1.2 emphasize their importance in gauge invariance restoration in systems far away from closed shells.

3.6 Halo pair addition and pygmy dipole modes: a new mechanism to break gauge invariance

Pairing correlations are intimately connected with particle number violation and thus spontaneous breaking of gauge invariance, as testified by the order parameter $\langle BCS | P^\dagger | BCS \rangle = \alpha_0$. In the nuclear case and, at variance with condensed matter, dynamical breaking of gauge symmetry is similarly important to that associated with static distortions. The fact that the average single-particle potential acts as an external field is one of the reasons for the existence of a critical value G_c of the pairing strength G to bind Cooper pairs in nuclei. Spatial quantization in finite systems at large and in nuclei in particular, is intimately connected with the paramount role the surface plays in these systems⁵⁰. Another consequence of this role is the fact that in nuclei an important fraction ($\approx 50\%$ in the case of nuclei lying along the stability valley and even more, up to 80% , for light halo nuclei) of Cooper pair binding is due to the exchange of collective vibrations between the pair partners⁵¹, the rest being associated with the bare NN -interaction in the 1S_0 channel (see Fig. 3.6.1) plus possible $3N$ corrections⁵².

The study of light exotic nuclei lying along the neutron drip line have revealed a novel aspect of the interplay between shell effects and induced pairing interaction. It has been found that there are situations in which spatial quantization screens, essentially completely, the bare nucleon-nucleon paring interaction. This happens

⁴⁸Cf. for example Flynn et al. (1972a), Lanford and McGrory (1973), Bortignon et al. (1978); Clark et al. (2006).

⁴⁹For details cf. Bès and Broglia (1977), Broglia et al. (1968), Bès et al. (1988), Barranco and Broglia (1987) Shimizu et al. (1989), Shimizu (2013), Vaquero et al. (2013) and references therein.

⁵⁰See Bohr and Mottelson (1975); see also Broglia (2002) and references therein.

⁵¹See Barranco et al. (1999), Terasaki et al. (2002a), Brink and Broglia (2005), Saperstein and Baldo (2013), Avdeenkov and Kamerdzhev (2013), Lombardo et al. (2013); Barranco et al. (2001), Potel et al. (2010), Pankratov et al. (2011), Barranco et al. (1999); cf. also Bohr and Mottelson (1975), p. 432.

⁵²See footnote 10 Ch. 1.

$$\begin{aligned}
P^\dagger &= \sum_{\nu>0} a_\nu^\dagger a_\nu^\dagger \\
x' &= \frac{2G\Omega'}{D} = GN(0) \\
&\quad \left| \begin{array}{l} x' > 1 \\ \alpha_0 = \langle P^\dagger \rangle = \frac{\Delta}{G} \approx 7 \end{array} \right. \quad \left| \begin{array}{l} x' > 1 \\ \alpha_{dyn} = \frac{\langle PP^\dagger \rangle^{1/2} + \langle P^\dagger P \rangle^{1/2}}{2} \\ \approx \frac{1}{2} \left(\frac{E_{corr}(A+2)}{G} + \frac{E_{corr}(A-2)}{G} \right) \approx 10 \end{array} \right. \\
&\quad \frac{\alpha_0}{\alpha_{dyn}} \approx 0.7 \\
&\quad \frac{\beta_2}{(\beta_2)_{dyn}} \approx 3 - 6
\end{aligned}$$

Figure 3.5.7: Relative importance of dynamic and static pairing distortion (α_{dyn} and α_0 respectively) associated with closed shell and open shell nuclei, calculated in terms of a two level model, as compared with similar quantities for the case of quadrupole surface degrees of freedom (β_2 -values). The parameter x' (product of the effective pairing strength $G' = Z_\omega^2(v_p^{bare} + v_p^{ind})$ and of the effective density of levels at the Fermi energy $N'(0) = Z_\omega^{-1}N(0) = Z_\omega^{-1}(2\Omega/D) = 2\Omega'/D = 2\Omega/D'; \Omega' = Z_\omega^{-1}\Omega, D' = Z_\omega D$), measures the relative importance of the single-particle gap $D' = Z_\omega D$ and of the pair correlation $G'\Omega$ (see Sect. 3.2; see also Brink and Broglia (2005) App. H, Sect. H.4). It is of notice that in referring to α_0 , one has actually in mind α'_0 ($\alpha_0 = e^{-2i\phi}\alpha'_0$, α'_0 being a real quantity). We nonetheless avoid the use of α'_0 in keeping with the use of primes to refer to renormalized quantities, both here and in Sect. 3.2.

in the case in which the nuclear valence orbitals are s, p -states at threshold⁵³. An example of situations of this type is provided by some of the $N = 7$ isotones, in particular ^3_3Li and, to some extent, ^4_4Be . Nuclei which display “parity inversion” in the sequence of single-particle levels $1p_{1/2}$ and $2s_{1/2}$ as compared to the Mayer-Jensen prediction (Fig. 1.1.2). In what follows we discuss the (unbound) nucleus ^{10}Li (see also Sect. 5.2.4), in connection with the (bound) two-neutron halo system ^{11}Li .

The $N = 7$ isotope ^{10}Li displays a virtual $s_{1/2}$ and a resonant $p_{1/2}$ state⁵⁴. Let us elaborate on this point. The binding provided by a contact pairing interaction $V_\delta(|\mathbf{r} - \mathbf{r}'|)$ (δ -force) to a pair of fermions moving in time-reversal states in a single j -shell⁵⁵ is given by the matrix element,

$$M_j = \langle j^2(0) | V_\delta | j^2(0) \rangle = -\frac{(2j+1)}{2} V_0 I(j) \approx -\frac{(2j+1)}{2} V_0 \frac{3}{R^3}. \quad (3.6.1)$$

It is of notice that $G = V_0 I(j)$ ($\approx 25/A$ MeV ≈ 2.3 MeV ($A = 11$)). The ratio of the above matrix element associated with the two halo neutrons of ^{11}Li and with an hypothetical normal nucleus of mass $A = 11$ is

$$r = \frac{(M_j)_{\text{halo}}}{(M_j)_{\text{core}}} = \frac{2}{(2j+1)} \left(\frac{R_0}{R} \right)^3. \quad (3.6.2)$$

The quantities $R_0 = 1.2A^{1/3}\text{fm} = 2.7\text{fm}$ ($A = 11$), and $R = \sqrt{\frac{5}{3}} \langle r^2 \rangle_{^{11}\text{Li}}^{1/2} = \sqrt{\frac{5}{3}} (3.55 \pm 0.1) \text{ fm} = (4.58 \pm 0.13) \text{ fm}$ are the radius of a nucleus of mass $A = 11$ (systematics), and the measured (mean square) radius of ^{11}Li , respectively. The quantity j is the effective angular momentum of a single j -shell which can accommodate 8 neutrons ($j \sim k_F R_0 \approx 1.36 \text{ fm}^{-1} \times 2.7 \text{ fm} \approx 3.7$). One thus obtains

$$r = 0.048. \quad (3.6.3)$$

Consequently, the bare NN -nucleon pairing interaction is expected to become strongly screened, the resulting effective G -value ($A \approx 11$)

$$G_{\text{scr}} = r \times G = 0.048 \times 25 \text{ MeV}/A \approx 1 \text{ MeV}/A \approx 0.1 \text{ MeV}, \quad (3.6.4)$$

becoming subcritical and thus unable to bind the halo Cooper pair ($2\tilde{\epsilon}_{s_{1/2}} = 0.3$ MeV, see Fig. 2.7.1) to the ^9Li core.

Further insight into this question can be shed making use of the multipole expansion of a general interaction

$$v(|\mathbf{r}_1 - \mathbf{r}_2|) = \sum_{\lambda} V_{\lambda}(r_1, r_2) P_{\lambda}(\cos \theta_{12}). \quad (3.6.5)$$

⁵³Pairing anti-halo effect; Bennaceur et al. (2000), Hamamoto and Mottelson (2003), Hamamoto and Mottelson (2004).

⁵⁴See Barranco et al. (2020), Moro et al. (2019) and refs. therein, as well as footnotes 30, 31 and 32 of Ch. 5.

⁵⁵cf. e.g. Eq. (2.12) Brink and Broglia (2005).

Because the function P_λ drops from its maximum at $\theta_{12} = 0$ in an angular distance $1/\lambda$, particles 1 and 2 interact through the component λ of the force, only if $r_{12} = |\mathbf{r}_1 - \mathbf{r}_2| < R/\lambda$, where R is the mean value of the radii \mathbf{r}_1 and \mathbf{r}_2 . Thus, as λ increases, the effective force range decreases. For a force of range much greater than the nuclear size, only the lowest λ (long wavelength) terms are important. At the other extreme, a δ -function force has coefficients $V_\lambda(r_1, r_2) \left(= \frac{(2\lambda+1)}{4\pi r_1^2} \delta(r_1 - r_2)\right)$ that increase with λ . In the case of $^{11}\text{Li}(\text{gs})$ one is thus confronted with the need to accept a long range, low λ pairing interaction, as responsible for the binding of the dineutron, halo Cooper pair to the ^9Li core. This is equivalent to saying, an induced pairing interaction arising from the exchange of vibrations with low λ -value.

3.6.1 Cooper pair binding: a novel embodiment of the Axel-Brink hypothesis.

In what follows we discuss a possible novel test of the Axel-Brink hypothesis⁵⁶. Within the s, p subspace, the most natural low multipolarity, long-wavelength vibration is the dipole mode. From systematics, the centroid of these vibrations is found at $\hbar\omega_{GDR} \approx 100 \text{ MeV}/R$, R being the nuclear radius⁵⁷. Thus, in the case of ^{11}Li , one expects the centroid of the *Giant Dipole Resonance* (GDR) carrying $\approx 100\%$ of the energy weighted sum rule (EWSR) at $\hbar\omega_{GDR} \approx 100 \text{ MeV}/4.6 \approx 22 \text{ MeV}$. Such a high frequency mode can hardly be expected to give rise to anything, but polarization effects. On the other hand, there exists experimental evidence which testifies to the presence in ^{11}Li , of a well defined dipole resonance with centroid at $\lesssim 1 \text{ MeV}$ and carrying $\approx 8\%$ of the EWSR⁵⁸. The existence of this “pygmy (dipole) resonance” (PDR) which can be viewed as a consequence of the existence of a low-lying particle-hole state associated with the transition $s_{1/2} \rightarrow p_{1/2}$ testifies, arguably, to the *coexistence*⁵⁹ of two states with rather different radii in the

⁵⁶The color of an object can be determined in two ways: by illuminating it with white light and see which wavelength it absorbs, or by heating it up and see the wavelength it emits. In both cases one is talking about dipole radiation. To describe the de-excitation process of hot nuclei requires the knowledge of the photon interactions with excited states. The common assumption, known as the Axel-Brink hypothesis, has been that each excited state of a nucleus carries a giant dipole resonance (GDR) on top of it, and that the properties of such resonances are unaffected by any excitation of the nucleus (Brink (1955) (unpublished), Lynn (1968) pag. 321, Axel (1962); cf. also Bertsch and Broglia (1986), Bortignon et al. (1998), Martin et al. (2017) and refs. therein).

⁵⁷See Bohr and Mottelson (1975) Bortignon et al. (1998) and Bertsch and Broglia (2005) and refs. therein.

⁵⁸See footnote 121 Ch. 2.

⁵⁹Within this context one can mention similar situations concerning the coexistence of spherical and quadrupole deformed states (cf. e.g. Wimmer et al. (2010), Federman and Talmi (1965), Federman and Talmi (1966), Dönau et al. (1967) and refs. therein; cf. also Bohr and Mottelson (1963 (unpublished)), typically of nuclei with $N \approx Z$. The fact that the associated inhomogeneous damping on the GDR has modest consequences concerning dipole strength at low energies as compared with radial (isotropic) deformations in ^{11}Li is understood in terms of the plasticity (non-Newtonian material) of the atomic nucleus regarding quadrupole deformations (low-lying collective 2^+ surface vibrations, fission, exotic decay (cf. Barranco et al. (1988), Barranco et al. (1989); Bertsch (1988),

ground state. One, closely connected with the ${}^9\text{Li}$ core, (≈ 2.5 fm), the second with the diffuse halo (≈ 4.6 fm), namely displaying a large radial (isotropic) deformation (neutron skin), and thus able to induce a conspicuous inhomogeneous damping to the dipole mode.

The importance of this mechanism is underscored by the fact that ${}^{11}\text{Li}$, displaying a neutron excess $(N - Z)/A \approx 0.45$ as compared to the value of 0.21 in the case of ${}^{208}\text{Pb}$, is able to bring down by tens of MeV a consistent fraction ($\approx 8\%$) of the TRK-sum rule. A consequence of the fact that the nucleus is most sensitive to changes in density (saturation phenomena). In the case of the halo of ${}^{11}\text{Li}$ we are confronted with nuclear structure phenomena in a medium displaying a density of $\approx 4\%$ of saturation density (Sect. 4.10, see also Eq. (4.3.6)).

Before proceeding, let us estimate the overlap O between the two “ground states”. Making use of a schematic expression for the single-particle radial wave-functions⁶⁰

$$\mathcal{R} = \sqrt{3/R_0^3} \Theta(r - R_0), \quad (3.6.6)$$

where

$$\Theta = 1 \quad (r \leq R_0); \quad 0 \quad (r > R_0),$$

leading to,

$$\int_0^\infty dr r^2 \mathcal{R}^2(r) = \frac{3}{R_0^3} \int_0^{R_0} dr r^3 / 3 = 1, \quad (3.6.7)$$

one can work out the overlap O between the two halo neutrons and the core nucleons. That is,

$$\begin{aligned} O &= |\langle \mathcal{R}_{\text{halo}} | \mathcal{R}_{\text{core}} \rangle|^2 = \left(\sqrt{\frac{3}{R_0^3}} \sqrt{\frac{3}{R^3}} \int_0^\infty dr r^2 \Theta(r - R) \Theta(r - R_0) \right)^2 \\ &= \left(\sqrt{\frac{3}{R_0^3}} \sqrt{\frac{3}{R^3}} \int_0^{R_0} dr r^3 / 3 \right)^2 = (R_0/R)^3 = 0.20, \end{aligned} \quad (3.6.8)$$

where use has been made of $\Theta(r - R)\Theta(r - R_0) = \Theta(r - R_0)$, $R_0 = 1.2A^{1/3}$ fm = 2.7 fm ($A = 11$) and $R = (4.58 \pm 0.013)$ fm. Because of the small value of this overlap, one can posit that the $E_x \lesssim 1$ MeV ($\Gamma \approx 0.5$ MeV) soft $E1$ -mode of ${}^{11}\text{Li}$ is a *bona fide* dipole resonance based on an exotic, unusually extended $|0,\rangle$ state of radial dimensions equivalent, according to systematics, to a system of effective A mass number about 5 times that of the actual system ($A \approx (4.6/1.2)^3 \approx 60$). Thus consistent with the connotation of PDR.

It is of notice that the small values of r and of O have essentially the same origin. On the other hand, they have apparently, rather different physical consequences. In fact, the first makes the bare pairing interaction strength G subcritical,

Bertsch et al. (1987)), and of the little tolerance to both compressibility and rarification displayed by the same system and connected with saturation properties (see also Broglia et al. (2019a) and Broglia et al. (2019b)).

⁶⁰Bohr and Mottelson (1969).

while the second one screens the repulsive symmetry potential $V_1(\approx +25 \text{ MeV})^{61}$, that is, the energy price one has to pay to separate protons from neutrons. This effect allows for a consistent fraction of the dipole Thomas–Reiche–Kuhn sum rule, that is of the $J^\pi = 1^-$ energy weighted sum rule (EWSR), to come low in energy ($p_{1/2} - s_{1/2}$ transition) from the value $E_{GDR} \approx (100/R) \text{ MeV}$ and, acting as an intermediate boson between the two halo neutrons, glue them to the ${}^9\text{Li}$ core. *Summing up, the halo anti-pairing effect $G_{scr} = r \times G \ll G < G_{crit}$ triggers ($OV_1 \ll V_1$) the virtual presence of a “gas” of dipole (pygmy) bosons which, exchanged between the two halo neutrons (cf. Fig. 3.6.2), overcompensates the reduction of the bare pairing interaction, leading to the binding of the halo Cooper pair to the core (anti-(halo anti-pairing effect)). It can thus be stated that the halo of ${}^{11}\text{Li}$ and the pygmy dipole resonance⁶² built on top of it constitute a pair of symbiotic states (see also Chapter 7).*

Let us further elaborate on these issues. Making use of the relation $\langle r^2 \rangle^{1/2} \approx (3/5)^{1/2} R$ between mean square radius and the radius, one may write

$$\langle r^2 \rangle_{{}^{11}\text{Li}} \approx \frac{3}{5} R_{eff}^2({}^{11}\text{Li}). \quad (3.6.9)$$

with

$$R_{eff}^2({}^{11}\text{Li}) = \left(\frac{9}{11} R_0^2({}^9\text{Li}) + \frac{2}{11} \left(\frac{\xi}{2} \right)^2 \right), \quad (3.6.10)$$

where

$$R_0({}^9\text{Li}) = 2.5 \text{ fm}, \quad (3.6.11)$$

is the ${}^9\text{Li}$ radius ($R_0 = r_0 A^{1/3}$, $r_0 = 1.2 \text{ fm}$), while ξ is the correlation length of the Cooper pair neutron halo. An estimate of this quantity is provided by the relation

$$\xi = \frac{\hbar v_F}{\pi E_{corr}} \approx 20 \text{ fm}, \quad (3.6.12)$$

in keeping with the fact that in ${}^{11}\text{Li}$, $(v_F/c) \approx 0.16$ and $E_{corr} \approx -0.5 \text{ MeV}$ (see App. 7.C). Consequently,

$$R_{eff}({}^{11}\text{Li}) \approx 4.8 \text{ fm}, \quad (3.6.13)$$

and $\langle r^2 \rangle_{{}^{11}\text{Li}}^{1/2} \approx 3.7 \text{ fm}$, to be compared with the experimental value⁶³ $\langle r^2 \rangle_{{}^{11}\text{Li}}^{1/2} = 3.55 \pm 0.1 \text{ fm}$. It is of notice that this experimental value implies the radius $R({}^{11}\text{Li}) = \sqrt{5/3 \langle r^2 \rangle_{{}^{11}\text{Li}}} = 4.58 \pm 0.13 \text{ fm}$.

We now proceed to the calculation of the centroid of the dipole pygmy resonance of ${}^{11}\text{Li}$ in RPA making use of the separable interaction

$$H_D = -\kappa_1 \vec{D} \cdot \vec{D} \quad (3.6.14)$$

⁶¹See e.g. Bortignon et al. (1998) Eq. (3.48) and refs. therein.

⁶²In which the two halo neutrons oscillate out of phase with respect to nucleons of the core ${}^9\text{Li}$ moving in phase (see Broglia et al. (2019a)).

⁶³Kobayashi et al. (1989).

where $\vec{D} = \vec{r}$ and

$$\kappa_1 = \frac{-5V_1}{AR^2}. \quad (3.6.15)$$

$V_1 = 25$ MeV being the symmetry potential energy. The RPA dispersion relation is⁶⁴

$$W(E) = \sum_{k,i} \frac{2(\epsilon_k - \epsilon_i)|\langle \tilde{i}|F|k\rangle|^2}{(\epsilon_k - \epsilon_i)^2 - E^2} = \frac{1}{\kappa_1}. \quad (3.6.16)$$

Making use of this relation as well as (see Fig. 2.7.1) $\epsilon_{\nu_k} - \epsilon_{\nu_i} = \tilde{\epsilon}_{p_{1/2}} - \tilde{\epsilon}_{s_{1/2}} \approx 0.45$ MeV, and that the EWSR associated with the ^{11}Li pygmy resonance is $\approx 8\%$ of the total Thomas-Reiche-Kuhn sum rule⁶⁵

$$\sum_n |\langle 0|F|n\rangle|^2(E_n - E_0) = \frac{\hbar^2}{2M} \int d\mathbf{r} |\vec{\nabla} F|^2 \rho(r), \quad (3.6.17)$$

which, for $F = r$ has the value⁶⁶ $\hbar^2 A / 2M$ one can write the numerator of Eq. (3.6.16) as,

$$2 \times 0.08 \times \frac{\hbar^2 A}{2M} = \frac{1}{\kappa_1} [(0.45 \text{ MeV})^2 - (\hbar\omega_{\text{pygmy}})^2], \quad (3.6.18)$$

and thus

$$(\hbar\omega_{\text{pygmy}})^2 = (0.45 \text{ MeV})^2 - 2 \times 0.08 \times \frac{\hbar^2 A}{2M} \kappa_1, \quad (3.6.19)$$

where⁶⁷

$$\kappa_1 = -\frac{5V_1}{A(\xi/2)^2} \left(\frac{2}{11} \right) = -\frac{125 \text{ MeV}}{A 100 \text{ fm}^2} \left(\frac{2}{11} \right) \approx -0.021 \text{ fm}^{-2} \text{ MeV}, \quad (3.6.20)$$

the ratio in parenthesis reflecting the fact that only 2 out of 11 nucleons, slosh back and forth in an extended configuration with little overlap with the other nucleons. The quantity,

$$\kappa_1^0 = -\frac{5V_1}{AR_{\text{eff}}^2(^{11}\text{Li})} \approx 0.49 \text{ MeV fm}^{-2}, \quad (3.6.21)$$

is the “standard” self consistent dipole strength⁶⁸. The screening factor $s = (\kappa_1/\kappa_1^0) = (R_{\text{eff}}^2 \times (2/11)) / (\xi/2)^2 = 0.043$ is very close in magnitude to the ratio $r (= 0.048,$

⁶⁴See (3.30) p.55 of Bortignon et al. (1998).

⁶⁵The Thomas-Reiche-Kuhn sum rule (Bohr and Mottelson (1975); Bortignon et al. (1998)) TRK = $\frac{9}{4\pi} \frac{\hbar^2 e^2}{2m} \frac{NZ}{A} = 14.8 \frac{NZ}{A} e^2 \text{ fm}^2 \text{ MeV}$ has a value of $32.3 e^2 \text{ MeV fm}^2$ for $^{11}\text{Li}_8$. Assuming a systematic behaviour, the centroid of the giant dipole resonance is expected at $\hbar\omega_D \approx 80/A^{1/3}$ MeV ≈ 36 MeV, leading to the ratio $32.3 e^2 \text{ MeV fm}^2 / (36 \text{ MeV}) \approx 0.9 e^2 \text{ fm}^2$. The $E1$ -single-particle (Weisskopf) unit can be written as (Bohr and Mottelson (1969) p. 389, Eq. (3C-38)) $B_W(E1) \approx ((1.2)^2/4\pi)(3/4)^2 A^{2/3} (e_{E1})^2 \text{ fm}^2 = 0.32 (e_{E1})^2 \text{ fm}^2$, e_{E1} being the effective dipole charge equal to $(N/A)e = 0.73e$ for the protons of ^{11}Li , and $-(Z/A)e = 0.27$ for the neutrons. Making use of the average value one can write $\bar{B}_W(E1) \approx 0.1 e^2 \text{ fm}^2$. Thus $8\% \text{TRK} / \bar{B}_W(E1) \approx 0.072/0.1 \approx 0.7$. In other words, about 1 single-particle unit is associated with the eventual γ -decay of the PDR of ^{11}Li .

⁶⁶cf. Bertsch and Broglia (2005) pag. 53.

⁶⁷see Bortignon et al. (1998).

⁶⁸cf. Bohr and Mottelson (1975).

Eq. (3.6.3)) and has a similar physical origin. It is of notice that $(V_1)_{scr} = sV_1 \approx 1$ MeV. Making use of (3.6.20) one obtains,

$$-2 \times 0.08 \frac{\hbar^2 A}{2M} \kappa_1 \approx 0.74 \text{MeV}^2 \approx (0.86 \text{MeV})^2. \quad (3.6.22)$$

Consequently

$$\hbar\omega_{pygmy} = \sqrt{(0.45)^2 + (0.86)^2} \text{MeV} \approx 1 \text{ MeV}, \quad (3.6.23)$$

in overall agreement with the experimental findings⁶⁹. It is of notice that the centroid of the pygmy resonance (microscopically) calculated in QRPA with the help of a separable dipole interaction is⁷⁰ $\approx (0.6 \text{ MeV} + 1.6 \text{ MeV})/2 \approx 1.1 \text{ MeV}$.

Let us now estimate the binding energy which the exchange of the pygmy resonance between the two neutrons of the halo Cooper pair of ^{11}Li can provide. The associated particle–vibration coupling⁷¹ is $\Lambda = (\partial W(E)/\partial E|_{\hbar\omega_{pygmy}})^{-1/2}$. Note the use in what follows of a dimensionless dipole single-particle field $F' = F/R_{eff}(^{11}\text{Li})$. This is in keeping with the fact that one wants to obtain a quantity with energy dimensions ($[\Lambda] = \text{MeV}$), and that κ_1 has been introduced through the Hamiltonian H_D with the self consistent value normalized in terms of $R_{eff}^2(^{11}\text{Li})$ (Eq. (3.6.21)). One then obtains

$$\begin{aligned} \Lambda &= \left\{ 2\hbar\omega_{pygmy} \frac{2 \times 0.08(\frac{\hbar^2 A}{2M})/R_{eff}^2}{[(\tilde{\epsilon}_{p_{1/2}} - \tilde{\epsilon}_{s_{1/2}})^2 - (\hbar\omega_{pygmy})^2]^2} \right\}^{-1/2}, \\ &= \left\{ 2 \text{MeV} \frac{0.16(\hbar^2 A/2M)(1/4.8)^2 \text{ fm}^2}{[(0.45 \text{ MeV})^2 - (1 \text{ MeV})^2]^2} \right\}^{-1/2}, \\ &= \left(\frac{3 \text{ MeV}^2}{(0.8)^2 \text{ MeV}^4} \right)^{-1/2} \approx 0.5 \text{ MeV}, \end{aligned} \quad (3.6.24)$$

The value of the induced interaction matrix elements is then given by (Fig. 3.6.2),

$$M_{ind} = \frac{2\Lambda^2}{DEN} \approx -\frac{2\Lambda^2}{\hbar\omega_{pygmy}} \approx -0.5 \text{ MeV}, \quad (3.6.25)$$

the factor of two arising from the two time ordering contributions. The resulting correlation energy is thus $E_{corr} = 2\tilde{\epsilon}_{s_{1/2}} - G_{scr} + M_{ind} = (0.3 - 0.1 - 0.5) \text{ MeV} \approx -0.3$ MeV, in overall agreement with the experimental⁷² findings (-0.380 MeV). It is of notice that in this estimate the (subcritical) effect of the screened bare pairing

⁶⁹See footnote 121 Ch. 2.

⁷⁰Barranco et al. (2001); in particular Fig. 2a where the doubled peaked ($\approx 0.6 \text{ MeV}$ and $\approx 1.6 \text{ MeV}$) $dB(E1)/dE$ strength function is displayed.

⁷¹cf. e.g. Brink and Broglia (2005) Eq. (8.42) p.189.

⁷²Bachelet et al. (2008), Smith et al. (2008).

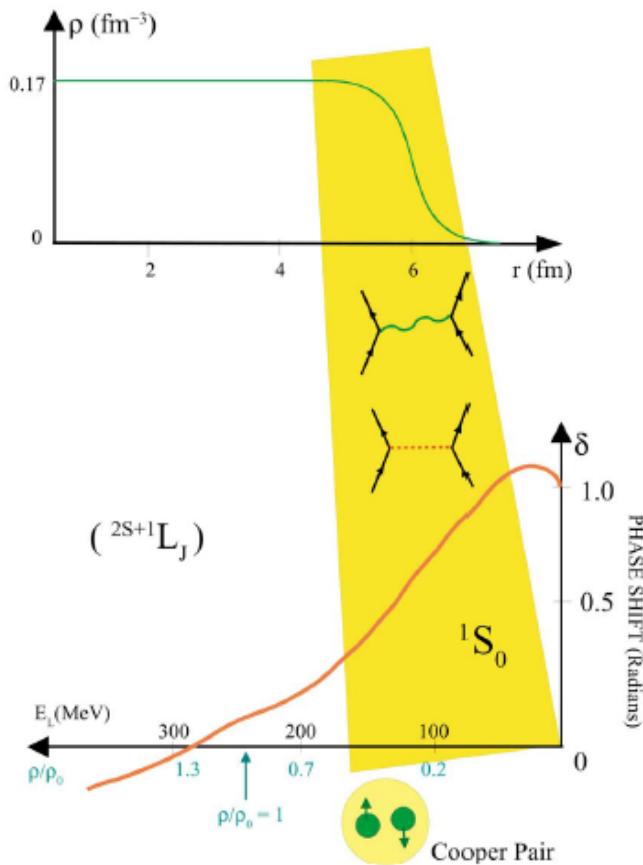


Figure 3.6.1: (top) Nuclear density ρ in units of fm $^{-3}$, plotted as a function of the distance r (in units of fm) from the center of a nucleus lying along the stability valley (for comparison with a bound unstable nucleus lying at the neutron drip line see Fig. 4.3.2). Saturation density correspond to ≈ 0.17 fm $^{-3}$, equivalent to 2.8×10^{14} g/cm 3 . Because of the short range of the nuclear force, the nuclear density changes from 90% of saturation density to 10% within 0.65 fm, i.e. within the nuclear diffusivity. (bottom) Phase shift parameter associated with the elastic scattering of two nucleons moving in states of time reversal, so called 1S_0 phase shift, in keeping with the fact that the system is in a singlet state of spin zero. Positive values of δ implies an attractive interaction, negative a repulsive one. For low relative velocities (kinetic energies E_L), i.e. for a situation similar to that found for nucleons moving around the nuclear surface where the density is low, the 1S_0 phase shift arising from the exchange of mesons (e.g. pions, represented by an horizontal dotted line) between nucleons (represented by upward pointing arrowed lines) is attractive. This mechanism provides about half of the glue to nucleons moving in time reversal states to form Cooper pairs. Cooper pair formation is further assisted by the exchange of collective surface vibrations (wavy curve in the scattering process) between the members of the pair.

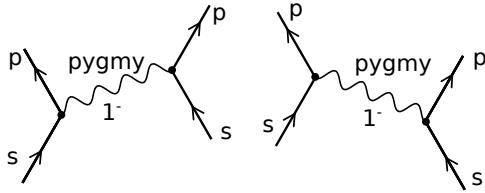


Figure 3.6.2: Diagrammatic representation of the exchange of a collective 1^- pygmy resonance between pairs of nucleons moving in the time-reversal configurations $s_{1/2}^2(0)$ and $p_{1/2}^2(0)$. It is of notice that both these configurations can act as initial states the figure showing only one of the two possibilities. Consequently, the energy denominator to be used in the simple estimate (3.6.25) is the average value $DEN = (DEN_1 + DEN_2)/2 = -\hbar\omega_{pygmy}$ where $DEN_1 = \Delta\epsilon - \hbar\omega_{pygmy}$ and $DEN_2 = -\Delta\epsilon - \hbar\omega_{pygmy}$, while $\Delta\epsilon = \epsilon_{s_{1/2}} - \epsilon_{p_{1/2}}$.

interaction has also been used (see Eq. (3.6.4))⁷³, as well as the theoretical value⁷⁴ $\tilde{\epsilon}_{1/2} = 0.15$ MeV (Fig. 2.7.1).

This schematic model⁷⁵ has been implemented with microscopic detail⁷⁶ within the framework of a field theoretical description of the interweaving of collective vibrations and single-particle motion, and is also discussed within the context of single-particle (Chapter 5) and two-particle (Chapter 7) transfer processes. Here we provide a summary of the theoretical findings.

In Fig. 3.6.3 (I), the lowest single-particle neutron virtual and resonant states of ^{10}Li are indicated⁷⁷. The position of the levels $s_{1/2}$ and $p_{1/2}$ determined making use of mean-field theory is shown (left hatched area and thin horizontal line, respectively). The coupling of a single-neutron (upward pointing arrowed line) to a vibration (wavy line) calculated making use of NFT Feynman diagrams (schematically depicted also in terms of either solid dots (neutron) or open circles (neutron hole) moving in a single-particle level around or in the ^9Li core (grey circle)), leads to conspicuous shifts in the energy centroid of the $s_{1/2}$ and $p_{1/2}$ resonances (shown by thick horizontal lines to the right) and eventually to an inversion in their sequence. In Fig. 3.6.3 (II) the processes contributing to binding the halo neutron system ^{11}Li are displayed. One starts with the dressed mean field picture in which two neutrons (solid dots) coupled to angular momentum and parity 0^+ move in time-reversal states around the core ^9Li (hatched area) in the $s_{1/2}$ virtual state leading to an unbound $s_{1/2}^2(0)$ configuration. The associated spatial structure

⁷³That new physics, namely a novel mechanism to (dynamically) violate gauge invariance finds, to express itself, a scenario of a barely bound Cooper pair at the drip line (half life 8.75 ms), seems to confirm a recurrent expectation. That truly new complex phenomena appear at the border between rigid order and randomness (see de Gennes (1994)).

⁷⁴See footnotes 30 and 31 of Ch. 5.

⁷⁵See also Broglia et al. (2019b).

⁷⁶cf. Barranco et al. (2001); see also Potel et al. (2010).

⁷⁷Barranco et al. (2001). For a more detailed (NFT) study of ^{10}Li see Barranco et al. (2020) (see also Sect. 5.2.4). See also Moro et al. (2019).

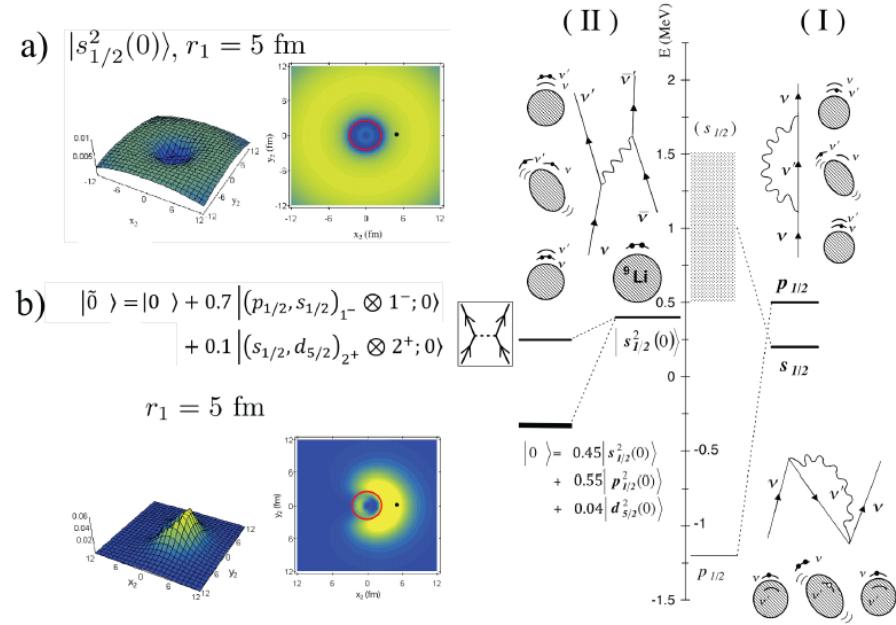


Figure 3.6.3: (Color online) In (I) and (II) the NFT processes renormalizing the single-particle motion (${}^{10}\text{Li}$) and leading to the effective interaction, sum of the bare (horizontal dotted line, inset) and induced (wavy line) interactions which bind the two-neutron halo to the core of ${}^9\text{Li}$ thus leading to the $|{}^{11}\text{Li}\rangle$ ground state are displayed. It is of notice that the odd $p_{3/2}(\pi)$ proton, considered as a spectator, is not shown. In a) and b) the spatial structure of the pure $|s_{1/2}^2(0)\rangle$ configuration and that of the two-neutron halo $|\tilde{0}\rangle$ Cooper pair is displayed. The modulus squared wave function $|\Psi_0(\mathbf{r}_1, \mathbf{r}_2)|^2 = |\langle \mathbf{r}_1, \mathbf{r}_2 | 0^+ \rangle|^2$ describing the motion of the two halo neutrons around the ${}^9\text{Li}$ core is shown as a function of the cartesian coordinates of particle 2, for fixed values of the position of particle 1 ($r_1 = 5$ fm) represented by a solid dot, while the core radius (${}^9\text{Li}$), is shown as a solid circle. The numbers appearing on the z -axis of the three-dimensional plots displayed on the left side of the figure are in units of fm^{-2} .

of the uncorrelated pair is shown in **a**). The exchange of vibrations between the two neutrons displayed in the upper part of the figure leads to an induced pairing interaction which, added to the bare nucleon-nucleon Argonne potential⁷⁸ (see boxed inset) which, as can be seen from the figure, is subcritical, correlates the two-neutron system leading to a bound state $|\tilde{0}\rangle$. The corresponding wavefunction is displayed in **b**), together with the spatial structure of the resulting Cooper pair. It is of notice that a large fraction of the induced interaction arises from the exchange of the pygmy dipole resonance between the two halo neutrons. Within this scenario one can posit that the ^{11}Li PDR can hardly be viewed but in symbiosis with the ^9Li halo neutron pair addition mode and vice versa. Furthermore, that the two halo neutrons of ^{11}Li , provide the first example of a Van der Waals Cooper pair, the first of its type in nuclei (Sect.. 3.7 and Fig. 3.7.1). For further details see Chapter 7 as well as⁷⁹.

3.7 Nuclear van der Waals Cooper pair

The atomic van der Waals (dispersive, retarded) interaction which, like gravitation, acts between all atoms and molecules, also non-polar, can be written for two systems placed at a distance R as (see App. 3.B),

$$\Delta E = -\frac{6 \times e^2 \times a_0^5}{R^6} = -\frac{16}{(R/a_0)^6} \left(\frac{e^2}{2a_0} \times 0.75 \right), \quad (3.7.1)$$

where a_0 is the Bohr radius. One recognizes in the term in parenthesis the transition energy $2P \rightarrow 1S$ of the hydrogen atom. A possible qualitative nuclear parallel (see Fig. 3.7.1) can be established making the correspondences of the term in parenthesis with $\tilde{\epsilon}_{p_{1/2}} - \tilde{\epsilon}_{s_{1/2}} \approx 0.45$ MeV, R with $R_{eff} \approx 4.6$ fm and a_0 with $R_0(^{11}\text{Li}) \approx 3$ fm, leading to $\Delta E \approx -0.6$ MeV ($= M_{ind}$). Thus,

$$E_{corr} = 2\tilde{\epsilon}_{s_{1/2}} - G_{scr} + \Delta E = 0.3 \text{ MeV} - 0.1 \text{ MeV} - 0.6 \text{ MeV} \approx -0.4 \text{ MeV}.$$

3.8 Renormalized coupling constants ^{11}Li : resumé

The fact that the screening factors r and s (see Eqs. (3.6.3), (3.6.2); also (3.6.21) and subsequent paragraph), essentially coincide within numerical approximations is in keeping with the fact that both quantities are closely related to the overlap⁸⁰

$$O = \left(\frac{R_0}{R} \right)^3 = \left(\frac{2.7 \text{ fm}}{4.6 \text{ fm}} \right)^3 = 0.2, \quad (3.8.1)$$

⁷⁸Wiringa et al. (1984).

⁷⁹Barranco et al. (2001).

⁸⁰One can equivalently use $(R_0/R_{eff})^3 \approx (2.7 \text{ fm}/(4.8 \text{ fm}))^3 \approx 0.18$.

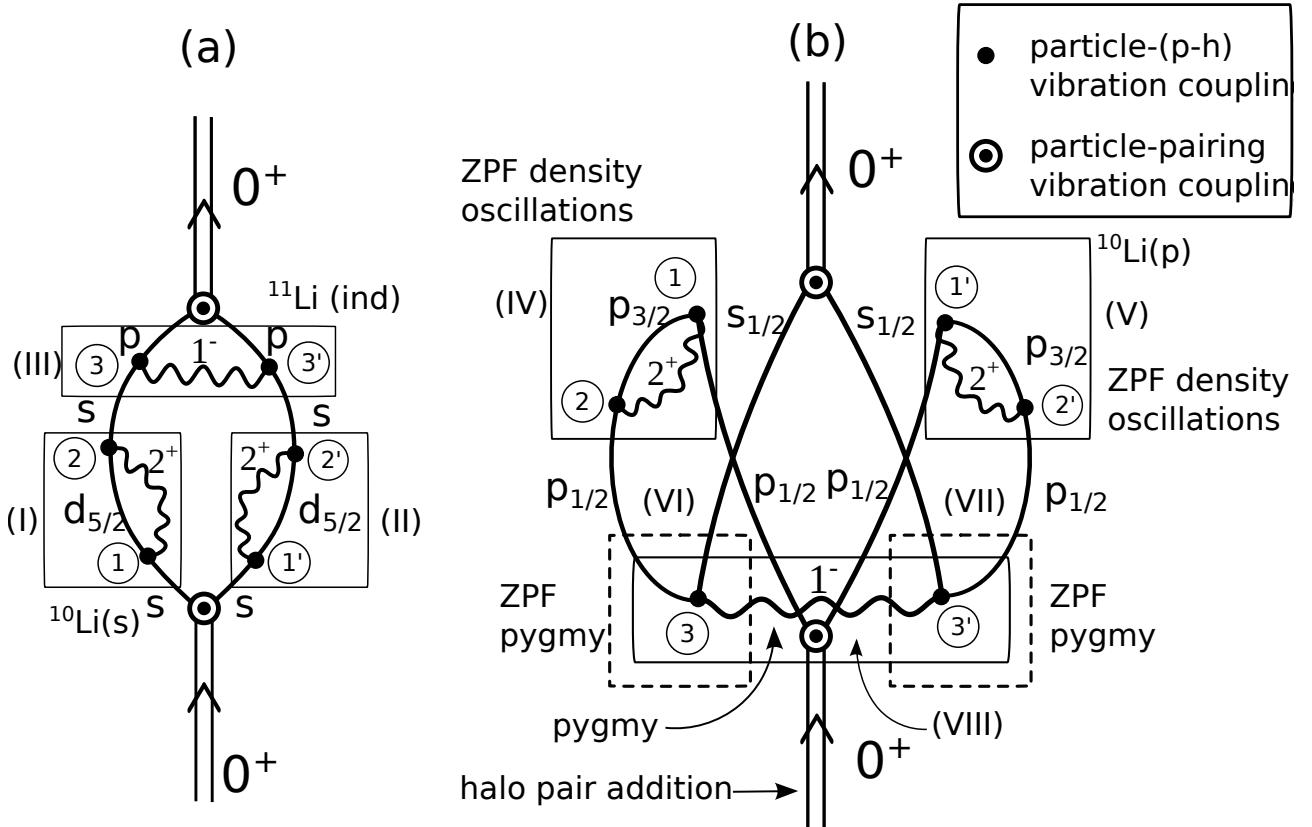


Figure 3.7.1: NFT Feynman diagrams describing the binding of the halo Cooper pair through pygmy. That is, producing the symbiotic mode involving the pair addition mode and the PDR. The single-particle states $s_{1/2}$ and $p_{1/2}$ are labeled in (a) s and p for simplicity. The different particle-vibration coupling vertices (either with the quadrupole (2^+) or with the pygmy (1^-) modes drawn as wavy lines) are denoted by a solid dot, and numbered in increasing time sequence so as to show that diagram (b) emerges from (a) through time ordering. The motion of the neutrons are drawn in terms of continuous solid curves. In keeping with the fact that the occupation of the single-particle states is neither 1 nor 0 we treat, for simplicity, these states as quasiparticle states. Thus no arrow is drawn on them. Diagram (a) emphasizes the self-energy renormalization of the state $s_{1/2}$ lying in the continuum and which through its clothing with the quadrupole mode is brought down becoming a virtual ($\tilde{\epsilon}_{s_{1/2}} \approx 0.2$ MeV) state (see (I) and (II)), while (III) contributes to the induced pairing interaction through pygmy (see also Fig. 3.6.2). The “eagle” diagram (b) contains ((IV) and (V)) Pauli corrections which push the bound state $p_{1/2}$ into a resonant state in the continuum ($\tilde{\epsilon}_{p_{1/2}} \approx 0.5$ MeV). In other words, processes (I), (II), (III), (IV) and (V) are at the basis of parity inversion, and of the appearance of the new magic number $N = 6$. Processes (VI) and (VII) are associated with the pygmy ZPF, while (VIII) contributes to the induced pairing interaction through pygmy (van der Waals-like process).

quantity which has two main effects concerning the mechanism which is at the basis of much of the nuclear structure of the halo exotic nucleus ^{11}Li : **1)** it makes subcritical the screened bare NN -pairing interaction $G_{scr} = r \times G < G_c$ ($G_{scr} = 1 \text{ MeV}/A$); **2)** it screens the symmetry potential drastically, reducing the energy price one has to pay to separate protons from delocalized neutrons, permitting a consistent chunk ($\approx 8\%$) of the TRK sum rule⁸¹ to become essentially degenerate with the ground state ($(V_1)_{scr} = 1 \text{ MeV}$). In other words, allowing for, allegedly, the first nuclear example of a van der Waals Cooper pair and a novel mechanism to break dynamically gauge invariance. Namely, dipole–dipole fluctuating fields associated with the exchange of the pygmy dipole resonance between the halo neutrons of ^{11}Li . As a result, a new, (composite) elementary mode of nuclear excitation joins the ranks of the previously known: the halo pair addition mode carrying on top of it, a low-lying collective pygmy resonance. This symbiotic mode can be studied through two-particle transfer reactions, eventually in coincidence with γ -decay. In particular, making use of the reaction $^9\text{Li}(t, p)^{11}\text{Li}(f)$ for

$|f\rangle$: ground state ($L = 0$) and pygmy ($L = 1; E_x \lesssim 1 \text{ MeV}$).

Similar, but in this case $^{10}\text{Be}(t, p)^{12}\text{Be}(f)$ for

$|f\rangle$: first excited 0^+ state ($E_x = 2.24 \text{ MeV}$),

as well as pygmy ($L = 1$) on top of it, arguably the 1^- state at $E_x = 2.70 \text{ MeV}$ being part of this mode⁸².

Appendix 3.A Lindemann criterion and connection with quantity parameter

The original Lindemann criterion⁸³ compares the atomic fluctuation amplitude $\langle \Delta r^2 \rangle^{1/2}$ with the lattice constant a of a crystal. If this ratio, which is defined as the disorder parameter Δ_L , reaches a certain value, fluctuations cannot increase without damaging or destroying the crystal lattice. The results of experiments and simulations show that the critical value of Δ_L for simple solids is in the range of 0.10 to 0.15, relatively independent of the type of substance, the nature of the interaction potential, and the crystal structure⁸⁴. Applications of this criterion to an inhomogeneous finite system like a protein in its native state (aperiodic crystal)⁸⁵, requires evaluation of the generalized Lindemann parameter⁸⁶

$$\Delta_L = \frac{\sqrt{\sum_i \langle \Delta r_i^2 \rangle / N}}{a'}, \quad (3.A.1)$$

⁸¹See footnote 121 Ch. 2.

⁸²Iwasaki et al. (2000).

⁸³Lindemann (1910).

⁸⁴Bilgram (1987); Löwen (1994); Stillinger (1995).

⁸⁵Schrödinger (1944).

⁸⁶Stillinger and Stillinger (1990).

where N is the number of atoms and a' the most probable non-bonded near-neighbor distance, \mathbf{r}_i being the position of atom i , $\Delta r_i^2 = (\mathbf{r}_i - \langle \mathbf{r}_i \rangle)^2$, and $\langle \rangle$ denotes configurational averages at the conditions of measurement or simulations (e.g. biological, in which case $T \approx 310$ K, $\text{PH} \approx 7$, etc.⁸⁷). The dynamics as a function of the distance from the geometric center of the protein is characterized by defining an interior (*int*) Lindemann parameter,

$$\Delta_L^{int}(r_{cut}) = \frac{\sqrt{\sum_{i,r_i < r_{cut}} \langle \Delta r_i^2 \rangle / N}}{a'}, \quad (3.A.2)$$

which is obtained by averaging over the atoms that are within a chosen cutoff distance, r_{cut} , from the center of mass of the protein.

Simulations and experimental data for a number of proteins, in particular Barnase, Myoglobin, Crambin and Ribonuclease A indicate 0.14 as the critical value distinguishing between solid-like and liquid-like behaviour, and $r_{cut} \approx 6$ Å. As can be seen from Table 3.A.1, the interior of a protein, under physiological conditions, is solid-like ($\Delta_L < 0.14$), while its surface is liquid-like ($\Delta_L > 0.14$). The beginning of thermal denaturation in the simulations appears to be related to the melting of its interior (i.e. $\Delta_L^{int} > 0.14$), so that the entire protein becomes liquid-like. This is also the situation of the denatured state of a protein under physiological conditions⁸⁸

	$\Delta_L(\Delta_L^{int}(6 \text{ \AA}))(300 \text{ K})$			
	MD simulations			X-ray data
Proteins	Barnase	Myoglobin	Crambin	Ribonuclease A
all atoms	0.21(0.12)	0.16(0.11)	0.16(0.09)	0.16(0.12)
backbone atoms only	0.16(0.10)	0.12(0.09)	0.12(0.08)	0.13(0.10)
side-chain atoms only	0.25(0.14)	0.18(0.12)	0.19(0.10)	0.19(0.13)

Table 3.A.1: The heavy-atom $\Delta_L(\Delta_L^{int})$ value, for four proteins at 300 K. After Zhou et al. (1999).

3.A.1 Lindemann (“disorder”) parameter for a nucleus

An estimate of $\sqrt{\sum_i \langle \Delta r_i^2 \rangle / A}$ in the case of nuclei considered as a sphere of nuclear matter of radius R_0 , is provided by the “spill out” of nucleons due to quantal effects. That is⁸⁹ $\sqrt{\dots} \approx 0.69 \times a_0$, where a_0 is of the order of the nuclear diffusivity (≈ 0.65 fm). The average internucleon distance can be estimated to be $2 \times r_0 \approx 2.4$ fm.

⁸⁷Fluctuations, classical (thermal) or quantal imply a probabilistic description. While one can only predict the odds for a given outcome of an experiment in quantum mechanics, probabilities themselves evolve in a deterministic fashion (Born (1948)).

⁸⁸see e.g. Rösner et al. (2017).

⁸⁹Bertsch and Broglia (2005), see e.g. Ch. 5. See also paragraph following Eq. (4.10.3).

Thus,

$$\Delta_L = \frac{0.69a_0}{2.4 \text{ fm}} \approx 0.19. \quad (3.A.3)$$

While it is difficult to compare among them crystals, aperiodic finite crystals and atomic nuclei, arguably, the above value indicates that a nucleus is liquid-like.

Appendix 3.B The van der Waals interaction

Historically one can distinguish two contributions to the van der Waals interaction⁹⁰:

1. **dispersive** retarded contribution⁹¹, emerging from the dynamical dipole-dipole, as well as from higher multipolarities, interaction associated with the quantum mechanical zero point fluctuations (ZPF) of the ground state of the two interacting atoms or molecules⁹².
2. **inductive** implying the polarization of one molecule in the permanent dipole or quadrupole field of the other molecule⁹³

It is only the first one which is a *bona fide* van der Waals interaction. In fact with the advent of quantum mechanics it was very early recognized⁹⁴ that for most molecules, interactions of type 2 are small compared with interactions of type 1. That is the interaction corresponding to the mutual polarization of one molecule in the rapidly changing field –due to the instantaneous configuration of electrons and nuclei– of the other molecule⁹⁵.

3.B.1 van der Waals interaction between two hydrogen atoms

For large values of the internuclear distance $r_{AB} = R$, the exchange phenomenon is unimportant (Pauli principle) and one can take as the unperturbed wavefunction for a system of two hydrogen atoms the simple product of two hydrogenlike wavefunctions,

$$\Psi^0 = u_{1sA}(1)u_{1sB}(2). \quad (3.B.1)$$

The perturbation for this wavefunction arises from the potential energy terms

$$H' = \frac{e^2}{r_{12}} + \frac{e^2}{r_{AB}} - \frac{e^2}{r_{A2}} - \frac{e^2}{r_{B1}}, \quad (3.B.2)$$

⁹⁰Let us mainly think of non polar (NP) molecules.

⁹¹dispersion: variation of a quantity, e.g. spatial separation of white light (rainbow), as a function of frequency (c.f. e.g. Israelachvili (1985), p.65).

⁹²These forces act between all atoms and molecules, even non-polar, totally neutral ones.

⁹³Debye (1920, 1921).

⁹⁴London (1930).

⁹⁵Pauling and Wilson Jr. (1963) p. 384, Born (1969) p. 471.

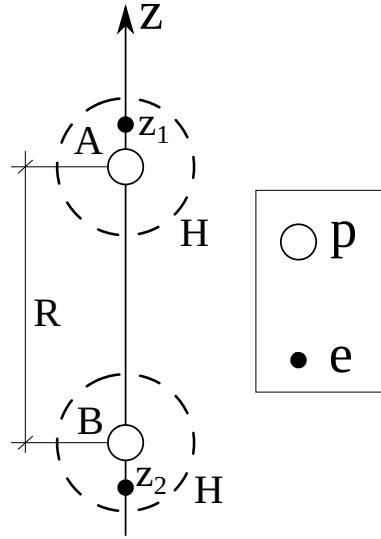


Figure 3.B.1: Planar configuration assumed for two hydrogen atoms at a relative distance R , where p stands for proton and e for electron.

corresponding to the variety of Coulomb interactions involving electrons and protons. Let us assume for simplicity that we are dealing with a one-dimensional problem, in which case one can write (Fig. 3.B.1)

$$\begin{aligned}\mathbf{r}_{12} &= (R + z_1 + z_2) \hat{\mathbf{z}}, \\ \mathbf{r}_{AB} &= R \hat{\mathbf{z}}, \\ \mathbf{r}_{A2} &= (R + z_2) \hat{\mathbf{z}}, \\ \mathbf{r}_{B1} &= (R + z_1) \hat{\mathbf{z}}.\end{aligned}\quad (3.B.3)$$

Because all these distances are much larger than the radius of the atom ($a_0 \approx 0.529 \text{ \AA}$, Bohr radius) the expression (3.B.2) can be calculated making use of a Taylor expansion. One obtains

$$r_{12}^2 = (R + z_1 + z_2)^2 = R^2 \left[1 + 2\frac{(z_1 + z_2)}{R} + \frac{(z_1 + z_2)^2}{R^2} \right], \quad (3.B.4)$$

which leads to

$$\frac{e^2}{r_{12}} = \frac{e^2}{R \left[1 + \frac{2(z_1 + z_2)}{R} + \frac{(z_1 + z_2)^2}{R^2} \right]^{1/2}} \approx \frac{e^2}{R} \left[1 - \frac{(z_1 + z_2)}{R} - \frac{(z_1 + z_2)^2}{2R^2} \right]. \quad (3.B.5)$$

Similarly

$$r_{A2}^2 = (R^2 + 2Rz_2 + z_2^2) = R^2 \left(1 + 2\frac{z_2}{R} + \frac{z_2^2}{R^2} \right), \quad (3.B.6)$$

and

$$r_{B1}^2 = R^2 \left(1 + 2 \frac{z_1}{R} + \frac{z_1^2}{R^2} \right), \quad (3.B.7)$$

leading to

$$-\frac{e^2}{r_{A2}} = -\frac{e^2}{R} \left(1 - \frac{z_2}{R} - \frac{z_2^2}{2R^2} \right), \quad (3.B.8)$$

and

$$-\frac{e^2}{r_{B1}} = -\frac{e^2}{R} \left(1 - \frac{z_1}{R} - \frac{z_1^2}{2R^2} \right). \quad (3.B.9)$$

Finally

$$\frac{e^2}{r_{AB}} = \frac{e^2}{R}. \quad (3.B.10)$$

With the exception of the cross term of (3.B.5) there is complete cancellation between the different contributions to (3.B.2). Thus

$$H' = -\frac{\mathbf{D}_1 \cdot \mathbf{D}_2}{R^3}, \quad (3.B.11)$$

where

$$\mathbf{D}_i = ez_i \hat{\mathbf{z}} \quad (3.B.12)$$

is the dipole moment operator associated with electron i . Because $R \gg z_i$, one can diagonalize the interaction Hamiltonian (3.B.11) perturbatively. In keeping with the fact that a single-particle quantum state displaying a given parity (and in the present case angular momentum $(-1)^\ell = \pi$) cannot sustain a permanent dipole moment, in particular

$$\int d\tau u_{1s}(z) \times z \times u_{1s}(z) = \int d\tau (u_{1s}(z))^2 z = \int d\tau \rho(z) z = 0, \quad (3.B.13)$$

the lowest perturbative correction to (3.B.1) is of second order. This is in keeping with the fact that we are dealing with interaction through virtual processes. The associated energy correction is given by the relation,

$$\Delta E_z^{(2)} = - \sum_{int} \frac{\langle 0 | H' | int \rangle \langle int | H' | 0 \rangle}{E_{int} - E_0}. \quad (3.B.14)$$

Because we are concerned with the $1s \rightarrow 2p$ transition (Fig. 3.B.2), the intermediate state is

$$\Psi^{int} = u_{2pA}(1)u_{2pB}(2), \quad (3.B.15)$$

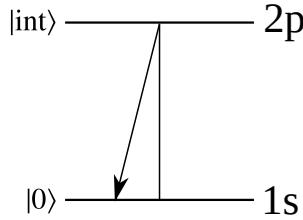


Figure 3.B.2: Schematic representation of the virtual process associated with (3.B.14), the intermediate, virtual state being the $2p$ state.

and thus

$$D_{en} = E_{int} - E_0. \quad (3.B.16)$$

One can then write

$$\begin{aligned} \Delta E_z^{(2)} &= -\frac{|\langle 0|H'|0\rangle|^2}{D_{en}} = -\frac{e^4}{R^6} \frac{|\langle 0|z_1^2 z_2^2|0\rangle|^2}{D_{en}}, \\ &\approx -\frac{e^4}{R^6} \frac{\int d\tau_1 d\tau_2 u_{1s}^2(1) u_{1s}^2(2) z_1^2 z_2^2}{D_{en}}, \\ &= -\frac{e^4}{R^6} \frac{\int d\tau_1 \rho(z_1) z_1^2 \int d\tau_2 \rho(z_2) z_2^2}{D_{en}}, \\ &= -\frac{e^4}{R^6} \frac{\bar{z}_1^2 \bar{z}_2^2}{D_{en}} \approx -\frac{e^4}{R^6} \frac{a_0^2 a_0^2}{\frac{e^2}{2a_0}} = -\frac{2e^2 a_0^5}{R^6}. \end{aligned} \quad (3.B.17)$$

This result corresponds to the z -degree of freedom of the system (two H atoms at a distance $R \gg a_0$). One has then to multiply the above result by 3 to take into account the x and y degrees of freedom. Thus

$$\Delta E^{(2)} = -\frac{6e^2 a_0^5}{R^6}. \quad (3.B.18)$$

Let us now calculate the van der Waals interaction between two H-atoms at a distance of the order of ten times the summed radii of the two atoms ($\approx 2a_0 \approx 1\text{\AA}$), that is for $R \approx 10\text{\AA}$,

$$\begin{aligned} \Delta E_{H-H}^{(2)}(10 \text{\AA}) &\approx -\frac{6 \times 14.4 \text{ eV \AA}(0.529 \text{\AA})^5}{(10 \text{\AA})^6} \\ &\approx -3.6 \times 10^{-6} \text{ eV} = -3.6 \mu\text{eV} \end{aligned} \quad (3.B.19)$$

Making use of the relation

$$1 \text{ eV} = 2.42 \times 10^{14} \text{ Hz}, \quad (1\text{Hz}=\text{s}^{-1}) \quad (3.B.20)$$

one obtains

$$|\Delta E_{H-H}^{(2)}(10 \text{ \AA})| \approx 3.6 \times 10^{-6} \times 2.42 \times 10^{14} \text{ Hz} \approx 9 \times 10^8 \text{ Hz} \approx 10^3 \text{ MHz}, \quad (3.B.21)$$

a quantity which can be compared with the Lamb shift (1058 MHz; see Fig. 5.2.4). It is of notice that $|\Delta E_{H-H}^{(2)}(2.5 \text{ \AA})| \approx 15 \text{ meV/part} \approx 0.35 \text{ kcal/mole}$, ($1 \text{ meV/part} \approx 0.02306 \text{ kcal/mole}$), a value of the order of $kT/2$. That is, one half of the thermal energy under biological conditions ($T \approx 300 \text{ K}$, $kT \approx 0.6 \text{ kcal/mole}$)⁹⁶, in keeping with the role played by the Van der Waals interaction in the folding and stability of proteins.

⁹⁶Huang (2005).

Chapter 4

Phase-coherent pair transfer

A superconductor has rather perfect internal phase order... The importance of the Josephson effect... is that... it can pin down the order parameter.

P. W. Anderson

4.1 Simultaneous versus successive Cooper pair transfer in nuclei

Cooper pair transfer is commonly thought to be tantamount to simultaneous transfer. In this process a nucleon goes over through the NN -interaction v , the second one does it making use of the correlations with its partner (cf. Figs. 4.1.1 and 6.5.1 (I)). Consequently, in the independent particle limit, simultaneous transfer should not be possible (see Sect. 6.5.1). Nonetheless, it remains operative. This is because, in this limit, the particle transferred through v does it together with a second one which profits from the non-orthogonality of the wavefunctions describing the single-particle motion in target and projectile (Figs. 4.1.2 and 6.5.1 (II)). This is the reason why this (non-orthogonality) transfer amplitude has to be treated on equal footing with the previous one representing, within the non-orthogonal basis employed, a natural contribution to simultaneous transfer. In fact $(T^{(1)} - T_{NO}^{(1)})$ is the correct, two-nucleon transfer amplitude to lowest order (first) in v . The resulting cancellation is quite conspicuous in actual nuclei (see Figs. 4.5.2 (b), 4.5.3, and Fig. 4.5.4 for examples in open shell (superfluid) and closed shell (normal) nuclei respectively). This is in keeping with the fact that Cooper pairs are weakly correlated systems ($2\Delta \ll \epsilon_F$), and the reason why the successive transfer process in which v (e.g. the mean field U in the post-post representation¹) acts twice, is

¹Bayman and Chen (1982); Potel et al. (2013a). See also Pinkston and Satchler (1982).

the dominant mechanism in pair transfer reactions. While this mechanism seems antithetical to the transfer of correlated fermions pairs, it probes, in the nuclear case, the same pairing correlations as simultaneous transfer does². This is because nuclear Cooper pairs (quasi-bosons) are quite extended objects, the two nucleons being (virtually) correlated over distances ξ larger than typical nuclear dimensions (see Fig. 4.1.3). In a two-nucleon transfer process this virtual property becomes real, in the sense that the presence of (normal) density over regions larger than that of the dimensions of each of the interacting nuclei allows for Cooper pair (abnormal density) manifestation over distances of the order of ξ .

Within this context, let us refer to the Josephson effect, associated with the Cooper pair tunneling across a thin barrier separating two metallic superconductors. Because the probability of one-electron-tunneling is small, (conventional) simultaneous tunneling associated with that probability quantity square would hardly be observed. Nonetheless, Josephson currents are standard measures in low temperature laboratories³. The same arguments related to the large value of the correlation length is operative in explaining the fact that Coulomb repulsion is rather weak between partners of Cooper pairs which are, in average, at a distance $\xi (\approx 10^4 \text{ \AA})$ much larger than the Wigner-Seitz radius r_s typical of metallic elements ($\approx 1 - 2 \text{ \AA}$). Consequently, it can be overcompensated and turned into an effective (weakly) attractive force by the long range electron-phonon pairing.

Similarly, in widely extended light halo nuclei, the short range bare pairing interaction may become subcritical due to the poor overlap between the halo neutrons and the core ones (cf. Sect. 3.6). The fact that such systems are nonetheless bound, although weakly, testifies to the important role the exchange of collective vibrations between halo nucleons have in binding the associated halo Cooper pair.

The above arguments are at the basis of the fact that second order DWBA theory which add both successive and non-orthogonality contributions to the simultaneous transfer amplitudes, are needed in a quantitative description of the experimental findings.

4.2 One- and two-nucleon transfer probabilities

In what follows we define transfer probabilities P_1 and P_2 for one- and two-nucleon transfer. For this purpose let us remind some useful relations. In particular that of the differential reaction cross section

$$\frac{d\sigma}{d\Omega} = |f(\theta)|^2, \quad (4.2.1)$$

²Within this context note however the possible interplay in reactions between heavy ions at energies below the Coulomb barrier, between the distance of closest approach and the correlation length ξ (see Sect. 7.3 as well as Eqs. (4.9.5)–(4.9.7)).

³cf. e.g. Rogalla and Kes (2012) and references therein.

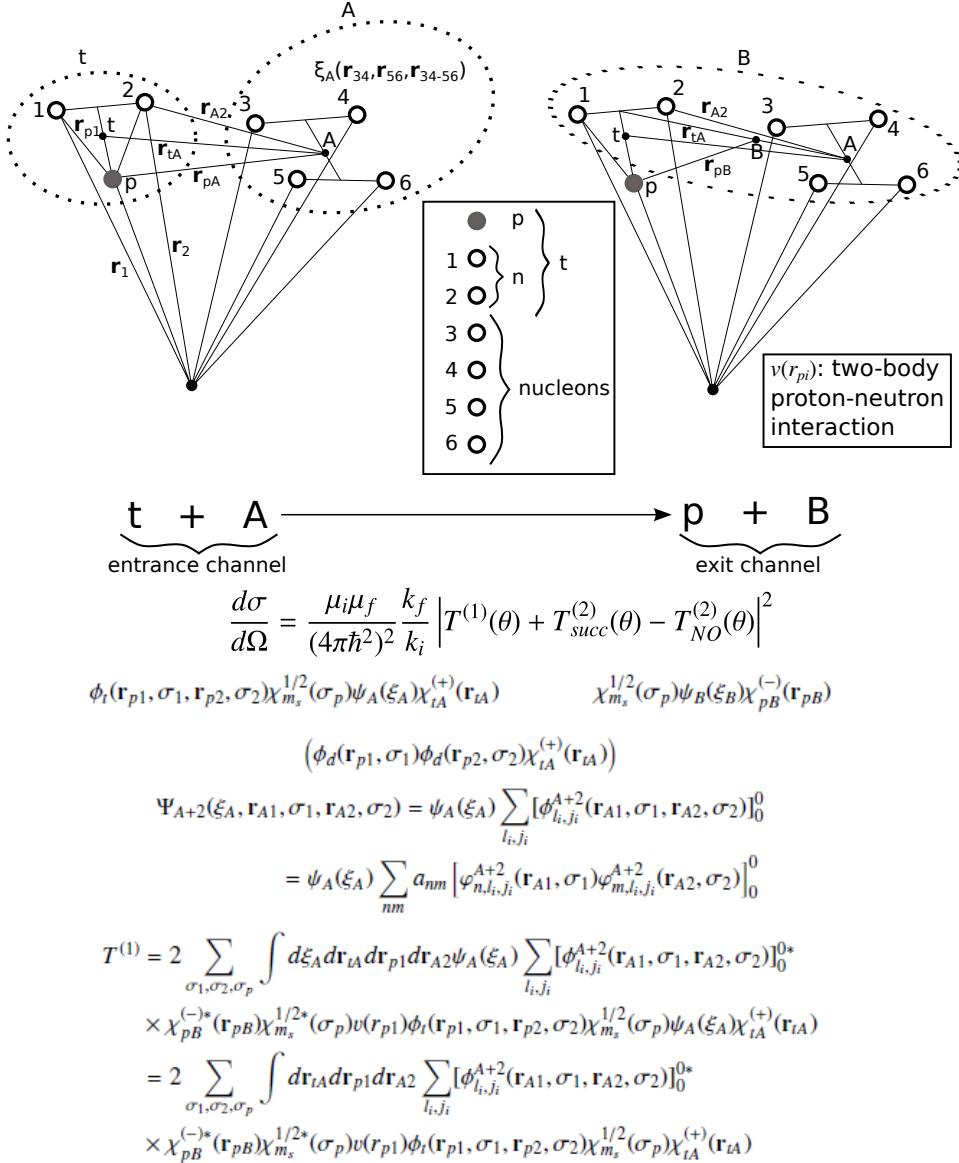
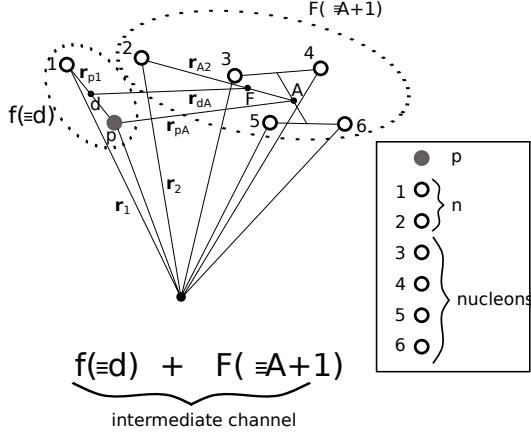


Figure 4.1.1: Contribution of simultaneous transfer, in first order DWBA, to the reaction $A(t, p)B(\equiv A + 2)$. The nucleus A is, for simplicity, assumed to contain four nucleons, the triton being composed of two neutrons and one proton. The set of coordinates used to describe the entrance and exit channels are shown in the upper part (bold face vectors represent the coordinates used to describe the relative motion, while the intrinsic coordinates ξ_A are \mathbf{r}_{34} , \mathbf{r}_{56} and \mathbf{r}_{34-56}). In the lower part of the figure, the simultaneous two-nucleon transfer amplitude is written (see Potel et al. (2013b)). It is of notice that of all the relative motion coordinates, only those describing the relative motion of (t, A) and of (p, B) have asymptotic values, being those associated with distorted waves.



$$\begin{aligned}
& \chi_{m_s}^{1/2}(\sigma_p) \phi_d(\mathbf{r}_{p1}, \sigma_1) \psi_A(\xi_A) \varphi_{l_f, j_f, m_f}^{A+1}(\mathbf{r}_{A2}, \sigma_2) \\
G(\mathbf{r}_{dF}, \mathbf{r}'_{dF}) &= i \sum_l \sqrt{2l+1} \frac{f_l(k_{dF}, r_<) g_l(k_{dF}, r_>)}{k_{dF} r_{dF} r'_{dF}} \left[Y^l(\hat{r}_{dF}) Y^l(\hat{r}'_{dF}) \right]_0^0 \\
T_{succ}^{(2)} &= 2 \sum_{l_i, j_i} \sum_{l_f, j_f, m_f} \sum_{\sigma_1 \sigma_2} \int d\xi_A d\mathbf{r}_{dF} d\mathbf{r}_{p1} d\mathbf{r}_{A2} \chi_{pB}^{(-)*}(\mathbf{r}_{pB}) \psi_B^*(\xi_B) v(\mathbf{r}_{p1}) \phi_d(\mathbf{r}_{p1}) \varphi_{l_f, j_f, m_f}^{A+1}(\mathbf{r}_{A2}, \sigma_2) \\
&\times \chi_{m_s}^{1/2}(\sigma_p) \Psi_A(\xi_A) \frac{2\mu_{dF}}{\hbar^2} \int d\xi'_A d\mathbf{r}'_{dF} d\mathbf{r}'_{p1} d\mathbf{r}'_{A2} G(\mathbf{r}_{dF}, \mathbf{r}'_{dF}) \\
&\times \chi_{tA}^{(+)}(\mathbf{r}_{tA}) \psi_A^*(\xi'_A) v(\mathbf{r}'_{p2}) \phi_d(\mathbf{r}'_{p1}) \varphi_{l_f, j_f, m_f}^{A+1}(\mathbf{r}'_{A2}, \sigma'_2) \\
&= 2 \sum_{l_i, j_i} \sum_{l_f, j_f, m_f} \sum_{\sigma_1 \sigma_2} \int d\mathbf{r}_{dF} d\mathbf{r}_{p1} d\mathbf{r}_{A2} \chi_{pB}^{(-)*}(\mathbf{r}_{pB}) v(\mathbf{r}_{p1}) \phi_d(\mathbf{r}_{p1}) \left[\varphi_{l_f, j_f, m_f}^{A+2}(\mathbf{r}_{A1}, \sigma_1, \mathbf{r}_{A2}, \sigma_2) \right]_0^0 \\
&\times \frac{2\mu_{dF}}{\hbar^2} \int d\mathbf{r}'_{dF} d\mathbf{r}'_{p1} d\mathbf{r}'_{A2} G(\mathbf{r}_{dF}, \mathbf{r}'_{dF}) \chi_{tA}^{(+)}(\mathbf{r}'_{tA}) v(\mathbf{r}'_{p2}) \phi_d(\mathbf{r}'_{p1}, \sigma'_1) \phi_d(\mathbf{r}'_{p2}, \sigma'_2) \varphi_{l_f, j_f, m_f}^{A+1}(\mathbf{r}'_{A2}, \sigma'_2) \\
T_{NO}^{(1)} &= 2 \sum_{l_i, j_i} \sum_{l_f, j_f, m_f} \sum_{\sigma_1 \sigma_2} \int d\xi_A d\mathbf{r}_{dF} d\mathbf{r}_{p1} d\mathbf{r}_{A2} \chi_{pB}^{(-)*}(\mathbf{r}_{pB}) \chi_B^*(\xi_B) v(\mathbf{r}_{p1}) \phi_d(\mathbf{r}_{p1}) \varphi_{l_f, j_f, m_f}^{A+1}(\mathbf{r}_{A2}, \sigma_2) \\
&\times \chi_{m_s}^{1/2}(\sigma_p) \Psi_A(\xi_A) \frac{2\mu_{dF}}{\hbar^2} \int d\xi'_A d\mathbf{r}'_{dF} d\mathbf{r}'_{p1} d\mathbf{r}'_{A2} \\
&\times \chi_{tA}^{(+)}(\mathbf{r}_{tA}) \psi_A^*(\xi'_A) \phi_d(\mathbf{r}'_{p1}) \varphi_{l_f, j_f, m_f}^{A+1}(\mathbf{r}'_{A2}, \sigma'_2) \\
&= 2 \sum_{l_i, j_i} \sum_{l_f, j_f, m_f} \sum_{\sigma_1 \sigma_2} \int d\mathbf{r}_{dF} d\mathbf{r}_{p1} d\mathbf{r}_{A2} \chi_{pB}^{(-)*}(\mathbf{r}_{pB}) v(\mathbf{r}_{p1}) \phi_d(\mathbf{r}_{p1}) \left[\varphi_{l_f, j_f, m_f}^{A+2}(\mathbf{r}_{A1}, \sigma_1, \mathbf{r}_{A2}, \sigma_2) \right]_0^0 \\
&\times \frac{2\mu_{dF}}{\hbar^2} \int d\mathbf{r}'_{dF} d\mathbf{r}'_{p1} d\mathbf{r}'_{A2} \chi_{tA}^{(+)}(\mathbf{r}'_{tA}) \phi_d(\mathbf{r}'_{p1}, \sigma'_1) \phi_d(\mathbf{r}'_{p2}, \sigma'_2) \varphi_{l_f, j_f, m_f}^{A+1}(\mathbf{r}'_{A2}, \sigma'_2)
\end{aligned}$$

Figure 4.1.2: Successive and non-orthogonality contributions to the amplitude describing two-nucleon transfer in second order DWBA, entering in the expression of the absolute differential cross section $d\sigma/d\Omega = \frac{\mu_i \mu_f}{(4\pi\hbar^2)^2} \frac{k_f}{k_i} |T^{(1)} + T_{succ}^{(2)} - T_{NO}^{(1)}|^2$. Concerning $T^{(1)}$ we refer to Fig. 4.1.1. In the upper part of the figure the coordinates used to describe the intermediate channel $d + F(\equiv A + 1)$ are given (bold face vectors represent the coordinates used to describe the relative motion, while the intrinsic coordinates ξ_A are \mathbf{r}_{34} , \mathbf{r}_{56} and \mathbf{r}_{34-56}). In the lower part of the figure, the expressions corresponding to the (t, p) process are displayed (Potel et al., 2013b). Schematically, the three contributions $T^{(1)}$, $T_{succ}^{(2)}$ and $T_{NO}^{(1)}$ to the transfer amplitude can be written as $\langle pB|v|tA\rangle$, $\sum \langle pB|v|dF\rangle \langle dF|v|tA\rangle$ and $\sum \langle pB|v|dF\rangle \langle dF|\mathbf{1}|tA\rangle$ respectively, where v is the proton–neutron interaction and $\mathbf{1}$ the unit operator. Within this context, while $T_{NO}^{(1)}$ receives contributions from the intermediate (virtual) closed $(d + F)$ channel as $T_{succ}^{(2)}$ does, it is first order in v as $T^{(1)}$ is.

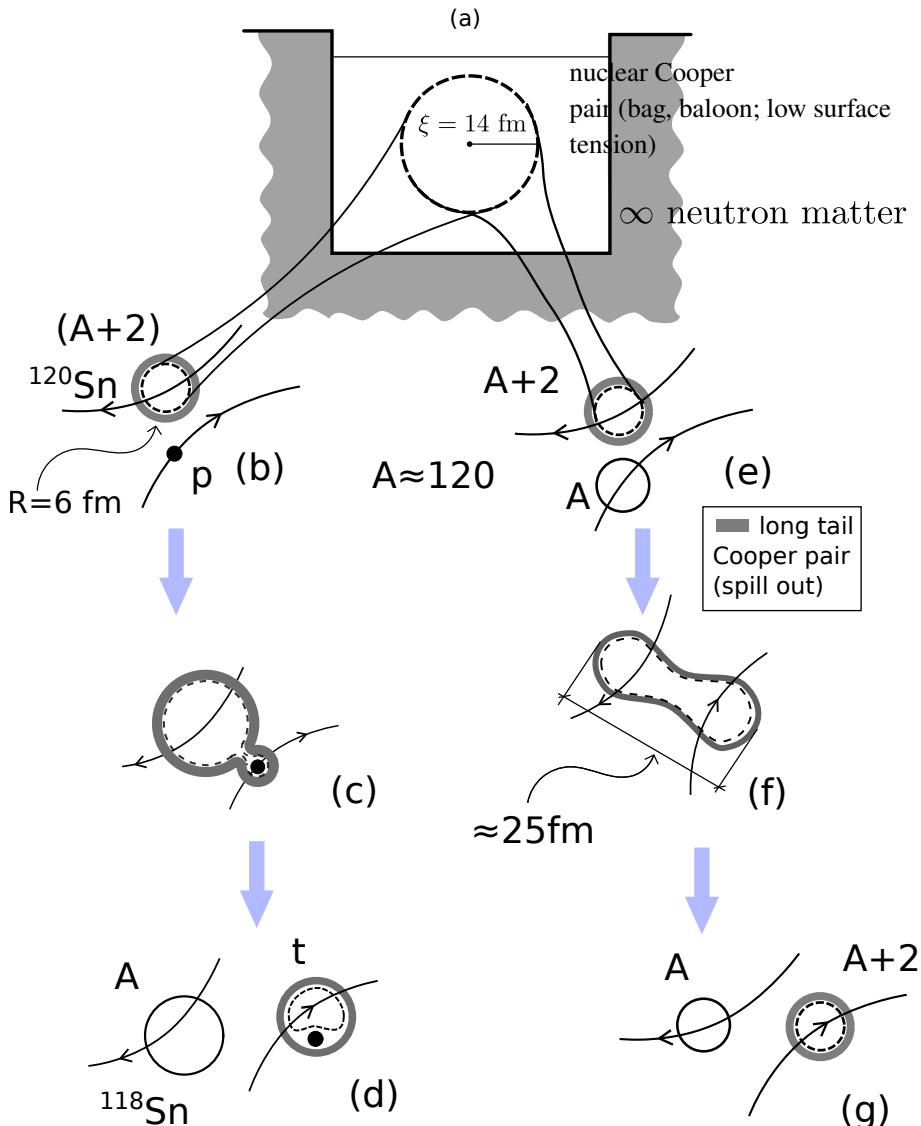


Figure 4.1.3: The correlation length associated with a nuclear Cooper pair is of the order of $\xi \approx \hbar v_F / (\pi \Delta) \approx 14 \text{ fm}$ ($\Delta \approx 1.2 \text{ MeV}$, $v_F/c \approx 0.27$). (a) in neutron matter at typical densities of the order of 0.5 saturation density, the $NN^{-1}S_0$ short range force, eventually renormalized by medium polarization effects, makes pairs of nucleons moving in time reversal states to correlate over distances larger than nuclear dimensions (dashed circle). How can one get evidence for such an extended object? (b) Hardly when the Cooper bag (balloon) is introduced in the mean field of a superfluid nucleus which, acting as a strong external field, leads to spatial quantization (related to angle correlations, see Sect. 4.11) and constrains the Cooper pair to be inside the nucleus with some spill out (tail of Cooper pair, grey, shaded area; see Bertsch and Broglia (2005) p. 88) as testified by (d) two-nucleon transfer process (e.g. (p, t) reaction), in which case the absolute cross section can change in a conspicuous fashion, in going from pure two-particle (uncorrelated configurations) to long tail Cooper pair spill outs. This effect is expected to become stronger by allowing pair transfer between similar superfluid nuclei (see Sect. 7.3), in which case the Cooper pair can move in the combined nuclear system resulting from the two heavy ions in weak contact, and at the same time profit of the same type of correlations (superfluidity) as resulting from similar pair mean fields (e), (f), (g) (see e.g. von Oertzen (2013); von Oertzen and Vitturi (2001), and references therein).

where

$$f(\theta) = \frac{1}{k} \sum_l (2l+1) e^{i\delta_l} \sin \delta_l P_l(\cos \theta), \quad (4.2.2)$$

δ_l being the partial wave l phase shift. Let us now use for simplicity the results associated with hard sphere scattering⁴ in the low and high energy limit. Making use of the fact that in the case under discussion the phase shifts δ_l are related to the regular and irregular spherical Bessel functions,

$$\tan \delta_l = \frac{j_l(kR)}{n_l(kR)}, \quad (4.2.3)$$

and that $\sin^2 \delta_l = \tan^2 \delta_l / (1 + \tan^2 \delta_l)$, one can write in the case in which $kR \ll 1$, i.e. in the low-energy, long wavelength, regime

$$\tan \delta_l \approx \frac{-(kR)^{2l+1}}{(2l+1)[(2l-1)!!]}, \quad (4.2.4)$$

implying that one can ignore essentially all δ_l with $l \neq 0$. Thus, $\tan \delta_0 = -kR$, $\sin^2 \delta_0 = (kR)^2$ and one can write,

$$\frac{d\sigma}{d\Omega} = \frac{\sin^2 \delta_0}{k^2} = R^2. \quad (4.2.5)$$

Consequently,

$$\sigma_{tot} = \int \frac{d\sigma}{d\Omega} d\Omega = 4\pi R^2 \quad (kR \ll 1), \quad (4.2.6)$$

a cross section which is four times the geometric cross section πR^2 , namely the area of the disc of radius R that blocks the propagation of the incoming (plane) wave, and has the same value as that of a hard sphere. Because $kR \ll 1$ implies long wavelength scattering, it is not surprising that quantal effects are important, so as to overwhelm the classical picture. Let us now consider the high energy limit $kR \gg 1$. The total cross section is in this case, given by

$$\begin{aligned} \sigma_{tot} &= \int |f_l(\theta)|^2 d\Omega = \frac{1}{k^2} \int_0^{2\pi} d\phi \int_{-1}^1 d(\cos \theta) \sum_{l=1}^{kR} \sum_{l'=1}^{kR} (2l+1)(2l'+1) \\ &\times e^{i\delta_l} \sin \delta_l \times e^{-i\delta_{l'}} \sin \delta_{l'} \times P_l P_{l'} = \frac{4\pi}{k^2} \sum_{l=1}^{kR} (2l+1) \sin^2 \delta_l. \end{aligned} \quad (4.2.7)$$

Making use of the relation

$$\sin^2 \delta_l = \frac{\tan^2 \delta_l}{1 + \tan^2 \delta_l} = \frac{[j_l(kR)]^2}{[j_l(kR)]^2 + [n_l(kR)]^2} \approx \sin^2 \left(kR - \frac{\pi l}{2} \right), \quad (4.2.8)$$

⁴cf. e.g. Sakurai (1994)

and the fact that so many l -values contribute to (4.2.7), one can replace $\sin^2 \delta_l$ by its average value $1/2$. Because the number of terms of the sum is roughly kR , the same being true for the average value of $(2l + 1)$, one can thus write

$$\sigma_{tot} = \frac{4\pi}{k^2} (kR)^2 \frac{1}{2} = 2\pi R^2, \quad (kR \gg 1) \quad (4.2.9)$$

which, in this short wavelength limit, is not the geometric cross section either. In fact, (4.2.9) can be split into two contributions each of value πR^2 . One due to reflection in which it can be shown that there is no interference amongst contributions from different l -values. A second one (coherent contribution in the forward direction) called shadow because for hard-sphere scattering at high energies, waves with impact parameter less than R must be deflected⁵. Consequently, behind the scatterer there must be zero probability for finding the scattered particle and a shadow must be generated⁶. The quantity

$$\lambda = \frac{\lambda}{2\pi} = \frac{h}{2\pi p} = \frac{\hbar}{p} = \frac{1}{k} = \frac{\hbar}{\sqrt{2mE}}, \quad (4.2.10)$$

is the reduced de Broglie wavelength for a massive particle ($E = p^2/2m$). For a proton of energy $E_p \approx 20$ MeV, typical of beams used in $^{120}\text{Sn}(p, t)^{118}\text{Sn(gs)}$ and $^{120}\text{Sn}(p, d)^{119}\text{Sn(j)}$ reactions⁷ $\lambda \approx 1$ fm, to be compared with the value $R \approx 6$ fm of the radius of ^{120}Sn . Consequently, we are in a situation of type (4.2.9), that is,

$$\sigma_{tot} = 2\pi(6)^2 \text{ fm}^2 \approx 2.3 \text{ b.} \quad (4.2.11)$$

⁵The original German word *Unbestimmtheitsprinzip* (“indefiniteness” or “indeterminacy principle”) is sometimes incorrectly translated as “uncertainty principle”. This is very misleading, since it suggests that the electron or the nucleon actually has a definite position and momentum of which one is uncertain. In fact, the quantum formalism simply does not allow the ascription of a definite position and momentum simultaneously (see Leggett (1987)).

⁶In terms of wave mechanics, this shadow is due to the destructive interference between the original wave (which would be there even if the scatterer was absent), and the newly scattered wave. Thus, one needs scattering in order to create a shadow. This contribution is intimately related to the optical theorem (Sakurai (1994))

$$\sigma_{tot} = \frac{4\pi}{k} \Im[f(\theta = 0, k)] = \frac{4\pi}{k} [f_{shad}(\theta = 0, k)] = \frac{4\pi}{k^2} \sum_l (2l + 1) \sin^2 \delta_l,$$

to which it provides its physical interpretation. In fact, there are two independent ways of measuring σ_{tot} , namely: i) by integrating the differential cross section $d\sigma/d\Omega = |f(\theta)|^2$ moving around the detector, ii) measuring the attenuation of the incoming beam. Both procedures should give the same result. One then identifies $(4\pi/k)f(\theta = 0, k)$ with the attenuation arising from the interference of the elastic wave with the incoming wave. Of notice that in (4.2.7) the factor $(\pi/k^2)(2l + 1) = \pi\lambda^2(2l + 1)$ is the area of a ring with radius $b = (l + 1/2)\lambda$ and width λ due to quantal indeterminacy. Thus

$$\sigma_{tot} = 2\pi(R + \lambda/2)^2 \quad (kR \gg 1).$$

⁷Of notice that the reduced wavelength of a photon ($p = E/c$) of the same energy ($E = 20$ MeV) is $\lambda(\lambda/2\pi = \hbar/p = \hbar c/E) \approx 10$ fm (cf. Table 2.1 p. 22 Satchler (1980)).

Because typical values of the absolute one-particle cross section associated with the (p, d) reaction mentioned above are few mb (see e.g. Fig. 5.2.3 right panel) one can use, for order of magnitude estimate purposes,

$$P_1 \approx \frac{5.35 \text{ mb}}{2.3 \text{ b}} \approx 10^{-3}, \quad (4.2.12)$$

as typical probability for such processes. *Consequently, one may argue that the probability for a pair of nucleons to simultaneously tunnel in e.g. the (p, t) process mentioned above is $(P_1)^2 \approx 10^{-6}$, as near impossible as no matter.* Within this context we note that the integrated $\text{gs} \rightarrow \text{gs}$ absolute cross section $\sigma(^{120}\text{Sn}(p, t)^{118}\text{Sn}(\text{gs})) \approx 2.25 \pm 0.338 \text{ mb}$, quantity representative of the superfluid Sn-isotopes ground state rotational band (see Figs. 3.1.3, 3.1.7 and 7.2.1)⁸. This fact implies that the empirical two-nucleon transfer probability is of the order of $P_2 \approx 10^{-3}$, i.e. $P_2 \approx P_1$.

It is to be noted that perplexities similar to the one expressed above, also emerged in connection with the study of Cooper pair tunneling between weakly coupled superconductors. In fact, objections were raised⁹ in connection with the prediction of Josephson¹⁰ that there should be a contribution to the current through an insulating barrier between two superconductors which would behave like direct tunneling of condensed pairs. This is in keeping with the fact that having the single electron a rather small probability of getting through the insulating junction, the “classical” estimate of simultaneous pair tunneling will not be observable. Let alone the fact that the (local) pairing interaction vanishes inside the barrier¹¹.

4.3 Phase correlation and enhancement factor

When one turns on, in an open shell atomic nucleus like e.g. $^{120}_{50}\text{Sn}_{70}$, a pairing interaction of strength larger than critical, the system moves into a Cooper pair regime¹². This fact has quantitative (but not qualitative) consequences concerning the one-particle transfer mechanism, and regards the size of the mismatch between the relative motion incoming ($p + ^{120}\text{Sn}(\text{gs})$) and outgoing ($d + ^{119}\text{Sn}(\text{gs})$) trajectories (Q -value and recoil effect). This is in keeping with the fact that one has to break a Cooper pair to populate a single quasiparticle state. From a structure point of view, the depletion of the occupation probability measured in a (p, d) process is correlated with the corresponding increase in occupation observed in (d, p) (U^2, V^2

⁸Concerning the marked fluctuations observed in these cross sections see Sect. 3.1.1, second paragraph.

⁹Bardeen (1962, 1961) (see also Pippard (2012); Cohen et al. (1962); McDonald (2001)).

¹⁰Josephson (1962), Anderson (1964b).

¹¹Anderson (1970); see also Broglia (2020) p. xxi-xxiii in connection with the nuclear case.

¹²Regime which is conditioned by the “external” mean field. In other words, regime (abnormal density) which express itself provided there is nucleon (normal) density available. It is of notice that pairing in turn may help extend the range over which normal nuclear density can reach, as in the case of the neutron halo nucleus ^{11}Li (lying at) defining the neutron drip line.

factors). Aside from the detailed quantitative values, this is quite similar to what is observed in dressed single-particle states in normal nuclei, the single-particle sum rule involving both the ($A - 1$) and ($A + 1$) aside from the A systems (see Sect. 5.8). Concerning the phase coherence of the pair correlated wavefunction, it has no consequence for one-particle transfer process, in keeping with the fact that $|e^{i\phi} \sqrt{P_1}|^2 = P_1$. A further reminder that not one, but two-particle transfer is the specific probe of pairing in nuclei.

In fact, qualitative differences are found concerning pair transfer when the Fermi system changes from the normal into the superconducting state. A result of the fact that in this phase transition, more or less the same set of (phase disordered) renormalized many-body states become superposed with a fixed phase relation described by $|BCS\rangle_{\mathcal{K}'} = \prod_{\nu>0} (U'_\nu + e^{-2i\phi} V'_\nu a_\nu^\dagger a_\nu^\dagger) |0\rangle$ state, which displays ODLRO¹³. As a consequence, in the calculation of the Cooper pair transfer probability, one has to add phased amplitudes before one takes modulus squared. That is

$$\begin{aligned} P_2 &= \left| \frac{1}{\sqrt{2}} (U'_\nu \sqrt{P_1} + e^{-2i\phi} V'_\nu \sqrt{P_1}) \right|^2 \\ &= \frac{P_1}{2} (1 + 2U'_\nu V'_\nu \cos 2\phi) \approx P_1 \end{aligned} \quad (4.3.1)$$

In other words, it is like interference in optics, with phase-coherent mixing. We note that above it was assumed $U'_\nu V'_\nu|_{\epsilon_F} \approx 1/2$ and $\cos 2\phi \approx 1$. Within this context see Sect. 5.5.1.

In connection with pair transfer experimental probes involving intrinsic (as opposite to relative motion) energy exchange $< 2\Delta$ and relative distances of closest approach $< \xi$, only Cooper pairs exist in the nuclear pairing correlated system. In this case, the partners nucleons are correlated over distances of the order of 15–20 fm (Fig. 4.3.1 (b)), the intrinsic motion being frozen because of the pairing gap. A situation quite different from that of a normal system characterized by independent nucleon motion, where the relative distance between nucleons is ≈ 2.4 fm is the characterizing feature (Fig. 4.3.1 (a)).

To the extent that the mean field acting as an “external” field allows particle density to be present, the properties of phase correlated, independent Cooper pair motion will explicit themselves. And thus it is a physical condition which is assumed fulfilled each time one makes use of the physical picture provided by Fig. 4.3.1 (b). In other words, inside ^{120}Sn all (6–8) Cooper pairs will be found within a volume of radius $R_0 \approx 6$ fm, in a similar way in which the transferred Cooper pair can be distributed over two nuclear volumes during the contact time in e.g. a Sn+Sn heavy ion reaction (Fig. 4.5.1). A phenomenon to which the long-range induced pairing interaction (exchange of phonons) contributes in the case of nuclei lying along the stability valley, let alone in the case of light halo nuclei like ^{11}Li

¹³See footnote 42 Ch. 1.

(Fig. 4.3.2)¹⁴.

The interest of the picture shown in Fig. 4.3.1 (b) can also be exemplified by referring to the fact that the moment of inertia of heavy deformed nuclei is considerably smaller than the rigid moment of inertia, but still larger than the irrotational one¹⁵ ($5\mathcal{J}_{irrot} \lesssim \mathcal{J} \lesssim \mathcal{J}_r/2$). Even confined within the mean field of the nucleus, the small but finite number of phase correlated pairs having the “intrinsic”, infinite-matter-like tendency will, to some extent, average out the different orientations of the rotating system and react to it in terms of an effective inertia smaller than the one related to the independent particle motion in a deformed potential (rigid moment of inertia). However, constrained by such a potential, they cannot fully profit of BCS condensation (ODLRO). To have irrotational flow, one has to have $\Delta \gg \hbar\omega_0$. Summing up, $\mathcal{J}_{irrot} = \delta^2 \mathcal{J}_{rig} \approx \delta^2 \times 2\mathcal{J}_{exp} \approx (2/10) \times \mathcal{J}_{exp} \approx \mathcal{J}_{exp}/5$, where¹⁶ $\delta \approx 0.95\beta_2$, typical values of β_2 being ≈ 0.3 .

Within this context we note that the (approximate) form of the (local) pair wavefunction can be written as¹⁷

$$F(r) \approx \Delta N(0) \frac{\sin k_F r}{k_F r} \exp\left(-\frac{\sqrt{2}r}{\xi}\right), \quad (4.3.2)$$

where $N(0)$ is the density of levels at the Fermi energy for one spin orientation. For $r \lesssim \xi$ the pair wavefunction is approximately proportional to that of two particles at the Fermi energy moving freely in a relative *s*-wave state. In a typical metallic superconductor ξ is of the order of $1 \mu\text{m}$ (10^4 \AA), much larger than the inter electron spacing ($\approx 2 \text{ \AA}$). Note that relative to the Fermi energy, the correlation energy ($E_{corr} = (-1/2)N(0)\Delta^2$) associated with Cooper pairing is very small, $\approx 10^{-7} - 10^{-8}$. Arguably, the most important consequence of this fact, is the exponentially large radius and thus very small value of the relative momentum associated with Cooper pairs. In other words, the typical scenario for a very small value of the quantum localization kinetic energy and thus of the generalized quantity parameter (see Sect. 3.5 and App. 7.C), implying that the two partners of the Cooper pair, are anchored (entangled) to each other. This phenomenon is at the basis of the emergence of new elementary modes of excitation. Pairing vibrations for single Cooper pairs, pairing rotations for few ones, supercurrents and Josephson currents for macroscopic amounts of them.

The situation of very extended Cooper pairs sounds, in principle, very different in the case of condensed matter (e.g. low-temperature superconductors) than in atomic nuclei, in keeping with the fact that nuclear Cooper pairs are, as a rule, subject to an overwhelming external (mean) field ($|E_{corr}| \approx 2\Delta \approx 2.4 \text{ MeV} \ll |U(r \approx R_0)| \approx |V_0/2| \approx 25 \text{ MeV}$). But even in this case one can posit that the transition from independent-particle to phase correlated, independent-pair motion,

¹⁴In this case $R(^{11}\text{Li})=4.58\pm0.13 \text{ fm}$, the “effective” Wigner-Seitz radius being $(r_0)_{eff} \approx 4.58 \text{ fm}/(11)^{1/3} \approx 2.1 \text{ fm}$.

¹⁵Bohr and Mottelson (1975); Belyaev (1959, 2013).

¹⁶See p. 47, Bohr and Mottelson (1975).

¹⁷cf. Leggett (2006) p. 185; for the non local nuclear version cf. e.g. Broglia and Winther (1983).

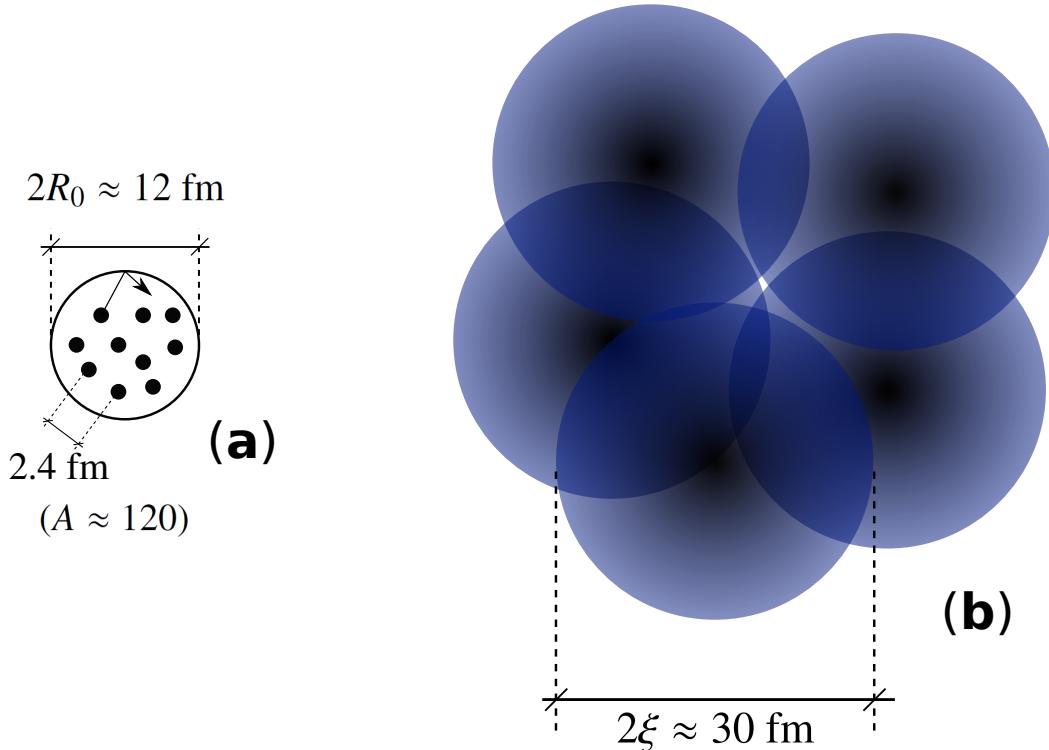


Figure 4.3.1: (a) Schematic representation of independent-particle motion and (b) independent-pair motion. In the first case nucleons (fermions) move independently of each other, their relative distance being $2r_0 (\approx 2.4 \text{ fm})$. Switching on the pairing interaction (bare plus induced) leads to Cooper pair formation in which the correlation length (mean square radius) is ξ . Pairs of nucleons moving in time reversal states close to the Fermi energy will tend to recede from each other lowering their relative momentum ($r_0 \rightarrow \xi$) thus boosting the stability of the system, provided that the external mean field allows for it. Or better, if there is (normal) nucleon density available to do so, something controlled to a large extent by the single-particle potential. From this point on, and at least for the levels lying close to the Fermi surface, one cannot talk about particles but about Cooper pairs (unless one does not intervene the system with an external field, e.g. (p, d) and provides the energy, angular and linear momentum needed to break a pair). The picture displayed in (b) is likely to be the one that is operative in the case of two nuclei of Sn at a relative (CM) distance somewhat larger than $2R_0 (\approx 12 \text{ fm})$, but still allowing for (weak) contact. The pair field associated with Cooper pairs will extend from one to the other partner of the heavy ions participating in the reaction through the weakly overlapping interaction region, allowing two nucleons to correlate over a distance ξ and, eventually, in a reaction like e.g. $\text{Sn} + \text{Sn} \rightarrow \text{Sn(gs)} + \text{Sn(gs)}$ allow for the successive transfer of two nucleons correlated over distances larger than nuclear dimensions.

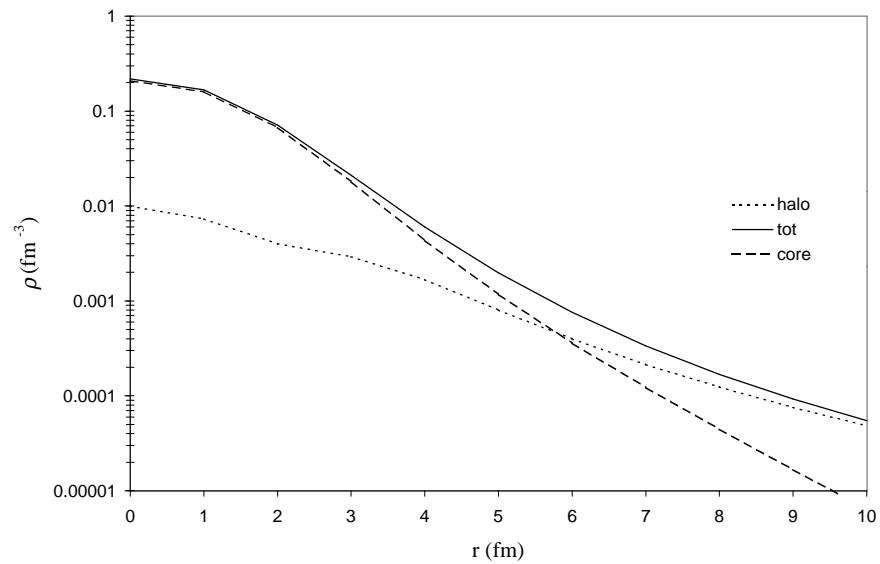


Figure 4.3.2: The nuclear density associated with ^{11}Li , as resulting from the microscopic NFT calculations which are at the basis of the results displayed in Figs. 3.6.3 and 7.1.2 (Barranco et al. (2001)). The contribution arising from the core (^9Li) is displayed with a dashed curve, while that associated with the two halo neutrons is shown in term of a dotted curve. The sum of these two contributions labeled tot (total) is drawn with a continuous curve.

implies that Cooper pair partners recede from each other. Let us clarify this point for the case of a single pair, e.g. $^{210}\text{Pb}(\text{gs})$. It is true that allowing the pair of neutrons to correlate in the valence orbitals leads to a pair wavefunction which is angle correlated ($\Omega_{12} \approx 0$), as compared to e.g. the pure $j^2(0)(j = g_{9/2})$ configuration¹⁸ (Sect. 4.11). On the other hand, the correlated pair addition mode (Tables 3.5.4 and 3.5.5) will display a sizeable spill out as compared to the pure two particle state, and thus a lower density and larger related average distance between Cooper pair partners. This is also the reason why close to $\approx 40\%$ of the pairing matrix elements is contributed by the induced pairing interaction resulting from the exchange of long wavelength, low-lying, collective modes, the other $\approx 60\%$ resulting from the bare nucleon-nucleon 1S_0 pairing interaction¹⁹.

The situation described above becomes likely clearer, even if extreme, in the case of ^{11}Li . In this case, the Fermi momentum is $k_F \approx 0.8 \text{ fm}^{-1}$, the radius $R \approx 4.58 \text{ fm}$ being much larger than $R_0 = 2.7 \text{ fm}$ expected from systematics. Furthermore essentially all of the correlation energy ($E_{corr} \approx -0.5 \text{ MeV}$)²⁰ is associated with the exchange of the pygmy dipole resonance between the halo neutrons²¹. It is of notice that in this case, as already stated, renormalization effects due to the clothing of single-particle states by vibrations, are as strong as mean field effects.

Again in this case $s_{1/2}^2(0)$ and $p_{1/2}^2(0)$ are not correlated in Ω_{12} , while the Cooper state probability density displays a clear angular correlation (see Fig. 3.6.3 (II) (a) and (b)). Nonetheless, the average distance between the partners of the neutron halo Cooper pair, is considerably larger than that associated with the ^9Li core nucleons, as testified by the following numerical estimates (see also Fig. 4.3.2):

$$\text{a)} \quad R(^{11}\text{Li}) = 4.58 \pm 0.13 \text{ fm} \quad (V = (4\pi/3)R^3 = 402.4 \text{ fm}^3) \quad (4.3.3)$$

$$\text{b)} \quad R_0(^{11}\text{Li}) = 2.7 \text{ fm} \quad (V = 82.4 \text{ fm}^3) \quad (4.3.4)$$

$$\text{c)} \quad R_0(^9\text{Li}) = 2.5 \text{ fm} \quad (V = 65.4 \text{ fm}^3), \quad (4.3.5)$$

and associated mean distance between nucleons,

$$\text{a)} \quad \left(\frac{337 \text{ fm}^3}{2} \right)^{1/3} \approx 5.5 \text{ fm}, \quad ((402.4 - 65.4) \text{ fm}^3) \quad (4.3.6)$$

$$\text{b)} \quad \left(\frac{82.4 \text{ fm}^3}{11} \right)^{1/3} \approx 1.96 \text{ fm}, \quad (4.3.7)$$

¹⁸Bertsch et al. (1967), Ferreira et al. (1984); Matsuo (2013) and references therein.

¹⁹In carrying out the above arguments the values of $(|E_{corr}|/\epsilon_F)^2 \approx \left(\frac{2.4 \text{ MeV}}{36 \text{ MeV}}\right)^2 \approx 10^{-3}$ and $\xi = \frac{\hbar v_F}{\pi \Delta} \approx 14 \text{ fm}$ ($(\frac{v_F}{c}) \approx (k_F)_{\text{fm}^{-1}}/5 \approx 0.27$; $\Delta \approx 1.2 \text{ MeV}$), typical for superfluid nuclei lying along the stability valley, were used.

²⁰ $(E_{corr}/\epsilon_F)^2 \approx (0.5/13)^2 \approx 10^{-3}$, $\xi \approx 20 \text{ fm}$ ($v_F/c \approx 0.2(k_F)_{\text{fm}^{-1}} \approx 0.16$; Eq. (7.C.6) and footnote 31 of Chapter 7).

²¹See Barranco et al. (2001); see also Broglia et al. (2019a).

$$\text{c)} \quad \left(\frac{65.4 \text{ fm}^3}{9} \right)^{1/3} \approx 1.94 \text{ fm}. \quad (4.3.8)$$

The above quantities are to be compared with the definition²²,

$$d = \left(\frac{\frac{4\pi}{3} R^3}{A} \right)^{1/3} = \left(\frac{4\pi}{3} \right)^{1/3} \times r_0 \approx 1.93 \text{ fm}, \quad (4.3.9)$$

consistent with the standard parametrization $R_0 = r_0 A^{1/3}$ of the nuclear radius. It is written in terms of the radius r_0 (=1.2 fm) of the Wigner-Seitz cell associated with each nucleon, and derived from systematics of stable nuclei lying along the stability valley.

4.3.1 Interplay between mean field and correlation length

In Fig. 4.3.1 one displays a schematic representation of the basis for two possible *gedankenexperiment* situations: (a) (*independent particle motion*), system which can be probed in a (p, t) reaction leading insight into non-interacting pairs of nucleons moving in time reversal states and confined by a mean field potential; (b) (*phase correlated independent pair motion*), target (or projectile) of two-nucleon transfer process induced by a heavy ion reaction between superfluid nuclei involving for example Sn-isotopes (see Sect. 7.3). In this case it is assumed that pairs of nucleons moving in time reversal states interact through an effective pairing interaction v_p^{eff} , sum of a bare $NN^{-1}S_0$ potential (v_p^{bare}) and an induced (v_p^{ind}) pairing component. As a result they become weakly bound with an energy 2Δ ($\ll \epsilon_F$), condensing in the BCS state and giving rise to the nuclear abnormal density in which the average distance between the pair partners is considerably larger than the relative distance associated with normal density ($\approx 2r_0$). Expressed differently and in keeping with the parallel with the BCS description of low temperature superconductivity found at the basis of nuclear BCS²³, one can assume that pairs of correlated nucleons moving in time reversal states close to the Fermi energy, will tend to recede from each other. For example, in the case of Sn, Cooper pair partners are expected to be at the antipodes²⁴, in keeping with the fact that $\xi \approx 14$ fm, and $R_0 \approx 6$ fm.

This expectation is not confirmed by studies of the Cooper pair wavefunction²⁵ or better, its modulus squared, which indicated that in going from situation (a) to situation (b), Cooper pair partners come close to each other, if nothing else because of angular correlation (see Sect. 4.11). A result which is also valid for systems with two nucleons outside closed shells.

²²Brink and Broglia (2005) App. C.

²³Bohr et al. (1958).

²⁴While not profiting completely for the latitude given by Cooper's mean square radius ξ , they do the best in the strong "external" mean field they are subject to, eventually displaying a non-negligible amount of spill out.

²⁵See footnote 18 of this Chapter.

In hindsight, one can posit that this result is interesting in itself in order to compare (nuclear structure) theory with theory, but less theory with experiment. At least not with the experiments associated with the specific probe²⁶ of Cooper pair correlations, namely two-nucleon transfer reactions, in which case the closest quantity to be observable is the two-nucleon transfer formfactor (Sect. 7.4). The soundness of picture (b) (Fig. 4.3.1) gets also support from the fact that one- and two-particle transfer absolute cross sections have similar values²⁷.

Within this context we note that the fact that ${}^9_3\text{Li}_6$ is well bound ($N = 6$ isotone parity-inverted closed shell), ${}^{10}_3\text{Li}_7$ is not while ${}^{11}_3\text{Li}_8$ is again bound, indicates that we are confronted with a pairing phenomenon. Allowing the two neutrons moving outside $N = 6$ closed shell to correlate in the configurations $j^2(0)(s_{1/2}^2, p_{1/2}^2, d_{5/2}^2 \dots)$ through a short range bare pairing interaction, e.g. the v_{14} Argonne NN -potential²⁸, does not lead to a bound state. The system lowers the relative momentum of the pair by exchanging the low-lying dipole vibration (soft $E1$ mode, pygmy resonance)²⁹ of the associated diffuse system becoming, eventually, bound, ever so weakly ($S_{2n} = 380$ MeV). The radius of the resulting system ($R({}^{11}\text{Li})=4.58 \pm 0.13$ fm) corresponds, in the parametrization $R_0 = 1.2A^{1/3}$ fm, to an effective mass number $A \approx 60$. So undoubtedly the system has swelled in moving from $A = 9$ to $A = 11$ in a manner that goes beyond the $1.2A^{1/3}$ (fm) expected dependence. Although the correlation length of the neutron Cooper is restricted to a spherical volume of diameter $2 \times R({}^{11}\text{Li}) \approx 9.2$ fm, half of the estimated value $\xi \approx 20$ fm (see App. 7.C), it is almost double as large as $2 \times R_0({}^{11}\text{Li}) \approx 5.4$ fm. Consequently, the function ($|\Psi_0(\mathbf{r}_1, \mathbf{r}_2)|^2$) displayed in Fig. 3.6.3 (II) b) should be read with care.

It will be surprising if the above mentioned bootstrap-like mechanism (see Fig. 2.7.1) namely that of profiting from low, unstable, nuclear densities to generate transient medium polarization effects to stabilize a Cooper pair halo system, was a unique property of ${}^{11}\text{Li}$. In fact, one can expect situations of s and p states at threshold eventually leading to a symbiotic halo Cooper pair with a small value of S_{2n} , also in connection with nuclear excited states.

Within the bootstrap ansatz of symbiotic Cooper pair binding, we introduce a straightforward generalization of the Axel-Brink³⁰ hypothesis based on well es-

²⁶On the other hand, if one is interested in the structure of the Cooper pair subject to the single-particle potential acting as a strong external field, and leading to conspicuous confinement effects, then the experiment to be carried out is e.g. (p, pn) knockout reactions at high ($\gg \epsilon_F$) energy/nucleon. This is also true concerning ${}^{11}\text{Li}$, not because of confinement, but because of spatial quantization, in keeping with the fact that the component $|s_{1/2}^2(0)\rangle$ and $|p_{1/2}^2(0)\rangle$ enter the wavefunction describing $|gs({}^{11}\text{Li})\rangle$ with almost equal amplitudes, and of the symmetry relations of the spherical harmonics entering the single-particle wavefunction, as discussed in Sect. 4.11, in particular in the text following Eq. (4.11.11) (within this scenario see ?? Kubota 2020).

²⁷ See footnote 26 Ch. 3.

²⁸Wiringa et al. (1984).

²⁹Broglia et al. (2019a).

³⁰Namely, the hypothesis that on top of each excited level of the nuclear spectrum is built a GDR equal to that built on the ground state; Axel (1962); Brink (1955 (unpublished)).

tablished experimental results. Namely the fact that the line shape and thus also the percentage of EWSR per energy interval as well as the decay properties of the GDR will reflect the static (splitting) and dynamic (motional narrowing) deformation properties of the state on which the GDR is built upon³¹.

In the case of halo nuclei, the above mentioned generalization is not only qualitative, but also quantitative. A sensible fraction of the TRK sum rule is found almost degenerate with the ground state. From the elastic antenna-like response typical of the high energy GDR ($\hbar\omega_{GDR} \approx 80\text{MeV}/A^{1/3}$) one is now confronted with a very low energy ($\lesssim 1\text{ MeV}$, $\Gamma \approx 0.5\text{ MeV}$) plastic dipole response pygmy dipole resonance (PDR), or soft $E1$ -mode³² where pairing, i.e. Cooper pairs play an important role. Regarding the consequences this phenomenon has for the $L = 1$ induced interaction between nucleons, one moves from dipole-dipole (static moment interactions, effective charges) to dispersive (retarded) contributions, emerging essentially from quantum mechanical ZPF³³. In other words, and making use of an analogy with atomic physics, one moves from an interaction between polar molecules, to a “purely” quantal interaction arising from the mutual polarization of one molecule in the rapidly changing field of the other (due to the instantaneous configuration of electrons and nuclei associated with ZPF) and viceversa, only one operative in the case of non-polar molecules. It is this second one which dominates the van der Waals interaction (Sect. 3.7) and, similarly, it is one which can lead to an almost resonant (PDR) gluing of Cooper pair halos (see Eq. 4.8.19). Resonant mechanism, phonon in this case, found at the basis of superconductivity in metals (see Eq. (4.8.62) and following discussion).

The energy centroid, the width and the percentage of the TRK sum rule (EWSR) of dipole resonances can be strongly affected by nuclear dynamical fluctuations and deformations which, in turn, depend on pairing, and thus on Cooper pairs. Because of angular momentum conservation, such phenomenon is restricted in lowest order, to fluctuations and deformations of quadrupole and monopole type. In the case of ^{11}Li , to an isotropic radial deformation, build up by the two less bound neutrons. This highly extended, low-momentum neutron halo, vibrating out of phase against the nucleons of the core (^9Li), gives rise to the soft $E1$ -mode (PDR)³⁴. It can be viewed as the taylored glue which, exchanged between the halo neutrons, bind the resulting neutron halo Cooper pair to the core³⁵.

The challenges faced in trying to learn about the physical basis of pairing in nuclei are comparable to those encountered to extract a collective vibration from a background much larger than the signal, as it was the case in the discovery of the

³¹Le Tourneau (no. 11, 1965); Bohr and Mottelson (1975); Bortignon et al. (1998) and refs. therein.

³²See footnote 121 of Ch. 2.

³³Within this context, see Bohr and Mottelson (1975), Sect. 6-5 f p. 432.

³⁴Broglia et al. (2019a) and refs. therein.

³⁵Barranco et al. (2001).

GDR in hot nuclei^{36,37}. In trying to observe pairing effects in nuclei close to the ground state, one has the advantage to start with the system at zero temperature for free. On the other hand one needs to subtract the very large, state dependent effects of the “external” mean field, a challenge not second to that faced by condensed matter practitioners to study low-temperature superconductivity in general, and the Josephson effect in particular. Within this context, ¹¹Li provides a textbook example of the fact that, given the possibility³⁸, nucleon pair partners recede from each other, lowering in the process the momentum of relative motion and thus the confinement kinetic energy³⁹, allowing to extend the limits of stability of nuclear species through a subtle long range pairing mechanism.

4.4 Correlations between nucleons in Cooper pair tunneling

Let us call x_1 and x_2 the coordinates of the Cooper pair partners. Let us furthermore assume they can only take two values: 0 when they are bound to the target nucleus, 1 when they have tunneled and become part of the outgoing particle (see Fig. 4.4.1).

The correlation between the two nucleons is measured by the value⁴⁰

$$\langle x_1 x_2 \rangle - \langle x_1 \rangle \langle x_2 \rangle = \int d\gamma P_2 \times 1 \times 1 - \int d\gamma P_1 \times 1 \int d\gamma' P'_1 \times 1 = P_2 - P_1 P'_1, \quad (4.4.1)$$

$d\gamma$ being the differential volume in phase space, normalized with respect to the corresponding standard deviations, that is, with respect to

$$\sigma_{x_1} \sigma_{x_2} = \left[(\langle x_1^2 \rangle - \langle x_1 \rangle^2) (\langle x_2^2 \rangle - \langle x_2 \rangle^2) \right]^{1/2}. \quad (4.4.2)$$

³⁶Snover (1986); Gaardhøje (1985); Gaardhøje et al. (1986). See also Bortignon et al. (1998) Figs. 1.4 and 6.8, and refs. therein.

³⁷A statement which becomes more than an analogy in connection with the nuclear analogue to the (ac) Josephson effect proposed in ?? (Potel Josephson γ), which can eventually be tested by extracting a dipole γ -signal of Josephson frequency $\nu_J = Q_{2n}/h$ (or better, that of the somehow higher peaked γ -strength function, energy shifted by the γ -phase space volume element), from a continuum plus discrete lines background.

³⁸Namely, the presence of normal density. A feature which in the present case goes hand in hand with the presence of abnormal density (single Cooper pair).

³⁹That is, from $\hbar^2/(m(2r_0)^2) \approx 7$ MeV ($r_0 = 1.2$ fm) to $\hbar^2/(m(2R(^{11}\text{Li}))^2) \approx 0.5$ MeV ($R(^{11}\text{Li}) \approx 4.6$ fm; within this context it is of notice that the overlap between radial wavefunctions of the halo neutrons and of the core neutrons is $O = (R_0/R)^3 \approx 0.2$ see Fig. 2.7.1; see also Broglia et al. (2019b)). While the system loses in this way a consistent fraction of the bare, short range pairing interaction through screening, it opens up for long-range pairing contributions which, ever so weak, can still bind the halo Cooper pair to the core. In other words, symbiotic binding by the confinement. In the case of superconducting metals, the correlation length is $\xi \approx 10^4$ Å, a distance to be compared to the Wigner-Seitz radius of ≈ 2 Å. This variation leads to a decrease of the Coulomb repulsion by almost 4 orders of magnitude. From $U_c = e^2/r = 14.4 \times \text{eV} \times \text{\AA}/(2\text{\AA}) = 7.2$ eV to $U_c = 14.4 \times \text{eV} \times \text{\AA}/(10^4\text{\AA}) = 1.4$ meV.

⁴⁰Basdevant and Dalibard (2005).

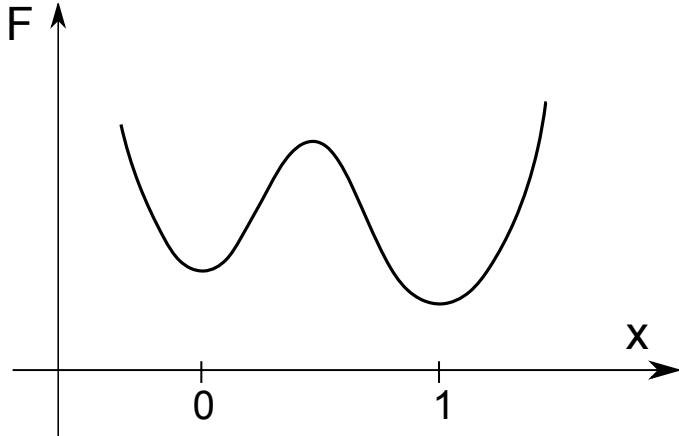


Figure 4.4.1: A schematic representation of nucleon tunneling between target and projectile. The free energy $F = U - TS$ which for the zero temperature situation under consideration (e.g. $^{120}\text{Sn}(p,t)^{118}\text{Sn}$) coincides with the potential energy as a function of the nucleon coordinate x . For $x = 0$ the nucleon is assumed to be bound to the target system. For $x = 1$ the nucleon has undergone tunneling becoming bound to the outgoing particle. In other words x_1 jumps from the value 0 to the value 1 in the tunneling process ($x, 0 \rightarrow 1$), the same for the coordinate of the second nucleon.

Making use of the fact that

$$\langle x_1^2 \rangle = \int d\gamma P_1 \times 1^2 = P_1, \quad (4.4.3)$$

and

$$\langle x_1 \rangle = \int d\gamma P_1 \times 1 = P_1, \quad (4.4.4)$$

one can calculate the function which measures the correlations between nucleons 1 and 2, namely,

$$Corr = \frac{\langle x_1 x_2 \rangle - \langle x_1 \rangle \langle x_2 \rangle}{\sqrt{(\langle x_1^2 \rangle - \langle x_1 \rangle^2)(\langle x_2^2 \rangle - \langle x_2 \rangle^2)}} = \frac{P_2 - P_1 P'_1}{\sqrt{(P_1 - P_1^2)(P'_1 - P'^2_1)}}. \quad (4.4.5)$$

Because both nucleons are identical and thus interchangeable, $P_1 = P'_1$. Thus

$$Corr = \frac{P_2 - P_1^2}{P_1 - P_1^2}. \quad (4.4.6)$$

Making use of the empirical values

$$P_1 \approx P_2 \approx 10^{-3} \quad (4.4.7)$$

l	p_l
0	1.02×10^{-3}
1	2.40×10^{-3}
2	1.26×10^{-2}
3	1.84×10^{-2}
4	6.13×10^{-3}
5	1.39×10^{-3}
6	2.89×10^{-4}
7	5.04×10^{-5}
8	6.51×10^{-6}
9	5.87×10^{-7}

Table 4.4.1: Probabilities p_l associated with the reaction ${}^1\text{H}({}^{11}\text{Li}, {}^{10}\text{Li(gs)}){}^2\text{H}$ calculated with the same bombarding conditions as those associated with ${}^1\text{H}({}^{11}\text{Li}, {}^9\text{Li(gs)}){}^3\text{H}$ (see also Table 7.1.2).

leads to,

$$\text{Corr} = \frac{10^{-3} - 10^{-6}}{10^{-3} - 10^{-6}} \approx 1. \quad (4.4.8)$$

In other words, within the Cooper pair motion regime, nucleon partners are solidly anchored (entangled) to each other: if one nucleon goes over, the other does it also. This is so in spite of the very liable and fragile (intrinsic) structure of the nuclear Cooper pairs ($2\Delta/\epsilon_F \ll 1$). An example of such a scenario is provided by ${}^{11}\text{Li}$. In fact, if one picks-up a neutron from ${}^{11}\text{Li}({}^{11}\text{Li}(p, d){}^{10}\text{Li})$, the other one breaks up essentially instantaneously, ${}^{10}\text{Li}$ being unbound (Sect. 5.2.4). In spite of this fact, the probability associated with the reaction ${}^1\text{H}({}^{11}\text{Li}, {}^9\text{Li(gs)}){}^3\text{H}$ is given by⁴¹

$$P_2 = \frac{(5.7 \pm 0.9) \text{ mb}}{2\pi((4.58 \pm 0.13) \text{ fm})^2} \approx (4.3 \pm 1)10^{-3}. \quad (4.4.9)$$

This value is of the order of that associated with the breakup process mentioned above⁴², and reported in Table 4.4.1.

The fact that $P_2 \approx P_1$ contained in the relation (4.3.1), is applicable both for static and dynamic pairing modes, is in keeping with the fact that in nuclei dynamic spontaneous breaking of gauge invariance, is of similar importance as the static one⁴³ (see Figs. 3.5.7 and 4.4.2).

⁴¹See Fig. 7.1.2, Sect. 7.1.2 and Table 7.1.2 concerning the experimental value of $\sigma({}^{11}\text{Li} \rightarrow {}^9\text{Li (gs)})$, and Kobayashi et al. (1989) concerning the experimental value of $R({}^{11}\text{Li})$. See also paragraph following Eq. (3.6.13).

⁴²It is of notice that the value given in Eq. (4.4.9) essentially coincides with that shown in the $l = 0$, column **1** of Table 7.1.2.

⁴³cf. Fig. 3.5.7, cf. also Sect. 4.12 and Fig. 4 of Potel et al. (2013b).

$$\text{Order parameter} \quad \left(\langle \tilde{0} | P P^\dagger | \tilde{0} \rangle \right)^{1/2} = \begin{cases} \alpha_0 = \sum_{\nu>0} U'_\nu V'_\nu \\ \alpha_{dyn} = \sum_{\nu>0} U_\nu^{eff} V_\nu^{eff} \end{cases}$$

pairing vibrations

$$(V_\nu^{eff})^2 = 2Y_a^2(j_\nu)/\Omega_\nu; \quad (U_\nu^{eff})^2 = 1 - (V_\nu^{eff})^2$$

$$\begin{cases} X_a(j_\nu) \\ Y_a(j_\nu) \end{cases} = \frac{(\sqrt{\Omega_j}/2)\Lambda_a}{2|E_j| \mp W_a}$$

pairing rotations

$$\begin{cases} U'_\nu \\ V'_\nu \end{cases} = \frac{1}{\sqrt{2}} \left(1 \pm \frac{\epsilon_\nu}{\sqrt{\epsilon_\nu^2 + \Delta^2}} \right)^{1/2}$$

Figure 4.4.2: Order parameter associated with static and dynamic pair correlations.

4.5 Pair transfer

In the semi-classical approximation, the second order two-nucleon transfer amplitude can be written as (see Eq. (6.5.3)),

$$a(t = +\infty) = a^{(1)}(\infty) - a^{(NO)}(\infty) + \tilde{a}^{(2)}(\infty), \quad (4.5.1)$$

where $a^{(1)}(\infty) = a_{sim}^{(1)}$ and $\tilde{a}^{(2)}(\infty) = a_{succ}^{(2)}$ describe the simultaneous and successive transfer amplitudes respectively. In the **independent particle limit**, these amplitudes fulfill the relations⁴⁴,

$$a_{sim}^{(1)} = a_{NO}^{(1)}, \quad (4.5.2)$$

and

$$a_{succ}^{(2)} = a_{one-part}^{(1)} \times a_{one-part}^{(1)}, \quad (4.5.3)$$

with

$$a + A \rightarrow f + F \rightarrow b + B, \quad (4.5.4)$$

corresponding to the product of two single nucleon transfer amplitudes. On the other hand, in the **strong correlation limit** one can write, making use of the post-prior representation

$$a_{succ}^{(2)} = \tilde{a}_{succ}^{(2)} - a_{NO}^{(1)}, \quad (4.5.5)$$

⁴⁴see Sect. 6.5, also Potel et al. (2013a).

relation which fulfills,

$$\lim_{E_{corr} \rightarrow \infty} a_{succ}^{(2)} = 0. \quad (4.5.6)$$

That is, the transfer process is, in this case, associated with simultaneous transfer. Actual nuclei are close to the independent particle limit ($E_{corr} \approx (2\text{--}3 \text{ MeV}) \ll \epsilon_F (\approx 36 \text{ MeV})$). Successive transfer is thus expected to be the major contribution to pair transfer processes. But successive transfer seems to break the pair *right? Wrong. Why?* let us see below.

4.5.1 Cooper pair dimensions

At the basis of the relations used to estimate the dimensions of a Cooper pair (correlation length; mean square radius) one finds

$$\delta x \delta p \geq \hbar \quad \delta \epsilon \approx |E_{corr}|, \quad (4.5.7)$$

where

$$\epsilon = \frac{p^2}{2m}; \quad \delta \epsilon = \frac{p \delta p}{m} \approx v_F \delta p, \quad (4.5.8)$$

and thus

$$\delta \epsilon \approx |E_{corr}| \approx v_F \delta p, \quad (4.5.9)$$

leading to

$$\xi \approx \delta x \approx \frac{\hbar}{\delta p} \approx \frac{\hbar v_F}{|E_{corr}|} \quad (\text{correlation length}). \quad (4.5.10)$$

In what follows we use, for normal systems,

$$\xi = \frac{\hbar v_F}{\pi |E_{corr}|}, \quad (4.5.11)$$

and, in the case of open shell superfluid nuclei⁴⁵

$$\xi = \frac{\hbar v_F}{\pi \Delta}. \quad (4.5.12)$$

For Sn-isotopes, a typical value of the pairing gap is $\Delta \approx 1.2 \text{ MeV}$. Making use of $k_F = 1.36 \text{ fm}^{-1}$ associated with nuclei lying along the stability valley, and thus ($v_F/c \approx 0.2(k_F)_{\text{fm}^{-1}} \approx 0.27$),

$$\xi \approx 14 \text{ fm}. \quad (4.5.13)$$

Consequently, successive and simultaneous transfer can feel equally well the pairing correlations giving rise to long range order⁴⁶. This virtual property can become real in e.g. a pair transfer between two superfluid nuclei⁴⁷, in particular two tin isotopes (see Figs. 4.1.3 and 4.5.1).

⁴⁵So called Pippard correlation length. See Schrieffer (1964) p. 18 and 34, Annott (2013) p. 62.

⁴⁶See footnote 42 Ch. 1; see also e.g. Potel et al. (2017) and references therein.

⁴⁷See von Oertzen (2013); von Oertzen and Vitturi (2001); Weiss (1979); for a more recent experimental and theoretical study see Montanari et al. (2014, 2016).

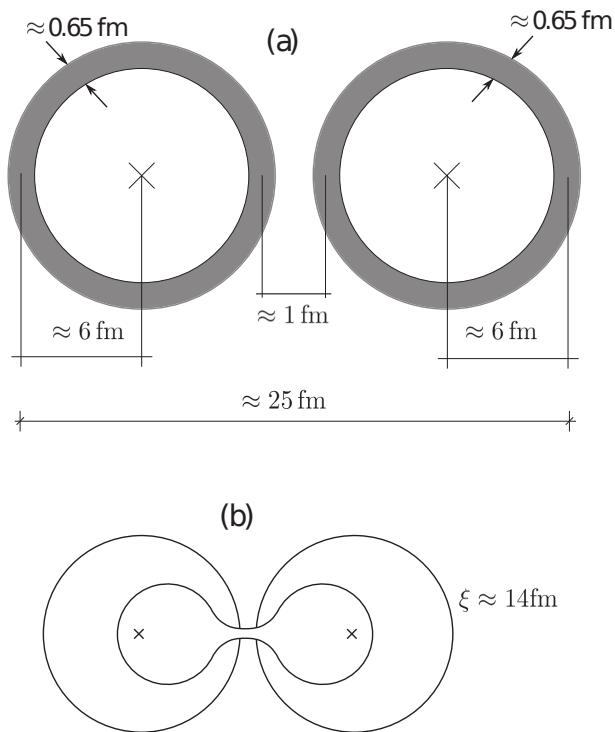


Figure 4.5.1: (a) Schematic representation of two Sn-isotopes (radius $R_0 \approx 6 \text{ fm}$) at the distance of closest approach in a heavy ion collision; (b) single Cooper pair in which each nucleon partner is in a different nucleus. The gray areas indicate the nuclear diffusivity.

Objection

What about $v_{pairing}(= G)$ becoming zero, e.g. between the two nuclei?

Answer

$$\frac{d\sigma(a(= b + 2) + A \rightarrow b + B(= A + 2))}{d\Omega} \sim |\alpha_0|^2, \quad (4.5.14)$$

$$\alpha_0 = \langle BCS(A + 2)|P^\dagger|BCS(A)\rangle = \sum_{\nu>0} U_\nu(A)V_\nu(A + 2). \quad (4.5.15)$$

That is, pair transfer does not depend explicitly but only functionally on G .

Objection

Relation (4.5.15) is only valid for simultaneous transfer, *right? Wrong.*

Answer

The order parameter can also be written as,

$$\begin{aligned} \alpha_0 &= \sum_{\nu, \nu' > 0} \langle BCS | a_\nu^\dagger | int(\nu') \rangle \langle int(\nu') | a_{\bar{\nu}}^\dagger | BCS \rangle \\ &\approx \sum_{\nu, \nu' > 0} \langle BCS(A + 2) | a_\nu^\dagger a_{\nu'}^\dagger | BCS(A + 1) \rangle \langle BCS(A + 1) | \alpha_{\nu'} a_{\bar{\nu}}^\dagger | BCS(A) \rangle \\ &= \sum_{\nu, \nu' > 0} \langle BCS(A + 2) | V_\nu(A + 2) \alpha_{\bar{\nu}} a_{\nu'}^\dagger | BCS(A + 1) \rangle \\ &\times \langle BCS(A + 1) | \alpha_{\nu'} U_\nu(A) a_{\bar{\nu}}^\dagger | BCS(A) \rangle = \sum_{\nu > 0} V_\nu(A + 2) U_\nu(A), \quad (4.5.16) \end{aligned}$$

where the (inverse) quasiparticle transformation relation $a_\nu^\dagger = U_\nu a_\nu^\dagger + V_\nu a_{\bar{\nu}}$ was used⁴⁸. Examples of two-nucleon spectroscopic amplitudes involving superfluid targets, namely those associated with the reactions $^{112}\text{Sn}(p, t)^{110}\text{Sn(gs)}$ and $^{124}\text{Sn}(p, t)^{122}\text{Sn(gs)}$ are given in Table 3.1.1 (see also Table 7.2.1). Making use of some of these amplitudes (first column of Table 3.1.1) and of global optical parameters (Table 7.2.2), the two-nucleon transfer absolute differential cross section of the reaction $^{112}\text{Sn}(p, t)^{110}\text{Sn(gs)}$ at a bombarding energy of $E_p = 26$ MeV, was calculated making use of the software COOPER based on second order DWBA and taking into account successive and simultaneous transfer properly corrected for non-orthogonality (cf. Chapter 6 and App. 7.A). It is compared with experimental data in Fig. 4.5.2 (a). The corresponding absolute integrated cross sections are $1310 \mu\text{b}$ and $1309 \pm 200 \mu\text{b}$ respectively. The largest contribution to the cross section arise from successive transfer, the cancellation between simultaneous and non-orthogonality amplitudes being important (Fig. 4.5.2 (b)). The above provides

⁴⁸See e.g. Brink and Broglia (2005) App. G.

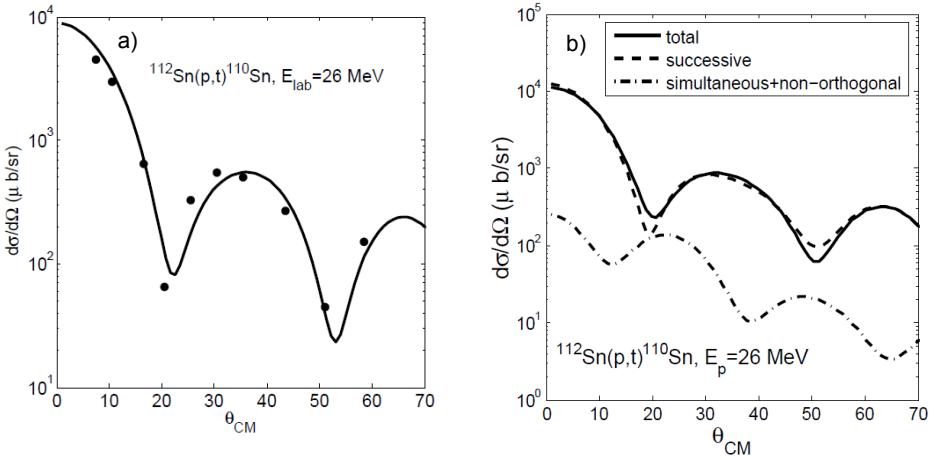


Figure 4.5.2: a) Absolute differential cross section associated with the reaction $^{112}\text{Sn}(p,t)^{110}\text{Sn}(\text{gs})$ calculated with the software COOPER in comparison with the experimental data (Guazzoni et al. (2006)). b) Details of the different contributions to the total absolute (p,t) differential cross section.

a typical example of results of a systematic study of two-nucleon transfer reactions in terms of absolute cross sections⁴⁹ (Figs. 3.1.7 and 7.2.1).

Making use of two-nucleon spectroscopic amplitudes worked out within the framework of an extended shell model calculation (Table 3.1.1, second column) one obtains very similar results to those displayed in Fig. 4.5.2 (a). In Fig. 4.5.3 we report results similar to those displayed in Fig. 4.5.2, but for the case of the reaction $^{124}\text{Sn}(p,t)^{122}\text{Sn}(\text{gs})$ calculated within second order DWBA making use of the BCS spectroscopic amplitudes (Table 3.1.1 third column).

Let us now discuss an example of two-nucleon transfer around a closed shell nucleus displaying well defined collective pairing vibrational modes. We refer, in particular, to the pair removal mode of ^{206}Pb , that is, to the reaction, $^{206}\text{Pb}(t,p)^{208}\text{Pb}(\text{gs})$. Making use of the spectroscopic amplitudes displayed in Tables 3.5.2 and 3.5.3 and of global optical parameters, the associated absolute differential cross section was calculated with the software COOPER. It is displayed in Fig. 4.5.4 in comparison with the experimental findings. In the same figure, the absolute total differential cross section is compared with that associated with the TD (Tamm-Dancoff) description of $^{206}\text{Pb}(\text{gs})$, that is, setting the pairing ground state correlations to zero ($\sum_i X_r^2(i) = 1$, $Y_r(k) \equiv 0$, see Table 3.5.2). In this case, theory underpredicts observation by about a factor of 2. Also given in Fig. 4.5.4 is the predicted cross section associated with the pure configuration $|p_{1/2}^{-2}(0)\rangle$.

It is of notice, that within the effective reaction mechanism described in Sect. 6.4 pairing correlations increase the value of $\Omega_0 (\approx 0.97)$. As a consequence the $l = n = 0$ two-neutron system gives a consistently larger contribution to the two-

⁴⁹Potel et al. (2013a), Potel et al. (2013b).

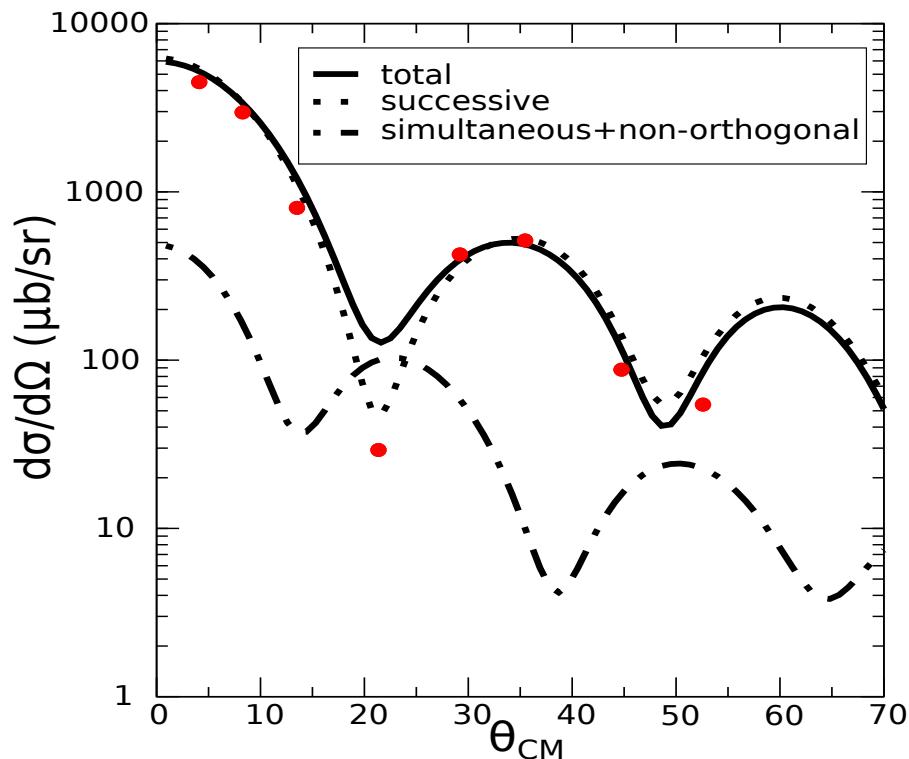


Figure 4.5.3: Absolute differential cross section associated with the reaction $^{124}\text{Sn}(p, t)^{122}\text{Sn}(\text{gs})$ calculated making use of second order DWBA taking into account non-orthogonality corrections and the two-nucleon spectroscopic amplitudes resulting from BCS (see Table 3.1.1, third column; for details see Potel et al. (2013a), Potel et al. (2013b)) in comparison with experimental data (Guazzoni et al. (2011)).

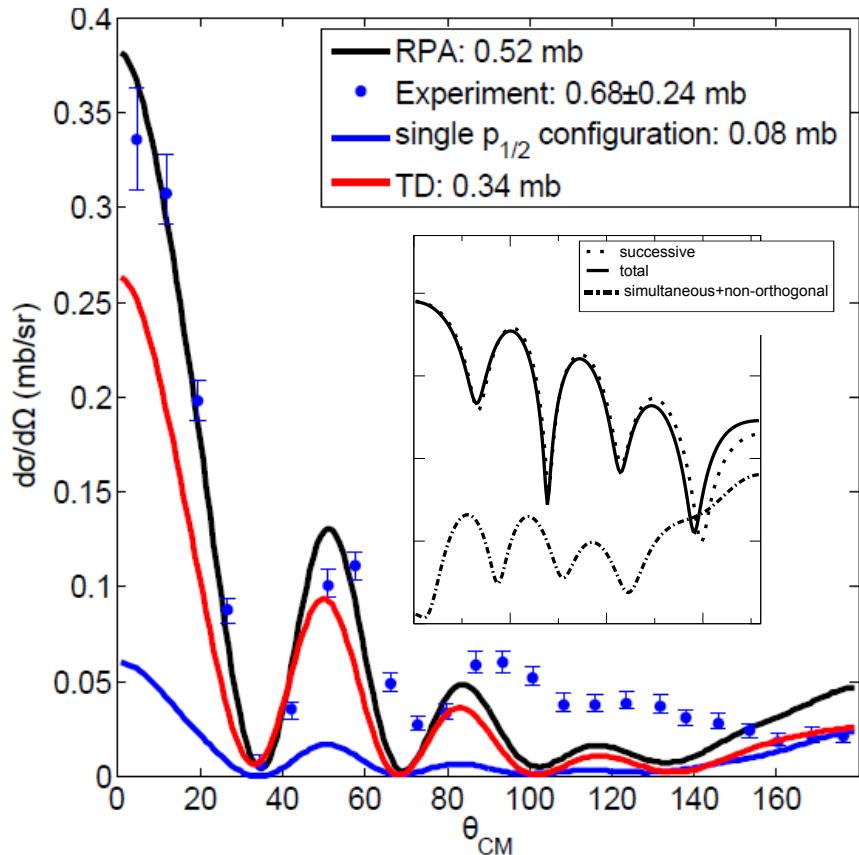


Figure 4.5.4: Absolute two-nucleon transfer differential cross section associated with the $^{206}\text{Pb}(t, p)^{208}\text{Pb}(\text{gs})$ transfer reaction, that is, the annihilation of the pair removal mode of ^{208}Pb in comparison with the data (Bjerregaard et al. (1966)). The theoretical cross sections were calculated making use of the spectroscopic amplitudes given in Tables 3.5.2 and 3.5.3 and of global optical parameters as reported in the reference above. Both RPA and TD amplitudes were used as well as a pure configuration $p_{1/2}^2(0)$. **Gregory control**

nucleon transfer process than those associated with $n = 1$ and 2, that is those proportional to Ω_1 and Ω_2 whose values are 0.25 and 0.06 respectively (cf. Eq. (6.4.13)). All these features boost the effective absolute two-nucleon transfer cross section. While the results displayed in Fig. 4.5.4 were calculated making use of the full formalism of second order DWBA (Figs. 4.1.1 and 4.1.2) the simplified expressions given in Eqs. (6.4.1–6.4.3) and (6.4.13–6.4.16) were useful to gain physical insight into the effective two-nucleon transfer formfactors $u_{LSJf}^{J_i J_f}(R)$. In fact, these functions, multiplied by the factors D_0 , provided a simple parametrization to account for the absolute two-nucleon transfer differential cross sections. *On the other hand, they were a source of misunderstanding concerning the reaction mechanism of two-nucleon transfer; let alone of the spatial structure of Cooper pairs*⁵⁰.

Collective surface vibrations of closed shell nuclei can be viewed as correlated particle-hole excitations. The phase coherence existing between the different RPA amplitudes of the corresponding wavefunction leads to a decrease of the average distance (angular correlation) between the particle and the hole, as compared with pure ph -configurations, similar to what happens in pair addition (subtraction) modes (concerning this parallel see Fig. 4.9.3; see also Fig. 7 of Barranco et al. (2019b)), in which case are the pp (hh) which approach each other. It has been argued that this is the reason why both pairing and surface vibrations of closed shell nuclei display enhanced (t, p) cross sections as compared to pure configurations⁵¹. The two neutrons lie rather close to each other also in the triton.

In the case of collective surface vibrations of closed shell nuclei, like e.g. the octupole vibration of ^{208}Pb ($E_x = 2.65$ MeV), the specific probe is not two-nucleon transfer, but Coulomb excitation or inelastic scattering, the wavelength of the γ -ray associated with the corresponding electromagnetic decay being $\lambda \approx 460$ fm. This is about two orders of magnitude larger than nuclear dimensions. Within this scenario, whether the particle lies closer to the hole or less in the correlated state, as compared to the pure $p - h$ state can hardly be of any relevance to explain the enhancement of the absolute transition probability ($B(E3) = 32$ Bsp) of the particle-hole correlated state, one of the largest of the whole mass table. Not only this, it sets a question mark on the validity of the argument as applied to Cooper pair transfer and pairing correlations. This question is taken up in Sects. 4.9–4.12.

4.6 Weak link between superconductors

Two-nucleon transfer reactions involving superfluid nuclei display some similarities with Cooper pair tunneling between weakly coupled superconductors, in particular when discussing heavy ion reactions, but not only⁵². Within this context it

⁵⁰For more (historic) details see Broglia (2019) (Overview).

⁵¹Bertsch et al. (1967)

⁵²Goldanskii and Larkin (1968); Dietrich (1970); Dietrich et al. (1971); Hara (1971); Kleber and Schmidt (1971); Weiss (1979), von Oertzen and Vitturi (2001); von Oertzen (2013); Broglia and

is useful to remind the basic elements of the pair tunneling which is at the basis of the Josephson effect. In this section we essentially reproduce the description of the tunneling of Cooper pairs between two weakly coupled superconductors to be found in⁵³, likely the best physical presentation of the Josephson effect⁵⁴.

One starts with the many-body Hamiltonian of⁵⁵

$$H = H_1 + H_2 + \sum_{kq} T_{kq} (a_{k\uparrow}^\dagger a_{q\uparrow} + a_{-q\downarrow}^\dagger a_{-k\downarrow}) + hc \quad (4.6.1)$$

where H_1 and H_2 are the separate Hamiltonians of the two superconductors on each side of the junction, T_{kq} being the (exponentially) small tunneling matrix element from state k on one side to state q on the other.

One can arrive to (4.6.1) by first finding sets of single-particle wavefunctions for each side separately, in the absence of the potential of the other system. *Then one eliminates the non-orthogonality effects by perturbation theory (cf. the similarity with the arguments used in Sect. 4.1 to introduce the two-nucleon transfer amplitudes in (2nd order) DWBA; see also App. 2.B, Eq. (2.B.23)).* A nuclear embodiment of such strategy⁵⁶ is worked out in Ch. 6 (see in particular Sect. 6.2.9 and Sect. 6.5) and was implemented in the code COOPER⁵⁷.

At the basis of the Josephson effect one finds $P_2 \approx P_1$ which, in the nuclear case implies $\sigma_{2p} \approx \sigma_{1p}$ (see Sect. 4.3; also footnote 26 Ch. 3). A remarkable finding valid also whether only one of the two interacting systems belongs to a closed shell pairing vibrational band, or is superfluid. This is because of the importance of pairing vibrations in nuclei as compared to superconductors. Similar results are found in the case in which both systems are superfluid, namely in the case of one- and two-nucleon transfer reactions in a heavy ion collision between superfluid nuclei (see Ch. 7)⁵⁸. It is likely that the results obtained in this case at bombarding energies of few MeV below the Coulomb barrier, come closest to the tunneling phenomena observed between two weakly coupled superconductors.

Let us now calculate the second order expression of (4.6.1) in the case in which the gaps of the two weakly linked superconductors are different. Making use of

Winther (2004).

⁵³Anderson (1964b)

⁵⁴Josephson (1962).

⁵⁵Cohen et al. (1962)

⁵⁶Because pairing vibrations in nuclei are quite collective, leading to effective U and V occupation factor (cf. Fig. 4.4.2) (see also Potel et al. (2013b)), the nuclear and the condensed matter expressions for Cooper tunneling processes are very similar. Of course no supercurrent is expected between nuclei. However, the systems $^{120}\text{Sn(gs)}$, $^{119}\text{Sn(j)}$, $^{118}\text{Sn(gs)}$ form an ensemble of weakly coupled Fermi superfluids, with different (average) number of particles ($N, N-1, N-2$), to which essentially all the BCS techniques, including those of the present Section can be applied.

⁵⁷Cf. App. 7.A; cf. also Broglia and Winther (2004).

⁵⁸Note that in the case of a heavy ion collision between two superfluid nuclei at energies below the Coulomb barrier, one compares gs → gs with gs → qp, summed over a number of qp states (see Sect. 7.3).

relations presented in⁵⁹ Sects. 1.4.2 and 4.5 one can write, for $T = 0$,

$$\Delta E_2 = -2 \sum_{kq} |T_{kq}|^2 \frac{|V_k U_q + V_q U_k|^2}{E_k + E_q}. \quad (4.6.2)$$

With the help of

$$2U_k V_k^* = \frac{\Delta_k}{E_k}, \quad 2U_q V_q^* = \frac{\Delta_q}{E_q}, \quad (4.6.3)$$

and

$$|U_k|^2 - |V_k|^2 = \frac{\epsilon_k}{E_k}, \quad |U_q|^2 - |V_q|^2 = \frac{\epsilon_q}{E_q}, \quad (4.6.4)$$

where

$$E = \sqrt{\epsilon^2 + \Delta^2} \quad (4.6.5)$$

and⁶⁰

$$\Delta_k = \Delta_1 e^{i\phi_1}, \quad \Delta_q = \Delta_2 e^{i\phi_2}, \quad (4.6.6)$$

one can write for the numerator of Eq. (4.6.2),

$$\begin{aligned} NUM &= (V_k U_q + V_q U_k)(V_k^* U_q^* + V_q^* U_k^*) \\ &= \{V_k^2 U_q^2 + V_q^2 U_k^2\} + [(U_k^* V_k)(U_q V_q^*) + (U_q^* V_q)(U_k V_k^*)]. \end{aligned} \quad (4.6.7)$$

It is of notice that, for simplicity, throughout this section

$$V^2 \equiv |V|^2. \quad (4.6.8)$$

With the help of (4.6.3) the expression in the squared bracket in (4.6.7) can be written as

$$[] = \frac{1}{4E_k E_q} (\Delta_k^* \Delta_q + (\Delta_k^* \Delta_q)^*) = \frac{1}{4E_k E_q} 2\Re(\Delta_k^* \Delta_q). \quad (4.6.9)$$

Making use of the relations

$$\begin{aligned} (U_k^2 - V_k^2)(U_q^2 - V_q^2) &= U_k^2 U_q^2 - U_k^2 V_q^2 - V_k^2 U_q^2 + V_k^2 V_q^2 \\ &= -(U_k^2 V_q^2 + V_k^2 U_q^2) + (U_k^2 U_q^2 + V_k^2 V_q^2), \end{aligned} \quad (4.6.10)$$

and

$$\begin{aligned} 1 &= (U_k^2 + V_k^2)(U_q^2 + V_q^2) = U_k^2 U_q^2 + U_k^2 V_q^2 + V_k^2 U_q^2 + V_k^2 V_q^2 \\ &= (U_k^2 V_q^2 + V_k^2 U_q^2) + (U_k^2 U_q^2 + V_k^2 V_q^2), \end{aligned} \quad (4.6.11)$$

⁵⁹See also App. G Brink and Broglia (2005).

⁶⁰Note the difference in gauge phasing used in this section to follow that of Anderson (1964b) as compared with that used in the rest of the monograph.

one obtains,

$$1 - (U_k^2 - V_k^2)(U_q^2 - V_q^2) = 2(U_k^2 V_q^2 + V_k^2 U_q^2), \quad (4.6.12)$$

that is, twice the expression written in curly brackets in (4.6.7). Consequently

$$\{ \} = \frac{1}{2}(1 - (U_k^2 - V_k^2)(U_q^2 - V_q^2)) = \frac{1}{2}\left(1 - \frac{\epsilon_k \epsilon_q}{E_k E_q}\right). \quad (4.6.13)$$

Thus, the sum of (4.6.9) and (4.6.13) leads to,

$$NUM = \frac{1}{2}\left(1 - \frac{\epsilon_k \epsilon_q}{E_k E_q} + \frac{\Re(\Delta_q^* \Delta_k)}{E_k E_q}\right) \quad (4.6.14)$$

and

$$\Delta E_2 = -\sum_{kq} \frac{|T_{kq}|^2}{E_k + E_q} \left(1 - \frac{\epsilon_k \epsilon_q}{E_k E_q} + \frac{\Re(\Delta_q^* \Delta_k)}{E_k E_q}\right). \quad (4.6.15)$$

In what follows, we concentrate in the gauge phase-dependent term.

With the help of (4.6.6) one can write

$$\Delta_k \Delta_q^* = \Delta_1 \Delta_2 e^{i(\phi_1 - \phi_2)} = \Delta_1 \Delta_2 (\cos(\phi_1 - \phi_2) + i \sin(\phi_1 - \phi_2)). \quad (4.6.16)$$

Thus

$$\Re \Delta_k \Delta_q^* = \Delta_1 \Delta_2 \cos(\phi_1 - \phi_2), \quad (4.6.17)$$

where \Re stands for real part. Making use of

$$\sum_k \rightarrow N_1 \int d\epsilon_1, \quad \sum_q \rightarrow N_2 \int d\epsilon_2 \quad (4.6.18)$$

where N_1 and N_2 are the density of levels of one spin orientation at the Fermi energy one obtains⁶¹

$$\begin{aligned} \Delta E_2 &\approx -N_1 N_2 \Delta_1 \Delta_2 \langle |T_{kq}|^2 \rangle \cos(\phi_1 - \phi_2) \int_{-\infty}^{\infty} \int_{-\infty}^{\infty} \frac{d\epsilon_1 d\epsilon_2}{E_1 E_2 (E_1 + E_2)} \\ &\approx -2\pi N_1 N_2 \langle |T_{kq}|^2 \rangle \cos(\phi_1 - \phi_2) \left(\pi \frac{\Delta_1 \Delta_2}{\Delta_1 + \Delta_2}\right). \end{aligned} \quad (4.6.19)$$

Replacing the phase difference in (4.6.19) by the gauge-invariant phase difference $\gamma = (\phi_1 - \phi_2) - \frac{2e}{\hbar} \int_1^2 \mathbf{A} \cdot d\ell$, where \mathbf{A} is the magnetic vector potential ($\nabla \times \mathbf{A} = \mathbf{B}$), and the line integral is carried across the barrier of thickness d ($\int_1^2 d\ell = d$), one can

⁶¹It is of notice that ΔE_2 is bilinear in the density of levels (see Potel et al. (2017)).

calculate the supercurrent⁶² whose maximum value for two identical superconductors is⁶³

$$J_0 = \left(\frac{\pi \Delta}{2e} \right) \times \frac{1}{R_b} = \frac{\pi}{4} \frac{V_{equiv}}{R_b}, \quad (4.6.20)$$

where

$$\frac{1}{R_b} = \frac{4\pi e^2}{\hbar} N^2 \langle |T|^2 \rangle. \quad (4.6.21)$$

The quantity R_b is the tunneling resistance per unit area of the junction (see footnote 62 of this Chapter) when both metals are in the normal state, and $V_{equiv} = 2\Delta/e$ is the minimum voltage at which a normal current, upon the breaking of Cooper pairs, starts to flow⁶⁴.

As it emerges from Fig. 4.6.1, in which a bias of value $V \gtrsim 2\Delta/e$ is applied to the barrier, Cooper pairs are broken and quasiparticle excitations created –thus the labeling (S-Q)– resulting in the flow of a normal (dissipative) current of carriers $q = e$. For $T = 0$, one is in presence of processes connecting a ground state (S) with ground and excited states (Q). The importance of this fact in connection with the Josephson-like junction transiently formed in heavy ion reactions becomes apparent below (Sect. 7.3).

For Pb and Sn, the gaps are 1.4 meV and 0.7 meV, leading to eV_{equiv} ($V_{equiv} \approx 2.8$ meV (2.8 mV) and 1.4 meV (1.4 mV) respectively. Assuming R_b to be of the order of 1Ω per unit area, implies maximum values of the Josephson supercurrent $J = J_0 \sin \gamma$, of the order of $J_0 \approx 2$ mA as experimentally observed.

It is suggestive that the expression (4.6.19) is formally similar to that of the ion-ion potential acting between two heavy ions in weak contact, namely at a distance a diffusivity away from the grazing distance r_g . In this case the role of the reduced gap is played by a quantity closely related to the reduced radius of curvature⁶⁵

$$U_{aA}^N(r_g + a) \sim -\gamma \frac{R_a R_A}{R_a + R_A} a. \quad (4.6.22)$$

In the above expression $\gamma \approx 0.9$ MeV/fm² is the surface tension, $a = 0.63$ fm the diffusivity of the potential, $R_i (= (1.233A^{1/3} - 0.98A^{-1/3})$ fm) being the radii of nuclei

⁶² $I = \frac{1}{d} (\delta(\Delta E_2)/\delta \mathbf{A}) = (2e/\hbar) N_a N_A \langle |T|^2 \rangle 2\pi (\Delta_a \Delta_A / (\Delta_a + \Delta_A)) \sin \left((\phi_1 - \phi_2) - \frac{2e}{\hbar} \int_1^2 \mathbf{A} \cdot d\mathbf{l} \right)$, where one has divided by the thickness of the barrier d , so as to obtain, in keeping with the Meissner effect (Meissner and Ochsenfeld (1933)), a supercurrent per unit area ($\langle |T|^2 \rangle$ is assumed to be an energy squared per unit area).

⁶³ Tinkham (1996) Ch. 6, Eq. (6-4) and subsequent discussion; see also Anderson (1964a) Eq. (11) and following discussion.

⁶⁴ Giaever (1973).

⁶⁵ Broglia and Winther (2004) p.114 Eq. (40), $U_{aA}^N(r) = -V_0/(1 + \exp(\frac{r-R_0}{a}))$, $V_0 = 16\pi\gamma R_a A a$, $R_0 = R_a + R_A + 0.29$ fm, which for two ¹²⁰Sn nuclei ($R = 5.883$ fm) leads to $R_0 \approx 12.1$ fm and $V_0 = 83.8$ MeV. For energies somewhat above the Coulomb barrier, the grazing distance (Eq. (25) p. 128) of the above reference is $r_g = r_B - \delta \approx 12.8$ fm ($r_B \approx 13.3$ fm, $\delta \approx 0.5$ fm). Thus $(1 + \exp(\frac{r_g+a-R_0}{a})) \approx 9.26$.

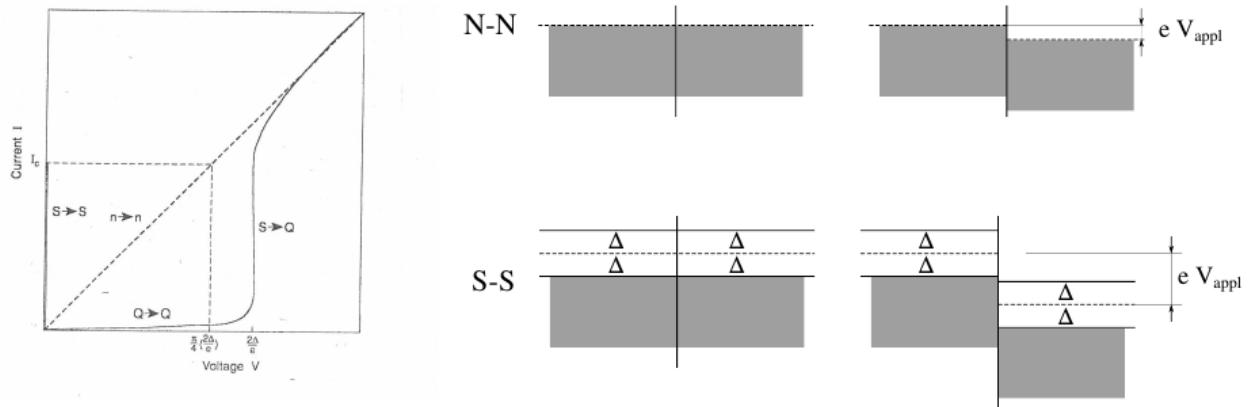


Figure 4.6.1: Current-voltage characteristic curves for tunneling between two metals across a barrier (thin oxide layer of 10–30 Å). It is assumed that both metals are equal and that they can become superconductors at low temperature (e.g. Al-Al₂O₃-Al, Sn-SnO_x-Sn). N-N indicates the current resulting from the *tunneling of single electrons* between the metals, both of them in the normal phase. S-Q indicates the single-electron tunneling between two superconductors (S-S) at finite temperature ($T < T_c$). It shows a small current flow for $0 < V < 2\Delta/e$, and the usual large flow for $V \gtrsim 2\Delta/e$, in which case Cooper pairs are broken and quasiparticles (Q) become excited. S-S labels the (super) current associated with the *tunneling of Cooper pairs* through the unbiased barrier ($V = 0$, no potential loss). That is, the (dc) Josephson effect. The critical (dc) Josephson current is equal to $\pi/4 (\approx 80\%)$ of the N-N state normal (single electron carrier) current at the voltage $V = 2\Delta/e$, 2Δ being the energy necessary to break Cooper pairs and excite quasiparticles.

$i = a, A$. For two identical nuclei $R_a = R_A = R$ and $V_0 = 8\pi\gamma Ra$. In the case in which the interacting nuclei are two ^{120}Sn systems $U_{aA}^N(r_g + a) = U_{aA}^N(13.43 \text{ fm}) \approx -9.1 \text{ MeV}$.

Being leptodermous systems, nuclei can be described at profit concerning a number of properties, with the help of the liquid drop. Because at the grazing distance the two leptodermous objects overlap, although weakly, two “unit” areas disappear. To reconstruct them one has to separate the two liquid drops until these areas are formed again. The energy needed to do so has to compensate for the value (4.6.22). Now, this quantity is a measure of the force that must be applied to surface molecules (nucleons) so that they experience the same situation as a molecule does when it is in the interior of the liquid (nucleus).

Similarly, the weak link (4.6.19) between the two superconductors 1 and 2, is associated with the situations in which each partner of the Cooper pair which tunnels is in a different superconductor, a kind of incipient (covalent) phenomenon, the Cooper pair being simultaneously shared by the two superconductors. The resulting tunneling interaction energy is proportional to the reduced pairing gap, that is $(\Delta_a\Delta_A)/(\Delta_a + \Delta_A)$.

4.7 Rotation in gauge space

The occurrence of rotation as a feature of the nuclear spectrum, e.g. of pairing rotational bands, originates in the phenomenon of spontaneous symmetry breaking of rotational invariance in the two dimensional gauge space. In other words, violation of particle number conservation, which introduces a deformation that makes it possible to specify an orientation of the system in gauge space.

The condensate in the nuclear superfluid system involve a deformation of the field that creates the fermion pairs. The process of addition or removal of a Cooper pair constitutes *a rotational mode in gauge space in which particle number plays the role of angular momentum*. Pairing rotational bands represents the collective mode associated with restoration of particle number conservation. Let us elaborate on the above points within the framework of a simple model which contains the basic physical features one is interested in discussing.

We consider N nucleons moving in a single j -shell⁶⁶ of energy ϵ_j and total angular momentum $(ls)_j$. The number of pairs moving in time reversal state which can be accommodated in the shell is $\Omega = (2j + 1)/2$. Consequently, the value of the BCS occupation parameters are $V = (N/2\Omega)^{1/2}$ and $U = (1 - N/2\Omega)^{1/2}$. The solution of the BCS number and gap equations associated with a pairing force of constant matrix elements G are

$$\lambda = -\frac{G}{2}(\Omega - N), \quad (4.7.1)$$

⁶⁶For details see e.g. App. H of Brink and Broglia (2005) and refs. therein.

and

$$\Delta = \frac{G}{2} \sqrt{N(2\Omega - N)}, \quad (4.7.2)$$

respectively.

The BCS ground state energy of the superfluid system is

$$U = 2 \sum_{\nu>0} (\epsilon_\nu - \lambda) V_\nu^2 - \frac{\Delta^2}{G}. \quad (4.7.3)$$

The fact that the single-particle energies are measured with respect to λ implies that the nucleons feel the Coriolis force ($\hbar\dot{\phi} = \lambda$) associated with rotation in gauge space (Eq. (4.7.11) below). In other words, the BCS solution is carried out in the intrinsic system of reference \mathcal{K}' .

We now calculate the ground-state energy in the laboratory frame \mathcal{K} (see Fig. 1.4.1),

$$E_0 = U + \lambda N = 2 \sum_{\nu>0} \epsilon_\nu V_\nu^2 - \frac{\Delta^2}{G}. \quad (4.7.4)$$

Because all m -substates of the single j -shell are degenerate $\epsilon_\nu = \epsilon$. Setting, for simplicity, $\epsilon_\nu = 0$ the above equation can be written as,

$$E_0 = -G \frac{\Omega}{2} N + \frac{G}{4} N^2. \quad (4.7.5)$$

Assuming $\Omega \gg N$ one obtains from Eq. (4.7.1) $\lambda \approx -G\Omega/2$, leading to

$$E_0 = \lambda N + \frac{\hbar^2}{2\mathcal{J}} N^2, \quad (4.7.6)$$

where

$$\frac{\hbar^2}{2\mathcal{J}} = \frac{G}{4}, \quad (4.7.7)$$

\mathcal{J} being the moment of inertia of the associated pairing rotational band⁶⁷.

4.7.1 Phase coherence

The phase of a wavefunction and the number of nucleons (electrons in condensed matter) are conjugate variables. In other words, gauge angle and particle number operators $\hat{\phi}$ and \hat{N} satisfy the commutation relation,

$$[\hat{\phi}, \hat{N}] = i. \quad (4.7.8)$$

⁶⁷See Bès and Broglia (1966), Broglia et al. (2000), and references therein.

In the particle number representation

$$\hat{N} = -i\partial/\partial\phi, \quad \hat{\phi} = \phi, \quad (4.7.9)$$

while

$$\hat{N} = N, \quad \hat{\phi} = i\partial/\partial N, \quad (4.7.10)$$

in the gauge angle representation. The time derivative of the gauge angle is given by the equation of motion⁶⁸,

$$\dot{\phi} = \frac{i}{\hbar} [H, \phi] = \frac{1}{\hbar} \frac{\partial H}{\partial N} = \frac{1}{\hbar} \lambda, \quad (4.7.11)$$

where λ is the chemical potential (Fermi energy).

Gauge invariance, i.e. invariance under phase changes, implies number of particle conservation in a similar way that rotational invariance implies angular momentum conservation.

Example: let us introduce the invariant many-body state⁶⁹,

$$|\Psi_N\rangle = a_1^\dagger a_2^\dagger \cdots a_N^\dagger |0\rangle, \quad (4.7.12)$$

and rotate it an angle ϕ with the help of the operator $\mathcal{G}(\phi) = e^{-i\hat{N}\phi}$. In other words, making use of

$$\mathcal{G}(\phi) a_\nu^\dagger \mathcal{G}^{-1}(\phi) = e^{-i\phi} a_\nu^\dagger = a'_\nu^\dagger, \quad (4.7.13)$$

and as a result of the relation $a_\nu^\dagger = e^{i\phi} a'^\dagger_\nu$, one obtains

$$|\Psi_N\rangle = e^{iN\phi} |\Psi'_N\rangle, \quad (4.7.14)$$

where

$$|\Psi'_N\rangle = a'_1^\dagger a'_2^\dagger \cdots a'_N^\dagger |0\rangle. \quad (4.7.15)$$

Thus,

$$\hat{N} |\Psi_N\rangle = -i \frac{\partial}{\partial \phi} |\Psi_N\rangle = N |\Psi_N\rangle. \quad (4.7.16)$$

A phase change for a gauge invariant function is just a trivial operation. Like to rotate a rotational invariant function. Quantum mechanically nothing happens rotating a spherical, symmetry conserving, system (in 3D-, gauge, etc.) space.

⁶⁸See e.g. Brink and Broglia (2005) App. I.

⁶⁹Anderson (1964b).

The situation is different in the case of the wavefunction

$$\begin{aligned}
 |\text{BCS}(\phi)\rangle_{\mathcal{K}} &= \prod_{\nu>0} (U_\nu + V_\nu a_\nu^\dagger a_{\bar{\nu}}^\dagger) |0\rangle, \\
 &= \prod_{\nu>0} (U'_\nu + e^{-2i\phi} V'_\nu a_\nu^\dagger a_{\bar{\nu}}^\dagger) |0\rangle, \\
 &= \prod_{\nu>0} (U'_\nu + V'_\nu a_\nu'^\dagger a_{\bar{\nu}}'^\dagger) |0\rangle, \\
 &= |\text{BCS}(\phi = 0)\rangle_{\mathcal{K}'}, \tag{4.7.17}
 \end{aligned}$$

where

$$U_\nu = |U_\nu| = U'_\nu; \quad V_\nu = e^{-2i\phi} V'_\nu (V'_\nu = |V_\nu|). \tag{4.7.18}$$

In fact,

$$|\text{BCS}(\phi)\rangle_{\mathcal{K}} = \left(\prod_{\nu>0} U'_\nu \right) \sum_{N \text{ even}} \frac{e^{-iN\phi}}{(N/2)!} \left(\sum_{\nu>0} c'_\nu P_\nu^\dagger \right)^{N/2} |0\rangle, \tag{4.7.19}$$

with

$$c'_\nu = \frac{V'_\nu}{U'_\nu}; \quad P_\nu^\dagger = a_\nu^\dagger a_{\bar{\nu}}^\dagger, \tag{4.7.20}$$

is a coherent state (see Eq. (1.4.39)) in particle number which can be written as,

$$|\text{BCS}(\phi)\rangle_{\mathcal{K}} = \sum_N f_N(\phi) |\Psi_N\rangle. \tag{4.7.21}$$

Let us now apply the gauge angle operator to it,

$$\hat{\phi} |\text{BCS}(\phi)\rangle_{\mathcal{K}} = \hat{\phi} \sum_N f_N(\phi) |\Psi_N\rangle, \tag{4.7.22}$$

that is,

$$i \frac{\partial}{\partial N} f_N(\phi) = \phi f_N(\phi). \tag{4.7.23}$$

Thus,

$$f_N(\phi) \sim e^{-iN\phi}, \tag{4.7.24}$$

and, making use of Eq. (4.7.14),

$$|\text{BCS}(\phi)\rangle_{\mathcal{K}} \sim \sum_N e^{-iN\phi} |\Psi_N\rangle = \sum_N |\Psi'_N\rangle \sim |\text{BCS}(\phi = 0)\rangle_{\mathcal{K}'}, \tag{4.7.25}$$

the state $|\text{BCS}(\phi = 0)\rangle_{\mathcal{K}'}$ being aligned in gauge space in which it defines a privileged orientation (z') with respect to the laboratory system.

An isolated nucleus will not remain long in this product type state. Due to the term⁷⁰ $(G/4) \left(\sum_{\nu>0} (U_\nu^2 + V_\nu^2) (\Gamma_\nu^\dagger - \Gamma_\nu) \right)^2$ in the residual quasiparticle Hamiltonian it will fluctuate, and decay into a state

$$|N\rangle \sim \int d\phi e^{iN\phi} |BCS(\phi)\rangle_{\mathcal{K}} \sim \sum_{N'} \int_0^{2\pi} d\phi e^{-i(N'-N)\phi} |\Psi_{N'}\rangle \sim |\Psi_N\rangle \quad (4.7.26)$$

in keeping with the fact that

$$\int_0^{2\pi} d\phi e^{-i(N'-N)\phi} = \begin{cases} \frac{i}{N'-N} e^{-i(N'-N)\phi} \Big|_0^{2\pi} = 0 & (N \neq N'), \\ 2\pi \delta(N, N') & (N = N'). \end{cases} \quad (4.7.27)$$

The state $|N\rangle$ is a member of a pairing rotational band.

In the example discussed in connection with Figs. 3.1.2 and 3.1.3, $|N\rangle$ is one of the ground states of the Sn-isotopes and $N_0 = 68$. Making use of the fact that $E_R = (\hbar^2/2I)(N - N_0)^2$ and that $\delta N \delta\phi \approx 1$, the coherent state (4.7.21) will decay into one of the states $|N\rangle$, likely the one corresponding to $N = \sum_{\nu>0} 2V_\nu^2$, in a time \hbar/E_R . The member of Cooper pairs participating in the nuclear condensate in the case of e.g. ^{120}Sn is $\alpha'_0 \approx 5 - 6$, the number of neutrons being $2\alpha'_0 \approx 10$. Because $\delta N \sim \sqrt{N} \sim 3$ ($\delta\phi \sim 0.3$ rad, that is⁷¹ $\delta\phi \sim 0.3/0.017 \approx 17^\circ$), the energy of rotation in gauge space of a state which is defined with an uncertainty of $\delta\phi \approx 0.3$ rad is $E_R \approx 0.092 \text{ MeV} \times (3)^2 \approx 1 \text{ MeV}$ ($\hbar^2/2I \approx G/4 \approx 25 \text{ MeV}/(4N_0) \approx 0.092 \text{ MeV}$, see Fig. 3.1.3). Consequently $\hbar/E_R \approx 10^{-21} \text{ s}$.

This is also the case for metallic superconductors. In fact, the state (4.7.17) even if prepared in isolation will dissipate because there is a term in the energy of the superconductor depending on N , namely the electrostatic energy $e^2(N - N_0)^2/2C$, where C is the electrostatic capacity⁷². Because the number of overlapping Cooper pairs contributing to superconductivity lying around the Fermi surface⁷³ is $\alpha'_0 \approx 10^6$ and thus the associated number of electrons $2\alpha'_0 \approx 2 \times 10^6$, $\delta N \sim \sqrt{N} \sim 10^3$, one can write $E_{el} = \frac{e^2}{2C} (\delta N)^2 \approx \frac{14.4 \text{ eV} \times 10^{-8} \text{ cm}}{2 \times 1 \text{ cm}} (10^3)^2 \approx 10^{-7} \text{ MeV}$, and $\hbar/E_{el} \approx 10^{-14} \text{ s}$. In this case $\delta\phi \sim 10^{-3}$ radians, that is $\delta\phi \sim 10^{-3}/0.017 \sim 1^\circ$.

⁷⁰Within BCS theory of pairing, there are two parameters which determines spontaneous symmetry breaking in gauge space. The probability amplitude with which a pair state ($\nu\bar{\nu}$) is occupied, and that with which it is empty. Namely, V_ν and U_ν respectively. As a consequence, there only two fields F which contribute, through terms of type FF^\dagger , to the residual interaction H_{res} acting among quasiparticles, which is neglected in the mean field solution of the pairing Hamiltonian. One which is antisymmetric with respect to the Fermi surface namely $(U_\nu^2 - V_\nu^2)$ and which leads to pairing vibrations of the gauge deformed state $|BCS\rangle$. The other one, $(U_\nu^2 + V_\nu^2)$ is symmetric with respect to ϵ_F and leads to fluctuations which diverge in the long wavelength limit ($W''_1 \rightarrow 0$) in precisely the right way to set $|BCS\rangle$ into rotation with a finite inertia, and restore gauge symmetry (see Eq. 1.4.57).

⁷¹ $\delta\phi \approx 0.3/(\pi/180) \approx 0.3/0.017 \approx 17^\circ$.

⁷²The capacity of a sphere is $C = R$. Below we use $R = 1\text{cm}$ (see Anderson (1964b)).

⁷³In the case of Pb, $\xi \approx 0.3 \times 10^4 \text{ \AA}$, $r_s = 1.22 \text{ \AA}$, $\Delta \approx 1.4 \text{ meV}$ and $\epsilon_F \approx 9.47 \text{ eV}$, are the correlations length, the Wigner-Seitz radius, the pairing gap and the Fermi energy respectively. Consequently $(\xi/r_s)^3 \times (\Delta/\epsilon_F) \approx 10^6$.

The opposite situation is that of the case in which one considers different parts of the same superconductor. In this case one can define relative variables $n = N_1 - N_2$ and $\phi = \phi_1 - \phi_2$ and again $n = -i\partial/\partial\phi$ and $\phi = i\partial/\partial n$. Thus, locally there is a superposition of different n states: ϕ is fixed so n is uncertain. There is a dividing line between these two behaviors, perfect phase coherence and negligible coherence, namely the Josephson effect.

Again, the total phase of the assembly is not physical. However, the relative phases can be given a meaning when one observes, as one does in e.g. junctions between metallic superconductors, that pairs of electrons can pass back and forth through the barrier, leading to the possibility of coherence between states in which the total number of electrons is not fixed *locally*. Under such conditions there is, for instance, a coherence between the state with $N/2$ electrons in one half of the block and $N/2$ in the other, and that with $(N/2) + 2$ on one side and $(N/2) - 2$ on the other.

4.8 Medium polarization effects and pairing

In many-body systems, medium polarization effects play an important role in renormalizing single-particle motion and the four-point vertices, namely Coulomb interaction in condensed matter and the bare NN -potential in nuclei. In what follows, special attention is paid to this mechanism in connection with the pairing interaction.

4.8.1 Nuclei

Elementary modes of excitation constitute a basis of states in which correlations, as found in observables, play an important role. As a consequence, it allows for an economic, low dimensionality solution of the nuclear many-body problem of structure and reaction.

As a result of their interweaving, the variety of elementary modes of excitation may break in a number of states, eventually acquiring a lifetime and, within a coarse grain approximation, a damping width (imaginary component of the self energy). Moving into the continuum, as for example in the case of direct reactions, one such component is the imaginary part of the optical potential⁷⁴ operating in the particular channel selected. It can, in principle, be calculated microscopically using similar techniques and elements as e.g. those used in the calculation of the damping width of giant resonances or of single-particle states. In this way, the consistency circle structure-reaction based on elementary modes and codified by NFT could be closed. The rich variety of emergent properties found along the way eventually acquiring a conspicuous level of physical validation. At that time it would be possible, arguably if there is one, to posit that the *ultima ratio* of structure

⁷⁴See footnote 23 Ch. 1.

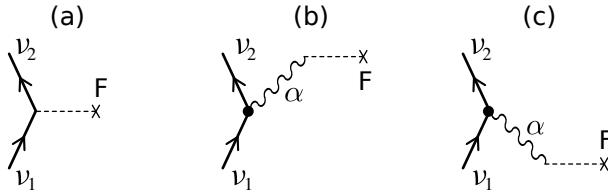


Figure 4.8.1: (a) F -moment of a single-particle and (b,c) renormalization effects induced by the collective vibration α .

and reactions, in any case that associated with pairing and Cooper pair transfer in nuclei, have been unveiled⁷⁵.

Effective moments

At the basis of the coupling between elementary modes of excitation, for example of single-particle motion and of collective vibrations, one finds the fact that, in describing the nuclear structure it is necessary to make reference to both of them simultaneously and in an unified way.

Within the harmonic approximation the above statement is economically embodied in e.g. the relation existing between the collective ($\hat{\alpha}$) and the single-particle (\hat{F}) representation of the operator creating e.g. a particle-hole excitation. That is⁷⁶,

$$\begin{aligned}
 \hat{F} &= \left\{ \langle k | F | i \rangle \Gamma_{ki}^\dagger + \langle i | F | k \rangle \Gamma_{ki} \right\} \\
 &= \sum_{k,i,\alpha'} X_{ki}^{\alpha'} \Gamma_{\alpha'}^\dagger - Y_{ki}^{\alpha'} \Gamma_{\alpha'} \\
 &= \sum_{\alpha'} \Lambda_{\alpha'} \sum_{ki} \frac{|\langle i | F | k \rangle|^2}{(\epsilon_k - \epsilon_i)^2 - (\hbar\omega_{\alpha'})^2} (\Gamma_{\alpha'}^\dagger + \Gamma_{\alpha'}) \\
 &= \sum_{\alpha'} \frac{\Lambda_{\alpha'}}{\kappa} (\Gamma_{\alpha'}^\dagger + \Gamma_{\alpha'}) = \sum_{\alpha'} \sqrt{\frac{\hbar\omega_{\alpha'}}{2C'_\alpha}} (\Gamma_{\alpha'}^\dagger + \Gamma_{\alpha'}) = \hat{\alpha}. \quad (4.8.1)
 \end{aligned}$$

This is a consequence of the self consistency relation

$$\delta U(r) = \int d\mathbf{r}' \delta\rho(r) v(|\mathbf{r}-\mathbf{r}'|), \quad (4.8.2)$$

⁷⁵In the above paragraph we allowed ourselves to paraphrase Jacques Monod writing in connection with biology and life: *L'ultima ratio de toutes les structures et performances téléonomiques des êtres vivants est donc enfermée dans les séquences des radicaux des fibres polypeptidiques "embryons" de ces démons de Maxwell biologiques que sont les protéines globulaires. En un sens, très réel, c'est à ce niveau d'organisation chimique qui gît, s'il y en a un, le secret de la vie. Et saurait-on non seulement décrire les séquences, mais énoncer la loi d'assemblage à laquelle obéissent, on pourrait dire que le secret est percé, l'ultima ratio découverte (Monod (1970)).*

⁷⁶cf. Bohr and Mottelson (1975), cf. also Brink and Broglia (2005) App. C.

existing between density (collective) and potential (single-particle) distortion, typical of normal modes of many-body systems. It is of notice that the assumption is made that F is dimensionless while κ (see Eq. (4.8.3)) has energy dimensions (see e.g. Eqs. (1.2.7) and (1.2.13)).

Relation (4.8.1) implies that at the basis of these normal modes one finds the (attractive $\kappa < 0$) separable interaction

$$H = \frac{\kappa}{2} \hat{F} \hat{F}, \quad (4.8.3)$$

but where now (Fig. 4.8.1 (a))

$$\hat{F} = \sum_{\nu_1, \nu_2} \langle \nu_1 | F | \nu_2 \rangle a_{\nu_1}^\dagger a_{\nu_2}, \quad (4.8.4)$$

is a general single-particle operator, while \hat{F} in Eq. (4.8.1) is its harmonic representation acting in the particle (k)–hole (i) space, Γ_{ki}^\dagger and Γ_{ki} being (quasi) bosons, i.e. respecting the commutation relation

$$[\Gamma_{ki}, \Gamma_{k'i'}^\dagger] = \delta(k, k') \delta(i, i'). \quad (4.8.5)$$

In other words, the representation (4.8.1), which is at the basis of the RPA (as well as QRPA), does not allow for scattering vertices, processes which become operative by rewriting (4.8.3) in terms of the particle–vibration coupling Hamiltonian

$$H_c = \kappa \hat{\alpha} \hat{F}. \quad (4.8.6)$$

It is of notice that κ is negative for an attractive field. Let us now calculate the effective single-particle moments (cf. Fig. 4.8.1 (b)),

$$\begin{aligned} \langle \nu_2 | \hat{F} | \nu_1 \rangle_{(b)} &= \frac{\langle \nu_2 | \hat{F} | \nu_2, n_\alpha = 1 \rangle \langle \nu_2, n_\alpha = 1 | H_c | \nu_1 \rangle}{(\epsilon_{\nu_1} - \epsilon_{\nu_2}) - \hbar\omega_\alpha}, \\ &= \frac{\langle 0 | \hat{\alpha} | n_\alpha = 1 \rangle \kappa \alpha \langle \nu_2 | F | \nu_1 \rangle}{(\epsilon_{\nu_1} - \epsilon_{\nu_2}) - \hbar\omega_\alpha}, \\ &= \kappa \alpha^2 \frac{\langle \nu_2 | F | \nu_1 \rangle}{(\epsilon_{\nu_1} - \epsilon_{\nu_2}) - \hbar\omega_\alpha}, \end{aligned} \quad (4.8.7)$$

and (Fig. 4.8.1 (c))⁷⁷

$$\begin{aligned} \langle \nu_2 | \hat{F} | \nu_1 \rangle_{(c)} &= \frac{\langle \nu_2 | H_c | \nu_1, n_\alpha = 1 \rangle \langle \nu_1, n_\alpha = 1 | F | \nu_1 \rangle}{\epsilon_{\nu_2} - (\epsilon_{\nu_1} + \hbar\omega_\alpha)}, \\ &= \kappa \alpha^2 \left(-\frac{\langle \nu_2 | F | \nu_1 \rangle}{(\epsilon_{\nu_1} - \epsilon_{\nu_2}) + \hbar\omega_\alpha} \right), \end{aligned} \quad (4.8.8)$$

⁷⁷In calculating the energy denominators one takes the difference between the energy of the initial and of the intermediate states. However, when an external field like \hat{F} acts on the system, before the PVC or four-point vertices operate, being equivalent to an observation, the energy denominator is to be calculated as the energy difference between the final and the intermediate states.

leading to

$$\begin{aligned} \langle v_2 | \hat{F} | v_1 \rangle_{(b)} + \langle v_2 | \hat{F} | v_1 \rangle_{(c)} &= \kappa \alpha^2 \frac{2\hbar\omega_\alpha \langle v_2 | F | v_1 \rangle}{(\epsilon_{v_1} - \epsilon_{v_2})^2 - (\hbar\omega_\alpha)^2}, \\ &= \frac{\kappa}{C_\alpha} \frac{(\hbar\omega_\alpha)^2 \langle v_2 | F | v_1 \rangle}{(\epsilon_{v_1} - \epsilon_{v_2})^2 - (\hbar\omega_\alpha)^2}. \end{aligned} \quad (4.8.9)$$

This is in keeping with the fact that the ZPF of the α -vibrational mode is,

$$\alpha = \sqrt{\frac{\hbar\omega_\alpha}{2C_\alpha}}. \quad (4.8.10)$$

The particle-vibration coupling strength used below is defined as,

$$\Lambda_\alpha = \kappa\alpha. \quad (4.8.11)$$

Together with $\langle v_2 | \hat{F} | v_1 \rangle_{(a)} = \langle v_2 | F | v_1 \rangle$ (see Fig. 4.8.1 (a)) one obtains

$$\langle v_2 | \hat{F} | v_1 \rangle = (1 + \chi(\omega)) \langle v_2 | F | v_1 \rangle, \quad (4.8.12)$$

where

$$\chi_\alpha(\omega) = \frac{\kappa}{C_\alpha} \frac{\omega_\alpha^2}{\omega^2 - \omega_\alpha^2} \quad (4.8.13)$$

is the polarizability coefficient while

$$\omega = |\epsilon_{v_1} - \epsilon_{v_2}|/\hbar. \quad (4.8.14)$$

In the static limit, e.g. in the case in which α is a giant resonance and $\omega_\alpha \gg \omega$ one obtains

$$\chi_\alpha(0) = -\frac{\kappa}{C_\alpha}. \quad (4.8.15)$$

The sign of $\chi_\alpha(0)$ is opposite to that of κ , since the static polarization effect produced by an attractive coupling ($\kappa < 0$) is in phase with the single-particle moment, while a repulsive coupling ($\kappa > 0$) implies opposite phases for the polarization effect and the one-particle moment⁷⁸.

Let us now calculate the two-body pairing induced interaction (Fig. 4.8.2) arising from the exchange of collective vibrations⁷⁹, summing over the two time

⁷⁸Bohr and Mottelson (1975); Mottelson (1962). Think in the first case about the effect the GQR ($\tau = 0$) plays in $(e(E2))_{eff}$, in the second that played by the GDR ($\tau = 1$) in $(e(E1))_{eff}$.

⁷⁹In the present discussion we do not consider spin modes. For details see e.g. Idini et al. (2015). See also Bortignon et al. (1983).

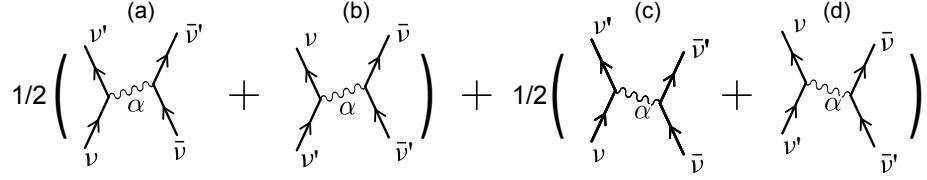


Figure 4.8.2: Diagrams associated with nuclear pairing induced interaction.

orderings and symmetrizing between initial and final states⁸⁰

$$\begin{aligned}
 v_{\nu\nu'}^{ind}(a) + v_{\nu\nu'}^{ind}(b) &= \kappa^2 \alpha^2 |\langle \nu' | F | \nu \rangle|^2 \left(\frac{1}{\epsilon_\nu - \epsilon_{\nu'} - \hbar\omega_\alpha} + \frac{1}{\epsilon_{\nu'} - \epsilon_\nu - \hbar\omega_\alpha} \right), \\
 &= \kappa^2 \alpha^2 |\langle \nu' | F | \nu \rangle|^2 \left(\frac{1}{(\epsilon_\nu - \epsilon_{\nu'}) - \hbar\omega_\alpha} - \frac{1}{(\epsilon_\nu - \epsilon_{\nu'}) + \hbar\omega_\alpha} \right), \\
 &= \Lambda_\alpha^2 |\langle \nu' | F | \nu \rangle|^2 \left(\frac{2\hbar\omega_\alpha}{(\epsilon_\nu - \epsilon_{\nu'})^2 - (\hbar\omega_\alpha)^2} \right), \\
 &= v_{\nu\nu'}^{ind}(c) + v_{\nu\nu'}^{ind}(d).
 \end{aligned} \tag{4.8.16}$$

Thus

$$\begin{aligned}
 v_{\nu\nu'}^{ind} &= \frac{1}{2} (v_{\nu\nu'}^{ind}(a) + v_{\nu\nu'}^{ind}(b)) + \frac{1}{2} (v_{\nu\nu'}^{ind}(c) + v_{\nu\nu'}^{ind}(d)) \\
 &= \Lambda_\alpha^2 |\langle \nu' | F | \nu \rangle|^2 \left(\frac{2\hbar\omega_\alpha}{(\hbar\omega)^2 - (\hbar\omega_\alpha)^2} \right).
 \end{aligned} \tag{4.8.17}$$

The diagonal matrix element,

$$v_{\nu\nu}^{ind} \equiv -\frac{2\Lambda_\alpha^2 |\langle \nu | F | \nu \rangle|^2}{\hbar\omega_\alpha},$$

testifies to the fact, for values of $\omega_\alpha \gtrsim \omega$, the induced pairing interaction is attractive. Summing to (4.8.17) the matrix element of the bare interaction (4.8.3), (Fig. 4.8.3 (b))⁸¹

$$v_{\nu\nu'}^{bare} = \kappa |\langle \nu' | F | \nu \rangle|^2, \tag{4.8.18}$$

⁸⁰Cf. Brink and Broglia (2005) p. 217.

⁸¹Within the framework of (4.8.3) and of its role in (4.8.19) one finds, in the case of superconductivity in metals to be discussed below, that the bare unscreened Coulomb interaction can be written as

$$U_c(r) = \frac{1}{2} \sum_{i,j} \frac{q_i q_j}{|\mathbf{r}_i - \mathbf{r}_j|},$$

i, j running over all particles (nuclei and electrons) and $q_i = -e$ for electrons and Ze for nuclei.

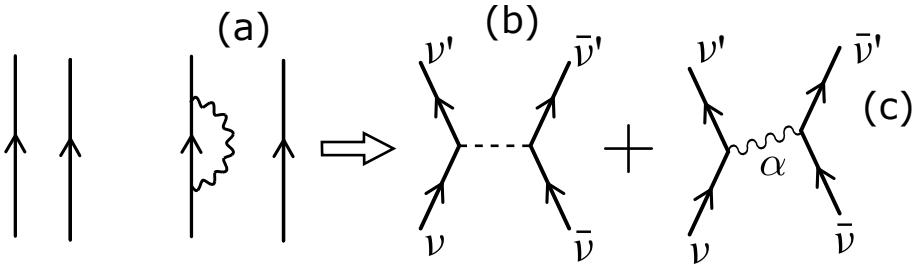


Figure 4.8.3: Starting with two bare nucleons moving around a closed shell system N_0 in Hartree-Fock orbitals (arrowed lines far left), a graphical (NFT) representation of (a) self energy processes and of (b) bare and (c) induced pairing interactions are displayed.

one obtains for the total pairing matrix element⁸²

$$\begin{aligned} v_{vv'} &= \kappa |\langle v'|F|v\rangle|^2 + \frac{\kappa^2}{C_\alpha} |\langle v'|F|v\rangle|^2 \frac{\omega_\alpha^2}{\omega^2 - \omega_\alpha^2} = v_{vv'}^{bare} + \frac{\kappa}{C_\alpha} v_{vv'}^{bare} \frac{\omega_\alpha^2}{\omega^2 - \omega_\alpha^2} \\ &= v_{vv'}^{bare} \left(1 + v_{vv'}^{bare} \Pi_{vv'}(\omega, \omega_\alpha) \right). \end{aligned} \quad (4.8.19)$$

where

$$\Pi_{vv'} = \begin{cases} \left(C_\alpha |\langle v'|F|v\rangle|^2 \right)^{-1} \frac{\omega_\alpha^2}{\omega^2 - \omega_\alpha^2}, \\ \left(D_\alpha |\langle v'|F|v\rangle|^2 \right)^{-1} \frac{1}{\omega^2 - \omega_\alpha^2}, \end{cases} \quad (4.8.20)$$

both expressions being equivalent in keeping with the fact that $\omega_\alpha = (C_\alpha/D_\alpha)^{1/2}$. In the second expression of $\Pi_{vv'}$ the inertia of the phonon appears in the denominator, similar to the factor (Z/AM) in (4.8.63) below. The units of both C_α and (\hbar^2/D_α) are MeV and, as a consequence, $\hbar\omega_\alpha = (\hbar^2 C_\alpha/D_\alpha)^{1/2}$ is also an energy. The units of $\Pi_{vv'}$ are MeV⁻¹ in keeping with the units of $v_{vv'}^{bare}$.

It is of notice that $v_{vv'}$ can display a resonant behaviour leading to attraction regardless whether $v_{vv'}^{bare}$ is attractive or repulsive ($\lim_{\omega \rightarrow \omega_\alpha^-} \Pi_{vv'} \rightarrow -\infty$), a situation of relevance in the case of dielectric polarization effects in metals. In this case, the resulting effective electron-electron interaction (phonon exchange) is attractive (see Eq. (4.8.62)), and at the basis of Cooper pair correlation and BCS superconductivity.

The importance of having available a bare interaction which allows for the calculation of HF and RPA solutions and thus of renormalized single-particle and of effective interactions on equal footing is apparent from Eq. (4.8.19) (see also the last paragraph of Sect. 4.8.3 in connection with pairing in metals).

Let us consider the nucleus ^{11}Li , in which case Cooper pair binding results mainly from the exchange of the low-lying dipole mode between the two halo nu-

⁸²It is of notice that $\Pi_{vv'}$ is closely related to Lindhard's function (Lindhard (no. 8, 1954)). See Eq. (4.8.63) below.

clei. Making use of the bare screened interaction⁸³ $G_{scr} \approx 0.05G \approx 0.1$ MeV ($G \approx 28/A$ MeV) ($v_{vv'}^{bare} = -G_{scr} = -0.1$ MeV), and of the induced pairing interaction $v_{vv'}^{ind}$ ($= M_{ind} \approx -0.6$ MeV), one can write for the last term in (4.8.19), $(0.1 \text{ MeV})^2 \Pi_{vv'}(\omega_\alpha, \omega) = -0.6 \text{ MeV}$ leading to $\Pi_{vv} = -60 \text{ MeV}^{-1}$. This large effect is not a strictly resonant phenomenon, although the energies $\hbar\omega$ ($= \tilde{\epsilon}_{p_{1/2}} - \tilde{\epsilon}_{s_{1/2}} \approx 0.5$ MeV) and $\hbar\omega_\alpha$ (≈ 0.7 MeV) are rather similar, in keeping also with the fact that the pygmy dipole resonance PDR has a width of the order of $\Gamma \approx 0.5$ MeV.

At the basis of the quite small value of the energy centroid of the ^{11}Li soft $E1$ -mode (pygmy dipole resonance⁸⁴), one finds the low energy dipole particle-hole excitation between the parity inverted levels (0.45 MeV) and the screened symmetry potential ($V_1 = 25$ MeV) which arising from the poor overlap between the core (^9Li) and the halo neutron wavefunctions, as in the case of G_{scr} can be assumed, for order of magnitude estimates, to be equal to it (i.e. ≈ 0.05).

Thus $(V_1)_{scr} \approx 1.3$ MeV. One can compare the parametrization⁸⁵ $\hbar\omega_{GDR} = \hbar\omega_0 (1 + \kappa_1/C^{(0)})^{1/2} = 41 \times A^{-1/3} \text{ MeV} (1 + 3.5V_1/(41 \text{ MeV}))^{1/2} \approx 41^{-1/3} \text{ MeV} \times (1 + 88/41)^{1/2} \approx 41^{-1/3} \text{ MeV} \times 1.8 \approx 73A^{-1/3} \text{ MeV}$, used to describe the centroid of the GDR with Eq. (3.6.23). That is, $\hbar\omega_{pygmy} \approx 0.45 \text{ MeV} (1 + 0.6(V_1)_{screened}/0.20 \text{ MeV})^{1/2} \approx 0.45 \text{ MeV} \times (1 + 0.8/0.2)^{1/2} \approx 0.45 \text{ MeV} \times 2.2$. Expression which leads to $\hbar\omega_{pygmy} \approx 1 \text{ MeV}$ while the above one gives, for the GDR of ^{11}Li to $\hbar\omega_{GDR} \approx 33 \text{ MeV}$.

Expressed it differently, the screened repulsive symmetry potential can hardly increase the energy of the dipole particle-hole excitation $\tilde{\epsilon}_{p_{1/2}} - \tilde{\epsilon}_{s_{1/2}} \approx 0.45$ MeV.

Let us now carry a simple estimate of the contribution of the induced pairing interaction to the (empirical) nuclear pairing gap for nuclei lying along the stability valley. For this purpose we introduce the quantity

$$\lambda = N(0)v_{vv'}^{ind} \quad (4.8.21)$$

where $N(0)$ is the density of levels of a single spin orientation at the Fermi energy. The above quantity is known as the nuclear mass enhancement factor. This is because of the role it plays in the nucleon ω -mass (see Sects. 5.4 and 5.8)

$$m_\omega = (1 + \lambda)m. \quad (4.8.22)$$

Systematic studies of this quantity, and of the related discontinuity displayed by the single-particle occupation number at the Fermi energy, namely $Z_\omega = (m/m_\omega)$ testifies to the fact that $\lambda \approx 0.4$, within an energy interval of the order of 10 MeV around the Fermi energy.

The BCS expressions of the pairing gap in terms of λ are

$$\Delta = \begin{cases} 2\hbar\omega_D e^{-1/\lambda}, & (\text{weak coupling } \lambda \ll 1) \\ \hbar\omega_D \lambda, & (\text{strong coupling } \lambda \geq 1) \end{cases} \quad (4.8.23)$$

⁸³See Broglia et al. (2019b).

⁸⁴Broglia et al. (2019a)

⁸⁵Bohr and Mottelson (1975) Eq. (6-315a). See also Eq. (6-27) of this reference.

where ω_D is the limiting frequency of the low-lying collective modes of nuclear excitation, typically of quadrupole and octupole vibrations.

In keeping with the fact that collective modes, in particular quadrupole vibrations are found at low energies, we choose $\hbar\omega_D \approx 5$ MeV which, together with $\lambda = 0.4$, intermediate between weak and strong coupling situations leads to,

$$\Delta \approx 0.8 \text{ MeV} \quad (4.8.24)$$

if one uses the first expression (Eq. (4.8.23)), and to

$$\Delta \approx 2.0 \text{ MeV}, \quad (4.8.25)$$

if one uses the second. Values to be compared with the empirical value

$$\Delta \approx 1.4 \text{ MeV}, \quad (4.8.26)$$

of superfluid medium heavy mass nuclei like for example in the Sn-isotopes.

While the relations (4.8.23) can hardly be relied to provide a quantitative number, they testify to the fact that induced pairing is expected to play an important role in nuclei. These expectations have been confirmed by detailed confrontation of theory with experiment⁸⁶.

Hindsight

Static polarization effects can be important in dressing single-particle states. For example, effective charges and induced interactions associated with moments induced by giant resonances⁸⁷. However, retarded ω -dependent self-energy effects and induced interactions are essential in describing structure and reactions of many-body systems. Examples are provided by the bootstrap binding of the halo neutrons (pair addition mode) to ^9Li , leading to the fragile $|^{11}\text{Li}(gs)\rangle$, displaying a $S_{2n} \approx 0.380$ MeV as compared to typical values of $S_{2n} \approx 16$ MeV as far as structure goes, and by the $^1\text{H}(^{11}\text{Li}, ^9\text{Li}(1/2^-; 2.69 \text{ MeV}))^3\text{H}$ population of the lowest member of the $(2^+ \times p_{3/2}(\pi))_{J^-}$ multiplet of ^9Li with a cross section $\sigma(1/2^-; 2.69 \text{ MeV}) \approx 1 \text{ mb}$, as far as reaction goes. If there was need for support coming from other fields of research, one can mention just two: the van der Waals force in e.g. folding of proteins, and the electron-phonon interaction in superconductivity.

It was recognized early in the study of dipole-dipole interaction in atomic systems that, of the variety of contributions to the van der Waals interaction, the retarded, fully quantal contribution, arising from (dipole) zero point fluctuations (ZPF) of the two interacting atoms or molecules, and the only active in the case of non-polar molecules⁸⁸, plays the most important role, static-induced interactions

⁸⁶See e.g. Barranco et al. (1999), Terasaki et al. (2002b), Barranco et al. (2005), and Idini et al. (2015).

⁸⁷See e.g. Bohr and Mottelson (1975), Eqs. (6-217) and (6-228).

⁸⁸In connection with quantal fluctuations and van der Waals forces, see London (1937).

being less important (App. 3.B). A consequence of this result is related to the size of folding domains of globular proteins⁸⁹.

Concerning superconductivity, the overscreening effect which weakly binds Cooper pairs stems from a delicate ω -dependent interaction which in the case of low temperature superconductors leads, eventually, to one of the first macroscopic manifestations of quantum mechanics, as e.g. “permanent” magnetic fields associated with persistent supercurrents.

The statement “*Life at the edge of chaos*” coined in connection with the study of emergent properties in biological molecules (e.g. protein evolution, folding and stability) reflects the idea, as expressed by de Gennes⁹⁰, that truly important new properties and results can emerge in systems lying at the border between rigid order and randomness, as testified by the marginal stability and conspicuous fluctuations characterizing, for example, nuclear Cooper pairs at the driplines, and that of the marginal stability of proteins of e.g. viral particles like the HIV-1– and HCV–proteases⁹¹.

Let us conclude by quoting again de Gennes but doing so with the hindsight of more than twenty years of nuclear research which have elapsed since “Les objets fragiles” was first published. The chapter entitled “Savoir s’arrêter, savoir changer” starting at p. 180 opens with the statement “En ce moment, la physique nucléaire (la science des noyaux atomiques) est une science qui, à mon avis, se trouve en fin de parcours... C’est une physique qui demande des moyens coûteux, et qui s’est constituée par ailleurs en un puissant lobby. Mais elle me semble naturellement exténuée... je suis tenté de dire: “Arrêtons”... mais ce serait aussi absurde que de vouloir arrêter un train à grande vitesse. Le mieux serait d’aguiller ce train sur une autre voie, plus nouvelle et plus utile à la collectivité.”

In a way, and even without knowing de Gennes remark, part of the nuclear physics community have followed them, capitalizing on the novel embodiment that concepts like elementary modes of excitation, spontaneous symmetry breaking and phase transitions have had in the paradigm of finite quantum many-body (FQMB) system the nucleus represents, where fluctuations can dominate over potential energy effects. The use of these concepts tainted by FQMB system effects as applied to proteins, in particular to the understanding of protein folding may, arguably, shed light on the possibility of designing leads to drugs which are less

⁸⁹It is of notice that similar arguments (cf. Sect. 3.6) are at the basis of the estimate (3.6.13) (see also Sect. 3.7) concerning the size of the halo nucleus ^{11}Li , a quantity which is influenced to a large extent by the maximum distance (correlation length) over which partners of a Cooper pair are virtually (of course only if particle, normal, density allows for it) but solidly anchored to each other (localized), and have to be seen as an extended (quasi) bosonic entity. The fact that Cooper pair transfer proceeds mainly in terms of successive transfer controlled by the single-particle mean field, reinforces the above physical picture of nuclear pairing. Even under the effect of an extremely large, as compared to the pair correlation energy, external single-particle fields, namely that of target and projectile, the Cooper pair field extends over the two nuclei, permeating the whole summed nuclear volume also through tiny (normal) density overlaps.

⁹⁰de Gennes (1994).

⁹¹See e.g. Broglia (2013).

prone to create resistance⁹², let alone all the parallels that one has been able to establish and the associated progress resulting from them concerning cluster and quantum dots physics, a particular example being the discovery of super shells⁹³.

4.8.2 Metals

Plasmons and phonons (jellium model)

The expression of the electron plasmon (ep) frequency of the antenna-like oscillations of the free, conduction electrons of mass m_e and charge $-e$, against the positive charged background (jellium model) is

$$\omega_{ep}^2 = \frac{4\pi n_e e^2}{m_e} = \frac{3e^2}{m_e r_s^3}, \quad (4.8.27)$$

where

$$n_e = \frac{3}{4\pi} \frac{1}{r_s^3}, \quad (4.8.28)$$

are the number of electrons per unit volume⁹⁴, r_s being

$$r_s = \left(\frac{3}{4\pi n_e} \right)^{1/3}, \quad (4.8.29)$$

that is, the radius of the Wigner-Seitz cell.

For⁹⁵ metallic Li

$$n_e = 4.70 \frac{10^{22}}{\text{cm}^3} = \frac{4.7 \times 10^{-2}}{\text{\AA}^3}, \quad (4.8.30)$$

while

$$r_s = \left(\frac{3\text{\AA}^3}{4\pi \times 4.7 \times 10^{-2}} \right)^{1/3} = 1.72\text{\AA}, \quad (4.8.31)$$

implying a value $(r_s/a_0) = 3.25$ in units of Bohr radius ($a_0 = 0.529\text{\AA}$). Making use of

$$\alpha = 7.2973 \times 10^{-3} = \frac{e^2}{\hbar c} \quad (4.8.32)$$

and

$$e^2 = 14.4 \text{ eV \AA}, \quad (4.8.33)$$

one obtains

$$\hbar c = \frac{14.4 \text{ eV \AA}}{7.2973 \times 10^{-3}} = 1973.3 \text{ eV \AA}. \quad (4.8.34)$$

⁹²See e.g. Broglia (2005); Rösner et al. (2017) and refs. therein.

⁹³Pedersen et al. (1991); de Heer et al. (1987); Brack (1993); Pacheco et al. (1991); Lipparini (2003); Martin et al. (1994); Bjornholm et al. (1994).

⁹⁴In the nuclear case one can write $\rho_n = 1/(4\pi r_0^3/3) = 0.14 \text{ fm}^{-3}$ ($r_0 = 1.2 \text{ fm}$), to be compared with the saturation density $\rho_0 = 0.17 \text{ fm}^{-3}$.

⁹⁵cf. page 5, table 1.1 of Ashcroft and Mermin (1987).

With the help of the above values and of

$$m_e c^2 = 0.511 \text{ MeV}, \quad (4.8.35)$$

one can write

$$\hbar^2 \omega_{ep}^2 = \frac{(\hbar c)^2}{m_e c^2} \frac{3e^2}{r_s^3} = \frac{(1973.3 \text{ eV } \text{\AA})}{0.511 \times 10^6 \text{ eV}} \frac{3 \times 14.4 \text{ eV } \text{\AA}}{(1.72 \text{ \AA})^3} = 64.7 \text{ eV}^2 \quad (4.8.36)$$

leading to⁹⁶

$$\hbar \omega_{ep} = 8.04 \text{ eV} \approx 1.94 \times 10^9 \text{ MHz} \quad (4.8.37)$$

For the case of metal clusters of Li, the Mie resonance frequency is⁹⁷

$$\hbar \omega_M = \frac{\hbar \omega_{ep}}{\sqrt{3}} = 4.6 \text{ eV}. \quad (4.8.38)$$

4.8.3 Elementary theory of phonon dispersion relation

Again, within the framework of the jellium model, one can estimate the long wavelength ionic plasma (ip) frequency introducing in (4.8.27) the substitution $e \rightarrow Ze$, $m_e \rightarrow AM$ ($A = N + Z$, mass number, M nucleon mass), $n_e \rightarrow n_i = n_e/Z$,

$$\omega_{ip}^2 = \frac{4\pi n_i (Ze)^2}{AM} = \frac{Zm_e}{AM} \omega_{ep}^2, \quad (4.8.39)$$

$AM(Ze)$ being the mass (charge) of the ions⁹⁸. For metallic Li, one obtains

$$\begin{aligned} \hbar \omega_{ip} &= \left(\frac{Zm_e}{AM} \right)^{1/2} \hbar \omega_{ep} = \left(\frac{3 \times 0.5}{9 \times 10^3} \right)^{1/2} \times 1.94 \times 10^{15} \text{ sec}^{-1} \\ &\approx 2.5 \times 10^{13} \text{ sec}^{-1} \approx 10^{13} \text{ sec}^{-1} \approx 1.04 \times 10^2 \text{ meV}. \end{aligned} \quad (4.8.40)$$

Now, both the relations (4.8.27) and (4.8.39), although being quite useful, are wrong from a many-body point of view: ω_{ep} because electrons appear as bare electrons not dressed by the phonons, neither by the plasmons; ω_{ip} because the static negative background does not allow for an exchange of electron plasmons between ions, exchange eventually leading to a screened, short-range ionic Coulomb repulsive field. Namely ions interact, in the approximation used above, in terms of the “bare” ion–ion Coulomb interaction. Being it infinite range it does not allow for a dispersion relation linear in k at long wavelengths (sound waves) but forces a finite “mass” also to the lattice phonons. Allowing for electron screening of the “bare” ion–ion Coulomb interaction, as embodied in the Thomas-Fermi electron gas dielectric function $1/\epsilon_{TF} = q^2/(q^2 + k_s^2)$, one obtains the dressed phonon frequency

$$\omega_q^2 = \frac{\omega_{ip}^2}{\epsilon_{TF}} = \frac{Zm_e}{AM} \frac{\omega_{ep}^2}{q^2 + k_s^2} q^2 = \frac{\omega_{ip}^2}{q^2 + k_s^2} q^2. \quad (4.8.41)$$

⁹⁶Kittel (1996) Table 2, p. 278.

⁹⁷See Bertsch and Broglia (2005), Sect. 5.1, also Table 5.1.

⁹⁸Ketterson and Song (1999), p. 230.

The quantity

$$k_s = \left(\frac{6\pi n_e e^2}{\epsilon_F} \right)^{1/2} = \left(\frac{4k_F}{\pi a_0} \right)^{1/2} = 0.82 k_F \left(\frac{r_s}{a_0} \right)^{1/2}, \quad (4.8.42)$$

is the Thomas-Fermi screening wave vector, a quantity which is of the order of the Fermi momentum, the associated screening length being then of the order of the Wigner-Seitz radius. In writing the above relations use has been made of

$$k_F = \left(\frac{9\pi}{4} \right)^{1/3} \frac{1}{r_s} = \frac{1.92}{r_s}, \quad (4.8.43)$$

and

$$\epsilon_F = \frac{e^2 a_0}{2} k_F^2 = \frac{50.1}{(r_s/a_0)^2}. \quad (4.8.44)$$

In the case of metallic Li ($k_F = 1.12 \text{ \AA}$, $r_s = 1.72 \text{ \AA}$),

$$k_s = 1.6 \text{ \AA}^{-1}. \quad (4.8.45)$$

and

$$k_F = 0.12 \text{ \AA}^{-1}, \quad \epsilon_F = 4.74 \text{ eV} \quad (r_s/a_0 = 3.25). \quad (4.8.46)$$

Let us return to (4.8.41), and take the long wavelength limit of ω_q . One finds

$$\lim_{q \rightarrow 0} \omega_q = c_s q \quad (4.8.47)$$

the sound velocity (squared) being

$$c_s^2 = \frac{Zm_e}{AM} \frac{4\pi n_e e^2}{m_e} \frac{\epsilon_F}{6\pi n_e e^2} = \frac{2Z}{3AM} \epsilon_F = \frac{Zm_e}{3AM} v_F^2, \quad (4.8.48)$$

With the help of (4.8.43), of the velocity of light and of the proton mass,

$$c = 3 \times 10^{10} \text{ cm/sec}; \quad Mc^2 = 938.87 \text{ MeV}, \quad (4.8.49)$$

one obtains,

$$\begin{aligned} v_F &= \left(\frac{\hbar}{m_e} \right) k_F = \left(\frac{\hbar c}{m_e c^2} \right) \times 3 \times 10^{10} \frac{\text{cm}}{\text{sec}} \frac{1.92}{r_s} \\ &= \left(\frac{1973.3 \text{ \AA eV}}{0.511 \times 10^6 \text{ eV}} \right) \times 3 \times 10^{10} \frac{\text{cm}}{\text{sec}} \frac{1.92}{1.72 \text{ \AA}} \approx 1.29 \times 10^8 \frac{\text{cm}}{\text{sec}}, \end{aligned} \quad (4.8.50)$$

and,

$$c_s^2 = \frac{1}{3} \frac{3m_e}{9M} v_F^2 \approx 6 \times 10^{-5} v_F^2, \quad (4.8.51)$$

that is,

$$c_s \approx 7.8 \times 10^{-3} v_F \approx 1.0 \times 10^6 \frac{\text{cm}}{\text{sec}}, \quad (4.8.52)$$

approximately equal to a hundredth of the Fermi velocity⁹⁹.

Let us now discuss the effective electron-electron interaction. Within the jellium model used above one can write it as

$$V(\mathbf{q}, \omega) = \frac{U_c(q)}{\epsilon(\mathbf{q}, \omega)}, \quad (4.8.53)$$

where the dielectric function

$$\frac{1}{\epsilon(\mathbf{q}, \omega)} = \frac{\omega^2 q^2}{\omega^2(q^2 + k_s^2) - \omega_{ip}^2 q^2} \quad (4.8.54)$$

contains the effects due to both the ions and the background electrons, while

$$U_c(q) = \frac{4\pi e^2}{q^2} \quad (4.8.55)$$

is the Fourier transform of the bare Coulomb interaction¹⁰⁰

$$U_c(r) = \frac{e^2}{r}. \quad (4.8.56)$$

For $\omega \gg \omega_{ip}$, $\epsilon(\mathbf{q}, \omega)^{-1} = q^2/(q^2 + k_s^2) = \epsilon_{TF}^{-1}$, and one obtains the screened Coulomb field,

$$U_c^{scr}(q) = \frac{U_c(q)}{\epsilon_{TF}(q)} = \frac{4\pi e^2}{q^2 + k_s^2}. \quad (4.8.57)$$

Its \mathbf{r} space Fourier transform is

$$U_c^{scr}(r) = \frac{e^2}{r} e^{-k_s r}. \quad (4.8.58)$$

A quantity that for large values of r falls off exponentially. Thus, in the high frequency limit, the electron-electron interaction, although strongly renormalized by the exchange of plasmons, as testified by the fact that (e.g. for Li),

$$U_c^{scr}(r = 5 \text{ \AA}) \approx U_c(r = 5 \text{ \AA}) e^{-1.6 \times 5} \approx 1 \text{ meV}, \quad (4.8.59)$$

as compared to $U_c(r = 5 \text{ \AA}) \approx 2.9 \text{ eV}$, is still repulsive.

Let us now consider frequencies $\omega \ll \omega_{ip}$ but for values of q of the order of a^{-1} , where a is the lattice constant ($a \approx 3 - 5 \text{ \AA}$, $a^{-1} \approx 0.25 \text{ \AA}^{-1}$) to be compared to $k_s \approx 1.6 \text{ \AA}^{-1}$ and $k_F \approx 1.12 \text{ \AA}^{-1}$ (metallic Li). In the case in which $\omega_{ip}^2/\omega^2 > (q^2 + k_s^2)/q^2$, V is attractive. This behavior explicitly involves the ions through ω_{ip} (electron-phonon coupling).

⁹⁹Ashcroft and Mermin (1987), p. 51, Ketterson and Song (1999) p. 234.

¹⁰⁰The units being $[U_c(r)] = \text{eV}$, $[U_c(q)] = \text{eV} \times \text{\AA}^3$.

The dispersion relation of the associated frequency collective modes follows from

$$\epsilon(\mathbf{q}, \omega) = 0. \quad (4.8.60)$$

In other words, making use of Eq. (4.8.54) one obtains the relation (4.8.41). One can now rewrite the reciprocal of the dielectric functions in terms of ω_q , that is,

$$\begin{aligned} \frac{1}{\epsilon(\mathbf{q}, \omega)} &= \frac{\omega^2 q^2}{\omega^2(q^2 + k_s^2) - \omega_{ip}^2 q^2} = \frac{q^2}{(q^2 + k_s^2)} \frac{\omega^2}{(\omega^2 - \omega_q^2)} \\ &= \frac{q^2}{q^2 + k_s^2} \left[1 + \frac{\omega_q^2}{\omega^2 - \omega_q^2} \right]. \end{aligned} \quad (4.8.61)$$

For $\omega \gg \omega_q$ one recovers the Thomas-Fermi dielectric function. For ω near, but smaller than ω_q the interaction eventually becomes attractive¹⁰¹. The effective electron-electron interaction can be then written as

$$\begin{aligned} V(\mathbf{q}, \omega) &= \frac{4\pi e^2}{q^2 + k_s^2} + \frac{4\pi e^2}{q^2 + k_s^2} \frac{\omega_q^2}{\omega^2 - \omega_q^2} \\ &= U_c^{scr}(q) + U_c^{scr}(q) \frac{\omega_q^2}{\omega^2 - \omega_q^2} \\ &= U_c^{scr}(q) (1 + U_c^{scr}(q) \Pi(\mathbf{q}, \omega)), \end{aligned} \quad (4.8.62)$$

where¹⁰²

$$\Pi(\mathbf{q}, \omega) = \left(\frac{Z}{AM} \right) \frac{n_e q^2}{\omega^2 - \omega_q^2}. \quad (4.8.63)$$

The units of $\Pi(\mathbf{q}, \omega)$ are $eV^{-1}\text{\AA}^{-3}$, in keeping with those of $U_c^{scr}(q)$.

In working out the last expression Eqs. (4.8.41) and (4.8.27) have been used. In other words $\omega_q^2 = (Zm_e/AM)(\omega_{ep}^2/(q^2 + k_s^2))q^2 = (Z/AM)(4\pi n_e e^2/(q^2 + k_s^2))q^2$. It is of notice the close relation of (4.8.62) with the expression (4.8.19) of the nuclear renormalized pairing interaction. The first term of $V(\mathbf{q}, \omega)$ contains the screened Coulomb field arising ultimately from the exchange of plasmons between electrons (see Fig. 4.8.4). The second term with the exchange of collective low frequency phonons calculated making use of the same screened interaction¹⁰³.

Concerning the parallels discussed above between pairing in nuclei and in metals, it is also important to point out the important differences. At least concerning the possibility of developing a unified theoretical working tool.

In metals the bare interaction emerges from the exchange of photons, a process which can be described at profit in terms of an instantaneous Coulomb interaction.

¹⁰¹Schrieffer (1964), Fig. 6-11, p. 152.

¹⁰²It is of notice that $\omega_q^2 = (Zm_e/AM)\omega_{ep}^2 \times q^2/(q^2 + k_s^2) = (Z/AM)4\pi e^2 n_e q^2/(q^2 + k_s^2) = U_c^{scr}(q) \times (Z/AM)n_e q^2$.

¹⁰³Lindhard (no. 8, 1954).

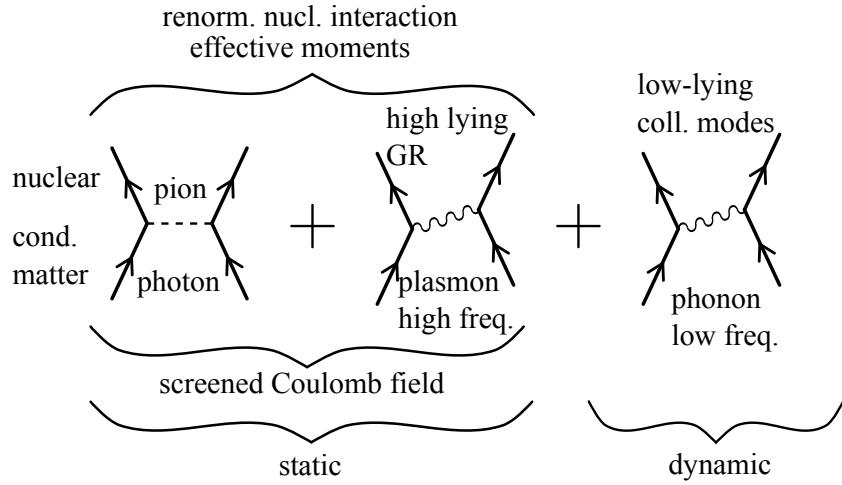


Figure 4.8.4: Schematic representation of the variety of contributions to the effective interaction in nuclei and in metals.

While major progress have been made concerning the bare nucleon interaction at large, and the NN -pairing interaction in particular, taking also into account three-body processes and carrying out ab initio calculations¹⁰⁴, we are not yet in possess of the equivalent to the Coulomb interaction. Let alone of a so called low- k version of such interaction, which could allow to work out on equal footing Hartree-Fock-Bogoliubov solutions, and QRPA microscopic calculations of low-lying collective modes in the variety of channels (density, spin, charge-exchange, etc.).

Concerning these phonons, they are built out of the same nucleon degrees of freedom which already exhaust the nuclear phase space. Double counting and Pauli principle violations have to be taken care of in using the collective modes as intermediate bosons. In metals, phonons are associated with lattice vibrations, that is degrees of freedom different from the electronic ones.

Nonetheless, in nuclei one can describe in terms of individual quantal states and of single Cooper pairs, if not identical, rather similar phenomena as those leading to some of the most remarkable and technically transferable quantum phenomena. In particular, the Josephson effect. This fact makes the nuclear pairing paradigm a unique laboratory of low-temperature (finite) many-body physics.

We conclude this section with a technical remark. Making use of (4.8.62) we calculate the mass enhancement factor¹⁰⁵

$$\lambda = N(0)V(\mathbf{q}, \omega) = N(0)U_c^{scr} (1 + U_c^{scr}\Pi). \quad (4.8.64)$$

¹⁰⁴See footnote 10 Ch. 1.

¹⁰⁵It is assumed that in this calculation we have multiplied (4.8.62) by n_e , as well as shifted the corresponding factor appearing in $\Pi(\mathbf{q}, \omega)$ to $U_c^{scr}(q)$. Schematically, it is assumed $n_e V = (n_e U_c^{scr}) \left(1 + (n_e U_c^{scr}) \left(\frac{Z}{AM} \right) \frac{q^2}{\omega^2 - \omega_q^2} \right)$.

In the weak coupling limit ($\lambda^2 \ll \lambda$)

$$\Delta = 2\omega_D e^{-1/\lambda}, \quad (4.8.65)$$

where ω_D is the Debye energy. Provided that one considers a situation in which ω is consistently different from ω_q ,

$$\frac{1}{\lambda} = \frac{1}{N(0)U_c^{scr}(1 + U_c^{scr}\Pi)} \approx \frac{1}{N(0)U_c^{scr}} (1 - U_c^{scr}\Pi), \quad (4.8.66)$$

Thus

$$\frac{1}{\lambda} = \frac{1}{N(0)U_c^{scr}} - \frac{\Pi}{N(0)}, \quad (4.8.67)$$

and

$$\Delta = \left(2\omega_D e^{\frac{\Pi}{N(0)}} \right) e^{-\frac{1}{N(0)U_c^{scr}}}. \quad (4.8.68)$$

Consequently, the renormalization effects of the pairing gap associated with phonon exchange are independent of the approximation used to calculate U_c^{scr} (Thomas–Fermi in the above discussion), provided one has used the same “bare” (screened) Coulomb interaction to calculate ω_q^2 . Otherwise, the error introduced through a resonant renormalization process entering the expression of e.g. the pairing gap, may be large.

4.8.4 Pairing condensation (correlation) energy

The condensation energy, namely the energy difference $W_N - W_S$ between the normal N – and superfluid S –state is defined as (Eq. (2-35) of ref¹⁰⁶)

$$W_{con} = W_N - W_S = \frac{1}{2} N(0) \Delta_0^2, \quad (4.8.69)$$

where $N(0)$ is the density of single-electron states of one-spin orientation evaluated at the Fermi surface, and Δ_0 is the pairing gap at $T = 0$.

The correlation energy E_{corr} introduced in equation (6-618) of ¹⁰⁷

$$E_{corr} = -\frac{1}{2d} \Delta^2 \quad (4.8.70)$$

to represent $W_S - W_N$ in the nuclear case, was calculated making use of a (single particle) spectrum of two-fold degenerate (Kramer degeneracy) equally spaced (spacing d) single-particle levels. Consequently, $2/d$ corresponds to the total level density, and $1/d = N(0)$. In keeping with the fact that a nucleus in the ground state (or in any single quantal state), is at zero temperature, Eq. (4.8.70) coincides with (4.8.69), taking into account the difference in sign in the definitions.

System	Δ_0		$N(0)$		W_{con}		E_{cohe}	BE/A	$\frac{W_{con}}{E_{cohe}}$	$\frac{W_{con}}{BE}$
	meV	MeV	$\frac{\text{meV}^{-1}}{\text{atom}}$	MeV^{-1}	$\frac{\text{meV}}{\text{atom}}$	MeV	$\frac{\text{meV}}{\text{atom}}$	$\frac{\text{MeV}}{A}$	10^{-7}	10^{-3}
Pb	^{120}Sn	1.4	1.46	0.276	4	3×10^{-4}	4.0	2030	8.5	

Table 4.8.1: Summary of the quantities entering the calculation of the condensation energy of superconducting lead (Pb), and of the single open shell superfluid nucleus ^{120}Sn .

Nuclei

The empirical value of the level density parameter for both states ($\nu, \bar{\nu}$) (Kramers degeneracy, both spin orientations) is $a \approx A/10 \text{ MeV}^{-1}$, $A = N + Z$ being the mass number. Thus, for neutrons one can write $a_N \approx N/10$ and $N_N(0) \approx N/20 \text{ MeV}^{-1}$. For $^{120}_{50}\text{Sn}$, $N_N(0) \approx 4 \text{ MeV}^{-1}$. Because $\Delta = 1.46 \text{ MeV}$, (Table 4.8.1)

$$W_{con} = \frac{1}{2} \times 4 \text{ MeV}^{-1} \times (1.46)^2 \text{ MeV}^2 \approx 4 \text{ MeV}. \quad (4.8.71)$$

The binding energy per nucleon is $BE/A = 8.504 \text{ MeV}$. Thus $BE = 120 \times 8.504 \text{ MeV} = 1.02 \times 10^3 \text{ MeV}$, and

$$\frac{W_{con}}{BE} \approx 4 \times 10^{-3}. \quad (4.8.72)$$

Superconducting lead

Making use of the value¹⁰⁸

$$N(0) = \frac{0.276 \text{ eV}^{-1}}{\text{atom}}, \quad (4.8.73)$$

and of $\Delta_0 = 1.4 \text{ meV}$, one obtains

$$W_{con} = 0.27 \times 10^{-6} \text{ eV/atom}. \quad (4.8.74)$$

In keeping with the fact that the cohesive energy of lead, namely the energy required to break all the bonds associated with one of its atoms is

$$E_{cohe} = 2.03 \frac{\text{eV}}{\text{atom}}, \quad (4.8.75)$$

one obtains

$$\frac{W_{con}}{E_{cohe}} \approx 1.3 \times 10^{-7}. \quad (4.8.76)$$

The different quantities are summarized in Table 4.8.1.

¹⁰⁶Schrieffer (1964).

¹⁰⁷Bohr and Mottelson (1975).

¹⁰⁸Beck and Claus (1970).

4.9 Cooper pair: radial dependence

The wavefunction of a Cooper pair can be written as

$$\Psi(\mathbf{r}_1\sigma_1; \mathbf{r}_2\sigma_2) = \phi_q(\mathbf{r})e^{i\mathbf{q}\cdot\mathbf{R}}\chi(\sigma_1, \sigma_2) \quad (4.9.1)$$

where $\mathbf{R} = (\mathbf{r}_1 + \mathbf{r}_2)/2$, $\mathbf{r} = \mathbf{r}_1 - \mathbf{r}_2$, and σ_1 and σ_2 denote the spins¹⁰⁹.

Let us consider the state with zero center of mass momentum ($q = 0$) and with zero spin, so that the two electrons carry equal and opposite momenta, aside of being in the singlet spin state state, with

$$\chi = \frac{1}{\sqrt{2}} \left[\begin{pmatrix} 1 \\ 0 \end{pmatrix} \begin{pmatrix} 0 \\ 1 \end{pmatrix} - \begin{pmatrix} 0 \\ 1 \end{pmatrix} \begin{pmatrix} 1 \\ 0 \end{pmatrix} \right]. \quad (4.9.2)$$

We have thus a pair of electrons moving in time reversal states and can write¹¹⁰,

$$\phi_0(\mathbf{r}) = \sum_{k>k_F} g(\mathbf{k}) e^{i\mathbf{k}\cdot\mathbf{r}} = \sum_{k>k_F} g(\mathbf{k}) e^{i\mathbf{k}\cdot\mathbf{r}_1} e^{-i\mathbf{k}\cdot\mathbf{r}_2}. \quad (4.9.3)$$

In the above wavefunction Pauli principle ($k > k_F$) and translational invariance (dependence on the relative coordinate \mathbf{r}) are apparent. Within this scenario, and in keeping with the fact that $|v_p^{eff}|/\epsilon_F \ll 1$, where $v_p^{eff} = v_p^{bare} + v_p^{ind}$, $\phi_0(\mathbf{r})$ consists mainly of waves of wavenumber k_F . Because the wavefunction of a Cooper pair represents a bound s -state, the motion it describes is a periodic back and forth movement of the two electrons in directions which are uniformly distributed, covering a relative distance $\approx \xi$, as schematically¹¹¹ shown in Fig. 4.9.1 (i). It is analogous to the motion of the two nucleon in a deuteron or the main ($L = 0$) component of the two neutrons in the triton. The hydrogen atom in s -state is also an example; in that case it is the electron that does most of the back and forth moving, whereas the proton only recoils slightly.

In keeping with the above arguments, $\phi_0(\mathbf{r})$ will look like $e^{i\mathbf{k}_F\cdot\mathbf{r}}$ for $r \ll \xi$, while for $r \gg \xi$ the waves $e^{i\mathbf{k}\cdot\mathbf{r}}$ weighted by $g(\mathbf{k})$ will destroy themselves by interference (Fig. 4.9.2). In other words, $\phi_0(\mathbf{r})$ will look like $e^{i\mathbf{k}_F\cdot\mathbf{r}}$ for $r \ll \xi$ while for $r \gtrsim \xi$ one can approximate the weighting function as,

$$g(\mathbf{k}) \sim \delta(\mathbf{k}, \mathbf{k}_F + i\hat{\mathbf{k}}_F/\xi), \quad (4.9.4)$$

where $\hat{\mathbf{k}}_F$ is a unit vector. One then obtains (see also Eq. (4.3.2)),

$$\phi_0(\mathbf{r}) \sim e^{-r/\xi} e^{ik_F r}. \quad (4.9.5)$$

¹⁰⁹In the limit $q \rightarrow 0$ the relative coordinate problem is spherically symmetric so that $\phi_0(\mathbf{r})$ is an eigenfunction of the angular momentum operator (Schrieffer (1964)).

¹¹⁰In other words, one expands the $l = 0$ wavefunction ϕ_0 in terms of s -states of relative momentum k and total momentum zero.

¹¹¹Weisskopf (1981).

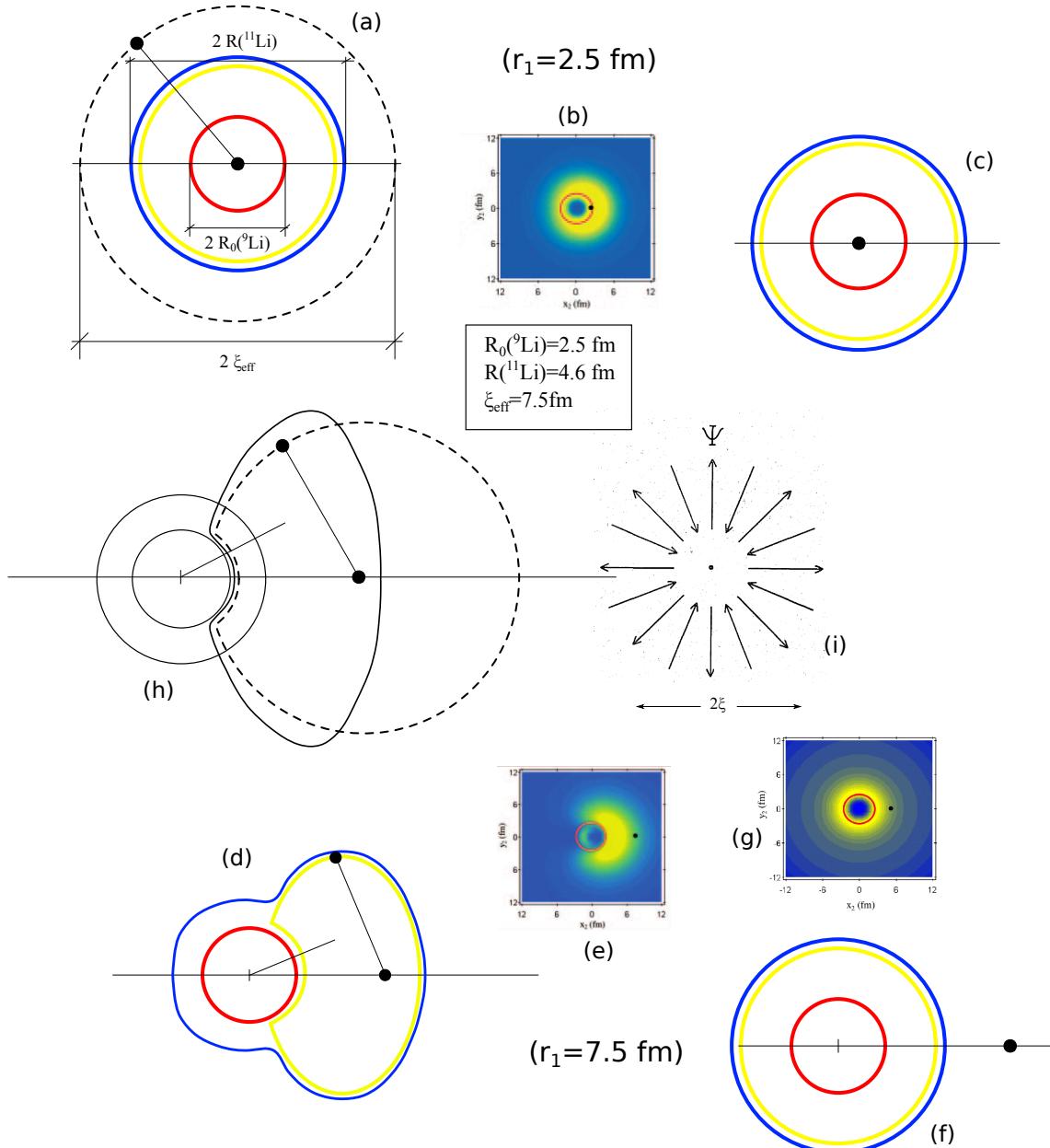


Figure 4.9.1: Aspects of the halo Cooper pair of the nucleus ^{11}Li .

Because we are dealing with a singlet state, and the total wavefunction has to be antisymmetric,

$$\phi_0(\mathbf{r}) \sim e^{-r/\xi} \cos k_F r, \quad (4.9.6)$$

A more proper solution of the Cooper pair problem leads to¹¹²

$$\phi_0(\mathbf{r}) \sim K_0(r/\pi\xi) \cos k_F r, \quad (4.9.7)$$

where K_0 is the zeroth-order modified Bessel function. For $x \gg 0$, $K_0(x) \sim (\pi/2x)^{1/2} \exp(-x)$, where $x = r/\pi\xi$.

A wavefunction which extends over distances much larger than the range of the binding potential is a well-known phenomenon when the binding energy is small. For example, in the case of the deuteron mentioned above. Be as it may, in the case of metallic superconductors the large size of the Cooper pair wavefunction also explains why the electrostatic repulsion between electron partners does not appreciably influence the binding.

Going back to Fig. 4.9.1, it is illustrative to elaborate on the two (NFT calculated) situations displayed in (b) and (e), concerning the relative distribution of the halo neutrons of ^{11}Li . Pairing correlations being, in particular in this case, mainly a surface phenomenon bring, for $r_1 = 7.5$ fm, the two nucleons close to each other as compared to the uncorrelated situation (diagram (g)). The fact that this result, which is not under discussion, is more subtle than just expressed, emerges by looking at (h), where the situations displayed in (b) and (e) (see also (a) and (d)), are schematically drawn in a single plot corresponding to: I) the (“free from external fields”) BCS-like, Cooper pair (dashed curve) which implies that the two neutrons recede from each other; II) the (“actual”), confined, situation in which the average potential $U(r) = \int d\mathbf{r}' \rho(\mathbf{r}') v(|\mathbf{r} - \mathbf{r}'|)$, acting as a strong external field which determines where normal ($\rho(r)$), and thus abnormal density can find themselves, distorts the halo Cooper pair, leading to the situation represented with the continuous (irregular) curve. Thus, to a situation in which the two halo neutrons have come closer to each other as compared to the uncorrelated configuration¹¹³.

In principle, both situations can be experimentally observed. That labelled II) in terms of high energy electron scattering or (p, pn) reactions, while that labelled I), making use of a two-particle pick up reaction ($^1\text{H}(^{11}\text{Li}, ^9\text{Li})^3\text{H}$). In fact, during the period of time the proton is around the grazing distance, and one of the halo neutrons joins it to form a virtual deuteron, the other neutron can be at the antipodes, essentially one diameter apart. The situation will essentially evolve into the density distribution represented with the dashed curve in (h), the two members of the halo Cooper pair being at a distance of the order of the correlation length and carrying a much lower relative momentum than that typical of the situation (e)

¹¹²Kadin (2007); see also van Witsen (2014).

¹¹³And to do so, the finite quantal system under discussion uses the mechanism discussed in Sect. 4.11 namely, spatial quantization.

(continuous irregular curve). It is then natural that successive transfer dominates the absolute differential cross section.

While it is true that both I) and II) can be the outcome of experiments, the specific probe of dynamic or static distortion in gauge space, and thus of Cooper pair structure is Cooper pair transfer.

Specific probe

From daily life experience one can state that, as a rule, macroscopic, condensed systems find themselves in the presence of external fields which fix the *order parameter* at some definite preferred value. Because of the long-range order, a very weak external field can pin down the *orientation* (violation of rotational invariance) and *position* (violation of translational invariance) of a crystal, or of a chair¹¹⁴. To determine the value of these order parameter values, one needs to have instruments, probes, experimental setups, or whatever one likes to call them, which violate themselves rotational and translational invariance. Again, daily life objects (instruments) do so.

In the case of superconductivity, the order parameter is α_0 , α'_0 being its magnitude and ϕ its phase. It is gauge invariance which is spontaneously violated, the (BCS) superconductor displaying a rather perfect internal gauge phase order. There are not many instruments which have such a property but another superconductor acting as a probe, weakly coupled¹¹⁵ to the probed superconductor. Said it differently, a Josephson junction and associated special effects. In particular direct currents of carriers of charge $2e$ and mass $2m_e$.

Returning to the nuclear scenario, pair transfer, eventually back and forth¹¹⁶ pair transfer, of nucleons correlated in time reversal states is the specific probe of pairing in atomic nuclei in general, and of halo nuclear Cooper pairs in particular.

To interact at profit through long wavelength medium polarization pairing, pairs of nucleons have to have low momentum. To do so they have to reduce the

¹¹⁴Quoting from Weinberg (1996b) Ch. 19: "...spontaneous symmetry breaking actually occurs only for idealized systems that are infinitely large. The appearance of broken symmetry for a chair arises because it has a macroscopic moment of inertia I , so that its ground state is part of a tower of rotationally excited states whose energies are separated by only tiny amounts, of the order of \hbar^2/I ... even very weak... rotationally asymmetric external fields will cause any... state of the chair with definite angular momentum rapidly to develop components with other angular momentum quantum numbers. States of the chair that are relatively stable... are... those with a definite orientation, in which the rotational symmetry of the underlying theory is broken". In connection with superconductivity, distortion in gauge space is measured by the number of Cooper pairs $\alpha'_0 (\approx 10^6)$ around the Fermi energy, the associated number of electrons taking the place of angular momentum in 3-D space. Now, is it 2×10^6 an infinitely large number of particles? And if yes, what about 12–16 nucleons ($\alpha'_0 \approx 6\text{--}8$ for Sn-isotopes)? An answer which can hardly be other than positive, in keeping with the observation of well defined pairing rotational bands, populated in two-nucleon transfer reactions. But if this is so, implying "angular momentum" fluctuations as large as $\approx 30\%$, why 2 nucleons ($\alpha'_0 = 1$, ^{11}Li) with associated fluctuations of $\approx 70\%$ can not display, in an incipient fashion, basic features of BCS pairing?

¹¹⁵Linear response. See footnote 1 Ch. 2.

¹¹⁶??.

effect of the strong external (mean) field by moving away from it, possible mechanisms being among others: halo (Fig. 4.9.1), transfer processes (see e.g. Fig. 4.5.1), exotic decay¹¹⁷ (see Fig. 4.9.3) and, if nothing else, some amount of spill out.

Caption Fig 4.9.1

Synthesis of the spatial structure of ^{11}Li neutron halo Cooper pair calculated in NFT (Barranco et al. (2001)). We use, for illustrative purposes $\xi_{\text{eff}} = 7.5$ fm, as an *effective Cooper pair mean square radius*¹¹⁸ (dashed circle in (a) and (h)). Diagrams (a) and (d) are the schematic representations of the modulus square $|\Psi_0(\mathbf{r}_1, \mathbf{r}_2)|^2 = |\langle \mathbf{r}_1, \mathbf{r}_2 | 0 \rangle|^2$ describing the motion of the two halo neutrons of ^{11}Li , moving around the ^9Li core as a function of the cartesian coordinates of neutron 2, for fixed (small and large as compared to $R(^{11}\text{Li}) \approx 4.6$ fm) values of the position r_1 of neutron 1 (for more details see Caption to Fig. 3.6.3). Diagrams (b), (e) and (g) are the results of NFT (see also Fig. 3.6.3 (II) a and b)). **(a)** The circles drawn with continuous lines correspond to the relative distance r at the radius of the ^9Li core and of ^{11}Li . The Cooper pair “intrinsic coordinate” r_{12} is also shown. Particle 1 of the Cooper pair is assumed to occupy the center of the nucleus ($r_1 = 0$). **(b)** Result of NFT for a situation similar to the above. **(c)** Schematic representation of an uncorrelated pair in a potential weakly binding the pure configuration $p_{1/2}^2(0)(r_1 = 0)$. **(d)** Same as (a) but for $r_1 = 7.5$ fm (note however that the dashed curve associated with ξ_{eff} is not displayed here). **(e)** Result of the NFT calculation for a similar setup ($r_1 = 7.5$ fm). **(f)** Schematic representation of a pure configuration $p_{1/2}^2(0)(r_1 = 7.5$ fm). **(g)** The result of the microscopic calculation for a weakly bound $p_{1/2}^2(0)$ configuration ($r_1 = 5$ fm). **(h)** The variety of situations found in (a) and (d) in comparison to each other in a single cartoon. **(i)** Schematic picture of the dynamics of the partners fermions in the quantum state ϕ_0 of the Cooper pair. It is a linear combination of motions away and towards one another. The fermions stay within a distance of the order ξ , root mean square radius of the Cooper pair (see Weisskopf (1981), Kadin (2007) and van Witsen (2014)).

¹¹⁷In Fig. 4.9.3, a parallel is made between correlation lengths associated with (pairing) particle-particle or hole-hole modes and particle-hole vibrations. These last modes also display a consistent spatial correlation (see e.g. Broglia et al. (1971c)).

¹¹⁸In connection with this figure we have estimated the Fermi momentum associated with the two halo neutrons of ^{11}Li as $k_F = (3\pi^2 \times 2 / ((4\pi/3)(4.6)^3 - (2.5)^3))^{1/3} \text{ fm}^{-1} \approx 0.56 \text{ fm}^{-1}$, the denominator being the volume associated with the halo. Thus $(v_F/c) \approx 0.2(k_F)_{\text{fm}^{-1}} \approx 0.1$ and $\xi = \hbar v_F / (\pi \times 0.5 \text{ MeV}) \approx 14 \text{ fm}$ (see also Eq. (4.10.5) as well as the caption to Fig. 4.9.3; see also App. 7.C).

It is of notice the uncertainties which can arise in connection with the estimate of k_F (and thus of v_F/c) when using a model like the Fermi gas model (see Eq. (7.C.6)) in the case of a nucleus which, most of it is surface, as compared with estimates based on the halo neutron shell (variety of volumes associated with core and halo nucleons, Eqs. (4.3.6)–(4.3.9)); see also footnotes 19 and 20 of this Chapter, as well as Eq. (4.10.5).

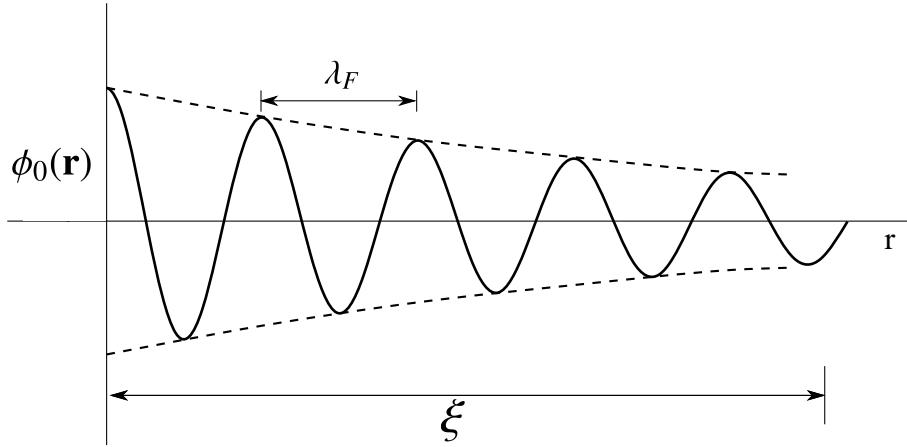


Figure 4.9.2: Schematic representation of the Cooper pair wavefunction. Indicated are the coherence length ξ and the Fermi wavelength $\lambda_F = h/p_F = 2\pi/k_F$. In the nuclear case, and for nuclei along the stability valley $\lambda_F \approx 4.6$ fm and $\xi \approx \hbar v_F/\pi\Delta \approx 14$ fm ($k_F \approx 1.36$ fm $^{-1}$, $v_F/c \approx 0.27$, $\Delta \approx 1.2$ MeV). Thus $\xi/\lambda_F \approx 3$.

4.9.1 Number of overlapping pairs

The correlation length for low temperature superconductors is of the order of 10^4 Å. In the case of e.g. bulk Pb, for which¹¹⁹ $\Delta = 1.4$ meV and $v_F = 1.83 \times 10^8$ cm/s one has $\xi \approx 0.3 \times 10^{-4}$ cm.

Since electrons in metals typically occupy a volume of the order of $(1 \times \text{Å})^3$ (Wigner-Seitz cell radius), there would be of the order of¹²⁰ $\xi^3/(\text{Å})^3 \approx 3 \times 10^{10}$ other Cooper pairs within a “coherence volume”. Eliminating the electrons deep within the Fermi sea as they behave essentially as if the metal was in the normal phase¹²¹, one gets¹²² $\approx 10^6$. *In other words, about a million of other Cooper pairs have their center of mass falling inside the coherence volume of a pair.* Thus, the isolated pair picture (Fig. 3.3.3) is not correct, but yes that displayed in Fig. 3.3.4.

In the nuclear case, the number of Cooper pairs participating in the condensate,

¹¹⁹ The experimental value of the critical temperature and of the pairing gap for Pb is $T_c = 7.193$ K ($k_B T_c = 0.62$ meV) and $\Delta_0(0) = 1.5$ meV respectively. Then $\Delta_0(0)/k_B T_c = 2.26$, to be compared with the BCS prediction of 1.76. Within this context, Pb is considered a strong-coupling superconductor (see Tinkham (1996), Fig. 2-1).

¹²⁰ Ketterson and Song (1999) p. 198.

¹²¹ The BCS ground state at $T = 0$ consists in two classes of electrons: those deep inside the Fermi sea, which behave essentially in the same way as those in the normal state, and those near the Fermi surface, which form the overlapping Cooper pairs. The latter electrons cannot scatter because they are in a coherent state, their intrinsic motion being frozen because of the pairing gap. The former electrons cannot either because blocked by Pauli principle. At $T = 0$, all electrons of both classes contribute to the lossless supercurrent (Waldrum (1996)).

¹²² Schrieffer (1964) p. 43. That is $3 \times 10^{10} \times (\Delta/\epsilon_F) \approx 3 \times 10^{10} \times (1.4 \text{ meV}/9.47 \text{ eV}) \approx 4 \times 10^6$.

namely

$$a'_0 = \langle BCS | P'^{\dagger} | BCS \rangle = \sum_j \frac{2j+1}{2} U'_j V'_j, \quad (4.9.8)$$

can be estimated with the help of the single j -shell model, in which case $V_j = (N/2\Omega)^{1/2}$ and $U_j = (1 - N/2\Omega)^{1/2}$, where $\Omega = (2j+1)/2$. For a half-filled shell ($N = \Omega$), $a'_0 = \Omega/2$. With the help of the approximate expression¹²³ $\Omega \approx (2/3)A^{2/3}$ one obtains, for ^{120}Sn ($N = 70$), $a'_0 \approx 6-8$.

In keeping with the fact that $\xi > R_0$, in the nuclear case one has a complete overlap between all Cooper pairs participating in the condensate. This, together with the fact that the nuclear Cooper pairs press against the nuclear surface in an attempt to expand and are forced to bounce elastically off from it, receive circumstantial evidence from the following experimental results:

1) while the moment of inertia of rotational bands¹²⁴ is $\mathcal{J}_r/2$ it is also $5\mathcal{J}_{irrot}$. In other words, while pairing in nuclei is important, its role is only partially exhausted, and certainly strongly distorted by the mean field;

2) One- and two-nucleon transfer reactions in pairing correlated nuclei have the same order of magnitude. For example $\sigma(^{120}\text{Sn}(p, d)^{119}\text{Sn}(5/2^+; 1.09 \text{ MeV})) = 5.35 \text{ mb}$ ($2^\circ < \theta_{cm} < 55^\circ$), while $\sigma(^{120}\text{Sn}(p, t)^{118}\text{Sn}(gs)) = 2.25 \text{ mb}$ ($7.6^\circ < \theta_{cm} < 59.7^\circ$) (see also footnote 26 of Ch. 3). Furthermore, in the Josephson-like junction, transiently established in the heavy ion collision $^{116}\text{Sn} + ^{60}\text{Ni}$ at bombarding energies around, but still few MeV below the Coulomb barrier ($\tau_{coll} \approx 10^{-21} \text{ s}$), one finds for distances of closest approach $D_0 \leq \xi \approx 13.5 \text{ fm}$, that σ_{2n} is of the order of σ_{1n} . However, while the Cooper pair channel $^{116}\text{Sn} + ^{60}\text{Ni} \rightarrow ^{114}\text{Sn} + ^{62}\text{Ni}$ is dominated by the ground-ground state transition, the single-particle transfer one is inclusive. In fact, the theoretical calculations¹²⁵ of the differential cross section associated with the channel $^{116}\text{Sn} + ^{60}\text{Ni} \rightarrow ^{115}\text{Sn} + ^{61}\text{Ni}$ indicates the incoherent contribution of a number of quasiparticle states of ^{61}Ni lying at energies $\lesssim 2.64 \text{ MeV}$. A value which is consistent with twice the pairing gap of Ni. In other words, we are in presence of superconducting-quasiparticle (S-Q) tunneling across the junction, as discussed in connection with Eq. (4.6.20) (see also Sect. 7.3).

3) The decay constant of the exotic decay $^{223}_{88}\text{Ra}_{135} \rightarrow ^{14}_6\text{C}_8 + ^{209}_{82}\text{Pb}_{127}$ has been measured to be $\lambda_{exp} = 4.3 \times 10^{-16} \text{ sec}^{-1}$. For theoretical purposes it can be written as $\lambda = PfT$, product of the formation probability P of ^{14}C in ^{223}Ra (saddle configuration, see bottom Fig. 4.9.3), the knocking rate f and the tunneling probability T . These two last quantities hardly depend on pairing. On the other hand P changes¹²⁶ from $\approx 2 \times 10^{-76}$ to 2.3×10^{-10} by introducing pairing and, consequently, the associated lifetimes from 10^{75} y to the “observed” (empirical) value of

¹²³Making use of the harmonic oscillator, one can write $\Omega = \frac{1}{2}(N+1)(N+2) \sim A^{2/3}$, where the proportionality constant has a value between 1/2 and 2/3.

¹²⁴Bohr and Mottelson (1975) p. 75.

¹²⁵Montanari et al. (2014); ?.

¹²⁶See Brink and Broglia (2005) Ch. 7 and refs. therein.

10^8 y allowing Cooper pairs to be correlated over distances which can be as large as 20 fm.

Within the above context, exotic halo nuclei open new possibilities to understand the physics at the basis of pairing in nuclei. In fact, one may be able to include in a bare 1S_0 NN -potential the $\approx 50\%$ contribution associated with the induced pairing interaction to the pairing gap in ${}^{120}\text{Sn}$. Hardly, essentially all of the binding energy ($S_{2n} \approx 380$ keV) of the neutron halo Cooper pair of ${}^{11}\text{Li}$ to the core ${}^9\text{Li}$. At the basis of the large magnitude of renormalization effects of elementary modes of excitation and of medium polarization contributions to the pairing interaction observed in light exotic halo nuclei like e.g. ${}^{11}\text{Li}$, one finds a fundamental parameter of NFT, namely the effective degeneracy $\Omega (\approx 2/3A^{2/3})$ of the single-particle phase space, $1/\Omega$ being the small expansion parameter. In the case of ${}^{11}\text{Li}$, Ω is rather small (≈ 3) as compared with heavy nuclei like e.g. ${}^{210}\text{Pb}$ ($\Omega \approx 24$). Furthermore, in the case of ${}^{11}\text{Li}$, the surface (S) to volume (V) ratio ($= aS/V$, a being the nuclear diffusivity) is much larger¹²⁷ (≈ 0.72) than in the case of heavy nuclei lying along the stability valley (≈ 0.27 in the case of ${}^{210}\text{Pb}$).

4.9.2 Coherence length and quantity parameter for (*ph*) vibrations

As schematically indicated in Fig. 2.2.1, both $\beta = 0$ and $\beta = \pm 2$ modes contribute to give rise to a sigmoidal shape to the single-particle occupation probabilities within an energy range $\delta\epsilon = |E_{corr}|$ around the Fermi energy. Concerning $\beta = 0$ see also Fig. 1.2.5. Consequently, the arguments used in Sect. 4.5.1 and leading to

$$\xi = \frac{\hbar v_F}{\pi |E_{corr}|}, \quad (4.9.9)$$

for pairing modes and resulting in $\xi = \hbar v_F / \pi \Delta$ for the case of superfluid nuclei, can be used equally well in connection with $\beta = 0$ modes. In other words, in a similar way in which pair addition and subtraction modes (Cooper pairs) can be viewed as correlated *pp* and *hh* modes over distances of the order ξ , $\beta = 0$ collective vibrations can be pictured as *ph* modes correlated again over distances inversely proportional to the correlation energy of e.g. the RPA collective roots. Summing up, ξ describes, both in the case of $\beta = \pm 2$ and $\beta = 0$ modes, a similar physical property of the vibration: the length over which pairs of fermions ((*ph*), (*pp*), (*hh*)) are correlated in normal and in superfluid nuclei¹²⁸ (see Fig. 4.9.3).

Caption Fig 4.9.3

Vibrations can be classified by the transfer quantum number β . Collective modes with $\beta = 0$ correspond to correlated particle-hole (*ph*) excitations. For example

¹²⁷In carrying out this estimate use was made of $a_{eff} = (R({}^{11}\text{Li})/R_0({}^{11}\text{Li}))a$ (within this context see Fig. 4.3.2).

¹²⁸See e.g. Barranco et al. (2019b) and refs. therein.

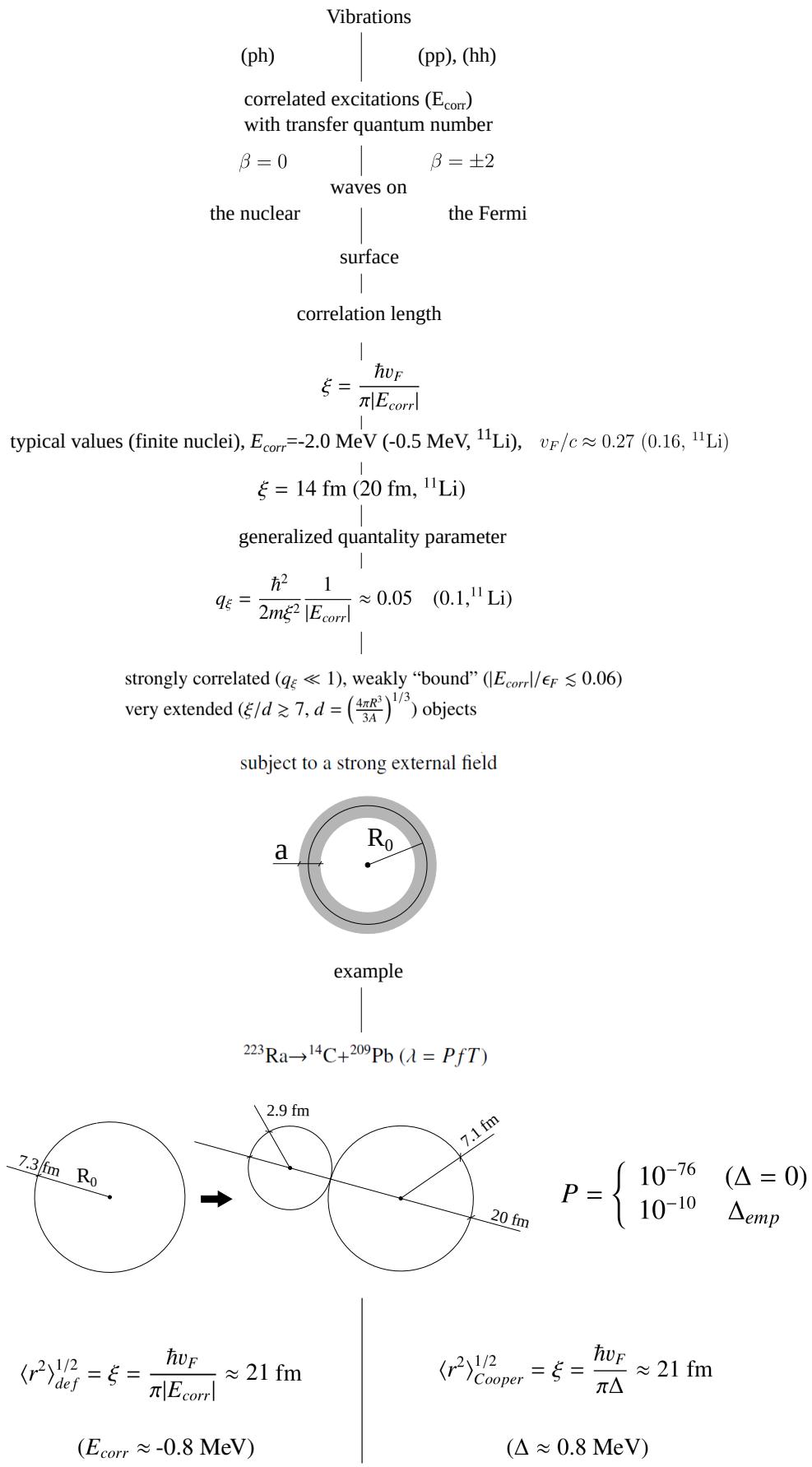


Figure 4.9.3

f		$\sigma(gs \rightarrow f)/\sigma(gs \rightarrow gs)$	Table A
J^π	E_x		
$0^+(gs)$	0	1	(hh) pair removal
3^-	2.62	0.21	(ph) collective mode
5^-	3.20	0.45	(ph) collective mode
0^+	4.87	0.45	(pp) pair addition

Table 4.9.1: Relative two-nucleon transfer cross sections $\sigma(^{206}\text{Pb}(t, p)^{208}\text{Pb}(f)/\sigma(^{206}\text{Pb}(t, p)^{208}\text{Pb}(gs))$ integrated in the range $5^\circ - 175^\circ$ of cm angles. (After Broglia et al. (1973), Table A. VIII b)

J^π	$\sigma(gs \rightarrow f)$ (mb)	$\sigma(gs \rightarrow f)/\sigma(gs \rightarrow gs)$
$0^+(gs)$	2250 ± 338	1
2^+	613 ± 92	0.27

Table 4.9.2: Absolute cross section associated with the reaction $^{120}\text{Sn}(p, t)^{118}\text{Sn}$ to the ground state and first excited state integrated in the range $7.6^\circ < \theta_{cm} < 69.7^\circ$. After Guazzoni et al. (2008).

low-lying quadrupole or octupole (surface) vibrations. Modes with $\beta = \pm 2$ correspond to correlated (pp) or (hh) modes, that is, pair addition and pair subtraction modes. Thinking of these modes propagating in uniform nuclear matter, the associated correlation length can be calculated making use of Eq. (4.9.9), the correlation energy being, as a rule estimated from the energy shift of the lowest root of the RPA solution for both, ph and pairing modes, from the lowest unperturbed configuration (pole). The (generalized) quantality parameter, ratio of the quantal kinetic energy of localization and the correlation energy, gives a measure of the tendency to independent particle ($q_\xi \approx 1$) or independent pp , hh , ph ($q_\xi \ll 1$) motion. A concrete example which testifies to the fact that (ph) excitations (large amplitude surface distortion) and phase correlated independent pair motion (superfluidity) are correlated over dimensions larger than typical nuclear dimensions, is provided by e.g. fission and exotic decay, in particular $^{223}\text{Ra} \rightarrow ^{14}\text{C} + ^{209}\text{Pb}$. In the related estimates use has been made of¹²⁹ $\Delta = 12/\sqrt{A}$ MeV; $C = 18.1$ MeV, $D/\hbar^2 = 29.1$ MeV⁻¹ and¹³⁰ $\hbar\omega = \hbar(C/D)^{1/2} \approx 0.8$ MeV; $E = 2\Delta$ (uncorrelated 2- qp states), $E_{corr} = -(E - \hbar\omega)$. In keeping with the uncertainties affecting the above simple estimates (factor 2 or π in the denominator of ξ , $\langle r^2 \rangle_{Cooper}^{1/2}$, $\sqrt{\frac{5}{3}}\langle r^2 \rangle_{Cooper}^{1/2}$, etc.), it seems fair to conclude that $15 \text{ fm} \lesssim \xi \lesssim 20 \text{ fm}$. Thus, one is likely faced with a situation in which $2 \lesssim \xi/R \lesssim 3$.

The parallel which can be traced between Cooper pairs and correlated particle-hole excitations is further supported by the fact that two-nucleon transfer reaction do excite quite strongly also the $\beta = 0$ modes (see Tables 4.9.1 and 4.9.2).

¹²⁹Bohr and Mottelson (1969).

¹³⁰Brink and Broglia (2005) Sect. 7.1.

Such parallel emerged already in the analysis¹³¹ of the $^{206}\text{Pb}(t, p)^{208}\text{Pb}$ which provided the experimental confirmation of the theoretical predictions concerning pairing vibrations. In particular concerning the population of the pair addition mode $|a\rangle \equiv |gs(^{210}\text{Pb})\rangle$ of the two-phonon pairing vibrational mode $|a\rangle \otimes |r\rangle$; see Fig. 3.1.1(b) transition denoted a (indicating addition mode) predicted at an excitation energy¹³² of 4.9 MeV (Figs. 3.5.1 and 1.3.2) in ^{208}Pb , and observed¹³³ at 4.87 MeV. Because the lowest excited state populated with a sizable cross section ($21 \pm 3\%$) of that associated with the ground state transition (and thus annihilating the pair removal mode ($|r\rangle \equiv |gs(^{206}\text{Pb})\rangle$)), see Fig. 3.1.1(b) transition denoted r) was observed at $E_x = 2.619$ MeV, much effort and time was dedicated to experimentally disentangle an alleged $J^\pi = 0^+$ essentially degenerated with the octupole vibration at $E_x = 2.62 \pm 0.01$ MeV. This was in keeping with the idea that around close shell nuclei, only pairing vibrations were strongly populated, this not being the case for $\beta = 0$ collective surface state.

The analysis of the data dispelled such believe, as testified already by the title of the paper reporting the results. In fact, collective $\beta = 0$ and $\beta = \pm 2$ are similarly strongly populated in two-nucleon transfer reactions, the main difference being the fact that while the Y -components (ground state correlations, see Table 3.5.3) contribute in the case of the pairing modes constructively coherent to the cross section, the similar components of the $\beta = 0$ modes give rise to destructive interference. Again an example of the competition for phase space between pairing and (dynamical) surface deformation (see e.g. Fig. 2.6.2)¹³⁴. These results originally found in Pb, were confirmed in other mass regions and for different $\beta = 0$ modes¹³⁵.

Within the above scenario one can look forward to the test of the $^9\text{Li}(t, p)^{11}\text{Li}$ predictions, concerning the population of both the ground state and the PDR ($E_x \lesssim 1$ MeV; see Fig. 4.12.1). Also to assess to which extent the $0^{*+} {^{12}\text{Be}}$ excited state and 1^- excitation on top of it, can be viewed as related elementary mode of excitation (see Fig. 4.12.1).

4.9.3 Tunneling probabilities

The state $|BCS(\phi)\rangle_{\mathcal{K}} = \prod_{\nu>0} (U'_\nu + V'_\nu e^{-2i\phi} a_\nu^\dagger a_\nu^\dagger) |0\rangle$ displays off-diagonal-long-range-order (ODLRO) because each pair is in a state $(U'_\nu + V'_\nu e^{-2i\phi} a_\nu^\dagger a_\nu^\dagger) |0\rangle$ with the same phase as all the others. In fact, the above wavefunction leads to a two-particle density matrix with the property $\lim_{\mathbf{r}_1, \mathbf{r}_2; \mathbf{r}_3, \mathbf{r}_4 \rightarrow \infty} \phi(\mathbf{r}_1, \mathbf{r}_2; \mathbf{r}_3, \mathbf{r}_4) \neq 0$ un-

¹³¹Broglia and Riedel (1967).

¹³²See Fig. 5 of Bès and Broglia (1966).

¹³³Bjerregaard et al. (1966).

¹³⁴But more important, a further example of the fact that two-nucleon transfer reaction is the specific tool to probe pairing vibrational modes. In fact, setting $\hbar\omega(\beta = 0) = 0$ as well as $\hbar\omega(\beta = 2) = \hbar\omega(\beta = -2) = 0$, in which case the (X, Y) -RPA amplitudes of the corresponding modes become equal, the two-nucleon absolute cross section associated with the $\beta = 0$ mode vanish, while that of the $\beta = \pm 2$ diverge.

¹³⁵Broglia et al. (1971c).

der the assumption that $r_{12}, r_{34} < \xi$, $(\mathbf{r}_1, \mathbf{r}_2)$ and $(\mathbf{r}_3, \mathbf{r}_4)$ being the coordinates of a Cooper pair, r_{ij} the modulus of the relative distance between the corresponding partner fermions, and ξ the coherence length¹³⁶. Quoting from Anderson (1996): “ODLRO is the idea that the density matrix contains a factorizable part $\rho(x, x') = f^*(x)f(x') + \text{remainder}$, with the factorizable part playing the role of a condensate wavefunction”.

Bringing the above argument into reaction implies, that $P_2 \approx P_1$ (see Eq. (4.3.1)). The importance of this result concerning the mechanism at the basis of Cooper pair transfer is connected with the fact that the probability of one-electron tunneling across a typical dioxide layer giving rise to a weak $S - S$ coupling¹³⁷ is likely quite small, e.g. of the order of 10^{-10} . Consequently, simultaneous pair transfer between two superconductors (S), with a probability $(10^{-10})^2$ would not be observed¹³⁸. Thus, the observed Josephson current of carriers of charge $2e$ results from the tunneling of a Cooper pair partner at a time, equally pairing correlated than when they are both in the same superconductor (S) than when each of them is, within a correlation length, in a different of the two (weakly coupled) S .

There is experimental evidence which testifies to the fact that through the Josephson-like junctions, transiently established in heavy ion collisions between superfluid nuclei at energies below the Coulomb barrier, one- and two-nucleon transfer processes display cross sections of the same order of magnitude, provided the distance of closest approach is¹³⁹ $D_0 \lesssim \xi$. More correctly Cooper pair transfer cross section, dominated by ground-ground state transition (i.e. exclusive transitions between members of the corresponding pairing rotational bands), are approximately equal to the incoherent contribution of single quasiparticle states up to an energy of about twice the Cooper pair binding energy ($\approx 2\Delta$). In other words, the parallel of the equality, within a factor ($\pi/4$) of the maximum Josephson current of an unbiased junction, and the normal (quasiparticle) single electron carrier current across a Josephson junction biased with the equivalent potential ($2\Delta/e$) needed to break Cooper pairs (see Eq. (4.6.20) and Fig. 4.6.1).

4.9.4 Nuclear correlation (condensation) energy

The pair-correlation energy is the difference between the energy with and without pairing. The energy including pair correlations is¹⁴⁰

$$E_p = 2 \sum_{\nu>0} |V_\nu|^2 \epsilon_\nu - G |\alpha_0|^2 \quad (4.9.10)$$

¹³⁶See footnote 42 of Ch. 1. See also Ambegaokar (1969), and Potel et al. (2017) and refs. therein.

¹³⁷Pippard (2012).

¹³⁸See e.g. McDonald (2001).

¹³⁹?

¹⁴⁰See Eq. (1.4.38).

while the energy without correlation is

$$E_0 = \sum_{\nu>0} |V_\nu^0|^2 \epsilon_\nu. \quad (4.9.11)$$

The occupation probabilities $|V_\nu^0|$ are unity below the Fermi energy level and zero above. In both Eqs. (4.9.10) and (4.9.11) the Fermi energy has to be chosen to give the correct number of particles. The pairing correlation energy is

$$E_{corr} = E_S - G |\alpha_0|^2, \quad (4.9.12)$$

where

$$E_S = \sum_{\nu>0} 2(|V_\nu|^2 - |V_\nu^0|^2) \epsilon_\nu. \quad (4.9.13)$$

The pairing energy $-G|\alpha_0|^2$ contribution is partially canceled by the first term describing the fact that, in the BCS ground state, particles moving in levels close to the Fermi energy are partially excited across the Fermi surface. This is because V_ν^2 changes smoothly from 1 to 0 around ϵ_F , being 1/2 at the Fermi energy.

In other words, the energy gain resulting from the potential energy term, where G is the pairing coupling constant while $|\alpha_0|$ measures the number of Cooper pairs is partially compensated by a quantal, zero point fluctuation like term. It can, in principle, be related to the Cooper pair kinetic energy of confinement $T_\xi = \frac{\hbar^2}{2m} \frac{1}{\xi^2}$ (already discussed in connection with the generalized quantality parameter (Sect. 3.4)), through the relation $2|\alpha_0|T_\xi$ (for one type of nucleons), in keeping with the fact that (4.9.13) is expressed in terms of single particle energies. Let us make a simple estimate which can help at providing a qualitative example of the above argument, and consider for the purpose the nucleus ^{223}Ra and $G \approx (22/A) \text{ MeV}$, $|\alpha_0| \approx 5$ and¹⁴¹ $\xi \approx 17 \text{ fm}$: $T_\xi \approx 0.1 \text{ MeV}$ while $2 \times (2 \times |\alpha_0| \times T_\xi) = 2 \text{ MeV}$, $2 \times (-G|\alpha_0|^2) = -5 \text{ MeV}$ is the pairing contribution¹⁴² (factor of 2, both protons and neutrons). The resulting pairing correlation energy thus being $E_{corr} = -3 \text{ MeV}$. This number can be compared with a “realistic” estimate provided by the relation¹⁴³

$$E_{corr} = -\frac{g\Delta^2}{4}, \quad (4.9.14)$$

where $g_n \approx N/10 \text{ MeV}^{-1}$ and $g_p \approx Z/10 \text{ MeV}^{-1}$ (see Bortignon et al. (1998) Eq. (5.2)). Taking into account both types of particles $g = g_n + g_p = A/10 \text{ MeV}^{-1}$ and

¹⁴¹This is in keeping with the number of protons and of neutrons outside closed shells. For the estimate of $\xi = \hbar c \times 0.27/(\pi\Delta)$ we have used $\Delta = 12/\sqrt{A} \text{ MeV} \approx 1 \text{ MeV}$.

¹⁴²This quantity, but divided by 2, i.e. -2.5 MeV can be compared with the effective pairing matrix element $v = \left(\frac{\Delta_n^2 + \Delta_p^2}{4G} \approx -2.9 \text{ MeV} \right)$, operative at level crossing in the calculation of the inertia of the exotic decay $^{223}\text{Ra} \rightarrow ^{14}\text{Ca} + ^{209}\text{Pb}$, cf. Brink and Broglia (2005) p.159 and refs. therein.

¹⁴³See e.g. Sect. 3.5 of Brink and Broglia (2005).

making use of $\Delta = 12/\sqrt{A}$ MeV, one obtains $E_{corr} = -\frac{144}{40}$ MeV ≈ -3.6 MeV. With the help of E_{corr} and T_ξ , one can estimate the generalized quantity parameter, $q_\xi = T_\xi/|E_{corr}| = 0.1/3.6 \approx 0.03$, as well as make a consistency check on the value of ξ used, namely $\hbar v_F/|E_{corr}| \approx 15$ fm.

4.10 Absolute Cooper pair tunneling cross section: quantitative novel physics at the edge between stability and chaos

In the study of many-body systems, in particular of finite quantum many-body systems (FQMBS) like the atomic nucleus, much can be learned from symmetries. However it is the texture of the associated emergent properties, concrete embodiment of spontaneously symmetry breaking (potential energy) and of its restoration (zero point fluctuations (collective modes), confinement kinetic energy), which provides insight into the eventual new physics. In fact, when one understands the many-body under study in terms of the detailed motion of single-particles (nucleons) and collective motion, taking properly into account their couplings, is that one may eventually reach a quantitative understanding of the problem and of its solutions. Furthermore, that these solutions are likely transferable, at profit, to the study of other FQMBS like e.g. metal clusters, fullerenes¹⁴⁴, quantum dots¹⁴⁵, and eventually soft matter, in particular proteins¹⁴⁶, let alone the fact that one can make predictions. Predictions which, in connection with the study of halo nuclei, in particular of pairing¹⁴⁷ involve novel physics¹⁴⁸. Within this context one can quote from Leon Cooper's contribution to the volume¹⁴⁹ BCS: 50 years: "It has become fashionable... to assert... that once gauge symmetry is broken the properties of superconductors follow... with no need to inquire into the mechanism by which the symmetry is broken¹⁵⁰. This is not... true, since broken gauge symmetry might lead to molecule-like and a Bose-Einstein rather than BCS condensation... in 1957... the major problem was to show... how... an order parameter or condensa-

¹⁴⁴See e.g. Gunnarsson (2004), Broglia et al. (2004) and refs. therein.

¹⁴⁵Lipparini (2003).

¹⁴⁶Broglia (2013).

¹⁴⁷Cf. e.g. Broglia and Zelevinsky (2013).

¹⁴⁸Cf. e.g. Barranco et al. (2001); Tanihata et al. (2008); Potel et al. (2010) and references therein.

¹⁴⁹Cooper (2011).

¹⁵⁰Quoting (Weinberg (2011)): "... In consequence of this spontaneous symmetry breaking, products of any even number of electron fields have non-vanishing expectation values in a superconductor, though a single electron field does not. All of the dramatic exact properties of superconductors – zero electric resistance, the expelling of the magnetic fields from superconductors known as the Meissner effect, the quantization of magnetic flux through a thick superconducting ring, and the Josephson formula for the frequency of the ac current at a junction between two superconductors with different voltages– follow from the assumption that electromagnetic gauge invariance is broken in this way, with no need to inquire into the mechanism by which the symmetry is broken." The above quotation is similar to saying that once the idea of a double DNA helix was thought, all about inheritance was solved and known, (cf. e.g. Stent (1980) and references therein).

tion in momentum space could come about... to show how... gauge-invariant symmetry of the Lagrangian could be spontaneously broken due to interactions which were themselves gauge invariant".

Nuclear physics has brought this quest a step further. This time in connection with the "extension" of the study of BCS condensation to its origin, a single Cooper pair in the rarified atmosphere resulting from the strong radial (isotropic) deformation observed in light exotic, halo nuclei in general, and in ^{11}Li in particular. During the last few years, the probing of this system in terms of absolute two-nucleon transfer (pick-up) reactions, has helped at making this field a quantitative one, errors below the %10 limit being the rule. This achievement which has its basis on much experimental work, in particular the remarkable experiments of Tanihata et al. (2008), is also the result of the combined effort made in treating the structure and reaction aspects of the subject, two sides of the same physics, on equal footing.

4.10.1 Saturation density, spill out and halo

In the incipit to the Section on bulk properties of nuclei of Bohr and Mottelson¹⁵¹ one reads: "The almost constant density of nuclear matter is associated with the finite range of nuclear forces; the range of the forces is r_0 (where r_0 enters the nuclear radius in the expression $R = r_0 A^{1/3}$) thus small compared to nuclear size. This "saturation" of nuclear matter is also reflected in the fact that the total binding energy of the nucleus is roughly proportional to A . In a minor way, these features are modified by surface effects and long-range Coulomb forces acting between the protons".

Electron scattering experiments (see the figure 2-1, p. 159 of the above reference) yield

$$\rho(0) = 0.17 \text{ fm}^{-3}. \quad (4.10.1)$$

Thus, one can posit that

$$\frac{4\pi}{3} R_0^3 \rho(0) = A, \quad (4.10.2)$$

leading to

$$r_0 = \left(\frac{3}{4\pi} \frac{1}{\rho(0)} \right)^{1/3} \approx 1.1 \text{ fm}. \quad (4.10.3)$$

Because the above relations imply a step function distribution, one should correct Eq. (4.10.3) by the nucleon spill out¹⁵². The relative correction being of the order of $(a_0/R_0) \ln 2 \approx (0.65/5.92) \ln 2 \approx 0.08$ ($A = 120$) associated with the fact that a

¹⁵¹Bohr and Mottelson (1969) p. 139.

¹⁵²Bertsch and Broglia (2005).

more realistic distribution is provided by a Fermi function of diffusivity $a_0 \approx 0.65$ fm. That is, an 8% correction leading to $r_0 = (1.12 + 0.08)$ fm ≈ 1.2 fm.

In the case of the nucleus ^{11}Li , observations¹⁵³ indicate a mean square (gyration radius¹⁵⁴) $\langle r^2 \rangle^{1/2} = 3.55 \pm 0.1$ fm. Thus

$$R(^{11}\text{Li}) = \sqrt{\frac{5}{3}} \langle r^2 \rangle^{1/2} \approx 4.58 \pm 0.13 \text{ fm.} \quad (4.10.4)$$

Making use of the relation $R_0 \approx 1.2A^{1/3}$ fm, the quantity (4.10.4) leads to $(4.58/1.2)^3 \approx 56$, an effective mass number larger five times the actual value $A = 11$. To be noted that the actual mass number predicts a “systematic” value of the nuclear radius $R_0(^{11}\text{Li}) \approx 2.7$ fm.

The above results testify to a very large “isotropic radial deformation”, or halo region (skin), in keeping with the fact¹⁵⁵ that $R(^{11}\text{Li}) - R_0(^9\text{Li}) = R_0(^9\text{Li}) \left(\frac{R(^{11}\text{Li})}{R_0(^9\text{Li})} - 1 \right) = 0.83R_0(^9\text{Li})$. In other words, ^{11}Li can be viewed as made out of a normal ^9Li core and of a skin made out of two neutrons extending over a shell of width of the order of that of the core radius. But even more important, it is the fashion in which the above mentioned “deformation” affects nuclear matter. Matter which is little compliant to undergo either compressions or, for that sake, “depressions”, without resulting in nuclear instability. In one case, through a mini supernova. In the second, by obliterating the effect of the short range strong force acting in the 1S_0 channel (screening effect of the bare pairing interaction).

In fact, in the case of the halo Cooper pair of ^{11}Li , that is of the last two weakly bound neutrons, one is dealing with a rarefied nuclear atmosphere of density

$$\rho \approx \frac{2}{\frac{4\pi}{3}(R^3(^{11}\text{Li}) - R_0^3(^9\text{Li}))} \approx 0.6 \times 10^{-2} \text{ fm}^{-3} \quad (4.10.5)$$

where the value $R_0(^9\text{Li}) \approx 2.5$ fm was used. That is, we are dealing with pairing in a nuclear system at a density which is only 4% of saturation density.

The quest for the long range pairing mechanism which is at the basis of the binding of the halo Cooper pair of ^{11}Li to the ^9Li core ($S_{2n} \approx 0.380$ keV, to be compared to typical systematic values of $S_{2n} \approx 16$ MeV), has lead to the discovery of what can be considered a novel nuclear mode of elementary excitation. The symbiotic halo pair addition mode, which has to carry its own source of binding (glue). A novel embodiment of the Axel–Brink scenario in which not only the line shape, but the main structure of the resonance depends on the state on which it is built, and to which it is deeply interwoven as to guarantee its stability¹⁵⁶ and thus its own existence. It also provides a novel realization of the Bardeen–Frölich–Pines¹⁵⁷

¹⁵³Kobayashi et al. (1989).

¹⁵⁴The radius of gyration R_g is a measure of an object of arbitrary shape, R_g^2 being the second moment in 3D space. In the case of a sphere of radius R , $R_g^2 = 3R^2/5$.

¹⁵⁵Let us parametrize the radius of ^{11}Li as (see Bohr and Mottelson (1975)), $R = R_0(1 + \alpha_{00}Y_{00}) = R_0(1 + \beta_0 \frac{1}{\sqrt{4\pi}})$. Thus $\beta_0 = \sqrt{4\pi}(\frac{R}{R_0} - 1) \approx 2.96$ which testifies to the extreme “exoticity” of the phenomenon.

¹⁵⁶Axel (1962); Brink (1955 (unpublished)).

¹⁵⁷Bardeen and Pines (1955); Fröhlich (1952).

microscopic mechanism to break gauge invariance: through the exchange of quite large dipole ZPF which ensures the same symmetries of the original Hamiltonian to a system displaying essentially a permanent dipole moment, as a consequence of the almost degeneracy of the pygmy dipole resonance (centroid ≈ 0.75 MeV) with the ground state¹⁵⁸. An example of a van der Waals Cooper pair (Sect. 3.7).

The NFT diagram shown in Fig. 3.7.1 describing this binding seems quite involved and high order. Thus unlikely to be at the basis of a new elementary mode of nuclear excitation, if nothing else because of the apparent lack of “elementarity”. This is not the case and, in fact, the physics at the basis of the process depicted by the oyster-like and eagle-like networks displayed in (a) and (b) is quite simple and recurrent throughout nuclear structure and reactions, let alone many-body theories and QED. In fact, it encompasses (see Fig. 3.7.1): (I,II) the changes in energy of single-particle levels as a function of quadrupole dynamical deformations leading to a Jahn-Teller like effect (III) the interaction between particles through the exchange of (quasi) bosons (vibrations), (IV,V) Pauli principle leading to a conspicuous Lamb shift-like effect, (VI,VII) the softening of collective modes due to ground state correlations ((ZPF)-components, QRPA),(VIII) the interaction between two non-polar systems through virtual, ZPF associated dynamical dipoles.

Referring to general many-degrees of freedom systems, (I,II) and (III) are at the basis of the fact that, in QED, the coupling between one- and two- photons is zero (Furry’s theorem). It is also related, through cancellation processes, to the small width displayed by (particle-hole) nuclear giant resonances, as well as of inhomogeneous damping of nuclear rotational motion at finite temperatures¹⁵⁹ – parallel to motional narrowing observed in NMR– as compared with the width of single-particles at similar excitation energies. Concerning (VIII), one can mention resonant interactions between fluctuating systems like e.g. two coupled harmonic oscillators¹⁶⁰.

In the case of halo Cooper pair binding by PDR in ^{11}Li , the system although not being on resonance it is not far from it, in keeping with the fact $\tilde{\epsilon}_{p_{1/2}} - \tilde{\epsilon}_{s_{1/2}} \approx 0.45$ MeV and $\hbar\omega_{PDR} \approx 0.75$ MeV with a width of ≈ 0.5 MeV. In other words the ^{10}Li inverted parity system is poised to acquire a permanent dipole moment or, almost equivalent, to display a large amplitude, dipole mode at very low energy as well as a collective $B(E1)$ to the halo ground state, of the order of a single-particle Weisskopf unit $B_W(E1)$. The PDR (see Fig. 4.12.1, see also Fig. 2.7.1) is in ^{11}Li screened from the GDR through the poor overlap between core and halo single-particle wavefunctions. In this way, is able to retain essentially all of its $E1$ -strength ($\approx 1 \times B_W$), and can likely be considered a new mode of excitation¹⁶¹. Said it differently, one is faced, already at the level of the single-particle spectrum (parity inversion), with the possibility of a plastic large amplitude dipole mode, as

¹⁵⁸Aumann (2019).

¹⁵⁹Broglia et al. (1987).

¹⁶⁰See e.g. Born (1969), App. XL p. 471.

¹⁶¹See Broglia et al. (2019a) and references therein.

can be seen from

$$\frac{dn}{d\beta_L} = \frac{1}{4} \sqrt{\frac{2L+1}{3\pi}} A. \quad (4.10.6)$$

Relation which defines the number of level crossings n , in terms of the isotropical radial deformation¹⁶² ($L = 0$ and $\beta_0 = 2.96$), leading to $n \approx 3$.

4.11 Pairing spatial correlation: simple estimate

Let us assume two equal nucleons above closed shell as the nuclear embodiment of Cooper's model. The two-particle wave function in configuration space can be written as,

$$\Psi(\mathbf{r}_1\sigma_1, \mathbf{r}_2\sigma_2) = \Psi_0(\mathbf{r}_1, \mathbf{r}_2)\chi_{S=0}(\sigma_1, \sigma_2) + [\Psi_1(\mathbf{r}_1, \mathbf{r}_2)\chi_{S=1}(\sigma_1\sigma_2)]_0, \quad (4.11.1)$$

where $\chi_{S=0}$ and $\chi_{S=1}$ are the singlet and triplet spin wavefunctions, respectively.

In what follows we shall consider a pairing interaction acting on pairs of particles moving in time reversal states. Consequently we shall concentrate in the spin singlet radial component of (4.11.1). In the Tamm-Dancoff approximation (in keeping with Cooper ansatz) one can write

$$\Psi_0(\mathbf{r}_1, \mathbf{r}_2) = \sum_{nn'lj} X_{nn'lj} R_{nl}(r_1) R_{n'l}(r_2) \sqrt{\frac{2j+1}{2(2l+1)}} [Y_l(\hat{r}_1) Y_l(\hat{r}_2)]_0. \quad (4.11.2)$$

This wave function can be rewritten as

$$\begin{aligned} \Psi_0(\mathbf{r}_1, \mathbf{r}_2) &= \Psi_0(|\mathbf{r}_1|, |\mathbf{r}_2|, \theta) \\ &= \sum_{n'n'lj} X_{nn'lj} R_{nl}(r_1) R_{n'l}(r_2) \sqrt{\frac{2j+1}{2}} \frac{1}{4\pi} P_l(\cos \theta), \end{aligned} \quad (4.11.3)$$

where $\theta = \widehat{r_1 r_2}$. A convenient way to display two-particle correlation is by plotting $|\Psi_0(|\mathbf{r}_1|, |\mathbf{r}_2|, \theta)|^2$ in the $x - z$ plane.

In the case of pure configurations $a \equiv nlj$,

$$\Psi_0(|\mathbf{r}_1|, |\mathbf{r}_2|, \theta) = R_{nl}(r_1) R_{nl}(r_2) \sqrt{\frac{2j+1}{2}} \frac{1}{4\pi} P_l(\cos \theta). \quad (4.11.4)$$

In keeping with the fact that the specific probe of pairing correlation is two-nucleon transfer, a phenomenon which takes place mainly, although not only, on the nuclear surface, we shall set $r_1 = r_2 = R_0$, and use the empirical relation¹⁶³

$$R_{nl}(R_0) = \left(\frac{1.4}{R_0^3} \right)^{1/2}, \quad (4.11.5)$$

¹⁶²Brink and Broglia (2005), Eq. (7.35) and refs. therein. See also footnote 155 of this chapter.

¹⁶³Bohr and Mottelson (1969).

Thus

$$\Psi_0(R_0, R_0, \theta) = \left(\frac{1.4}{R_0^3} \right) \sqrt{\frac{2j+1}{2}} \frac{1}{4\pi} P_l(\cos \theta). \quad (4.11.6)$$

and

$$|\Psi_0(R_0, R_0, \theta)|^2 \sim |P_l(\cos \theta)|^2. \quad (4.11.7)$$

It is seen that the two particles have the same probability to be on top of each other ($\theta = 0^\circ; P_l(1) = 1$), or on opposite sides of the nucleus ($\theta = 180^\circ; P_l(-1) = (-1)^l$). Taking into account the actual radial dependence of $R_{nl}^2(r_1)$ for $r_1 = R_0$, the width of the two probability peaks is found to be ≈ 2 fm, i.e. $d \approx \left(\frac{4\pi}{3} R^3 / A \right)^{1/3}$.

Let us now consider the general expression (4.11.3), and assume, aside from $n = n'$, that the two nucleons are allowed to correlate in a phase space composed of N single-particle levels, and that all amplitudes are equal,

$$X \approx \frac{1}{\sqrt{N}}. \quad (4.11.8)$$

Thus

$$\Psi_0(R_0, R_0, \theta) = \left(\frac{1.4}{R_0^3} \right) \frac{1}{\sqrt{N}} \frac{1}{4\pi} \sqrt{\frac{2j+1}{2}} \sum_l P_l(\cos \theta), \quad (4.11.9)$$

where again (4.11.5) have been used. One can then write

$$|\Psi_0(R_0, R_0, \theta)|^2 \sim \left| \sum_l P_l(\cos \theta) \right|^2. \quad (4.11.10)$$

Assuming the closed shell nucleus to be ^{208}Pb and the N single-particle levels the neutron valence orbitals $2g_{9/2}, 1i_{11/2}, 1j_{15/2}, 3d_{5/2}, 4s_{1/2}, 2g_{7/2}$ and $3d_{3/2}$, one obtains

$$\frac{|\Psi_0(R_0, R_0, \theta = 0^\circ)|^2}{|\Psi_0(R_0, R_0, \theta = 180^\circ)|^2} \approx \left(\frac{7}{5} \right)^2 \approx 2, \quad (4.11.11)$$

in keeping with the fact that there is only a single state of opposite parity (intruder $j_{15/2}$ state).

Making use of an extended basis, containing a similar amount of positive and negative natural parity states, that is taking into account a large number of major shells ($\pi = (-1)^N, N$ principal quantum number), one can reduce in a consistent fashion the value of $|\Psi_0(R_0, R_0, \theta = 180^\circ)|^2$. This of course materializes already within the basis of valence states in e.g. ^{11}Li , in keeping with the fact that in this case $s_{1/2}$ and $p_{1/2}$ play a similar role.

Summing up, the above results have something to do with the Cooper pair problem, but much more with the peculiarities of spatial quantization associated with the nuclear self-bound many-body system at the level of mean field. That is, the nuclear Cooper pair phenomenon is to be expressed under the influence of a very strong external field which imposes not only confinement, but also spatial quantization with strong spin orbit effects resulting, among other things, in intruder states and thus parity mixing, let alone parity inversion due to quadrupole fluctuation of the mean field.

4.12 Hindsight

The formulation of superconductivity (BCS theory) described by Gor'kov¹⁶⁴ allows, among other things for a simple visualization of spatial dependences. In this formulation $F(\mathbf{x}, \mathbf{x}')$ is the amplitude for two Fermions (electrons) at \mathbf{x}, \mathbf{x}' , to belong to the Cooper pair (within the framework of nuclear physics cf. e.g. Fig. 3.6.3 $\Psi_0(\mathbf{r}_1, \mathbf{r}_2)$; see also Sect. 4.9). The phase of F is closely related to the angular orientation of the spin variable in Anderson's quasispin formulation of BCS theory¹⁶⁵. The gap function $\Delta(x)$ is given by $V(\mathbf{x})F(\mathbf{x}, \mathbf{x})$ where $V(\mathbf{x})$ is the local two-body interaction at the point \mathbf{x} . In the insulating barrier between the two superconductors of a Josephson junction, $V(\mathbf{x})$ is zero and thus $\Delta(x)$ is also zero.

The crucial point is that vanishing¹⁶⁶ $\Delta(x)$ does not imply vanishing F (abnormal density), provided, of course, that one has within the junction, a non-zero particle (electron) density, resulting from the overlap of densities from right (R) and left (L) superconductors. Now, these junctions are such that they allow for one-electron-tunneling with a probability P_1 quite small, but finite and, consequently, the above requirement is fulfilled. Nonetheless, conventional simultaneous pair transfer, with a probability P_1^2 will not be observed¹⁶⁷. *But because one electron at a time can tunnel profiting of the small, but finite electron density within the junction layer, $F(\mathbf{x}, \mathbf{x}')$ can have a sizable amplitude for Cooper pairs with partners electrons one on each side of the barrier (i.e. $\mathbf{x} \in L$ and $\mathbf{x}' \in R$), separated by distances $|\mathbf{x} - \mathbf{x}'|$ up to the coherence length.* Hence, for barriers thin compared with the coherence length, two electrons on opposite sides of the barrier can still be correlated and the pair current be consistent. An evaluation of its value shows that, at zero temperature, the pair current is equal to $\pi/4$ times the normal, single (quasiparticle) electron carrier current, at an equivalent voltage¹⁶⁸ $2\Delta/e$. Evalua-

¹⁶⁴Gor'kov (1958, 1959).

¹⁶⁵Anderson (1958); within the framework of nuclear physics cf. e.g. Bohr and Ulfbeck (1988), Potel et al. (2013b) and references therein.

¹⁶⁶This point was likely misunderstood by Bardeen who writes "...In my view, virtual pair excitations do not extend across the layer...", see McDonald (2001), see also Bardeen (1961) and Bardeen (1962).

¹⁶⁷Pippard (2012).

¹⁶⁸In the case of Pb at low temperatures (≈ 7.19 K (0.62 meV)) this voltage is ≈ 1 meV/e = 1 mV leading to ≈ 2 mA current for a barrier resistance of $R \sim 1\Omega$ (Ambegaokar and Baratoff (1963);

tion which has been experimentally confirmed¹⁶⁹.

The translation of the above parlance to the language of nuclear physics has to come to terms with the basic fact that nuclei are self-bound, finite many-body systems in which the surface, as well as the closely related property of spatial quantization, play an important role both as a static element of confinement, as well as a dynamic source for renormalization effects^{170,171}. Under the influence of the average potential which can be viewed as very strong external field ($|V_0| \approx 50$ MeV), Cooper pairs ($|E_{corr}| \approx 2$ MeV) will become constrained within its boundaries with some amount of spill out.

Let us now consider a two nucleon transfer reaction in the collision Sn+Sn assuming a distance of closest approach of ≈ 14 fm, in which the two nuclear surfaces are separated by ≈ 2 fm (Fig 4.5.1). In keeping with the fact that this distance is about $3 \times a$, the heavy ion system will display a few percent overlap of (normal, saturation) density. Ever so small this overlap and so narrow the surface hole between the two leptodermic systems resulting from it, Cooper pairs can now extend over the two volumes, in a similar way as electron Cooper pairs could partially be found in the R and L superconductors in a Josephson junction. If this is the case, Cooper pair partners can, in principle, be at distances of the order of the correlation length (see Sects. 4.9.3 and 7.3).

An example of the fact that Cooper pairs will “expand” if the external mean field allows for it, is provided by ^{11}Li in which case, profiting of the weak binding (≈ 380 keV), the extension of the constrained Cooper pair (≈ 4.58 fm ± 0.13 fm) is similar to that expected in a nucleus of mass number $A \approx 60$, assuming a standard radial behavior, i.e. $r_0 A^{1/3}$ fm. In keeping with this scenario, it could be expected that moving from one neutron pair addition 0^+ mode of the $N = 6$ isotones¹⁷² to another one ($|^{11}\text{Li(gs)}\rangle$, $|^{12}\text{Be(gs)}\rangle$ and $|^{12}\text{Be}(0^{+*}; 2.24 \text{ MeV})\rangle$) one would see the system expanding, contracting and expanding again, respectively, in keeping with the fact that the external (mean) field is weak, strong, weak respectively, as testified by S_{2n} (380 keV, 3672 keV, 1432 keV). Within this context, in Fig. 4.12.1 a schematic overall view of the pairing vibrational modes associated with $N = 6$ parity inverted closed shell isotones, together with low-energy $E1$ -strength modes is given. The possible candidates to the role of neutron halo pair addition modes

McDonald (2001); Tinkham (1996)).

¹⁶⁹Rogalla and Kes (2012).

¹⁷⁰Within this context it is of notice that the liquid drop model is a very successful nuclear model, able to accurately describe not only large amplitude motion (fission, exotic decay, low-lying collective surface vibrations, cf. e.g. Bohr and Wheeler (1939), Bertsch (1988), Barranco et al. (1990), and references therein), but also the masses of nuclides (see e.g. Møller et al. (1995)), provided the superfluid inertia and shell corrections respectively, are properly considered. Thus, it is likely that in the quest of more predictive theoretical tools of the global nuclear properties one should develop on equal footing ever more “accurate” bare NN -potentials as well as effective forces, and methods to deal with the long wavelength, renormalization effects and induced interactions. Within this context it is illuminating the contribution of Anderson and Thouless (1962) concerning the nuclear surface diffuseness, at the time of the development of Brueckner theory (Brueckner et al. (1961)).

¹⁷¹Broglia (2002).

¹⁷²See e.g. Gori et al. (2004).

and symbiotic state are explicitly indicated (boxed levels).

Appendix 4.A Coherent state

The BCS ground state can be written as,

$$\begin{aligned} |BCS(\phi)\rangle_{\mathcal{K}} &= \prod_{v>0} \left(U_v + V_v a_v^\dagger a_{\bar{v}}^\dagger \right) |0\rangle = \prod_{v>0} U_v \left(1 + \frac{V_v}{U_v} a_v^\dagger a_{\bar{v}}^\dagger \right) |0\rangle \\ &= \left(\prod_{v>0} U_v \right) \left(\prod_{v>0} \left(1 + c_v P^\dagger \right) \right) |0\rangle, \end{aligned} \quad (4.A.1)$$

where

$$c_v = \frac{V_v}{U_v} \quad \text{and} \quad P^\dagger = a_v^\dagger a_{\bar{v}}^\dagger. \quad (4.A.2)$$

In what follows, we work out a couple of simple examples:

a) $v=1,2$ (two pairs),

$$\begin{aligned} \prod_{v>0} \left(1 + c_v P^\dagger \right) &= \left(1 + c_1 P_1^\dagger \right) \left(1 + c_2 P_2^\dagger \right) = 1 + c_1 P_1^\dagger + c_2 P_2^\dagger \\ &\quad + c_1 c_2 P_1^\dagger P_2^\dagger = 1 + \sum_{v>0} c_v P_v^\dagger + \frac{1}{2!} \left(\sum_{v>0} c_v P_v^\dagger \right)^2, \end{aligned} \quad (4.A.3)$$

where use has been made of

$$\left(c_1 P_1^\dagger + c_2 P_2^\dagger \right)^2 = 2 c_1 c_2 P_1^\dagger P_2^\dagger, \quad (4.A.4)$$

in keeping with the fact that

$$\left(P_1^\dagger \right)^2 = \left(P_2^\dagger \right)^2 = 0. \quad (4.A.5)$$

The above expression underscores the fact that Cooper pairs are not bosons. **b) $v=1,2,3$ (three pairs):**

$$\begin{aligned} \prod_{v>0} \left(1 + c_v P^\dagger \right) &= \left(1 + c_3 P_3^\dagger \right) \left(1 + c_2 P_2^\dagger \right) \left(1 + c_1 P_1^\dagger \right) \\ &= \left(1 + c_3 P_3^\dagger \right) \left(1 + c_1 P_1^\dagger + c_2 P_2^\dagger + c_1 c_2 P_1^\dagger P_2^\dagger \right) \\ &= 1 + \left(c_1 P_1^\dagger + c_2 P_2^\dagger + c_3 P_3^\dagger \right) + \left(c_1 c_2 P_1^\dagger P_2^\dagger + c_1 c_3 P_1^\dagger P_3^\dagger + c_2 c_3 P_2^\dagger P_3^\dagger \right) \\ &\quad + c_1 c_2 c_3 P_1^\dagger P_2^\dagger P_3^\dagger = 1 + \sum_{v>0} c_v P_v^\dagger + \frac{1}{2!} \left(\sum_{v>0} c_v P_v^\dagger \right)^2 + \frac{1}{3!} \left(\sum_{v>0} c_v P_v^\dagger \right)^3, \end{aligned} \quad (4.A.6)$$

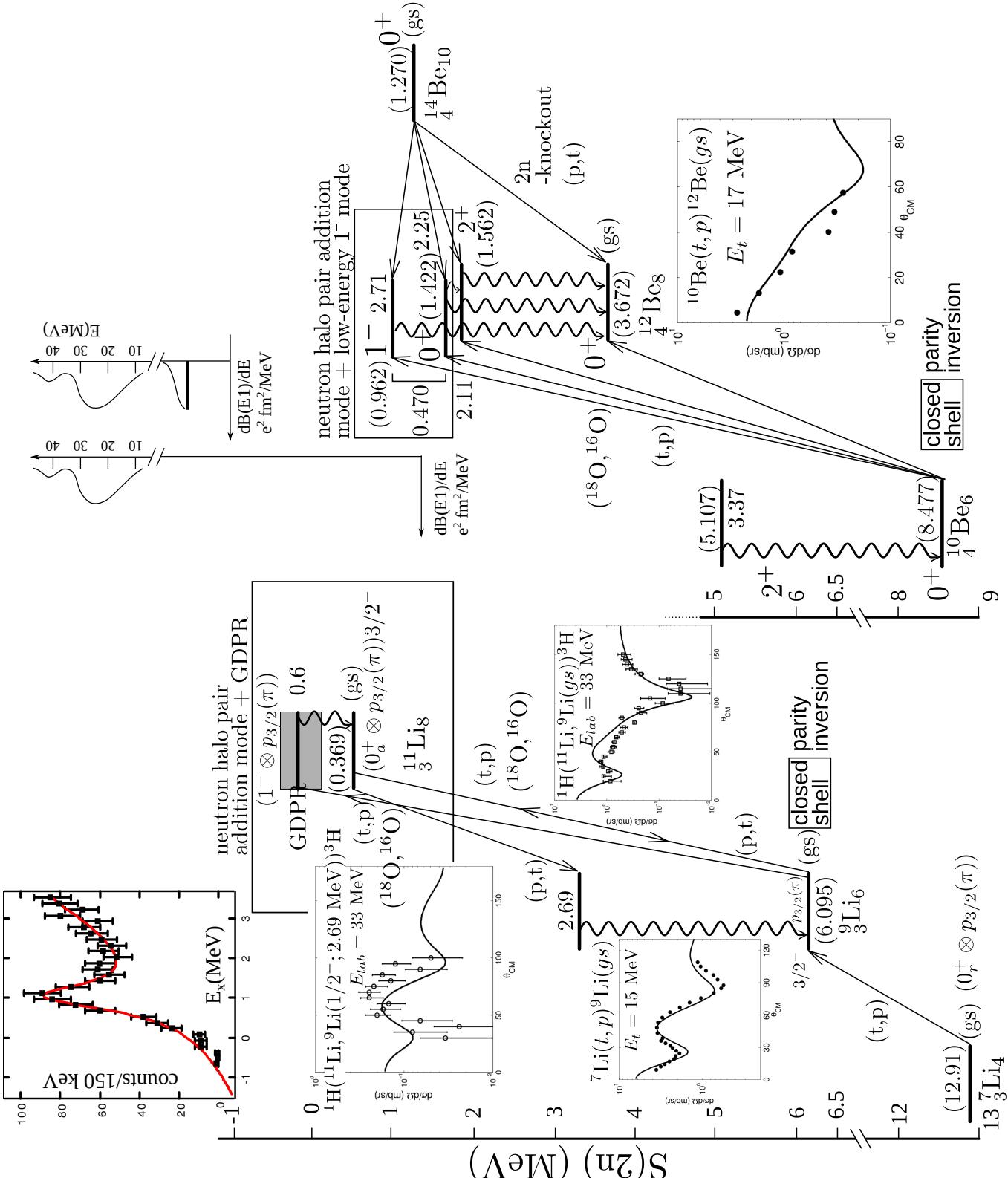


Figure 4.12.1: Monopole pairing vibrational modes associated with $N = 6$ parity inverted closed shell isotones, together with low-energy E1-strength modes. The levels are displayed as a function of the two-neutron separation energies $S(2n)$. These quantities are shown in parenthesis, the excitation energies with respect to the ground state are quoted in MeV. Absolute differential cross sections from selected (t, p) and (p, t) reactions calculated as described in the text (cf. Potel et al. (2010, 2014)), in comparison with the experimental data (Young and Stokes (1971); Fortune et al. (1994)).

where use has been made of the relations (4.A.5) and of,

$$\begin{aligned} [(a+b+c)(a+b+c)] &= ab + ac + ba + bc + ca + cb = 2ab + 2ac + 2bc, \\ a^2 = b^2 = c^2 &= 0, \end{aligned}$$

together with

$$(a+b+c)[(a+b+c)(a+b+c)] = 2abc + 2bac + 2cab = 6abc. \quad (4.A.7)$$

Thus,

$$\left(\sum_{\nu>0} c_{\nu} P_{\nu}^{\dagger} \right)^3 = 6c_1 c_2 c_3 P_1^{\dagger} P_2^{\dagger} P_3^{\dagger} = 3! c_1 c_2 c_3 P_1^{\dagger} P_2^{\dagger} P_3^{\dagger} \quad (4.A.8)$$

Making use of

$$e^x = 1 + x + \frac{x^2}{2!} + \frac{x^3}{3!} + \dots, \quad (4.A.9)$$

one can write

$$\begin{aligned} |BCS(\phi)\rangle_{\mathcal{K}} &= \left(\prod_{\nu>0} U_{\nu} \right) \left\{ 1 + \frac{1}{1!} \left(\sum_{\nu>0} c_{\nu} P_{\nu}^{\dagger} \right) + \frac{1}{2!} \left(\sum_{\nu>0} c_{\nu} P_{\nu}^{\dagger} \right)^2 + \frac{1}{3!} \left(\sum_{\nu>0} c_{\nu} P_{\nu}^{\dagger} \right)^3 + \dots \right\} |0\rangle, \\ &= \left(\prod_{\nu>0} U'_{\nu} \right) \left\{ 1 + \frac{e^{-2i\phi}}{1!} \left(\sum_{\nu>0} c'_{\nu} P_{\nu}^{\dagger} \right) + \frac{e^{-4i\phi}}{2!} \left(\sum_{\nu>0} c'_{\nu} P_{\nu}^{\dagger} \right)^2 + \frac{e^{-6i\phi}}{3!} \left(\sum_{\nu>0} c'_{\nu} P_{\nu}^{\dagger} \right)^3 + \dots \right\} |0\rangle, \end{aligned} \quad (4.A.10)$$

where

$$c_{\nu} = e^{-2i\phi} c'_{\nu}, \quad c'_{\nu} = V'_{\nu} / U'_{\nu}. \quad (4.A.11)$$

Thus,

$$\begin{aligned} |BCS(\phi)\rangle_{\mathcal{K}} &= \left(\prod_{\nu>0} U_{\nu} \right) \exp \left(\sum_{\nu>0} c_{\nu} P_{\nu}^{\dagger} \right) |0\rangle = \left(\prod_{\nu>0} U_{\nu} \right) \sum_{n=0,1,2,\dots} \frac{\left(\sum_{\nu>0} c_{\nu} P_{\nu}^{\dagger} \right)^n}{n!} |0\rangle \\ &= \left(\prod_{\nu>0} U'_{\nu} \right) \sum_{N \text{ even}} \frac{e^{-iN\phi}}{(N/2)!} \left(\sum_{\nu>0} c'_{\nu} P_{\nu}^{\dagger} \right)^{N/2} |0\rangle, \end{aligned} \quad (4.A.12)$$

(see Eq. (4.7.27)) and

$$\begin{aligned} |N_0\rangle &= \int d\phi e^{iN_0\phi} |BCS(\phi)\rangle_{\mathcal{K}} \\ &= \left(\prod_{\nu>0} U'_{\nu} \right) \sum_{N \text{ even}} \int d\phi e^{iN_0\phi} \frac{e^{-iN\phi}}{(N/2)!} \left(\sum_{\nu>0} c'_{\nu} P_{\nu}^{\dagger} \right)^{N/2} |0\rangle \sim \left(\sum_{\nu>0} c'_{\nu} P_{\nu}^{\dagger} \right)^{N_0/2} |0\rangle, \end{aligned} \quad (4.A.13)$$

is the member with N_0 particles of the pairing rotational band, while

$$\left(\sum_{\nu>0} c'_\nu P_\nu^\dagger \right) |0\rangle, \quad (4.A.14)$$

is the Cooper pair state. Because $U'_\nu \rightarrow 0$ for $\epsilon \ll \epsilon_F$, (4.A.14) is to be interpreted to be valid for values of ϵ_ν close to ϵ_F . In expression (4.A.12), $n = N/2$ is the number of Cooper pairs.

Making use of the single j -shell model

$$V' = \sqrt{\frac{N}{2\Omega}}, \quad U' = \sqrt{1 - \frac{N}{2\Omega}}, \quad (4.A.15)$$

and

$$\frac{V'}{U'} = \sqrt{\frac{N}{2\Omega - N}} \approx U' V' \left(1 + O\left(\frac{N}{2\Omega}\right) \right), \quad (4.A.16)$$

for a number of particles considerably smaller than the full degeneracy of the single-particle subspace in which nucleons can correlate, that is for $N \ll 2\Omega$. Consequently, one can write

$$|\tilde{0}\rangle \approx \frac{1}{\sqrt{N_{norm}}} \sum_{\nu>0} (\alpha'_0)_\nu P_\nu^\dagger |0\rangle, \quad (4.A.17)$$

where

$$(\alpha'_0)_\nu = \langle BCS | P_\nu^\dagger | BCS \rangle = U'_\nu V'_\nu, \quad (4.A.18)$$

and

$$N_{norm} = \sum_{\nu>0} (\alpha'_0)_\nu^2. \quad (4.A.19)$$

Chapter 5

One-particle transfer

... physics is an experimental science: it is concerned only with those statements which in some sense can be verified by an experiment. The purpose of the theory is to provide an unification, a codification, or however you want to say it, of those results which can be tested by means of some experiment.

J. Schwinger

In what follows we present a derivation of the one-particle transfer differential cross section within the framework of the distorted wave Born approximation (DWBA)¹.

The structure input in the calculations are, as a rule, single-particle states dressed, within the formalism of nuclear field theory, through the coupling with the variety of collective, (quasi-) bosonic vibrations, leading to renormalized energies, particle content and radial wavefunctions. With the help of the associated modified form-factors², and of global optical potentials, one can calculate the absolute differential cross sections, quantities which can be directly compared with the experimental findings.

In this way one avoids to introduce, let alone use spectroscopic factors, quantities which are rather elusive to calculate consistently³. This is in keeping with the fact that as a nucleon moves through the nucleus it feels the presence of the other

¹See Tobocman (1961), Austern (1963), Jackson (1970) Satchler (1980), Satchler (1983), Austern (1970), Glendenning (2004), Thompson and Nunes (2009), and refs. therein.

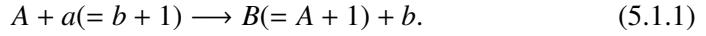
²See also Pinkston and Satchler (1982); Vaagen et al. (1979); Bang et al. (1980); Hamamoto (1970a) and refs. therein. See also Barranco et al. (2017).

³Duguet and Hagen (2012); Jenning (2011); Dickhoff and Barbieri (2004); Dickhoff and Van Neck (2005), and refs. therein. See also Feshbach (1992); Tamura (1974); Tamura et al. (1980).

nucleons whose configurations change as the motion progresses. It takes time for this information to be feed back on the nucleon. This renders the average potential nonlocal in time⁴. A time-dependent operator can always be transformed into an energy-dependent operator, implying an ω -dependence of the properties which are usually adscribed to particles like (effective) mass, charge, etc. Furthermore, due to Pauli principle, the average potential is also non local in space. Consequently, one is forced to deal with nucleons which carry around a cloud of (quasi) bosons, aside from exchanging its position with that of the other nucleons, and thus with renormalized energies, single-particle amplitude content and radial wavefunctions (formfactors) which eventually result in a *dynamical shell model*. It is of notice that the above mentioned phenomena are not only found in nuclear physics, but are universal within the framework of many-body systems as well as of field theories like quantum electrodynamics (QED). In fact, a basic result of such theories is that nothing is really free⁵. Within this context, in Section 5.2.1 we provide examples of one-particle transfer processes in nuclei lying along the stability valley (Sn-isotopes), where strongly renormalized quasiparticle states are populated, while in Sections 5.2.2–5.2.4 we give examples of exotic, halo nuclei. Namely of the isotones ^{11}Be and ^{10}Li , displaying also strong renormalization effects, as testified by the phenomenon of parity inversion.

5.1 General derivation

We now proceed to derive the transition amplitude for the reaction (Fig. 5.1.1).



For a simplified version –no recoil plus plane wave limit (see also Sect. 5.1.2)– we refer to Sect. 5.6, while for an alternative derivation within the framework of one-particle knock-out reactions we refer to Sect. 5.7.

Let us assume that the nucleon, bound initially to the core b is in a single-particle state with orbital and total angular momentum l_i and j_i respectively, and that the nucleon in the final state (bound to core A) is in the l_f, j_f state. The total spin and magnetic quantum numbers of nuclei A, a, B, b are $\{J_A, M_A\}, \{J_a, M_a\}, \{J_B, M_B\}, \{J_b, M_b\}$ respectively. Denoting ξ_A and ξ_b the intrinsic coordinates of the wavefunctions describing the structure of nuclei A and b , and \mathbf{r}_{An} and \mathbf{r}_{bn} the relative coordinates of the transferred nucleon with respect to the CM of nuclei A and b respectively, one can write the “intrinsic” wavefunctions of the colliding nuclei A, a as

$$\begin{aligned} & \phi_{M_A}^{J_A}(\xi_A), \\ & \Psi(\xi_b, \mathbf{r}_{b1}) = \sum_{m_i} \langle J_b \ j_i \ M_b \ m_i | J_a \ M_a \rangle \phi_{M_b}^{J_b}(\xi_b) \psi_{m_i}^{j_i}(\mathbf{r}_{bn}, \sigma), \end{aligned} \quad (5.1.2)$$

⁴See Mahaux et al. (1985) and references therein, see also Sect. 5.8.

⁵Feynman (1975).

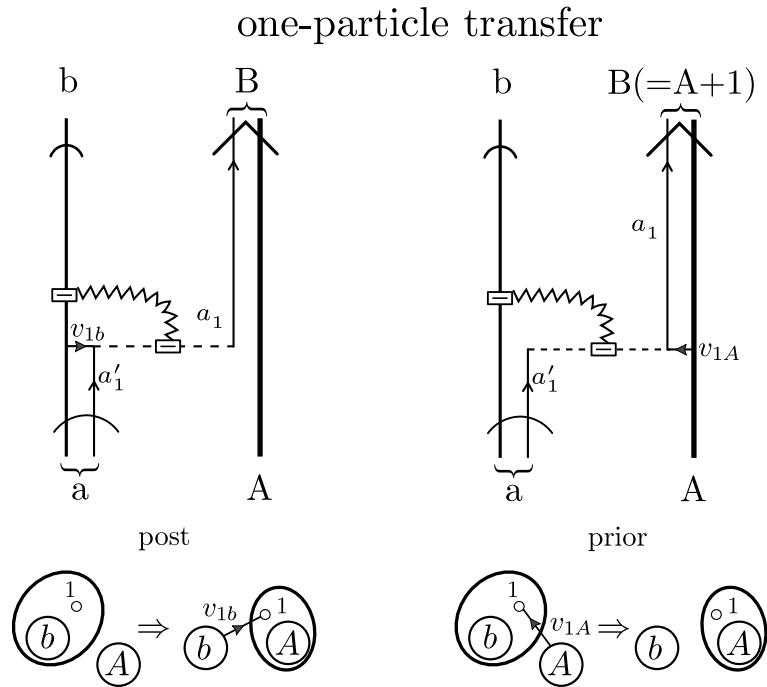


Figure 5.1.1: NFT graphical representation of the one-particle transfer reaction $a(=b+1)+A \rightarrow b+B(=A+1)$. The time arrow is assumed to point upwards. The quantum numbers characterizing the states in which the transferred nucleon moves in projectile and target are denoted a'_1 and a_1 respectively. The interaction inducing the nucleon to be transferred can act either in the entrance channel $((a, A); v_{1A}$, prior representation) or in the exit channel $((b, B); v_{1b}$, post representation), in keeping with energy conservation. In the transfer process, the nucleon changes orbital at the same time that a change in the mass partition takes place. The corresponding relative motion mismatch is known as the recoil process, and is represented by a jagged curve (this is the recoil elementary mode, mode which couples to the particle degrees of freedom through a Galilean transformation operator). The recoil mode provides information on the evolution of r_{1A} (r_{1b}). In other words, on the coupling between structure and reaction (relative motion) degrees of freedom.

while the “intrinsic” wavefunctions describing the structure of nuclei B and b are

$$\begin{aligned} \phi_{M_b}^{J_b}(\xi_b), \\ \Psi(\xi_A, \mathbf{r}_{A1}) = \sum_{m_f} \langle J_A \ j_f \ M_A \ m_f | J_B \ M_B \rangle \phi_{M_A}^{J_A}(\xi_A) \psi_{m_f}^{j_f}(\mathbf{r}_{An}, \sigma). \end{aligned} \quad (5.1.3)$$

For an unpolarized incident beam (sum over M_A, M_a and divide by $(2J_A + 1), (2J_a + 1)$), and assuming that one does not detect the final polarization (sum over M_B, M_b), the differential cross section in the DWBA can be written as

$$\begin{aligned} \frac{d\sigma}{d\Omega} = \frac{k_f}{k_i} \frac{\mu_i \mu_f}{4\pi^2 \hbar^4} \frac{1}{(2J_A + 1)(2J_a + 1)} \\ \times \sum_{\substack{M_A, M_a \\ M_B, M_b}} \left| \sum_{m_i, m_f} \langle J_b \ j_i \ M_b \ m_i | J_a \ M_a \rangle \langle J_A \ j_f \ M_A \ m_f | J_B \ M_B \rangle T_{m_i, m_f} \right|^2, \end{aligned} \quad (5.1.4)$$

where k_i and k_f are the relative motion linear momentum in both initial and final channels (flux), while μ_i and μ_f are the corresponding relative masses. The two quantities within $\langle \rangle$ brackets are Clebsch–Gordan coefficients taking care of angular momentum conservation⁶.

The transition amplitude T_{m_i, m_f} is

$$T_{m_i, m_f} = \sum_{\sigma} \int d\mathbf{r}_f d\mathbf{r}_{bn} \chi^{(-)*}(\mathbf{r}_f) \psi_{m_f}^{j_f*}(\mathbf{r}_{An}, \sigma) V(r_{bn}) \psi_{m_i}^{j_i}(\mathbf{r}_{bn}, \sigma) \chi^{(+)}(\mathbf{r}_i), \quad (5.1.5)$$

where

$$\psi_{m_i}^{j_i}(\mathbf{r}_{An}, \sigma) = u_{j_i}(r_{bn}) \left[Y^{l_i}(\hat{r}_i) \chi(\sigma) \right]_{j_i m_i}, \quad (5.1.6)$$

is the single-particle wavefunction describing the motion of the nucleon to be transferred, when in the initial state, u , Y and χ being the radial, angular (spherical harmonics) and spin components. Similarly for $\psi_{m_f}^{j_f}$. The distorted waves describing the relative motion of the incoming projectile and of the target nucleus and of the outgoing system and the residual nucleus are,

$$\chi^{(+)}(\mathbf{k}_i, \mathbf{r}_i) = \frac{4\pi}{k_i r_i} \sum_l i^l e^{i\sigma_i^l} g_l(\hat{r}_i) \left[Y^l(\hat{r}_i) Y^l(\hat{k}_i) \right]_0^0, \quad (5.1.7)$$

and

$$\chi^{(-)*}(\mathbf{k}_f, \mathbf{r}_f) = \frac{4\pi}{k_f r_f} \sum_l i^{-l} e^{i\sigma_f^l} f_l(\hat{r}_f) \left[Y^l(\hat{r}_f) Y^l(\hat{k}_f) \right]_0^0, \quad (5.1.8)$$

respectively. In the above relations f and g are, respectively, the solutions of the radial Schrödinger equation describing the relative motion associated with the corresponding optical potential (“elastic” scattering) in entrance and exit channel. Let

⁶Brink and Satchler (1968) and Edmonds (1960), also Bohr and Mottelson (1969).

us now discuss the angular components involved in the reaction process, starting with the relation

$$\begin{aligned} \left[Y^l(\hat{r}_f) Y^l(\hat{k}_f) \right]_0^0 \left[Y^{l'}(\hat{r}_i) Y^{l'}(\hat{k}_i) \right]_0^0 &= \sum_K ((ll)_0(l'l')_0 | (ll')_K (ll')_K)_0 \\ &\times \left\{ \left[Y^l(\hat{r}_f) Y^{l'}(\hat{r}_i) \right]^K \left[Y^l(\hat{k}_f) Y^{l'}(\hat{k}_i) \right]^K \right\}_0^0. \end{aligned} \quad (5.1.9)$$

The 9*j*-symbol can be explicitly evaluated to give,

$$((ll)_0(l'l')_0 | (ll')_K (ll')_K)_0 = \sqrt{\frac{2K+1}{(2l+1)(2l'+1)}}, \quad (5.1.10)$$

while the coupled expression can be written as

$$\begin{aligned} \left\{ \left[Y^l(\hat{r}_f) Y^{l'}(\hat{r}_i) \right]^K \left[Y^l(\hat{k}_f) Y^{l'}(\hat{k}_i) \right]^K \right\}_0^0 &= \sum_M \langle K \ K \ M \ -M | 0 \ 0 \rangle \left[Y^l(\hat{r}_f) Y^{l'}(\hat{r}_i) \right]_M^K \\ &\times \left[Y^l(\hat{k}_f) Y^{l'}(\hat{k}_i) \right]_{-M}^K = \sum_M \frac{(-1)^{K+M}}{\sqrt{2K+1}} \left[Y^l(\hat{r}_f) Y^{l'}(\hat{r}_i) \right]_M^K \left[Y^l(\hat{k}_f) Y^{l'}(\hat{k}_i) \right]_{-M}^K. \end{aligned} \quad (5.1.11)$$

Thus,

$$\begin{aligned} \left[Y^l(\hat{r}_f) Y^l(\hat{k}_f) \right]_0^0 \left[Y^{l'}(\hat{r}_i) Y^{l'}(\hat{k}_i) \right]_0^0 &= \sum_{K,M} \frac{(-1)^{K+M}}{\sqrt{(2l+1)(2l'+1)}} \left[Y^l(\hat{r}_f) Y^{l'}(\hat{r}_i) \right]_M^K \left[Y^l(\hat{k}_f) Y^{l'}(\hat{k}_i) \right]_{-M}^K. \end{aligned} \quad (5.1.12)$$

For the angular integral to be different from zero, the integrand must be coupled to zero angular momentum (scalar). Noting that the only variables over which one integrates in the above expression are \hat{r}_i, \hat{r}_f , we have to couple the remaining functions of the angular variables, namely the wavefunctions $\psi_{m_f}^{j_f*}(\mathbf{r}_{An}, \sigma) = (-1)^{j_f-m_f} \psi_{-m_f}^{j_f}(\mathbf{r}_{An}, -\sigma)$ and $\psi_{m_i}^{j_i}(\mathbf{r}_{bn}, \sigma)$ to angular momentum K , as well as to fulfill $M = m_f - m_i$. Let us then consider

$$\begin{aligned} (-1)^{j_f-m_f} \psi_{-m_f}^{j_f}(\mathbf{r}_{An}, -\sigma) \psi_{m_i}^{j_i}(\mathbf{r}_{bn}, \sigma) &= (-1)^{j_f-m_f} u_{j_f}(r_{An}) u_{j_i}(r_{bn}) \\ &\times \sum_P \langle j_f \ j_i \ -m_f \ m_i | P \ m_i - m_f \rangle \left\{ \left[Y^{l_f}(\hat{r}_{An}) \chi^{1/2}(-\sigma) \right]^{j_f} \left[Y^{l_i}(\hat{r}_{bn}) \chi^{1/2}(\sigma) \right]^{j_i} \right\}_{m_i-m_f}^P. \end{aligned} \quad (5.1.13)$$

Recoupling the spherical harmonics to angular momentum K and the spinors to $S = 0$, only one term survives the angular integral in (5.1.5), namely

$$\begin{aligned} & (-1)^{j_f - m_f} u_{j_f}(r_{An}) u_{j_i}(r_{bn}) ((l_f \frac{1}{2})_{j_f} (l_i \frac{1}{2})_{j_i} | (l_f l_i)_K (\frac{1}{2} \frac{1}{2})_0)_K \\ & \times \langle j_f \ j_i \ -m_f \ m_i | K \ m_i - m_f \rangle [Y^{l_f}(\hat{r}_{An}) Y^{l_i}(\hat{r}_{bn})]_{m_i - m_f}^K [\chi(-\sigma) \chi(\sigma)]_0^0. \end{aligned} \quad (5.1.14)$$

Making use of the fact that the sum over spins yields a factor $-\sqrt{2}$, and in keeping with the fact that $M = m_f - m_i$, one obtains,

$$\begin{aligned} T_{m_i, m_f} = & (-1)^{j_f - m_f} \frac{-16 \sqrt{2} \pi^2}{k_f k_i} \sum_{l'} i'^{-l} e^{\sigma'_f + \sigma'_i} \sum_K ((l_f \frac{1}{2})_{j_f} (l_i \frac{1}{2})_{j_i} | (l_f l_i)_K (\frac{1}{2} \frac{1}{2})_0)_K \\ & \times \langle j_f \ j_i \ -m_f \ m_i | K \ m_i - m_f \rangle [Y^l(\hat{k}_f) Y^{l'}(\hat{k}_i)]_{m_i - m_f}^K \int d\mathbf{r}_f d\mathbf{r}_{bn} \frac{f_l(r_f) g_{l'}(r_i)}{r_f r_i} \\ & \times u_{j_f}(r_{An}) u_{j_i}(r_{bn}) V(r_{bn}) (-1)^{K+m_f-m_i} [Y^l(\hat{r}_f) Y^{l'}(\hat{r}_i)]_{m_f - m_i}^K [Y^{l_f}(\hat{r}_{An}) Y^{l_i}(\hat{r}_{bn})]_{m_i - m_f}^K. \end{aligned} \quad (5.1.15)$$

Again, the only term of the expression

$$\begin{aligned} & (-1)^{K+m_f-m_i} [Y^l(\hat{r}_f) Y^{l'}(\hat{r}_i)]_{m_f - m_i}^K [Y^{l_f}(\hat{r}_{An}) Y^{l_i}(\hat{r}_{bn})]_{m_i - m_f}^K = \\ & (-1)^{K+m_f-m_i} \sum_P \langle K \ K \ m_f - m_i \ m_i - m_f | P \ 0 \rangle \left\{ [Y^l(\hat{r}_f) Y^{l'}(\hat{r}_i)]^K [Y^{l_f}(\hat{r}_{An}) Y^{l_i}(\hat{r}_{bn})]^K \right\}_0^P \end{aligned}$$

which survives after angular integration is the one with $P = 0$, that is,

$$\begin{aligned} & \frac{1}{\sqrt{(2K+1)}} \left\{ [Y^l(\hat{r}_f) Y^{l'}(\hat{r}_i)]^K [Y^{l_f}(\hat{r}_{An}) Y^{l_i}(\hat{r}_{bn})]^K \right\}_0^0 \\ & = \frac{1}{\sqrt{(2K+1)}} \sum_{M_K} \langle K \ K \ M_K \ -M_K | 0 \ 0 \rangle [Y^l(\hat{r}_f) Y^{l'}(\hat{r}_i)]_{M_K}^K \\ & \times [Y^{l_f}(\hat{r}_{An}) Y^{l_i}(\hat{r}_{bn})]_{-M_K}^K = \frac{1}{\sqrt{(2K+1)}} \sum_{M_K} \frac{(-1)^{K+M_K}}{\sqrt{(2K+1)}} [Y^l(\hat{r}_f) Y^{l'}(\hat{r}_i)]_{M_K}^K \\ & \times [Y^{l_f}(\hat{r}_{An}) Y^{l_i}(\hat{r}_{bn})]_{-M_K}^K \\ & = \frac{1}{2K+1} \sum_{M_K} (-1)^{K+M_K} [Y^l(\hat{r}_f) Y^{l'}(\hat{r}_i)]_{M_K}^K [Y^{l_f}(\hat{r}_{An}) Y^{l_i}(\hat{r}_{bn})]_{-M_K}^K, \end{aligned}$$

an expression which is spherically symmetric. One can evaluate it for a particular configuration, for example setting $\hat{r}_f = \hat{z}$ and the center of mass A, b, n in the $x - z$ plane (see Fig. 5.1.2). Once the orientation in space of this “standard” configuration is specified (through, for example, a rotation $0 \leq \alpha \leq 2\pi$ around \hat{z} , a rotation $0 \leq \beta \leq \pi$ around the new x axis and a rotation $0 \leq \gamma \leq 2\pi$ around

\hat{r}_{bB}), the only remaining angular coordinate is θ , while the integral over the other three angles yields $8\pi^2$. Setting $\hat{r}_f = \hat{z}$ one obtains

$$\left[Y^l(\hat{r}_f) Y^{l'}(\hat{r}_i) \right]_{M_K}^K = \langle l' l' 0 M_K | K M_K \rangle \sqrt{\frac{2l+1}{4\pi}} Y_{M_K}^{l'}(\hat{r}_i). \quad (5.1.16)$$

Because of $M = m_i - m_f$, and $m = m_f$, $T_{m_i, m_f} \equiv T_{m, M}$ where

$$\begin{aligned} T_{m, M} &= (-1)^{j_f - m} \frac{-64\sqrt{2}\pi^{7/2}}{k_f k_i} \sum_{ll'} i^{l'-l} e^{\sigma_f^l + \sigma_i^{l'}} \sqrt{2l+1} \sum_K \frac{(-1)^K}{2K+1} ((l_f \frac{1}{2})_{j_f} (l_i \frac{1}{2})_{j_i} | (l_f l_i)_K (\frac{1}{2} \frac{1}{2})_0)_K \\ &\times \langle j_f j_i - m M + m | K M \rangle \left[Y^l(\hat{k}_f) Y^{l'}(\hat{k}_i) \right]_M^K \int d\mathbf{r}_f d\mathbf{r}_{bn} \frac{f_l(r_f) g_{l'}(r_i)}{r_f r_i} \\ &\times u_{j_f}(r_{An}) u_{j_i}(r_{bn}) V(r_{bn}) \sum_{M_K} (-1)^{M_K} \langle l' l' 0 M_K | K M_K \rangle \left[Y^{l_f}(\hat{r}_{An}) Y^{l_i}(\hat{r}_{bn}) \right]_{-M_K}^K Y_{M_K}^{l'}(\hat{r}_i). \end{aligned} \quad (5.1.17)$$

We now turn our attention to the sum

$$\sum_{\substack{M_A, M_a \\ M_B, M_b}} \left| \sum_{m, M} \langle J_b j_i M_b m | J_a M_a \rangle \langle J_A j_f M_A M | J_B M_B \rangle T_{m, M} \right|^2, \quad (5.1.18)$$

appearing in the expression for the differential cross section (5.1.4). For any given value m', M' of m, M , the sum will be

$$\begin{aligned} &\sum_{M_a, M_b} \left| \langle J_b j_i M_b m' | J_a M_a \rangle \right|^2 \sum_{M_A, M_B} \left| \langle J_A j_f M_A M' | J_B M_B \rangle \right|^2 \left| T_{m', M'} \right|^2 \\ &= \frac{(2J_a + 1)(2J_B + 1)}{(2j_i + 1)(2j_f + 1)} \sum_{M_a, M_b} \left| \langle J_b J_a M_b - M_a | j_i m' \rangle \right|^2 \\ &\times \sum_{M_A, M_B} \left| \langle J_A J_B M_A - M_B | j_f M' \rangle \right|^2 \left| T_{m', M'} \right|^2, \end{aligned} \quad (5.1.19)$$

by virtue of the symmetry property of Clebsch–Gordan coefficients

$$\langle J_b j_i M_b m | J_a M_a \rangle = (-1)^{J_b - M_b} \sqrt{\frac{(2J_a + 1)}{(2j_i + 1)}} \langle J_b J_a M_b - M_a | j_i m \rangle. \quad (5.1.20)$$

The sum over the Clebsch–Gordan coefficients in (5.1.19) is equal to 1, so (5.1.18) becomes

$$\frac{(2J_a + 1)(2J_B + 1)}{(2j_i + 1)(2j_f + 1)} \sum_{m, M} \left| T_{m, M} \right|^2, \quad (5.1.21)$$

and the differential cross section can be written as,

$$\frac{d\sigma}{d\Omega} = \frac{k_f}{k_i} \frac{\mu_i \mu_f}{4\pi^2 \hbar^4} \frac{(2J_B + 1)}{(2j_i + 1)(2j_f + 1)(2J_A + 1)} \sum_{m, M} \left| T_{m, M} \right|^2. \quad (5.1.22)$$

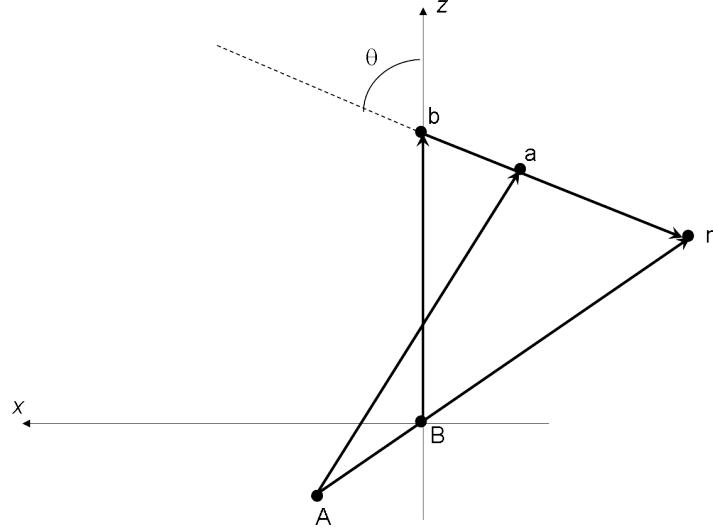


Figure 5.1.2: Coordinate system in the “standard” configuration. Note that $\mathbf{r}_f \equiv \mathbf{r}_{Bb}$, and $\mathbf{r}_i \equiv \mathbf{r}_{Aa}$.

where

$$T_{m,M} = \sum_{Kll'} (-1)^{-m} \langle j_f j_i - m | M + m | K M \rangle \left[Y^l(\hat{k}_f) Y^{l'}(\hat{k}_i) \right]_M^K t_{ll'}^K. \quad (5.1.23)$$

Orienting \hat{k}_i along the incident z -direction leads to,

$$\left[Y^l(\hat{k}_f) Y^{l'}(\hat{k}_i) \right]_M^K = \langle l l' M 0 | K M \rangle \sqrt{\frac{2l'+1}{4\pi}} Y_M^l(\hat{k}_f), \quad (5.1.24)$$

and

$$T_{m,M} = \sum_{Kll'} (-1)^{-m} \langle l l' M 0 | K M \rangle \langle j_f j_i - m | M + m | K M \rangle Y_M^l(\hat{k}_f) t_{ll'}^K, \quad (5.1.25)$$

with

$$\begin{aligned} t_{ll'}^K &= (-1)^{K+j_f} \frac{-32\sqrt{2}\pi^3}{k_f k_i} i^{l'-l} e^{\sigma_f^l + \sigma_i^{l'}} \frac{\sqrt{(2l+1)(2l'+1)}}{2K+1} ((l_f \frac{1}{2})_{j_f} (l_i \frac{1}{2})_{j_i} | (l_f l_i)_K (\frac{1}{2} \frac{1}{2})_0)_K \\ &\times \int dr_f dr_{bn} d\theta r_{bn}^2 \sin \theta r_f \frac{f_l(r_f) g_{l'}(r_i)}{r_i} u_{j_f}(r_{An}) u_{j_i}(r_{bn}) V(r_{bn}) \\ &\times \sum_{M_K} (-1)^{M_K} \langle l l' 0 | M_K | K M_K \rangle \left[Y^{l_f}(\hat{r}_{An}) Y^{l_i}(\hat{r}_{bn}) \right]_{-M_K}^K Y_{M_K}^{l'}(\hat{r}_i). \end{aligned} \quad (5.1.26)$$

5.1.1 Coordinates

To perform the integral in (5.1.26), one needs the expression of $r_i, r_{An}, \hat{r}_{An}, \hat{r}_{bn}, \hat{r}_i$ in term of the integration variables r_f, r_{bn}, θ . Because one is interested in evaluating these quantities in the particular configuration depicted in Fig. 5.1.2, one has

$$\mathbf{r}_f = r_f \hat{z}, \quad (5.1.27)$$

$$\mathbf{r}_{bn} = -r_{bn}(\sin \theta \hat{x} + \cos \theta \hat{z}), \quad (5.1.28)$$

$$\mathbf{r}_{Bn} = \mathbf{r}_f + \mathbf{r}_{bn} = -r_{bn} \sin \theta \hat{x} + (r_f - r_{bn} \cos \theta) \hat{z}. \quad (5.1.29)$$

One can then write

$$\mathbf{r}_{An} = \frac{A+1}{A} \mathbf{r}_{Bn} = -\frac{A+1}{A} r_{bn} \sin \theta \hat{x} + \frac{A+1}{A} (r_f - r_{bn} \cos \theta) \hat{z}, \quad (5.1.30)$$

$$\mathbf{r}_{an} = \frac{b}{b+1} \mathbf{r}_{bn} = -\frac{b}{b+1} r_{bn} (\sin \theta \hat{x} + \cos \theta \hat{z}), \quad (5.1.31)$$

and

$$\mathbf{r}_i = \mathbf{r}_{An} - \mathbf{r}_{an} = -\frac{2A+1}{(A+1)A} r_{bn} \sin \theta \hat{x} + \left(\frac{A+1}{A} r_f - \frac{2A+1}{(A+1)A} r_{bn} \cos \theta \right) \hat{z}, \quad (5.1.32)$$

where A, b are the number of nucleons of nuclei A and b respectively.

5.1.2 Zero-range approximation

In the zero range approximation,

$$\int dr_{bn} r_{bn}^2 u_{j_i}(r_{bn}) V(r_{bn}) = D_0; \quad u_{j_i}(r_{bn}) V(r_{bn}) = \delta(r_{bn})/r_{bn}^2. \quad (5.1.33)$$

It can be shown (see Fig. 5.1.2) that for $r_{bn} = 0$

$$\mathbf{r}_{An} = \frac{m_A + 1}{m_A} \mathbf{r}_f, \quad \mathbf{r}_i = \frac{m_A + 1}{m_A} \mathbf{r}_f. \quad (5.1.34)$$

One then obtains

$$t_{ll'}^K = \frac{-16\sqrt{2}\pi^2}{k_f k_i} (-1)^K \frac{D_0}{\alpha} i'^{-l} e^{\sigma_f^l + \sigma_i'^l} \frac{\sqrt{(2l+1)(2l'+1)(2l_i+1)(2l_f+1)}}{2K+1} ((l_f \frac{1}{2})_{j_f} (l_i \frac{1}{2})_{j_i} | (l_f l_i)_K (\frac{1}{2} \frac{1}{2})_0)_K \\ \times \langle l' l' 0 0 | K 0 \rangle \langle l_f l_i 0 0 | K 0 \rangle \int dr_f f_l(r_f) g_{l'}(\alpha r_f) u_{j_f}(\alpha r_f), \quad (5.1.35)$$

with

$$\alpha = \frac{A+1}{A}. \quad (5.1.36)$$

5.2 Examples and Applications

In the calculation of absolute reaction cross sections two elements melt together: reaction and structure. In the case of weakly coupled probes like, as a rule, direct one-particle transfer processes are, the first element can be further divided into two essentially separated components: elastic scattering (optical potentials), and transfer amplitudes connecting entrance and exit channels. In other words, the habitat of DWBA.

5.2.1 $^{120}\text{Sn}(p, d)^{119}\text{Sn}$ and $^{120}\text{Sn}(d, p)^{121}\text{Sn}$ reactions⁷.

The NFT calculation of the structure properties of the Sn-isotopes probed through (d, p) and (p, d) reactions, were carried out making use of an effective Skyrme interaction⁸, SLy4, to determine the mean field (HFB) and the collective modes (QRPA), and a $v_{14}(^1S_0) (\equiv v_p^{bare})$ Argonne pairing interaction⁹. Hartree-Fock-Bogoliubov (HFB) results provide the bare quasiparticle spectrum, while QRPA a realization of density ($J^\pi = 2^+, 3^-, 4^+, 5^-$) and spin ($2^\pm, 3^\pm, 4^\pm, 5^\pm$) modes.

Taking into account renormalization processes (self-energy, vertex corrections) in terms of the PVC mechanism, the dressed particles and the induced pairing interaction v_p^{ind} were calculated. Adding it to v_p^{bare} the induced pairing interaction v_p^{ind} , the total pairing interaction v_p^{eff} was determined. It is found that v_p^{bare} and v_p^{ind} contribute about equally to the state dependent pairing gap.

Making use of the above input, the lowest order NFT renormalizing diagrams (self-energy and vertex correction) were propagated to infinite order making use of Nambu-Gorkov equation¹⁰. In Fig. 5.2.1, the absolute differential cross section associated with the population of the low-lying state $|^{119}\text{Sn}(7/2^+; 788\text{keV})\rangle$ in the one-particle pick-up process $^{120}\text{Sn}(p, d)^{119}\text{Sn}$ and worked out with the help of the software ONE¹¹, of global optical parameters¹² and of NFT spectroscopic amplitudes, is compared with the experimental data.

Similar calculations have been carried for the reaction $^{120}\text{Sn}(d, p)^{121}\text{Sn} (J^\pi; E_x)$ in connection with the population of the $|3/2^+; \text{gs}\rangle$ and $|11/2^-; E_x \approx 0 \text{ MeV}\rangle$ states. In the stripping experiment¹³ the ground state and the $11/2^-$ state were not resolved in energy. This is the reason why theory and experiment are compared to the data for the summed $l = 2 + 5$ differential cross section (see Fig. 5.2.2 (a)), the separate theoretical predictions been displayed in Figs. 5.2.2 (b) and (c).

Let us now turn to the most fragmented low-lying quasiparticle state around ^{120}Sn , namely that associated with the $d_{5/2}$ orbital¹⁴. As shown in the right panel of

⁷Idini et al. (2015).

⁸Chabanat et al. (1997).

⁹Wiringa et al. (1984).

¹⁰For details see Idini et al. (2015).

¹¹Potel (2012a); see also App. 7.A.

¹²Dickey et al. (1982).

¹³Bechara and Dietzsch (1975).

¹⁴cf. Idini (2013), Idini et al. (2015).

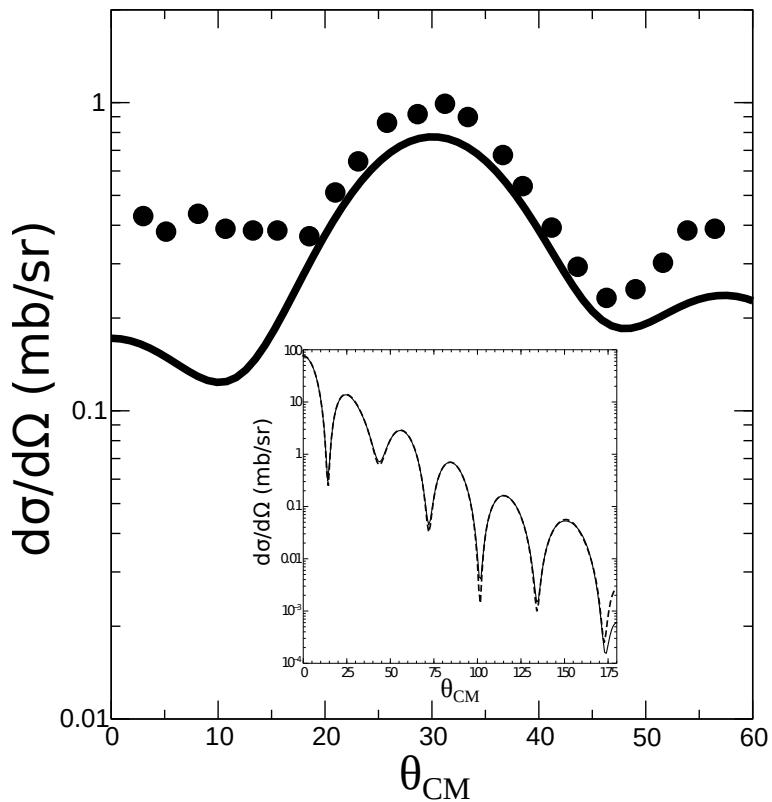


Figure 5.2.1: The absolute differential cross section $^{120}\text{Sn}(p, d)^{119}\text{Sn}(j^\pi)$ associated with the state $j^\pi = 7/2^+$. The theoretical prediction (continuous curve) is displayed in comparison with the experimental data (solid dots, Dickey et al. (1982)). The corresponding integrated cross sections are 5.0 and 5.2 ± 0.6 mb respectively. In the inset, and for the sake of accuracy control, the absolute differential cross sections associated with the reaction $^{124}\text{Sn}(p, d)^{123}\text{Sn}(\text{gs})$ calculated making use of the codes **ONE** (bold dashed) and **FRESCO** (thin continuous line (Thompson (1988))) are displayed.

Fig. 5.2.3 five low-lying $5/2^+$ states have been populated in the reaction $^{120}\text{Sn}(p, d)^{119}\text{Sn}$ with a summed cross section¹⁵ $\sum_{i=1}^5 \sigma(2^\circ - 25^\circ) \approx 8 \text{ mb} \pm 2 \text{ mb}$ while four are theoretically predicted with $\sum_{i=1}^4 \sigma(2^\circ - 25^\circ) = 6.2 \text{ mb}$. Within the present context, namely that of probing the single-particle content of an elementary excitation (coupling to doorway states¹⁶), the study of the $5/2^+$ quasiparticle strength is a rather trying situation, providing a measure of the limitations encountered in such studies.

Analysis of the type presented above allows one to posit that structure and reactions are but just two aspects of the same physics. If one adds to this picture the fact that the optical potential –that is, the energy and momentum dependent nuclear dielectric function describing the medium where direct nuclear chemistry reactions take place– can be calculated microscopically¹⁷ in terms of the same elements entering structure calculations (i.e. spectroscopic amplitudes, renormalized single-particle wavefunctions and transition densities and thus effective formfactors), the structure-reaction circle closes itself. Allowing halo nuclei to be part of the daily nuclear structure paradigm, the equivalence between structure and reactions becomes even stronger, in keeping with the central role the continuum plays in the structure of these nuclei¹⁸.

It seems then fair to state that the importance of the coupled channels approach to reactions¹⁹ is not so much, or at least not only, that it is able to handle situations like for example one-particle transfer to members of a quadrupole rotational band, alas at the expenses of including effects to all orders which can be treated in lowest one²⁰, but that it reminds us how intimately connected probed and probe are in nuclei.

On the other hand, for most of the situations dealt with in the present monograph, it is transparent the power, also to reflect the physics, of the approach based in perturbative DWBA (e.g. 1st order for one-nucleon transfer and 2nd for Cooper pair tunneling), coupled with NFT elementary modes of nuclear excitation. To which extent a FRESCO like software built on a NFT basis will ever be attempted is an open question. Note in any case the important attempts made at incorporating so called core excitations within the FRESCO framework²¹.

We conclude this section by recalling the fact that the dressing of single particles with pairing vibrations plays also a central role in the structure properties of

¹⁵Dickey et al. (1982).

¹⁶Feshbach (1958), Rawitscher (1987), Bortignon and Broglia (1981), Bertsch et al. (1983).

¹⁷See footnote 23 Ch. 1.

¹⁸For example, in the case of ^{10}Li (see Sect. 5.2.4) in which one is confronted with continuum spectroscopy, the self-energy functional describing the interweaving of the unbound neutron with vibrations of the core (^9Li), contribute both to the nuclear structure and nuclear reaction dielectric function.

¹⁹Thompson (1988), Thompson (2013), Tamura et al. (1970), Asciutto and Glendenning (1969), Asciutto and Glendenning (1970), Asciutto et al. (1971), Asciutto and Sørensen (1972); cf. also Fernández-García et al. (2010a), Fernández-García et al. (2010b).

²⁰Within this context see e.g. the couple channel Born approximation treatment of $^{208}\text{Pb}(t, p)^{210}\text{Pb}(3^-)$ carried out in Flynn et al. (1972b).

²¹Fernández-García et al. (2010a), Fernández-García et al. (2010b).

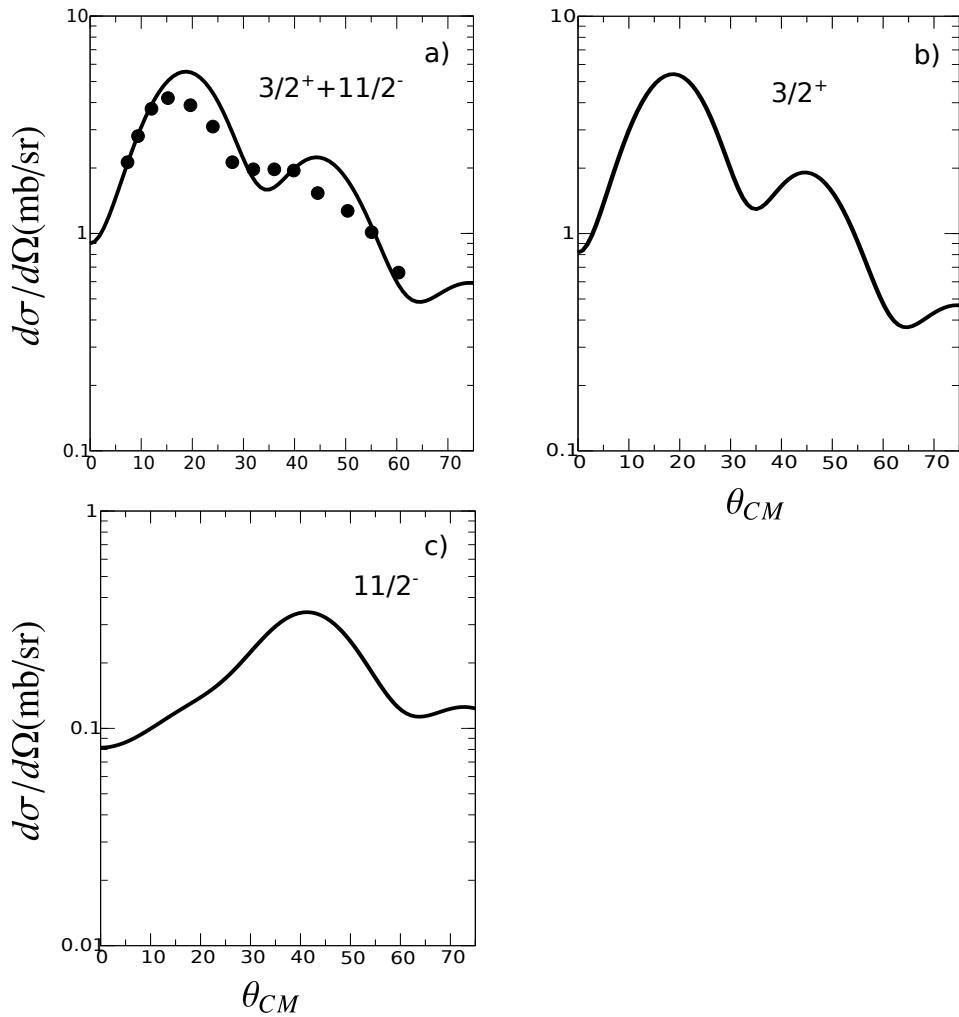


Figure 5.2.2: The theoretical absolute differential cross section (continuous curve) associated with the reaction $^{120}\text{Sn}(d, p)^{121}\text{Sn}$ populating the low-lying states $3/2^+$ and $11/2^-$ are shown in b) and c), while the incoherent summed differential cross section is displayed in a) in comparison with the data (Bechara and Dietzsche (1975)).

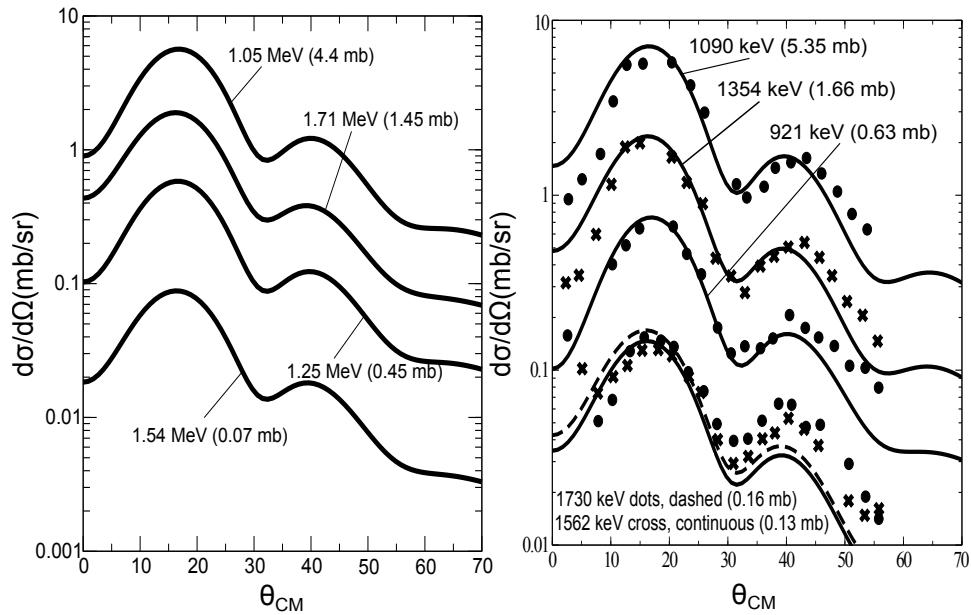


Figure 5.2.3: $^{120}\text{Sn}(p, d)^{119}\text{Sn}(5/2^+)$ absolute experimental cross sections (solid dots, Dickey et al. (1982)), together with the DWBA fit carried out in the analysis of the data (right panel) in comparison with the finite range, full recoil DWBA calculations carried out with the same global optical potentials making use of NFT structure inputs as explained in the text (after Idini et al. (2015)) and of the software ONE (Potel (2012a)); see also App. 7.A.

nuclei²².

5.2.2 NFT of ^{11}Be : one-particle transfer in halo nuclei

The nucleus $^{11}_{\text{4}}\text{Be}_7$ constitutes an example of one-neutron halo system, namely a halo neutron outside the $N = 6$ closed shell resulting from the phenomenon of parity inversion²³

Outlook²⁴

In the core of ^{11}Be , namely $^{10}_{\text{4}}\text{Be}_6$, six neutrons occupy the $1s_{1/2}$ and $1p_{3/2}$ levels (Fig. 5.2.4). The dominant ZPF is of quadrupole type, the main neutron component being the $((p_{1/2}, p_{3/2}^{-1}) \otimes 2^+)_0^+$ one. Because $\epsilon_{p1/2} - \epsilon_{p3/2} \approx 3.35$ MeV and $\hbar\omega_{2^+} = 3.368$ MeV, the largest amplitude of the quadrupole mode is associated with the particle-hole excitation $(p_{1/2}, p_{3/2}^{-1})_{2^+}$. The repulsion felt by a neutron moving in the $p_{1/2}$ state due to Pauli principle correction is ≈ 2.8 MeV (Fig. 5.2.4 inset (A)). The clothing of the $2s_{1/2}$ bare level by the quadrupole mode (Fig. 5.2.4 inset (B)) makes it heavier, lowering its energy by almost one-half MeV (0.43 MeV). The result of the two processes is parity inversion and the appearance of the $N = 6$ new magic number together with the melting away of the $N = 8$ standard one. In a similar way in which the Lamb shift (Fig. 5.2.4, inset (C)) provides a measure of the fluctuations of the QED vacuum²⁵, parity inversion measures ZPF of the nuclear vacuum (ground) state.

Calculations

In the calculations one has simultaneously dealt with the $1p_{3/2}$, $1p_{1/2}$, $2s_{1/2}$ and $1d_{5/2}$ valence single-particle states, treating their interweaving with the low-lying quadrupole collective vibration of the ^{10}Be core and the mixing between bound and continuum states. The bare energies of the single-particle orbitals were determined by freely varying the depth, diffusivity, radius and spin-orbit strength of a Woods-Saxon potential so that, making use of an effective radial dependent effective mass ($m_k(r=0) = 0.7m$, $m_k(r=\infty) = 0.91m$), the fully dressed, renormalized energies best reproduce the experimental findings²⁶.

²²Barranco et al. (1987), Bès et al. (1988), Baroni et al. (2004).

²³This nucleus has been extensively studied both experimentally (see Iwasaki et al. (2000); Fortier et al. (1999); Winfield et al. (2001); Auton (1970); Zwieglinski et al. (1979); Schmitt et al. (2013); Nörterhäuser et al. (2009); Kwan et al. (2014) and references therein) and theoretically (see Talmi and Unna (1960); Otsuka et al. (1993); Sagawa et al. (1993); Vinh Mau (1995); Gori et al. (2004); Nunes et al. (1996); Fos; Hamamoto and Shimoura (2007); Kanada-En'yo and Horiuchi (2002); Calci et al. (2016); Krieger et al. (2012); Timofeyuk and Johnson (1999); Keeley et al. (2004); Deltuva (2009, 2013); Lay et al. (2014); de Diego et al. (2014) and references therein).

²⁴Barranco et al. (2017).

²⁵Pais (1986) p. 451.

²⁶Namely, parallel to mass renormalization in QED (see Sect. 5.3.3).

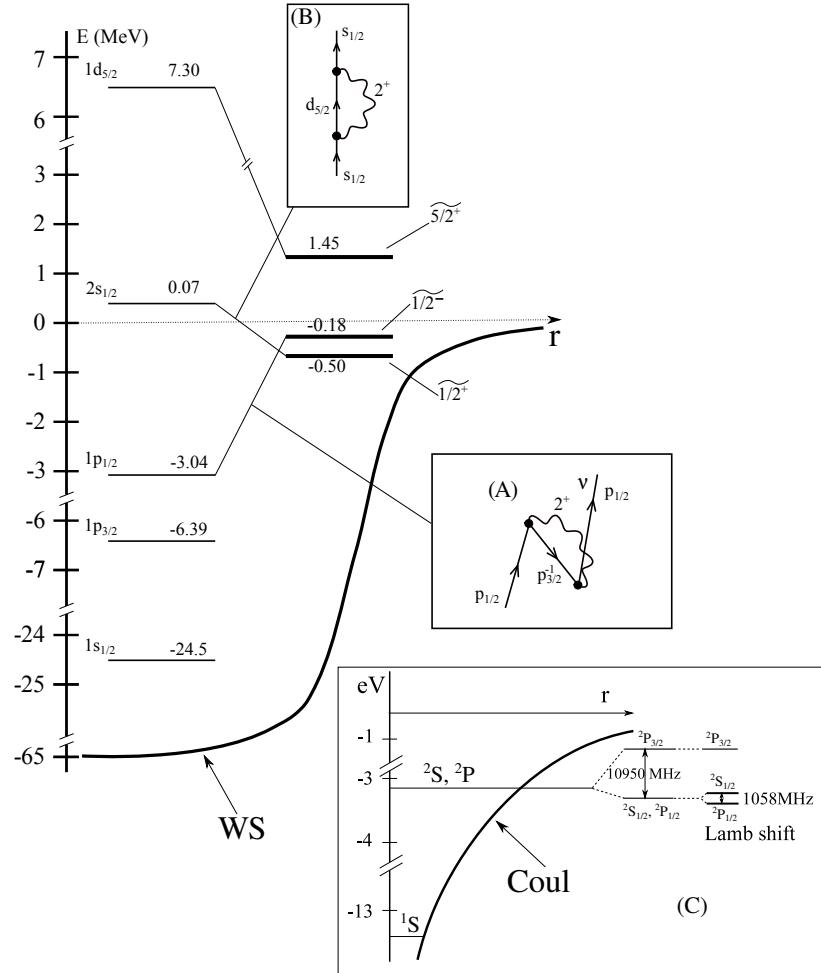


Figure 5.2.4: Bare ϵ_j (upper left thin horizontal lines) and dressed $\tilde{\epsilon}_j$ (bold face) single-particle levels of ^{11}Be . Due to the dressing of neutron motion with quadrupole vibrations of the core ^{10}Be (insets (A), (B) and (D)) inversion in sequence between the $2s_{1/2}$ and $1p_{1/2}$ levels (parity inversion) is observed. The numbers are energies in MeV. The Woods-Saxon (WS) bare mean field is indicated. In inset (C), the lowest energy levels of hydrogen are indicated, the Coulomb potential (Coul) is also schematically shown. The effects of the spin-orbit coupling and Lamb shift associated with the splitting of the $^2S_{1/2}$ and $^2P_{1/2}$ levels are displayed.

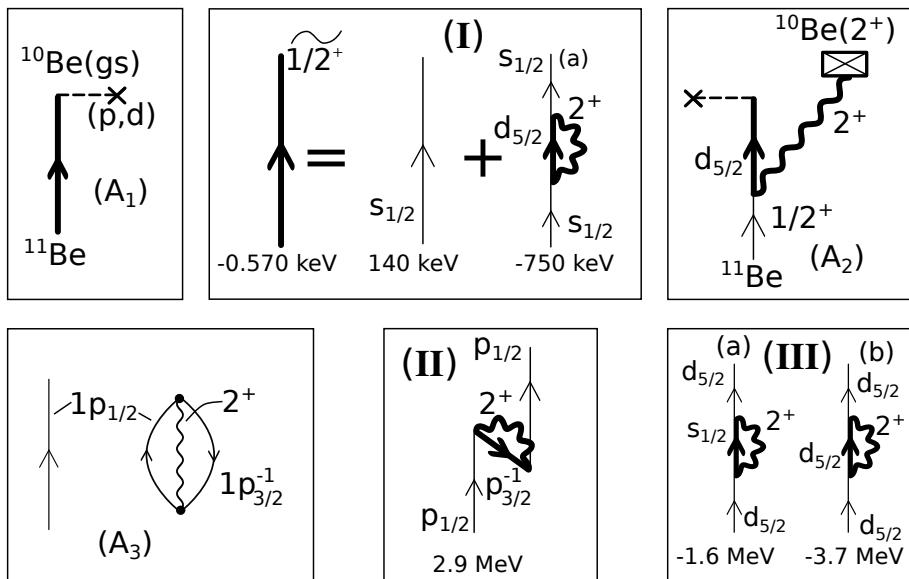


Figure 5.2.5: $(NFT)_{ren}$ diagrams describing the renormalization processes responsible for the different components of the clothed states (Eqs. (5.2.1)-(5.2.3)) associated with (I)-(III)) and the stripping and pickup processes populating the ground state of ^{11}Be (A_1) and the first excited 2^+ state (A_2) of ^{10}Be respectively. (A_3) Valence nucleon in presence of a virtual zero point fluctuation of the core ^{10}Be . Bold (thin) arrowed lines pointing upwards (downwards), describe dressed (bare) particle (hole) states. The wavy line represents the quadrupole vibration. A cross (crossed circle) followed by a horizontal dashed line stands for an external one-neutron pickup (p, d) (stripping (d, p)) field. A crossed box indicates a γ -detector, eventually revealing the decay of the quadrupole vibrational state of ^{10}Be .

The variety of self energy diagrams, renormalizing selfconsistently the motion of the odd neutron of ^{11}Be in both configuration– (Fig. 5.2.5) and conformational 3D-space (Fig. 5.2.6, form factors), through the coupling to quadrupole vibrations²⁷, have been worked out²⁸. The resulting states can be written as

$$|\widetilde{1/2^+}\rangle = \sqrt{0.80}|s_{1/2}\rangle + \sqrt{0.20}|(d_{5/2} \otimes 2^+)_{1/2^+}\rangle, \quad (5.2.1)$$

$$|\widetilde{1/2^-}\rangle = \sqrt{0.84}|(p_{1/2})\rangle + \sqrt{0.16}|(p_{1/2}, p_{3/2}^{-1})_{2^+} \otimes 2^+_{0+}, p_{1/2}\rangle, \quad (5.2.2)$$

and

$$|\widetilde{5/2^+}\rangle = \sqrt{0.49}|d_{5/2}\rangle + \sqrt{0.23}|(s_{1/2} \otimes 2^+)_{5/2^+}\rangle + \sqrt{0.28}|(d_{5/2} \otimes 2^+)_{5/2^+}\rangle. \quad (5.2.3)$$

The bare energies ϵ_j and the $((NFT)_{ren})$ values $\tilde{\epsilon}_j$ associated with the renormalised single-particle states are shown in Fig. 5.2.4. These last quantities reproduce quite accurately the experimental findings (Fig. 5.2.6, upper center). Also shown in this figure, are the corresponding wavefunctions $\phi_j(r)$ and $\tilde{\phi}_j(r)$. The form factors $\tilde{\phi}_j(r)$ were used, together with global optical potentials²⁹, to calculate the one-nucleon stripping and pickup absolute differential cross sections of the reactions $^{10}\text{Be}(d, p)^{11}\text{Be}(1/2^+, 1/2^-, \text{ and } 5/2^+)$ and $^{11}\text{Be}(p, d)^{10}\text{Be}(2^+)$. The results provide an overall account of the experimental findings.

We remark that the stripping process shown in inset (A₁) of Fig. 5.2.5 and populating ^{11}Be ground state implies the action of the external (d, p) field on the left hand side of the graphical representation of Dyson equation shown in Fig. 5.2.5(I), and involves, at the same time, the use of the corresponding radial wavefunction as form factor (Fig. 5.2.6 (a)). In the case of the population of the first 2^+ excited state of ^{10}Be (inset A₂), a (p, d) field acts on the $(d_{5/2} \otimes 2^+)_{1/2^+}$ virtual state of the second graph of the right hand side of Dyson equation (Fig. 5.2.5 (I) (a)), *involving this time the radial wave function $\tilde{\phi}_{1/2^+}^{(2+)}(r)$, namely the odd neutron moving around the quadrupole excited ^{10}Be core, as form factor* (Fig. 5.2.6 (d)).

Insets (A₁) and (A₂) and diagrams (I) of Fig. 5.2.5 testify to the subtle effects resulting from the unification of $(NFT)_{ren}$ of structure and reactions, and operative in the cross sections also shown in Fig. 5.2.6, as a result of the simultaneous and self consistent treatment of correlations in configuration and 3D-space. The bold face drawn state $|(d_{5/2} \otimes 2^+)_{1/2^+}\rangle$ shown in Fig. 5.2.5(I)(a) and the radial wave function $(NFT)_{ren}$ displayed with a continuous curve in Fig. 5.2.6 (d), can

²⁷It is of notice that the $d_{5/2}$ plays, through its coupling to the 2^+ state, an essential role in the parity inversion phenomenon, by lowering the energy of the $s_{1/2}$ state in a consistent way. The numbers quoted below also contain important contributions from Pauli principle correcting diagrams see Fig. 5.2.6 (IV) associated with the presence of two-phonon quadrupole states in intermediate, virtual configurations.

²⁸The results, obtained taking also into account the coupling to the octupole vibration and the pair removal mode of the core ^{10}Be is shown in Fig. 5.2.6. There are rather similar to the ones discussed in this Section (for details see Barranco et al. (2017), supplemental material).

²⁹Han et al. (2006); Koning and Delaroche (2003).

be viewed as *on par renormalized structure and reaction intermediate (virtual) elements of the quantal process* $^{11}\text{Be}(p,d)^{10}\text{Be}(2^+)$.

It is difficult, if not impossible, to talk about single-particle motion without also referring to collective vibrational states, concerning both structure and reactions. Within this context, the difficulties of defining spectroscopic factors are apparent.

5.2.3 Summary

In Figure 5.2.6 a “complete” $(\text{NFT})_{\text{ren(s+r)}}$ description of the single-neutron outside closed shell halo nucleus ^{11}Be is given, in terms of the reactions $^2\text{H}(\text{ ^{10}Be , ^{11}Be })^1\text{H}$ populating the $1/2^+$, $1/2^-$ and $5/2^+$ states and of the $^1\text{H}(\text{ ^{11}Be , ^{10}Be })^2\text{H}$ process populating the 2^+ mode. Also shown, in comparison with the data, are the $E1$ -transition between the parity inverted states $1/2^+$, $1/2^-$ and the isotopic shift of the charge radius of ^{10}Be . Repeating the last paragraph of Sect. 2.9 one can again state that *this, in a very real sense, is a nucleus*.

5.2.4 Continuum structure and reactions: ^{10}Li

As already stated, while the N -even isotopes $^9\text{Li}_6$ and $^{11}\text{Li}_8$ are bound, the N -odd system $^{10}\text{Li}_7$ is not. Based on the large corpus of experimental³⁰ and theoretical³¹ evidence concerning the fact that ^{10}Li also displays, as its bound isotope $^{11}\text{Be}_7$, parity inversion, the barely bound, very extended, ^{11}Li system can be viewed as two-weakly interacting fermions outside a closed shell ($N = 6$).

The consistency of the above picture was challenged by a $^9\text{Li}(d,p)^{10}\text{Li}$ experiment³² which found no $\ell = 0$ strength at low energies. In what follows an analysis of the corresponding results within the framework of continuum $(\text{NFT})_{\text{ren(s+r)}}$ is discussed.

The $^9\text{Li}(d,p)^{10}\text{Li}_7$ reaction

The NFT structure calculations were carried out as already explained in connection with the single halo isotope $^{11}\text{Be}_7$. In fact, in the determination of the bare potential parameters also the normal Mayer-Jensen level sequence, $N = 7$ isotones ^5_5B and ^6_6C , were included.

³⁰Wilcox et al. (1975); Young et al. (1994); Bohlen et al. (1997); Caggiano et al. (1999); Santi et al. (2003); Jeppesen et al. (2006); Smith et al. (2015); Sanetullaev et al. (2016); Kryger et al. (1993); Zinser et al. (1995); Kobayashi et al. (1997); Thoennessen et al. (1999); Chartier et al. (2001); Simon et al. (2007); Aksyutina et al. (2013); Amelin et al. (1990); Gornov et al. (1998); Chernysev et al. (2015); Kanungo et al. (2015); Tanihata et al. (2013); Fortune (2018); Tanihata et al. (2008).

³¹Tanihata et al. (2008); Thompson and Zhukov (1994); Bertsch et al. (1998); Blanchon et al. (2007); Mau and Pacheco (1996); Barker and Hickey (1977); Poppelier et al. (1993); Kitagawa and Sagawa (1993); Descouvemont (1997); Wurzer and Hofmann (1996); Kato and Ikeda (1993); Garrido et al. (2002, 2003); Potel et al. (2010); Casal et al. (2017); Vinh Mau (1995); Nunes et al. (1996); Sagawa et al. (1993); T. Myo et al. (2008); Barranco et al. (2001); Orrigo and Lenske (2009); Moro et al. (2019); Barranco et al. (2020).

³²Cavallaro et al. (2017).

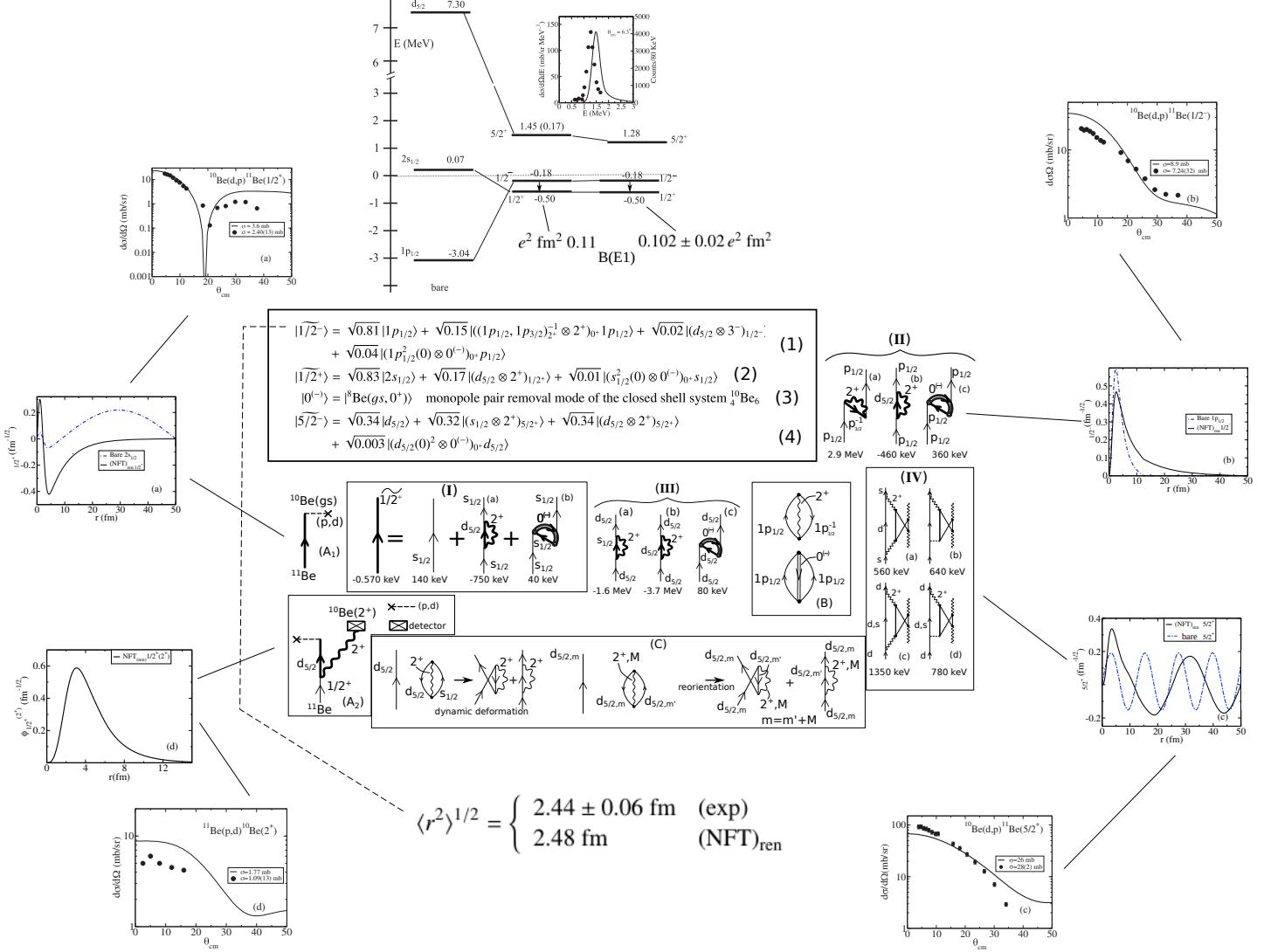


Figure 5.2.6: The clothing of the bare nucleons (single arrowed lines) with quadrupole and octupole particle-hole and monopole pair removal vibrations of the ^{10}Be core following the rules of renormalized nuclear field theory, give rise to values of the (renormalized) energies $\tilde{\epsilon}_j$ and renormalized single-particle wavefunctions $\tilde{\phi}_j$ which used as formfactors in connection with global optical potentials were used to calculate the absolute one-nucleon stripping and pickup differential cross sections (continuous curve (a), (b), (c), (d) (see also the corresponding form factors), displayed in comparison with the experimental data (solid dots)). The same is true concerning the $B(E1)$ transition between the parity inverted $1/2^+$, $1/2^-$ states and the isotopic shift of the charge radius.

The self energy matrix $\sum_{\nu_1 \nu_2}^{\mathbf{a}}(E)$ ($\mathbf{a} \equiv \{lj\}$) for the single-particle levels ν_1, ν_2 of ^{10}Li fulfilling $50 \text{ MeV} > \epsilon_{\nu_1}, \epsilon_{\nu_2} > \epsilon_F$, was calculated in a spherical box of 60 fm (continuum discretization). The dressed single particle states were determined by diagonalizing $\sum_{\nu_1 \nu_2}^{\mathbf{a}}(E)$. Up to about 2 MeV the spectrum is essentially determined by the $\widetilde{1/2}^+$ and $\widetilde{1/2}^-$ waves.

The $1/2^+$ scattering length was found to be³³ $\alpha = -8 \text{ fm}$, corresponding to an energy $\epsilon_{\widetilde{1/2}^+} = \hbar^2(1/\alpha)^2/2m \approx 0.3 \text{ MeV}$. The eigenfunction of a state lying close to this energy and, thus, representative of this virtual state is:

$$|\widetilde{1/2}^+\rangle = \sqrt{0.98}|s_{1/2}\rangle + \sqrt{0.02}|(d_{5/2} \otimes 2^+)_{1/2^+}\rangle. \quad (5.2.4)$$

Similarly, the resonant $\widetilde{1/2}^-$ state can be written as

$$|\widetilde{1/2}^-\rangle = \sqrt{0.94}|p_{1/2}\rangle + \sqrt{0.07}\left|((p_{1/2}, p_{3/2}^{-1})_{2^+} \otimes 2^+)_{0^+} p_{1/2}; 1/2^-\right\rangle. \quad (5.2.5)$$

The centroid and width of the resonance are $\epsilon_{\widetilde{1/2}^-} = 0.5 \text{ MeV}$, and $\Gamma_{\widetilde{1/2}^-} = 0.35 \text{ MeV}$ respectively³⁴. Theory predicts also the existence of a $5/2^+$ resonance with a centroid at $\approx 3.5 \text{ MeV}$. This state has also a clear many-body character, coupling to the quadrupole vibrations of the core through itself and through the $s_{1/2}$ state, not only in terms of virtual states containing one phonon ($(d_{5/2} \otimes 2^+)_{5/2^+}$, $((s_{1/2} \otimes 2^+)_{5/2^+})$) but also many (non crossing) phonon states which can be summed to infinite order (rainbow series)³⁵. Anharmonicities associated with states of two or more quadrupole phonons, resulting essentially from Pauli principle violation (see inset (IV) Fig. 5.2.6), are found to play an important role. Larger in the present case than in the case of ^{11}Be , in keeping with the fact that in ^{10}Li the $\widetilde{5/2}^+$ resonance lies closer in energy to the $\widetilde{1/2}^+ \otimes 2^+$ state than in ^{11}Be .

The coupling of the dressed neutron to the odd $1p_{3/2}^{-1}(\pi)$ proton hole leads to the doublets $(\widetilde{1/2}^+ \otimes p_{3/2}^{-1}(\pi))_{1^-, 2^-}$ and $(\widetilde{1/2}^- \otimes p_{3/2}^{-1}(\pi))_{1^+, 2^+}$ and to the quadruplet $(\widetilde{5/2}^+ \otimes p_{3/2}^{-1}(\pi))_{1^-, 2^-, 3^-, 4^-}$. The scattering length of the resulting $2^-, 1^-$ states lead to $\epsilon_{2^-} \approx 0.05 \text{ MeV}$ and $\epsilon_{1^-} = 0.8 \text{ MeV}$. For the positive-parity doublet, one finds $\epsilon_{1^+} \approx 0.3 \text{ MeV}$ and $\epsilon_{2^+} \approx 0.6 \text{ MeV}$, while the quadruplet spans the energy interval 2–6 MeV. Theory also predict a $3/2^-$ state which splits into four states $(\widetilde{3/2}^- \otimes p_{3/2}^{-1}(\pi))_{0^+, 1^+, 2^+, 3^+}$ with energies within the range 3–6 MeV.

³³This value provides circumstantial evidence on how extended the system described by the virtual state $|virt.\rangle = |\widetilde{1/2}^+(\kappa = 1/\alpha)\rangle$ is. In particular, as compared to the closed shell $N = 6$ core ^9Li of radius $R_0(^9\text{Li}) \approx 2.6 \text{ fm}$. A rough estimate of the overlap between $|virt.\rangle$ and single-particle states of ^9Li can be attempted through the ratio $(R_0/|\alpha|)^3 \approx 3.4 \times 10^{-2}$. Within this scenario, one may better assess the importance the many-body mixing amplitude $\langle(d_{5/2} \otimes 2^+)_{1/2^+} | \widetilde{1/2}^+ \rangle (= \sqrt{0.02})$ has in determine the properties of the renormalized $|\widetilde{1/2}^+\rangle$ state.

³⁴For details see Barranco et al. (2020), also supplemental material.

³⁵It is of notice that those many-phonon diagrams containing crossings are associated with vertex corrections. Because such diagrams imply recoupling coefficients of varied complexity and somewhat random phases, their summed contributions are expected to be considerably smaller than that of the rainbow series which displays coherence regarding the different contributions (see Barranco et al. (2019a)).

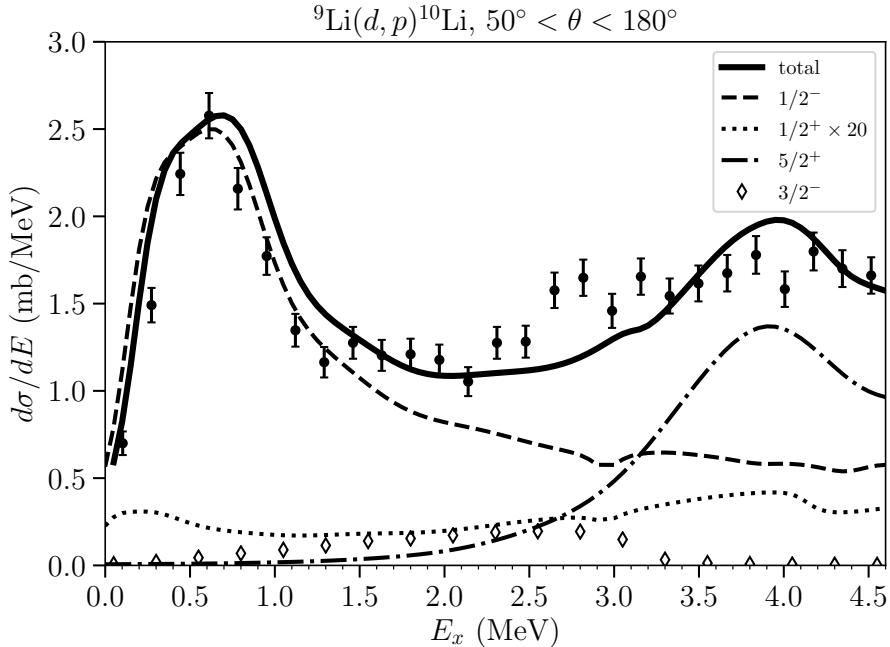


Figure 5.2.7: Theoretical predictions (continuous solid curve) of the absolute ^{10}Li strength function associated with the ${}^9\text{Li}(d, p){}^{10}\text{Li}$ reaction at 100 MeV incident energy and integrated over the angular range $5.5^\circ \leq \theta_{cm} \leq 16.5^\circ$, in comparison with experimental data (solid dots with errors, Cavallaro et al. (2017)). The partial contributions are the incoherent sums of the strength functions of the multiplets (\widetilde{J}^π , $1p_{3/2}(\pi)$). That is: (dashed line) $\widetilde{J}^\pi = \widetilde{1/2}^- (1^+, 2^+)$; (dash-dotted line) $\widetilde{5/2}^+ (1^- - 4^-)$; (diamonds) $\widetilde{3/2}^- (0^+ - 3^+)$; (dotted line) $\widetilde{1/2}^+ (1^-, 2^-)$ **multiplied by a factor of 20**.

Making use of the non-local self-energy matrices $\Sigma^a(r, r'; E)$, whose configuration space representation correspond to $\Sigma_{\nu_1 \nu_2}^a(E)$, together with global optical parameters³⁶ the absolute double differential cross sections $d\sigma/dEd\Omega$ containing the information the (d, p) reaction in question can provide on the structure of ${}^{10}\text{Li}$, was calculated.

One can now put the following question: recording the outgoing particle in the range of angles $5.5^\circ \leq \theta_{cm} \leq 16.5^\circ$, what the probability of observing the transferred neutron in the $\widetilde{1/2}^+$ final state is? The answer which arises from the strength function $d\sigma_{\widetilde{1/2}^+}/dE\Big|_{5.5^\circ-16.5^\circ}$ calculated by integrating $d\sigma/dEd\Omega$ in the chosen range of angles is univocal and reads *negligible*, as it emerges from Fig. 5.2.7, where theory is confronted with the experimental data to satisfactory agreement.

As seen from Fig. 5.2.8 the situation changes radically if *one puts the same question but changing the angular range from $5.5^\circ \leq \theta_{cm} \leq 16.5^\circ$ into $50^\circ \leq \theta_{cm} \leq$*

³⁶Schmitt et al. (2013).

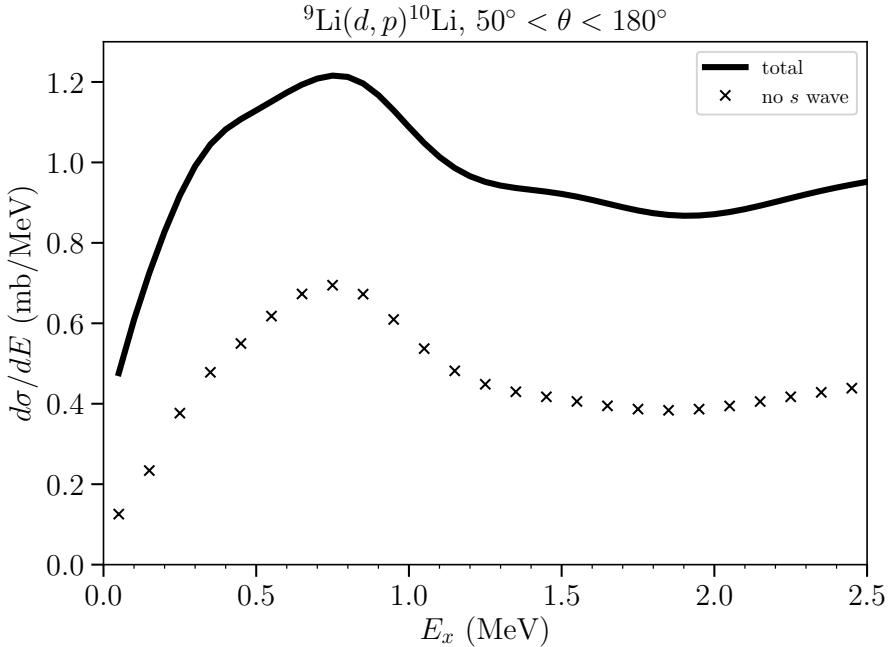


Figure 5.2.8: Predicted strength function obtained by integrating the absolute double differential cross section $d\sigma/dEd\Omega$ over the angular range $50^\circ \leq \theta_{cm} \leq 180^\circ$.

180° . Equally univocally the answer now reads *large*.

The reason at the basis of the essential difference between the two questions, and thus answers, is provided by the absolute differential cross section³⁷ $d\sigma(\theta)\widetilde{_{1/2}^+}/d\Omega$ obtained integrating $d\sigma\widetilde{_{1/2}^+}/dEd\Omega$ over the energy interval 0–0.2 MeV. As observed from Fig. 5.2.9 it results from a blend of structure, related to the single-particle content of the state (5.2.4), and the interference pattern displayed by the outgoing distorted waves which shows a minimum close to $\theta_{cm} = 16.5^\circ$. Returning to Fig. 5.2.7, and still within the above scenario, one can posit that the overall agreement between theory and experiment observed for $d\sigma/dE)_{5^\circ-16.5^\circ}$ within the energy range 3–4.5 MeV provides quantitative, although indirect, evidence of the presence of a robust $1/2^+$ state at low energy. This is in keeping with the strong coupling found in the NFT results between the $s_{1/2}$ and $d_{5/2}$ virtual and resonant states respectively, through the quadrupole vibration of the core ${}^9\text{Li}$. In fact, due to this strong mixing either one reproduces both the $d\sigma\widetilde{_{1/2}^+}/dE$ and the $d\sigma\widetilde{_{5/2}^+}/dE$ strength functions or likely none of them³⁸. Because these couplings renormalize single-particle content, energies, and widths, as well as the radial dependence of the wavefunctions (form factors), all these elements have to be calculated self-

³⁷The consistency with the ${}^9\text{Li}(d, p){}^{10}\text{Li}$ data of Jeppesen et al. (2006), recorded in the angular range $98^\circ \leq \theta_{cm} \leq 134^\circ$ is apparent. For details see Barranco et al. (2020). See also Moro et al. (2019).

³⁸Within this context see Barranco et al. (2020). See also Moro et al. (2019).

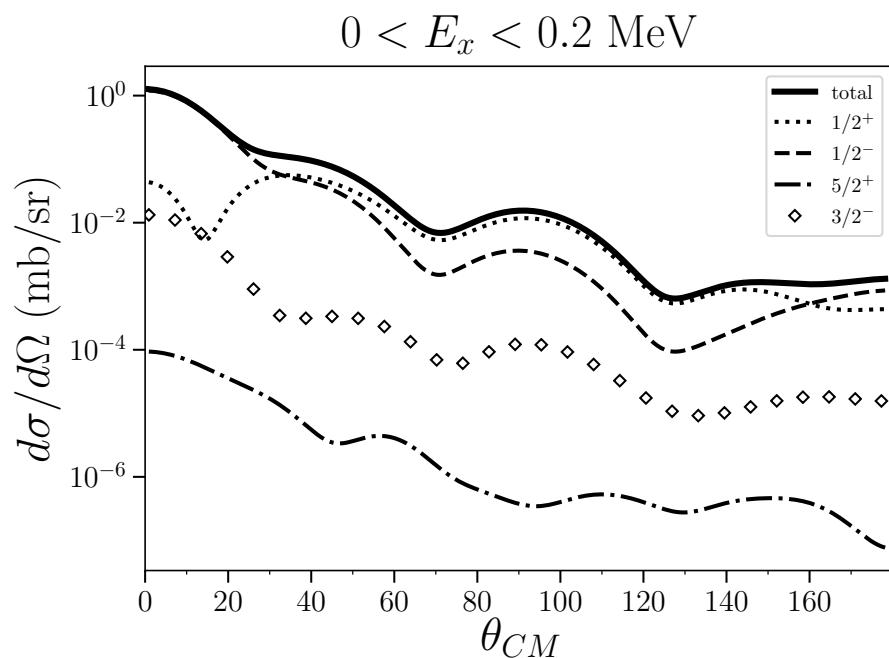


Figure 5.2.9: Predicted angular distributions ($d\sigma/d\Omega$) obtained by integrating the absolute double differential cross section $d\sigma/dEd\Omega$ in the energy interval 0–0.2 MeV. In the present case, and at variance with Fig. 5.2.7, no multiplicative factor was introduced in connection with $d\sigma_{1/2^+}/d\Omega$.

consistently. The role played by such requirements is likely emphasized in the case like the present one, of continuous spectroscopy, in which structure and reaction are, in a very real sense, just two aspects of the same physics.

Before concluding the present section, it may be useful to remind us what, within the framework of quantum mechanics, one can learn from a reaction experiment. It is not “what is the state after the collision” but “how probable is a given effect of the collision”³⁹.

5.3 Virtual states forced to become real through transfer reactions

One of the main subjects which has been discussed in the present monograph concerns the melting of *structure and reactions* and of *bare, virtual and renormalized states* into a higher unity –likely as it is in nature– which can be described in terms of Feynman diagrams where particles come in and the same or other particles come out and eventually interact with detectors whose clicks, can be translated into absolute cross sections and lifetimes. Thus conveying to observations the properties of renormalized, dressed, physical elementary modes of excitation.

The initial and final asymptotic states of the $(NFT)_{ren(s+r)}$ Feynman diagrams are anchored to the laboratory (incoming beams and target-outgoing particles, including γ -rays and detectors). The intermediate states break open the incoming systems –e.g. a beam of the halo nucleus ^{11}Li , of half life of 8.75 ms, which impinges on a hydrogen target (proton) in an inverse kinematic experiment– to allow the elementary modes of excitation which are probed by the experiment to become active, interact (melt) with other modes, and eventually pass the information concerning the properties of the physical modes to the outgoing particles.

Among these diagrams, one finds those associated with reaction processes in which a virtual state is acted upon and forced to become real, eventually reaching the detectors. A sort of Hawking radiation⁴⁰ to the extent that one concentrates on the (virtual) \rightarrow (real phenomenon) process, with the proviso of viewing the action of the external field as the (no-return) event horizon of the black hole. An example of such diagrams which describes a possible event associated with the reaction $^1\text{H}(^{11}\text{Li}, ^9\text{Li}(1/2^-))^3\text{H}$ is displayed in Fig. 2.7.3 (b): a virtual quadrupole vibration which, in the process of renormalizing the $s_{1/2}$ state, or of being exchanged

³⁹Born (1926a). If there is a lesson to be learned from the above discussion is the fact that, in dealing with a specific feature of a quantal many-body system, e.g. single-particle motion in nuclei (structure) and one-particle transfer process (reaction), one can hardly avoid to talk about other elementary modes of excitation and reaction channels, respectively. Within the scenario of the chosen example, this is because a nucleon which, in first approximation is in a mean field stationary state, can actually be viewed as a fermion moving through a gas of quadrupole phonons (ZPF) to which it couples, becoming eventually heavier because dressed by them (case of $s_{1/2}$). But also, because of Pauli principle the nucleon is forced to exchange role with the nucleons of the composite (particle-hole) phonons undergoing repulsion (case of $p_{1/2}$).

⁴⁰Barranco et al. (2019b).

between the partners of the neutron halo Cooper pair has been caught in the act by the pair transfer field produced by the ISAAC-2 facility at TRIUMF, forced to become a real final state and to bring this information to the active detector MAYA (see also Fig. 2.7.3 (d)).

5.3.1 Empirical renormalization⁴¹

Most if not all theories able to provide a complete description of structure and reactions of atomic nuclei can hardly avoid the separation between explicit and virtual configuration (phase) spaces⁴². Once this is recognized, a physical choice concerning the first type of degrees of freedom is: elementary modes of excitation. That is, the response to the variety of external probes of the nuclear system.

But here empiricism ends⁴³. Because the *bare* elementary modes of excitation which potentially contains all of the physics that experiment can eventually provide, are not observables. To become so, they have to loose their elementarity and become mixed, dressed, renormalized, and melt together into effective fields. In fact, what we call a physical nucleon moving inside the nucleus is only partially to be associated with that nucleon field alone. It is also partially to be associated with the vibrational fields, because the two are in interaction. On the other hand, vibrational modes participating in virtual states which intervened upon with an appropriate external field can become on-shell⁴⁴, have to be dressed, fully renormalized modes, poised to be forced to become observable on short call⁴⁵.

As shown in Fig. 5.3.2 (see also Fig. 5.3.1), the reaction $^{11}\text{Be}(p, d)^{10}\text{Be}(2^+)$ provides an embodiment of such processes. In fact, this reaction gives information concerning the most important process clothing of the $1/2^+$ parity inverted ground state of $^{11}_4\text{Be}_7$ through the coupling to the low-lying quadrupole vibration of the core $^{10}_4\text{Be}_6$ (see Fig. 5.3.2 I (a)). A schematic representation of the pickup of the neutron moving around a $N = 6$ closed shell and populating the low-lying

⁴¹The practical usefulness of the Feynman rules and diagrams made them one of the most essential elements of the scientific training of every theoretical physicist (Mehra (1996) p. 290).

⁴²While configuration space traditionally refers to structure and phase space to reactions (see e.g. Feshbach (1962)) these connotations tend to melt into a unity in the case of unstable, exotic nuclei like ^{11}Li ($\tau_{1/2} \approx 8$ ms), let alone ^{10}Li , transiently observed in terms of virtual and resonant states in a one-neutron transfer processes (see Sect. 5.2.4 and footnote 31 Ch. 5; see also Barranco et al. (2020) and refs. therein). As a consequence, both structure and reactions are to be worked out in the continuum, discretization being a connotation of methodological, rather than of physical approach. Think only on the central role the tail of the neutron halo Cooper pair, bound by only 380 keV to the core (^9Li), play in stabilizing ^{11}Li , and in connection with two-nucleon transfer process (formfactors). Within this scenario, inverse kinematics in e.g. the case of $^2\text{H}(^9\text{Li}, ^{10}\text{Li})^1\text{H}$ provides, in the laboratory set up, a large but finite box for both the probe (deuteron) and the probed (^{10}Li).

⁴³To the extent of attributing the ancient Greek meaning of “find” and “discover” to the word heuristic (*ευρισκω*) and of “serving to discover” of the Oxford dictionary, one can connote the above empirical protocol, as heuristic.

⁴⁴The intervened mode has an on-shell energy consistent with the Q -value of the reaction, and thus may not correspond to that of an actual final state in absence of that probe.

⁴⁵As a rule, the collective modes participating in these virtual state are calculated making use of empirical input. Thus, empirical renormalization.

quadrupole vibrational state of this core, in coincidence with the corresponding γ -decay is shown in **I(b)**. More detailed structure and reaction NFT diagrams, are shown in **I(d)** and **I(e)**, (the jagged line represents a graphic mnemonic of the recoil effect), together with a cartoon representation in **(f)**. The predicted (continuous curve) and experimental (solid dots) absolute differential cross sections are displayed in **I(c)**. Protons and neutrons are labeled π and ν respectively, d stands for deuteron while γ -detectors are represented by a crossed rectangle. Curved arrows indicate motion in the continuum (reaction). Normal arrowed lines, motion inside target or projectile (structure).

At the basis of the spontaneous γ -decay displayed in **I(b)** and **I(e)** we find the processes shown in the lower part of Fig. 5.3.2, as it is explained in what follows. **(II)** Interaction of protons in a nucleus with nuclear vibrations (solid dot, PVC vertex $\beta_L R_0 \partial U / \partial r Y_{LM}^*(\hat{r})$; β_L , dynamical distortion parameter; $U(r)$ central potential) and photons (normal vertex, electromagnetic interaction⁴⁶ $e \int d^4x J_\mu(x) A^\mu(x)$, the operator A^μ being the vector potential, and J_μ the current density ($\mu = 1, \dots, 4$)). While the variety of diagrams shown have general validity, we have assumed we are dealing with the low-lying correlated particle-hole quadrupole vibration ($L = 2$) of $^{10}_4\text{Be}_6$ lying at 3.368 MeV, $B(E2; 0^+ \rightarrow 2^+) = 0.0052 e^2 b^2$ being associated with $\beta_2 \approx 0.9$. An arrowed line pointing upward (downward) describes a proton (proton hole) moving in the $p_{1/2}$ ($1p_{3/2}$) orbital. Zero point fluctuations of the nuclear ground state associated with : **(a)** the nuclear quadrupole vibration, **(b)** the electromagnetic field associated with the corresponding spontaneous γ -decay. **(c)** Pauli principle correction to the simultaneous presence of the above two ZPF processes. **(d)** Intervening the virtual excitation of the nuclear vibrations (graph (c)) with an external (inelastic) field (cross followed by a dashed line), in coincidence with the γ -decay (γ -detector, crossed box), the virtual process (c) becomes real. **(e)**, **(f)** Time ordering of the above process correspond to the RPA contributions through backwardsgoing and forwardsgoing amplitudes and subsequent γ -decay.

The above is an example of empirical renormalization (use of dressed elementary modes of excitation in both structure and reaction or better, in the present case, decay channels), namely $((\text{NFT})_{\text{ren}(s+r)})^{47}$.

5.3.2 On-shell energy

It was stated that intermediate states are, aside from their energy, fully dressed, renormalized physically observable elementary modes of excitation. Concerning the eventual value of the on-shell-energy, it will naturally depend on the reaction process taking place. For example, in the reaction $^{11}\text{Be}(p, d)^{10}\text{Be}(2^+; 3.368 \text{ MeV})$ (Fig. 5.3.2 **(I)** (b) and (e)) populating the quadrupole vibration of ^{10}Be which, in the virtual $(\widetilde{5/2}^+ \otimes 2^+)_{1/2^+}$ state dress the $|1/2^+\rangle$ ground state of ^{11}Be (Fig. 5.3.2 **(I)** (a)), the on-shell-energy coincides with that which is observed in

⁴⁶Holstein (1989).

⁴⁷Broglia et al. (2016).

the inelastic process ${}^{10}\text{Be}(\text{gs} \rightarrow 2^+)$. On the other hand, the $5/2^+$ pickup neutron, while displaying the structural properties of the $5/2^+$ resonance (Fig. 5.3.1) ($\tilde{\epsilon}_{5/2^+} = 1.45$ MeV), it has a binding energy equal to $\tilde{\epsilon}_{1/2^+} - \hbar\omega_{2^+} = -3.868$ MeV ($\tilde{\epsilon}_{1/2^+} = -0.5$ MeV, $\hbar\omega_{2^+} = 3.368$ MeV), unrelated to the experimental energy of +1.28 MeV (continuum, unbound), recorded in the reaction ${}^{10}\text{Be}(d, p){}^{11}\text{Be}(5/2^+)$. This is a natural outcome of $(\text{NFT})_{\text{ren}}$ which, through the PVC and the Pauli mechanism, provides the proper clothing of the $d_{5/2}$ orbital so as to allow it to be able “to exist” inside the $|s_{1/2}\rangle$ state as a virtual, intermediate configuration. The asymptotic r -behavior results from the coherent superposition of many continuum states.

We now consider the process in which an inelastic field acting on $|{}^{11}\text{Be}(\text{gs})\rangle$ populates the $5/2^+$ resonance (see Fig. 5.3.1). An important contribution to this excitation results from the action of the external field on the quadrupole vibration of the virtual state $(\tilde{d}_{5/2} \otimes 2^+)_{1/2^+}$ which renormalizes the $1/2^+$ state (Fig. 5.3.1 (a)) leading to diagram (b) of the same figure. In this case, the external field has to provide an energy equal to $|\tilde{\epsilon}_{1/2} - \tilde{\epsilon}_{5/2^+}| = 1.95$ MeV, again unrelated to the on-shell energy of $|{}^{10}\text{Be}(2^+; 3.868 \text{ MeV})\rangle$ observed in the inelastic process ${}^{10}\text{Be}(\text{gs} \rightarrow 2^+)$, but needed to dress the bare single-particle valence orbital.

5.3.3 Perturbation and beyond

In Q.E.D. at the one-loop diagrams level⁴⁸, the renormalized physical electron mass m_e , that is the observed mass (≈ 0.511 MeV), is related to the parameter m in the Dirac equation, the bare electron mass (not observable), according to $m_e = m \left(1 + \frac{3\alpha}{4\pi} \log \left(\frac{\Lambda_{\text{cut}}}{m} \right)^2 \right)$, Λ_{cut} being the cutoff of the divergent integrals⁴⁹. This quantity appears inside the log, and in front of it one has the fine structure constant α which is small ($\approx 1/137$). Therefore, even pushing the cutoff to the Planck scale, $\Lambda_{\text{cut}} \sim 10^{19}$ GeV, with $m \sim \text{MeV}$ one has $\delta m_e/m \approx \frac{3\alpha}{4\pi} \log \left(\frac{\Lambda_{\text{cut}}}{m} \right)^2 \approx 0.1$. So δm_e is really a small correction.

Within this context one can compare the value of the four parameters (plus the effective mass functional) which determine the mean field (Saxon–Woods) potential used to calculate the bare single-particle energies of ${}^{11}\text{Be}$ –imposing the condition that the dressed levels best reproduce the experimental findings– with the value of the corresponding parameters of the standard, global Woods–Saxon potential (and of effective mass equal to the observed mass m). The results displayed in Table 5.3.1 testify to the fact that renormalization in nuclear physics, in particular in the case of halo, exotic parity inverted nuclei, is less “perturbative” than in the case of Q.E.D. This fact becomes even clearer if one compares the overall centroid of the valence orbitals as well as the density of levels associated with the two potentials. Within this context it is of notice that the Lamb (–like) shift taking place between the $s_{1/2}$ and $p_{1/2}$ valence levels of ${}^{11}\text{Be}$ has a value of

⁴⁸Diagram not allowed by rule III) of NFT in keeping with the fact that the nuclear bosons are (RPA) “composite particles”.

⁴⁹See Bjorken and Drell (1998) pp 162–166; Milloni (1994) p. 404.

approximately 10% ($(\Delta\epsilon_{1/2^+,1/2^-})_{bare} - (\Delta\epsilon_{1/2^+,1/2^-})_{ren} \approx 3.11 + 0.32\text{MeV}$) of that of the Fermi energy ($\approx 36\text{ MeV}$). This result can be compared with the ratio of the hydrogen $^2S_{1/2} - ^2P_{1/2}$ Lamb shift ($1058\text{ MHz} \approx 4.3 \times 10^{-9}\text{ eV}$) and the Rydberg constant ($R_H = 13.6\text{ eV}$), i.e. $\approx 10^{-10}$, a result which underscores the strong coupled situation one is confronted with in trying to describe the structure of light halo exotic nuclei (Fig. 5.2.4).

5.3.4 One-particle transfer (s+r)-dielectric function

The fact that in spite of this non-perturbative situation, $(\text{NFT})_{ren(r+s)}$ can provide an overall account of an essentially complete set of experimental data which characterizes ^{11}Be , within a 10% error, testifies to the power and flexibility Feynman version of Q.E.D. has. It can be used as paradigm to construct a field theory for both structure and reactions of a strongly interacting finite many-body system like the atomic nucleus. As already mentioned, examples of $(\text{NFT})_{ren(r+s)}$ diagrams, aside from that discussed in detail in connection with Fig. 5.3.2, are displayed in Figs. 2.7.3 and 2.7.2. The first describes the process $^1\text{H}(^{11}\text{Li}, ^9\text{Li})^3\text{H}$ providing a quantitative account of the data and evidence of phonon induced pairing in nuclei. The second shows in (a), one of the most important channels contributing to the optical potential needed to describe the elastic scattering process $^1\text{H}(^{11}\text{Li}, ^{11}\text{Li})^1\text{H}$. It is of notice that in the case of $^9\text{Li}(d, p)^{10}\text{Li}$ (self-energy) and reaction (optical potentials) dielectric functions melt together. The calculation of the optical potential constitutes, within the scheme of $(\text{NFT})_{ren(r+s)}$ a major challenge lying ahead.

5.3.5 $(\text{NFT})_{ren(s+r)}$ diagrams, S-matrix

Landau felt that Feynman diagrams, although usually derived from conventional field theory have an independent basic importance⁵⁰ This work is closely connected with both Heisenberg statement that in quantum theory, one should only introduce quantities⁵¹ which are, in principle, observable and with the special role the scattering matrix plays in such a program. As the wavefunctions themselves cannot be observed and because the Hamiltonian formalism is intimately connected with wavefunctions, it may not be the most appropriate one to describe quantal systems. The quantities to be studied are the scattering amplitudes where particles go in and another set, including eventually γ -rays, come out and directly determine the cross sections of the different physical processes. Within this context we refer to diagrams (b) and (e) of Fig. 5.3.2 (I) where the $(\text{NFT})_{ren(r+s)}$ diagram describing the becoming of a concrete virtual process⁵² real⁵³, is displayed.

⁵⁰Landau (1959).

⁵¹Heisenberg (1925); see english translation W. Heisenberg, Quantum-theoretical re-interpretation of kinematical and mechanical relations, in Van der Waerden (1967) p 261.

⁵²Dressing of the $s_{1/2}$ ground state of ^{11}Be , diagram (a), see also Table 5.3.1 for the parameters of the bare potential.

⁵³Through the action of an external field namely $^{11}\text{Be}(p, d)^{10}\text{Be}(2^+)$.

	$V_0(\text{MeV})$	$V_l(\text{MeV})$	$R_0(\text{fm})$	$a(\text{fm})$
Standard ^{a)}	-50	17	2.7 ^{c)}	0.65
bare ^{b)}	-68.9	14.47	2.15 ^{d)}	0.77

Table 5.3.1: Parametrization of the standard (Bohr and Mottelson (1969)) and of the bare mean field potential associated with ^{11}Be (Barranco et al. (2017), see also supplemental material). Changes of the order of 20–30% are observed. Within this context, in the case of Q.E.D. $\delta m_e/m = 0.1$.

^{a)} $m^* = m$

^{b)} $m^*(r=0) = 0.7m$, $m^*(r=\infty) = (10/11) \times m$

^{c)} $R_0 = r_0 A^{1/3}$, $r_0 = 1.2 \text{ fm}$

^{d)} $r_0 \approx 1.03 \text{ fm}$

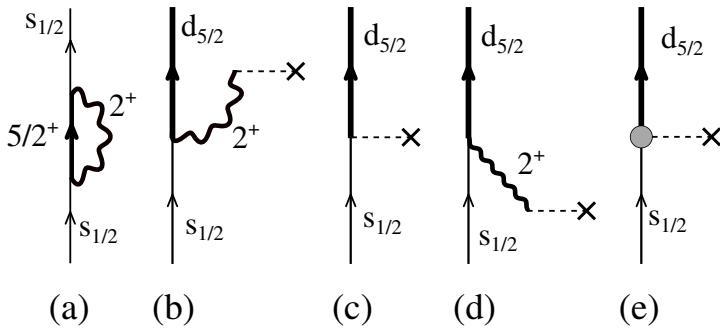


Figure 5.3.1: Inelastic population of the $|^{11}\text{Be}(5/2^+; 1.45 \text{ MeV})\rangle$ calculated resonance. (a) Main process dressing the neutron moving in the $s_{1/2}$ orbital and leading to the ^{11}Be ground state. The virtual state is made out of the fully renormalized (real, empirical) quadrupole vibration of the core ^{10}Be and of the low-lying $5/2^+$ resonance of ^{11}Be , and is indicated by using bold face symbols. (b) Intervening process (a) with an external inelastic hadron field (e.g. (p, p')) of quadrupole character (dashed horizontal line starting at a cross), one can excite the $5/2^+$ resonance. This can also happen if the external field acts on the $1/2^+$ state as in (c), or if this field first excites the quadrupole vibration of ^{10}Be which eventually couples to the $1/2^+$ state as in (d). (e) If the vibration was a high-lying quadrupole giant resonance the summed contribution of processes (b)–(d) could be replaced at profit by a single graph, namely the equivalent of (c) but with an effective charge (hatched circle). For low-lying modes like the 3.368 MeV quadrupole vibration of ^{10}Be , retardation, ω -dependent effects are to be explicitly taken into account.

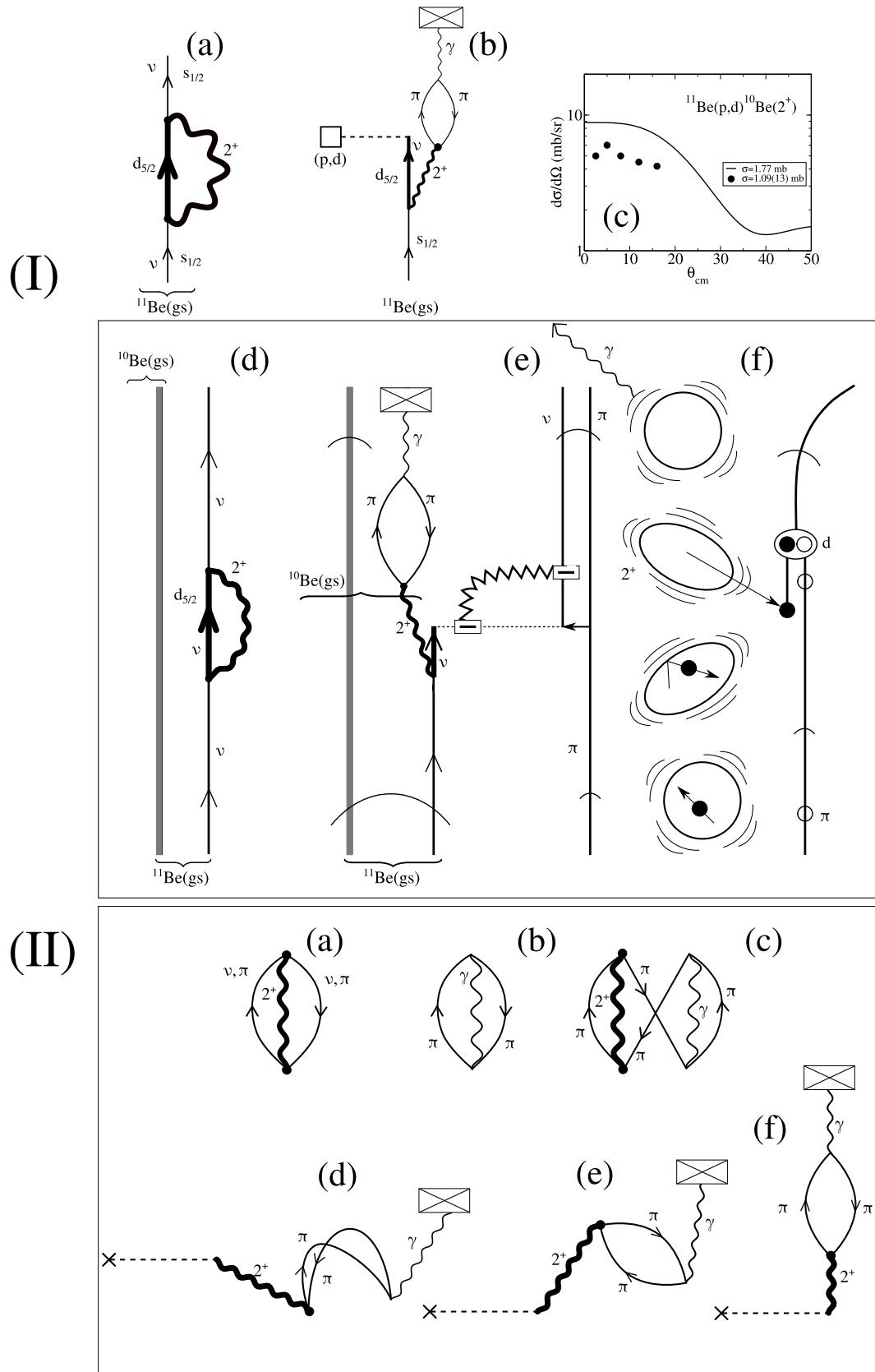


Figure 5.3.2: The reaction $^{11}\text{Be}(p, d)^{10}\text{Be}(2^+)$. A virtual process, namely the self-energy contribution of the $s_{1/2}$ ground state of ^{11}Be through the coupling to the low-lying collective quadrupole vibration of the ^{10}Be becomes real through the action of a one-particle pickup external field.

5.4 Minimal requirements for a consistent mean field theory

As seen from Figs. 1.2.2–1.2.4 the minimum requirements of selfconsistency to be imposed upon single-particle motion requires both non-locality in space (HF) and in time (TDHF)

$$i\hbar \frac{\partial \varphi_\nu}{\partial t} = -\frac{\hbar^2}{2m} \nabla^2 \varphi_\nu(x, t) + \int dx' dt' U(x - x', t - t') \varphi_\nu(x', t'). \quad (5.4.1)$$

Assuming, for simplicity, infinite nuclear matter (confined by a constant potential of depth V_0), and thus plane wave solutions, the above time-dependent Schrödinger equation leads to the quasiparticle dispersion relation

$$\hbar\omega = \frac{\hbar^2 k^2}{2m^*} + \frac{m}{m^*} V_0, \quad (5.4.2)$$

where the effective mass

$$m^* = \frac{m_k m_\omega}{m}, \quad (5.4.3)$$

is the product of the k -mass

$$m_k = m \left(1 + \frac{m}{\hbar^2 k} \frac{\partial U}{\partial k} \right)^{-1}, \quad (5.4.4)$$

closely connected with the Pauli principle ($\frac{\partial U}{\partial k} \approx \frac{\partial U_x}{\partial k}$), while the ω -mass

$$m_\omega = m \left(1 - \frac{\partial U}{\partial \hbar\omega} \right), \quad (5.4.5)$$

results from the dressing of the nucleon through the coupling with the (quasi) bosons.

The occupancy of levels around ε_F is related to Z_ω , a quantity which measures the discontinuity at the Fermi energy and which is equal to m/m_ω . This is in keeping with the fact that the time the nucleon is coupled to the vibrations it cannot behave as a single-particle and can thus not contribute to e.g. the single-particle pickup cross section. *The particle-vibration coupling not only renormalizes energies and single-particle content but also the radial dependence of the wavefunction (formfactors).*

It is of notice that the self-consistence requirements for the iterative solution of Eq. (5.4.1) reminds very much those associated with the solution of the Kohn-Sham equations in finite systems,

$$H^{KS} \varphi_\gamma(\mathbf{r}) = \lambda_\gamma \varphi_\gamma(\mathbf{r}), \quad (5.4.6)$$

where

$$H^{KS} = -\frac{\hbar^2}{2m_e} \nabla^2 + U_H(\mathbf{r}) + V_{ext}(\mathbf{r}) + U_{xc}(\mathbf{r}), \quad (5.4.7)$$

H^{KS} being known as the Kohn-Sham Hamiltonian, $V_{ext}(\mathbf{r})$ being the field created by the ions and acting on the electrons. Both the Hartree and the exchange-correlation potentials $U_H(\mathbf{r})$ and $U_{xc}(\mathbf{r})$ depend on the (local) density, hence on the whole set of wavefunctions $\varphi_\gamma(\mathbf{r})$. Thus, the set of KS -equations must be solved selfconsistently⁵⁴.

5.4.1 Density of levels

Making use of Eq. (5.4.2) ($E = \hbar\omega$), one can calculate dE/dk for a single nucleon and one spin orientation⁵⁵. The inverse of this expression is⁵⁶

$$\frac{dk}{dE} = \frac{m^*}{\hbar^2 k}, \quad (5.4.8)$$

which testifies to the fact that the energy spacing between levels, i.e. the density of levels (see below), changes as m^* does.

One can then calculate the average value over the Fermi distribution, obtaining

$$\left\langle \frac{dk}{dE} \right\rangle = \frac{m^*}{\hbar^2 (2/3) k_F}. \quad (5.4.9)$$

Let us now take into account all nucleons, both spin orientations and eliminate the unit (inverse) length, i.e.

$$\frac{2A}{k_F} \left\langle \frac{dk}{dE} \right\rangle = 3A \frac{m^*}{\hbar^2 k_F^2} = 3A \frac{m^*}{2m\epsilon_F}. \quad (5.4.10)$$

Assuming $m^* = m_\omega m_k/m \approx m$, where m is the experimental nucleon mass one obtains

$$\frac{3}{2} \frac{A}{\epsilon_F}, \quad (5.4.11)$$

a value which coincides with the Fermi gas model estimate for the one-particle level density g_0 ⁵⁷. Taking properly into account the geometry of the system, one obtains

$$a = \frac{\pi^2}{6} \frac{3}{2} \frac{A}{\epsilon_F} \quad (5.4.12)$$

for the prefactor of the exponential of the Fermi expression of the total density of single particle levels. Making use of $\epsilon_F = 36$ MeV leads to,

$$a \approx \frac{A}{14} \text{MeV}^{-1}. \quad (5.4.13)$$

⁵⁴See e.g. Broglia et al. (2004) and refs. therein.

⁵⁵Mahaux et al. (1985), p. 17.

⁵⁶P. F. Bortignon, private communication.

⁵⁷See e.g. Bohr and Mottelson (1969) Eq. (2-48).

In keeping with the fact that one can interpret dE/dk as the rate of change in energy when the momentum changes or, equivalently, when the number of nodes per unit length changes, and this can be used to label the single-particle states, (5.4.13) can be confronted at profit with the average degeneracy per unit energy of valence orbitals (see Table 5.4.1).

Within this context it is of notice that an estimate of the quantity a based on the harmonic oscillator, leads to $a \approx \frac{\pi^2}{6} \frac{(N_{max}+3/2)^2}{\hbar\omega_0}$ i.e. an expression inversely proportional to the (constant) energy separation of levels⁵⁸. In an attempt to bridge the gap between the nuclear matter expressions discussed above and finite nuclei, i.e. potential wells of finite range, we consider

$$H = \frac{p^2}{2D} + \frac{C}{2}x^2, \quad (5.4.14)$$

($p = D\dot{x}$) which leads to a constant level spacing,

$$\hbar\omega_0 = \hbar \sqrt{\frac{C}{D}}, \quad (5.4.15)$$

and implies that the density of states is proportional to the square root of the particle inertia (mass). However, this result follows from the assumption that the potential remains unchanged if the bare mass (in which case D is, for example, set equal to the HF k -mass, m_k) is replaced by an effective mass m^* (e.g. $m_k m_\omega / m$, Eq. (5.4.3)). In this case the ground state wavefunction

$$\varphi_0 \sim \exp\left(-\frac{x^2}{2b^2}\right), \quad (5.4.16)$$

with,

$$b = \sqrt{\frac{\hbar}{m^*\omega_0}}, \quad (5.4.17)$$

in the case of a dressed nucleon of effective mass $m^* > D$ will shrink in space compared with the one of mass D and, consequently, the mean square radius of the system

$$\langle r^2 \rangle = \frac{\hbar}{m^*\omega_0} \left(N + \frac{3}{2} \right) = b^2 \left(N + \frac{3}{2} \right) \quad (5.4.18)$$

will decrease. This is not correct, and one has to impose the condition $b^2=\text{constant}$. A condition which implies that the energy difference between levels is inversely proportional to the effective mass of the nucleon⁵⁹, that is,

$$\hbar\omega_0 = \frac{\hbar^2}{m^*b^2}. \quad (5.4.19)$$

⁵⁸See Bohr and Mottelson (1969) p. 188, Eq (2-125a).

⁵⁹See Sect. 3.2, Eq. (3.2.2); also Fig. 2.6.1 inset (c).

	MeV ⁻¹	
	empirical	<i>a</i>
²⁰⁸ Pb ₁₂₆ 82	17(10 ^a + 7 ^b)	15(9 ^a + 6 ^b)
¹²⁰ Sn ₁₂₆ 50	4 ^a)	5 ^a)

a) neutrons
b) protons

Table 5.4.1: Comparison of the factor (5.4.13) ($a = n/14\text{MeV}^{-1}$) corresponding to ²⁰⁸Pb for both $n = N$ and $n = Z$ and for ¹²⁰Sn in the case of $n = N$, in comparison with the empirical value associated with the valence orbitals of these nuclei. That is ($h_{9/2}, f_{7/2}, i_{13/2}, p_{3/2}, f_{5/2}, p_{1/2}, g_{9/2}, i_{11/2}, d_{5/2}, j_{15/2}, s_{1/2}, g_{7/2}, d_{3/2}$) and ($g_{7/2}, d_{5/2}, h_{11/2}, d_{3/2}, s_{1/2}, h_{9/2}, f_{7/2}, i_{13/2}$) for neutrons and protons in ²⁰⁸Pb, leading to $\sum_j^N (2j+1)/\Delta E_N \approx 102/10 \text{ MeV} \approx 10 \text{ MeV}^{-1}$ and $\sum_j^Z (2j+1)/\Delta E_Z = 64/(9 \text{ MeV}) \approx 7 \text{ MeV}^{-1}$, where ΔE_N is the experimental energy interval over which the valence orbitals are distributed (see e.g. Bohr and Mottelson (1969) p. 325, Fig. 3-3). In the case of neutrons of ¹²⁰Sn, use is made of the dressed valence orbitals $d_{5/2}, g_{7/2}, s_{1/2}, d_{3/2}, h_{11/2}$ resulting from the renormalization of HF-Sly4 levels through the coupling of collective modes making use of nuclear field theory plus Nambu-Gorkov techniques ((NFT)+(NG); for details see Idini et al. (2015) and Table I of Potel et al. (2017)). The result, taking into account the breaking of the single-particle strength, in particular that of the $d_{5/2}$ orbital is $\sum_j^N (2j+1)/\Delta E_N = 32/8 \text{ MeV} \approx 4 \text{ MeV}^{-1}$.

Let us conclude this Section with a remark concerning the dimensions of the parameters D and C entering Eq. (5.4.14). Because the variable x has dimensions of length ([x]=fm), the dimensions of the inertia and restoring force parameters are

$$[C] = \text{MeV fm}^{-2} \quad \text{and} \quad [D] = \text{MeV fm}^{-2}\text{s}^2.$$

Consequently the associated zero point fluctuations

$$\sqrt{\frac{\hbar\omega_0}{2C}} = \sqrt{\frac{\hbar^2}{2D} \frac{1}{\hbar\omega_0}} \quad (5.4.20)$$

have dimensions of fm. It is of notice that in the case of the harmonic oscillator Hamiltonian (1.1.9), the associated ZPF is a *c*-number, in keeping with the fact that the (dynamic) deformation parameter α is dimensionless and thus $[D]=\text{MeVs}^2$ and $[C]=\text{MeV}$.

5.5 Self-energy and vertex corrections

In Fig. 5.5.1 an example of the fact that in field theories (e.g. QED or NFT), nothing is really free and that e.g., the bare mass of a fermion (electron or nucleon),

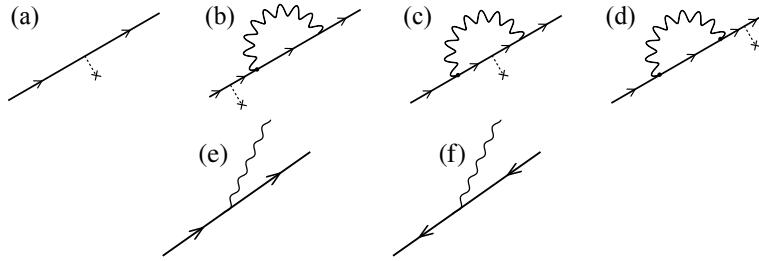


Figure 5.5.1: Self energy (effective-mass-like) processes. The result of the probing with an external field (dotted line started with a cross, observer) of the properties (mass, charge, single-particle content, etc) of a fermion (e.g. an electron or a nucleon, arrowed line) dressed through the coupling to bosons (photons or collective vibrations, wavy line), results from the modulus squared of the sum of the amplitudes associated with each of the four diagrams (a)–(d) (cf. (Feynman, 1975)). In graphs (e) and (f) the amplitudes eventually leading to the self-energy processes (b)–(d) are shown.

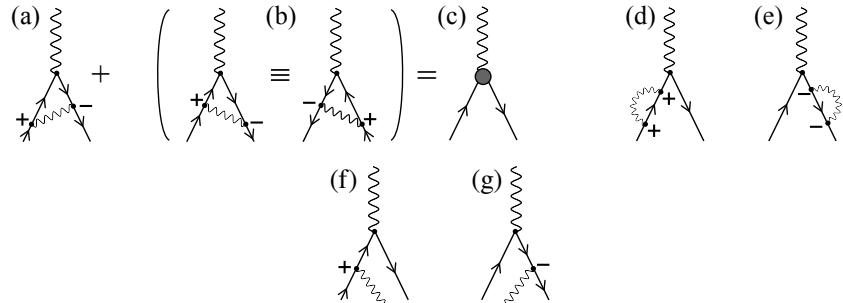


Figure 5.5.2: (a), (b) Vertex corrections. These are triple-interaction diagrams (phonon, particle and hole lines) in which none of the incoming lines can be detached from either of the other two by cutting one line. In connection with condensed matter Migdal's theorem (Migdal (1958)) states that for phonons, (Bardeen and Pines (1955), Fröhlich (1952)) vertex corrections can be neglected (cf. also Anderson (1964a)). Vertex corrections are, as a rule, important in the nuclear case where they lead to conspicuous cancellations of the self-energy contributions (d) and (e) (cf. e.g. Bortignon et al. (1983), cf. also Anderson (1964a)). The solid grey circle in (c) represents the effective, renormalized vertex. In graphs (f) and (g) the amplitudes describing the coupling of a $p - h$ vibrational mode with the $2p - 2h$ doorway states are shown, and eventually leading to the vertex correction and self energy processes (a)–(e) are schematically displayed.

is the parameter one adjusts so that the renormalized particle displays properties which agree with experimental findings, for example the observed mass (single particle energy). In Fig. 5.5.2, lowest order diagrams associated with the renormalization of the fermion–boson interaction (vertex corrections) are given. The sum of contributions (a) and (b) can, in principle, be represented by a renormalized vertex (see, diagram (c)). There is, as a rule, conspicuous interference (e.g. cancellation) in the nuclear case between vertex and self-energy contributions (see diagrams (a) and (d)+(e) of Fig. 5.5.2, a phenomenon closely related with conservation laws⁶⁰. In particular, cancellation in the case in which the bosonic modes are of isoscalar character⁶¹. *Consequently, one has to sum explicitly the different amplitudes with the corresponding phases and eventually take the modulus squared of the result to eventually obtain the quantities to be compared with the data, a fact that call attention to the use of an effective, ω -independent (renormalized) vertex.*

Within the framework of QED the above mentioned cancellations are exact implying that the interaction between one- and two-photon states vanishes (Furry theorem). The physics at the basis of the cancellation found in the nuclear case can be exemplified by looking at a spherical nucleus displaying a low-lying collective octupole vibration, like e.g. in ^{208}Pb where $|3^-; 2.615 \text{ MeV}\rangle$ is the first excited state, its $B(E3)$ being equal to $32 B_{sp}$. The associated zero point fluctuations (ZPF) lead to time dependent shapes with varied instantaneous values of the octupole moment. In other words, a component of the ground state wavefunction $(|j_p \otimes j_h^{-1} 3^- \otimes 3^-; 0^+\rangle)$, which can be viewed as a gas of octupole (quasi) bosons promoting a nucleon across the Fermi energy (particle-hole excitation) will lead to fermionic states which behave as having a positive (particle) and a negative (hole) effective octupole moment, in keeping with the fact that the closed shell system is spherical, and thus has zero octupole moment.

5.5.1 Phase coherence and damping of vibrational modes

Paradigmatic examples of vibrational modes are provided by giant resonances. In particular by the giant dipole resonance (GDR). This collective mode is made out of particle-hole excitations. An estimate of the associated damping width can be obtained in terms of the sum of the individual widths of the particle and of the hole. With the help of the relations (2.3.20), (2.3.21) valid for both particles and holes, one obtains

$$\Gamma_{GDR}(\hbar\omega_{GDR}) = \Gamma_p(\hbar\omega_{GDR}/2) + \Gamma_h(\hbar\omega_{GDR}/2) \approx 0.5 \hbar\omega_{GDR}, \quad (5.5.1)$$

where the assumption was made that the excitation energy $\hbar\omega_{GDR}$ is divided equally between the particle and the hole. Making use of the empirical expression $\hbar\omega_{GDR} \approx$

⁶⁰Schrieffer (1964)); and refs. Bortignon and Broglia (1981); Bertsch et al. (1983) and Bortignon et al. (1998) pp. 82–86.

⁶¹Bortignon et al. (1983).

$80/A^{1/2}$ MeV (Eq. (1.5.2) and following text), one can write

$$\Gamma_{GDR} \approx \frac{40}{A^{1/3}} \text{ MeV}, \quad (5.5.2)$$

which, in the case of ^{208}Pb leads to $\Gamma_{GDR} \approx 7$ MeV. Number to be compared with the experimental value of ≈ 4 MeV⁶². This overprediction of a factor of 2 is due to the fact of having taken the modulus squared of the amplitudes associated with graphs (f) and (g) of Fig. 5.5.2 separately, instead of doing so after having summed them with the proper phase. In other words,

$$\Gamma_{GDR} \approx \left| \frac{\sqrt{\Gamma_p} + e^{-i\alpha} \sqrt{\Gamma_h}}{\sqrt{2}} \right|^2 \approx \Gamma_p \left| \frac{1 + e^{-i\alpha}}{\sqrt{2}} \right|^2 = \Gamma_p (1 + \cos 2\alpha), \quad (5.5.3)$$

where

$$\alpha = (1 - c)\pi/2, \quad (5.5.4)$$

c being defined by the relation $(\tau \hat{F} \tau^{-1})^\dagger = -c \hat{F}$, where \hat{F} is the single-particle operator entering in the particle-vibration coupling Hamiltonian (see Eqs. (1.2.3)–(1.2.6)). In the case in which the (quasi-) bosonic mode is a low-lying surface vibration (see Eq. (1.2.7))⁶³, $c = -1$. As a result, a conspicuous cancellation between the particle- and hole-contributions⁶⁴ takes place (see diagrams (a) and (b) of Fig. 5.5.3), where typical amplitudes contributing to the damping width Γ_{GDR} of ^{208}Pb are displayed). It is of notice that in the case of the giant pairing vibration, built out of two correlated particles (or holes), the corresponding contributions (see diagrams (c) and (d), Fig. 5.5.3) have the same sign. That is, $c = +1$, and $\Gamma_{GPV} \approx 2\Gamma_p(\hbar\omega_{GPV}/2) \approx \hbar\omega_{GPV}$. As a result one would be confronted with a resonant state displaying a damping width about equal to the value of its centroid energy.

5.6 Single-nucleon transfer for pedestrians

In what follows we discuss some aspects of the relations existing between nuclear structure and one-particle transfer cross sections. To do so, we repeat some of the steps carried out in Sect. 5.1, but this time in a simpler way, essentially ignoring the implications associated with the spin carried out by the particles, the spin-orbit dependence of the optical model potential, the recoil effect, etc.

⁶²Bertrand (1981).

⁶³The precise formulation of this result is that in the case in which the field \hat{F} is spin- and isospin-independent, $c = -1$ (see Bortignon et al. (1983), see also Bortignon and Broglia (1981)).

⁶⁴The fact that expression (5.5.3) leads to complete cancellation is in keeping with the approximation that the single-particle and -hole subspaces are identical (i.e. $\Gamma_p = \Gamma_h$). In fact, the finite value of Γ_{GR} in general and of Γ_{GDR} in particular, is associated with the asymmetry existing between particle- and hole- subspaces.

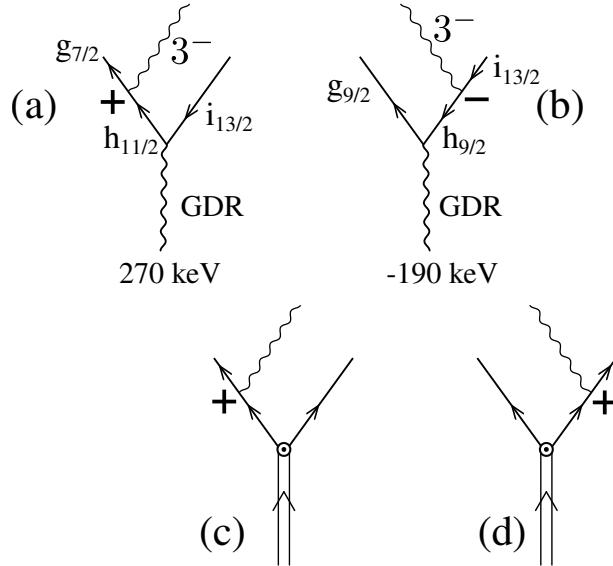


Figure 5.5.3: (a), (b) Example of cancellation between the particle- and hole- doorway contributions to the damping width Γ_{GDR} (Bortignon et al. (1983)); (c), (d) similar diagrams this time associated with a pair addition mode in general, and with the GPV in particular.

We consider the case of $A(d, p)A + 1$ reaction, namely that of neutron stripping. The intrinsic wave functions ψ_α and ψ_β , where $\alpha = (A, d)$ and $\beta = ((A + 1), p)$,

$$\psi_\alpha = \psi_{M_A}^{I_A}(\xi_A)\phi_d(\vec{r}_{np}), \quad (5.6.1a)$$

$$\begin{aligned} \psi_\beta &= \psi_{M_{A+1}}^{I_{A+1}}(\xi_{A+1}) \\ &= \sum_{l, I'_A} (I'_A; l | I_{A+1}) [\psi^{I_A}(\xi_A)\phi^l(\vec{r}_n)]_{M_{A+1}-M_A}^{I_{A+1}}, \end{aligned} \quad (5.6.1b)$$

where $(I'_A; l | I_{A+1})$ is a generalized fractional parentage coefficient. It is of notice that, as a rule, $(I'_A; l | I_{A+1})\phi^l(\vec{r}_n)$ should be able to describe a dressed quasiparticle state containing only a fraction of the single particle strength. Although ignoring possible radial renormalization we assume, for simplicity, the expansion to be operative. To further simplify the derivation we assume we are dealing with spinless particles. This is the reason why no “intrinsic” proton wavefunction appears in Eq. (5.6.1b). The variable \vec{r}_{np} is the relative coordinate of the proton and the neutron (see Fig. 5.6.1).

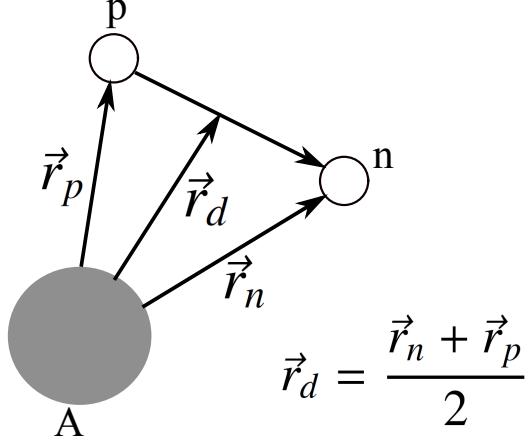


Figure 5.6.1: Coordinates used in the description of the $A(d, p)(A + 1)$ stripping process.

The transition matrix element can now be written as

$$\begin{aligned}
 T_{d,p} &= \langle \psi_{M_{A+1}}^{I_{A+1}}(\xi_{A+1}) \chi_p^{(-)}(\mathbf{k}_p, \mathbf{r}_p), V'_\beta \psi_{M_A}^{I_A}(\xi_A) \chi_d^{(+)}(\mathbf{k}_d, \mathbf{r}_d) \rangle \\
 &= \sum_{\substack{l, l'_A \\ M'_A}} (I'_A; l | I_{A+1}) (I'_A M'_A l M_{A+1} - M'_A | I_{A+1} M_{A+1}) \\
 &\quad \times \int d\mathbf{r}_n d\mathbf{r}_p \chi_p^{*(-)}(\mathbf{k}_p, \mathbf{r}_p) \phi_{M_{A+1}-M'_A}^{*l}(\mathbf{r}_n) (\psi_{M_A}^{I_A}(\xi_A), V'_\beta \psi_{M'_A}^{I'_A}(\xi_A)) \\
 &\quad \times \phi_d(\mathbf{r}_{np}) \chi_d^{(+)}(\mathbf{k}_d, \mathbf{r}_d) \delta_{I'_A, I_A} \delta_{M'_A, M_A}.
 \end{aligned} \tag{5.6.2}$$

In the stripping approximation

$$\begin{aligned}
 V'_\beta &= V_\beta(\xi, \mathbf{r}_\beta) - \bar{U}_\beta(r_\beta) \\
 &= V_\beta(\xi_A, \mathbf{r}_{pA}) + V_\beta(\mathbf{r}_{pn}) - \bar{U}_\beta(r_{pA}).
 \end{aligned} \tag{5.6.3}$$

Then

$$\begin{aligned}
 (\psi_{M_A}^{I_A}(\xi_A), V'_\beta \psi_{M_A}^{I_A}(\xi_A)) &= (\psi_{M_A}^{I_A}(\xi_A), V_\beta(\xi_A, \mathbf{r}_{pA}) \psi_{M_A}^{I_A}(\xi_A)) \\
 &\quad + (\psi_{M_A}^{I_A}(\xi_A), V_\beta(\mathbf{r}_{pn}) \psi_{M_A}^{I_A}(\xi_A)) - \bar{U}_\beta(r_{pA}).
 \end{aligned} \tag{5.6.4}$$

We assume

$$U_\beta(r_{pA}) = (\psi_{M_A}^{I_A}(\xi_A), V_\beta(\xi_A, \mathbf{r}_{pA}) \psi_{M_A}^{I_A}(\xi_A)). \tag{5.6.5}$$

Then

$$(\psi_{M_A}^{I_A}(\xi_A), V'_\beta \psi_{M_A}^{I_A}(\xi_A)) = V_{np}(\mathbf{r}_{pn}). \tag{5.6.6}$$

Inserting eq. (5.6.6) into eq. (5.6.2) we obtain

$$\begin{aligned} T_{d,p} &= \sum_l (I_A; l | I_{A+1}) (I_A M_A l M_{A+1} - M_A | I_{A+1} M_{A+1}) \\ &\times \int d\mathbf{r}_n d\mathbf{r}_p \chi_p^{*(-)}(\mathbf{k}_p, \mathbf{r}_p) \phi_{M_{A+1}-M_A}^{*l}(\mathbf{r}_n) V(\mathbf{r}_{pn}) \phi_d(\mathbf{r}_{np}) \chi_d^{(+)}(\mathbf{k}_d, \mathbf{r}_d). \end{aligned} \quad (5.6.7)$$

The differential cross section is then equal to

$$\frac{d\sigma}{d\Omega} = \frac{2}{3} \frac{\mu_p \mu_d}{(2\pi\hbar^2)^2} \frac{(2I_{A+1} + 1)}{(2I_A + 1)} \frac{k_p}{k_d} \sum_{l,m_l} \frac{(I_A; l | I_{A+1})^2}{2l + 1} |B_{m_l}^l|^2, \quad (5.6.8)$$

where

$$B_{m_l}^l(\theta) = \int d\mathbf{r}_n d\mathbf{r}_p \chi_p^{*(-)}(\mathbf{k}_p, \mathbf{r}_p) Y_m^l(\hat{\mathbf{r}}_n) u_{nl}(r_n) V(\mathbf{r}_{pn}) \phi_d(\mathbf{r}_{np}) \chi_d^{(+)}(\mathbf{k}_d, \mathbf{r}_d) \quad (5.6.9)$$

and

$$\phi_m^l(\mathbf{r}_n) = u_{nl}(r_n) Y_m^l(\hat{\mathbf{r}}_n), \quad (5.6.10)$$

is the single-particle wave function of a neutron bound to the core A. For simplicity, the radial wave function $u_{nl}(r_n)$ can be assumed to be a solution of a Saxon-Woods potential of parameters $V_0 \approx 50$ MeV, $a = 0.65$ fm and $r_0 = 1.25$ fm.

The relation (5.6.8) gives the cross section for the stripping from the projectile of a neutron that would correspond to the n^{th} valence neutron in the nucleus $(A+1)$. If we now want the cross section for stripping any of the valence neutrons of the final nucleus from the projectile, we must multiply eq. (5.6.8) by n . A more careful treatment of the antisymmetry with respect to the neutrons shows this to be the correct answer.

Finally we get

$$\frac{d\sigma}{d\Omega} = \frac{(2I_{A+1} + 1)}{(2I_A + 1)} \sum_l S_l \sigma_l(\theta), \quad (5.6.11)$$

where

$$S_l = n(I_A; l | I_{A+1})^2, \quad (5.6.12)$$

and

$$\sigma_l(\theta) = \frac{2}{3} \frac{\mu_p \mu_d}{(2\pi\hbar^2)^2} \frac{k_p}{k_d} \frac{1}{2l + 1} \sum_m |B_m^l|^2 \quad (5.6.13)$$

Distorted wave codes evaluate numerically the quantity $B_{m_l}^l(\theta)$, using for the wave functions $\chi^{(-)}$ and $\chi^{(+)}$ the solution of the optical potentials that fit the elastic scattering, i.e.

$$(-\nabla^2 + \bar{U} - k^2)\chi = 0, \quad (5.6.14)$$

Note that if the target nucleus is even-even, $I_A = 0$, $l = I_{A+1}$. That is, only one l value contributes in Eq. (5.6.8), and the angular distribution is uniquely given by $\sum_m |B_m^l|^2$. The l -dependence of the angular distributions helps to identify I_{A+1} . The

factor S_l needed to normalize the calculated function to the data yields (assuming a good fit to the angular distribution), is known in the literature as the spectroscopic factor. In the early stages of studies of nuclear structure with one-particle transfer reactions it was assumed, not only that such quantity could be accurately defined, but also that it contained all the nuclear structure information (aside from that associated with the angular distribution) which could be extracted from single-particle transfer. In other words, that it was the bridge directly connecting theory with experiment. *Because nucleons are never bare, but are dressed by the coupling to collective modes, coupling which renormalizes not only the single-particle content but also the single-particle wavefunctions (formfactors), the spectroscopic factor approximation is not straightforward to justify.*

There is a fundamental problem which makes the handling of integrals like that of (5.6.9) difficult to handle from a structure point of view, if not numerically at least conceptually. This difficulty is connected with the so called recoil effect⁶⁵, namely the fact that the center of mass of the two interacting particles in entrance ($\mathbf{r}_\alpha : \alpha = a + A$) and exit ($\mathbf{r}_\beta : \beta = b + B$) channels is different. This is at variance with what one is accustomed to deal with in nuclear structure calculations, in which the Hartree potential depends on a single coordinate, as well as in the case of elastic and inelastic reactions, situations in which $\mathbf{r}_\alpha = \mathbf{r}_\beta$. When $\mathbf{r}_\alpha \neq \mathbf{r}_\beta$ we enter a rather more complex many-body problem, in particular if continuum states are to be considered, than nuclear structure practitioners are accustomed to.

Of notice that similar difficulties have been faced in connection with the non-local Fock (exchange) potential. As a rule, the corresponding (HF) mean field equations are rendered local making use of the k -mass approximation or within the framework of Local Density Functional Theory (DFT).

Returning to the subject of the present Section, it is in general useful to be able to introduce approximations which can help the physics which is at the basis of the phenomenon under discussion (single-particle motion), to emerge in a natural way, if not to compare in detail with the experimental data. Within this context, to reduce the integral (5.6.9) one can assume that the proton-neutron interaction V_{np} has zero-range, i.e.

$$V_{np}(\mathbf{r}_{np})\phi_d(\mathbf{r}_{np}) = D_0\delta(\mathbf{r}_{np}) \quad (5.6.15)$$

so that B_m^l becomes equal to

$$B_m^l(\theta) = D_0 \int d\mathbf{r} \chi_p^{*(-)}(\mathbf{k}_p, \mathbf{r}) Y_{m_l}^{*l}(\hat{\mathbf{r}}) u_l(r) \chi_d^{(+)}(\mathbf{k}_d, \mathbf{r}), \quad (5.6.16)$$

which is a three dimensional integral, but in fact essentially a one-dimensional integral, as the integration over the angles can be worked out analytically.

⁶⁵While this effect could be treated in a cavalier fashion in the case of light ion reactions ($m_a/m_A \ll 1$), this was not possible in the case of heavy ion reactions, as the change in momenta involved was always sizeable (cf. Broglia and Winther (2004) and refs. therein).

5.6.1 Plane-wave limit

If in Eq. (5.6.14) one sets $\bar{U} = 0$, the distorted waves become plane waves i.e.

$$\chi_d^{(+)}(\mathbf{k}_d, \mathbf{r}) = e^{i\mathbf{k}_d \cdot \mathbf{r}}, \quad (5.6.17a)$$

$$\chi_d^{(-)}(\mathbf{k}_p, \mathbf{r}) = e^{-i\mathbf{k}_p \cdot \mathbf{r}}. \quad (5.6.17b)$$

Equation (5.6.16) can now be written as

$$B_m^l = D_0 \int d\mathbf{r} e^{i(\mathbf{k}_d - \mathbf{k}_p) \cdot \mathbf{r}} Y_m^{*l}(\hat{\mathbf{r}}) u_l(r). \quad (5.6.18)$$

The linear momentum transferred to the nucleus is $\mathbf{k}_d - \mathbf{k}_p = \mathbf{q}$. Let us expand $e^{i\mathbf{q} \cdot \mathbf{r}}$ in spherical harmonics, i.e.

$$\begin{aligned} e^{i\mathbf{q} \cdot \mathbf{r}} &= \sum_l l! j_l(qr) (2l+1) P_l(\hat{\mathbf{q}} \cdot \hat{\mathbf{r}}) \\ &= 4\pi \sum_l l! j_l(qr) \sum_m Y_m^{*l}(\hat{\mathbf{q}}) Y_m^l(\hat{\mathbf{r}}), \end{aligned} \quad (5.6.19)$$

so

$$\int d\hat{\mathbf{r}} e^{i\mathbf{q} \cdot \mathbf{r}} Y_m^l(\hat{\mathbf{r}}) = 4\pi l! j_l(qr) Y_m^{*l}(\hat{\mathbf{q}}). \quad (5.6.20)$$

Then

$$\begin{aligned} \sum_m |B_m^l|^2 &= \sum_m |Y_m^l(\hat{\mathbf{q}})|^2 D_0^2 16\pi^2 \\ &\times \left| \int r^2 dr j_l(qr) u_l(r) \right|^2 \\ &= \frac{2l+1}{4\pi} D_0^2 16\pi^2 \left| \int r^2 dr j_l(qr) u_l(r) \right|^2. \end{aligned} \quad (5.6.21)$$

Thus, the angular distribution is given by the integral $\left| \int r^2 dr j_l(qr) u_l(r) \right|^2$. If one assumes that the process takes place mostly on the surface, the angular distribution will be given by $|j_l(qR_0)|^2$, where R_0 is the nuclear radius.

We then have

$$\begin{aligned} q^2 &= k_d^2 + k_p^2 - 2k_d k_p \cos(\theta) \\ &= (k_d^2 + k_p^2 - 2k_d k_p) + 2k_d k_p (1 - \cos(\theta)) \\ &= (k_d - k_p)^2 + 4k_d k_p (\sin(\theta/2))^2 \\ &\approx 4k_d k_p (\sin(\theta/2))^2, \end{aligned} \quad (5.6.22)$$

since $k_d \approx k_p$ for stripping reactions at typical bombarding energies. Thus the angular distribution has a diffraction-like structure given by

$$|j_l(qR_0)|^2 = j_l^2(2R_0 \sqrt{k_d k_p} \sin(\theta/2)). \quad (5.6.23)$$

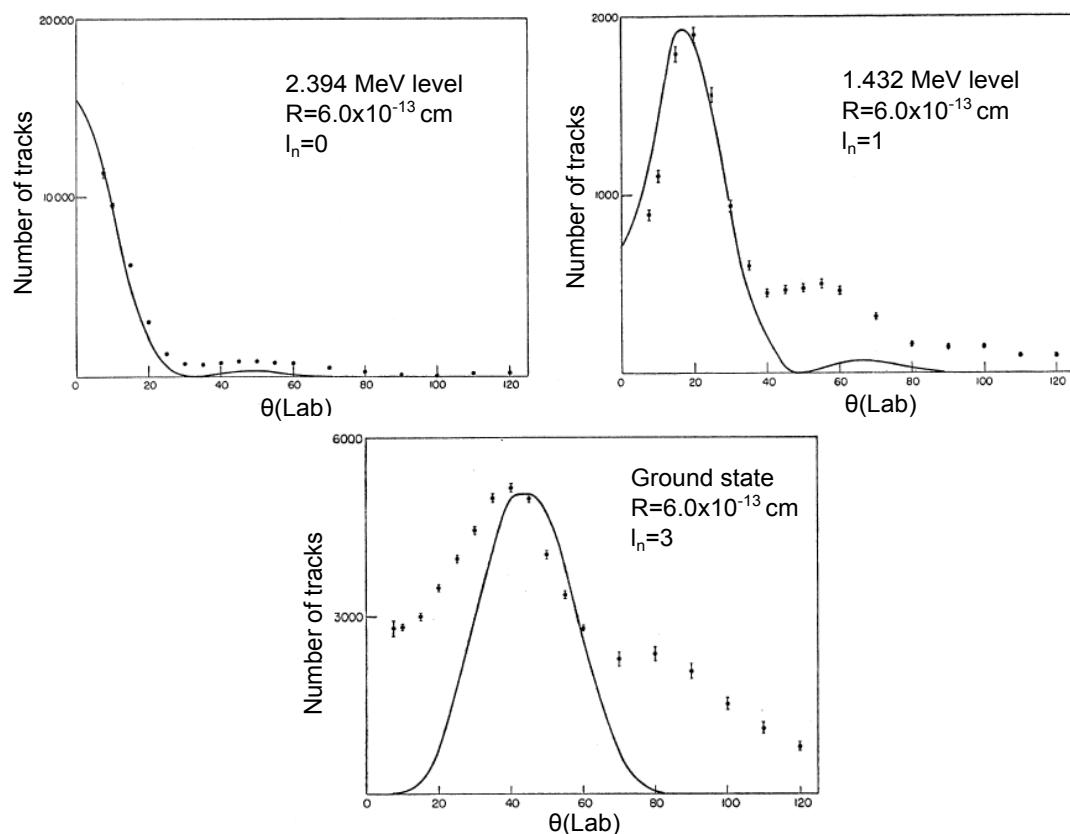


Figure 5.6.2: Plane wave approximation analysis of three $^{44}\text{Ca}(\text{d},\text{p})^{45}\text{Ca}$ differential cross sections leading to the ground state ($l_n = 3$) and to the 1.9 MeV ($l_n = 1$) and 2.4 MeV ($l_n = 0$) excited states, i.e. $f_{9/2}$, $p_{1/2}$ and $s_{1/2}$ states (Cobb and Guthe, 1957).

The function $j_l(x)$ has its first maximum at $x = l$, i.e. where

$$\sin(\theta/2) = \frac{l}{2R_0 k}, \quad (k_p \approx k_d = k), \quad (5.6.24)$$

Examples of the above relation are provided in Fig. 5.6.2.

5.7 One-particle knockout within DWBA

5.7.1 Spinless particles

We consider the reaction $A + a \rightarrow a + b + c$, in which particle b (group of particles, cluster) is knocked out from the nucleus $A (= c+b)$. Cluster b is thus initially bound, while the final states of a, b and the initial state of a are all in the continuum, and can be described with distorted waves defined as scattering solutions of an optical potential. A schematic depiction of the situation is shown in Fig. 5.7.1. While the derivation presented below is quite general, special emphasis is set to one-particle knock-out processes.

Transition amplitude

A first derivation will be given in which, for simplicity, all the “particles” (nuclei) involved in the reaction process are spinless and inert. Use is made of central, complex optical potentials ($U(r_{aA}), U(r_{cb}), U(r_{ac})$) without a spin-orbit term. In addition, the interaction $v(r_{ab})$ between a and b is taken to be a function of the distance r_{ab} . Within this scenario, the transition amplitude which is at the basis of the evaluation of the multi-differential cross section is the 6-dimensional integral

$$T_{mb} = \int d\mathbf{r}_{aA} d\mathbf{r}_{bc} \chi^{(-)*}(\mathbf{r}_{ac}) \chi^{(-)*}(\mathbf{r}_{bc}) v(r_{ab}) \chi^{(+)}(\mathbf{r}_{aA}) u_{lb}(r_{bc}) Y_m^{l_b}(\hat{\mathbf{r}}_{bc}). \quad (5.7.1)$$

Coordinates

The vectors $\mathbf{r}_{ab}, \mathbf{r}_{ac}$ can easily be written in function of the integration variables $\mathbf{r}_{aA}, \mathbf{r}_{bc}$ (see Fig. 5.7.1), namely

$$\begin{aligned} \mathbf{r}_{ac} &= \mathbf{r}_{aA} + \frac{b}{A} \mathbf{r}_{bc}, \\ \mathbf{r}_{ab} &= \mathbf{r}_{aA} - \frac{c}{A} \mathbf{r}_{bc}, \end{aligned} \quad (5.7.2)$$

where b, c, A stand for the number of nucleons of the species b, c and A respectively.

Distorted waves in the continuum

A standard way to reduce the dimensionality of the integral (5.7.1) consists in expanding the continuum wave functions $\chi^{(+)}(\mathbf{r}_{aA}), \chi^{(-)*}(\mathbf{r}_{ac}), \chi^{(-)*}(\mathbf{r}_{bc})$ in a basis

of eigenstates of the angular momentum operator (partial waves). Then one can exploit the transformation properties of these eigenstates under rotations to conveniently carry out the angular integrations. Making use of time-reversed phasing, that is

$$Y_m^l(\theta, \phi) = i^l \sqrt{\frac{2l+1}{4\pi} \frac{(l-m)!}{(l+m)!}} P_l^m(\cos \theta) e^{im\phi}, \quad (5.7.3)$$

the general form of these expansions is

$$\chi^{(+)}(\mathbf{k}, \mathbf{r}) = \sum_l \frac{4\pi}{kr} i^l \sqrt{2l+1} e^{i\sigma_l} F_l(r) [Y^l(\hat{\mathbf{r}}) Y^l(\hat{\mathbf{k}})]_0^0, \quad (5.7.4)$$

and

$$\chi^{(-)*}(\mathbf{k}, \mathbf{r}) = \sum_l \frac{4\pi}{kr} i^{-l} \sqrt{2l+1} e^{i\sigma_l} F_l(r) [Y^l(\hat{\mathbf{r}}) Y^l(\hat{\mathbf{k}})]_0^0, \quad (5.7.5)$$

σ_l being the Coulomb phase shift. The radial functions $F_l(r)$ are regular (finite at $r = 0$) solutions of the one-dimensional Schrödinger equation with an effective potential $U(r) + \frac{\hbar^2 l(l+1)}{2\mu r^2}$ and suitable asymptotic behaviour at $r \rightarrow \infty$ as boundary conditions. Thus, the distorted waves appearing in (5.7.1) are,

$$\chi^{(+)}(\mathbf{k}_a, \mathbf{r}_{aA}) = \sum_{l_a} \frac{4\pi}{k_a r_{aA}} i^{l_a} \sqrt{2l_a+1} e^{i\sigma^{l_a}} F_{l_a}(r_{aA}) [Y^{l_a}(\hat{\mathbf{r}}_{aA}) Y^{l_a}(\hat{\mathbf{k}}_a)]_0^0, \quad (5.7.6)$$

describing the relative motion of A and a in the entrance channel as determined by the complex optical potential $U(r_{aA})$,

$$\chi^{(-)*}(\mathbf{k}'_a, \mathbf{r}_{ac}) = \sum_{l'_a} \frac{4\pi}{k'_a r_{ac}} i^{-l'_a} \sqrt{2l'_a+1} e^{i\sigma^{l'_a}} F_{l'_a}(r_{ac}) [Y^{l'_a}(\hat{\mathbf{r}}_{ac}) Y^{l'_a}(\hat{\mathbf{k}}'_a)]_0^0, \quad (5.7.7)$$

which describes the relative motion of c and a , in the final channel controlled by the complex optical potential $U(r_{ac})$, and finally

$$\chi^{(-)*}(\mathbf{k}'_b, \mathbf{r}_{bc}) = \sum_{l'_b} \frac{4\pi}{k'_b r_{bc}} i^{-l'_b} \sqrt{2l'_b+1} e^{i\sigma^{l'_b}} F_{l'_b}(r_{bc}) [Y^{l'_b}(\hat{\mathbf{r}}_{bc}) Y^{l'_b}(\hat{\mathbf{k}}'_b)]_0^0, \quad (5.7.8)$$

final channel wavefunction describing the relative motion of b and c , as defined by the complex optical potential $U(r_{bc})$.

Recoupling of angular momenta

One now proceeds to the evaluation of the 6-dimensional integral

$$\begin{aligned} & \frac{64\pi^3}{k_a k'_a k'_b} \int d\mathbf{r}_{aA} d\mathbf{r}_{bc} u_{l_b}(r_{cb}) v(r_{ab}) \sum_{l_a, l'_a, l'_b} \sqrt{(2l_a+1)(2l'_a+1)(2l'_b+1)} \\ & \times e^{i(\sigma^{l_a} + \sigma^{l'_a} + \sigma^{l'_b})} \frac{F_{l_a}(r_{aA}) F_{l'_a}(r_{ac}) F_{l'_b}(r_{bc})}{r_{ac} r_{aA} r_{bc}} \\ & \times [Y^{l_a}(\hat{\mathbf{r}}_{aA}) Y^{l_a}(\hat{\mathbf{k}}_a)]_0^0 [Y^{l'_a}(\hat{\mathbf{r}}_{ac}) Y^{l'_a}(\hat{\mathbf{k}}'_a)]_0^0 [Y^{l'_b}(\hat{\mathbf{r}}_{bc}) Y^{l'_b}(\hat{\mathbf{k}}'_b)]_0^0 Y_{m_b}^{l_b}(\hat{\mathbf{r}}_{bc}), \end{aligned} \quad (5.7.9)$$

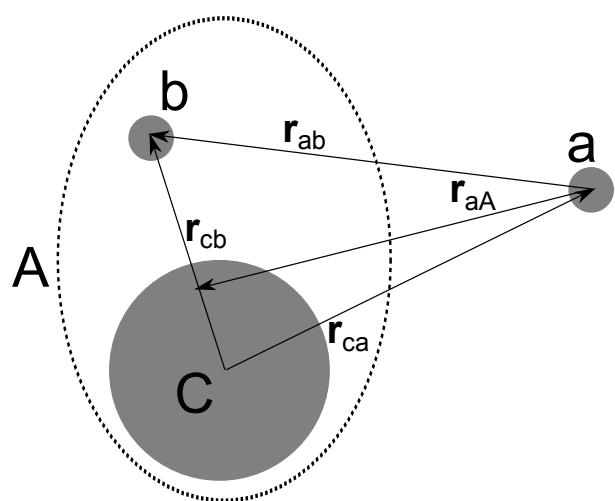


Figure 5.7.1: System of coordinates used to describe the reaction $A + a \rightarrow a + b + c$. The nucleus A is viewed as an inert cluster b bounded to an inert core c .

an expression which explicitly depends on the asymptotic kinetic energies and scattering angles $(\hat{\mathbf{k}}_a, \hat{\mathbf{k}}'_a, \hat{\mathbf{k}}'_b)$ of a, b as determined by k_a, k'_a, k'_b and $\hat{\mathbf{k}}_a, \hat{\mathbf{k}}'_a, \hat{\mathbf{k}}'_b$ respectively. In what follows we will take advantage of the partial wave expansion to reduce the dimensionality of the integral from 6 to 3. A possible strategy to follow is that of recoupling together all the terms that depend on the integration variables to a global angular momentum and retain only the term coupled to 0 as the only one surviving the integration. Let us start to separately couple the terms corresponding to particles a and b . For particle a we write

$$\begin{aligned} [Y^{l_a}(\hat{\mathbf{r}}_{aA}) Y^{l_a}(\hat{\mathbf{k}}_a)]_0^0 [Y'^{l_a}(\hat{\mathbf{r}}_{ac}) Y'^{l_a}(\hat{\mathbf{k}}'_a)]_0^0 &= \sum_K ((l_a l_a)_0 (l'_a l'_a)_0 | (l_a l'_a)_K (l_a l'_a)_K)_0 \\ &\times \left\{ [Y^{l_a}(\hat{\mathbf{r}}_{aA}) Y^{l_a}(\hat{\mathbf{r}}_{ac})]^K [Y^{l_a}(\hat{\mathbf{k}}_a) Y'^{l_a}(\hat{\mathbf{k}}'_a)]_0^K \right\}_0. \end{aligned} \quad (5.7.10)$$

We can now evaluate the $9j$ -symbol,

$$((l_a l_a)_0 (l'_a l'_a)_0 | (l_a l'_a)_K (l_a l'_a)_K)_0 = \sqrt{\frac{2K+1}{(2l'_a+1)(2l_a+1)}}, \quad (5.7.11)$$

and expand the coupling,

$$\begin{aligned} \left\{ [Y^{l_a}(\hat{\mathbf{r}}_{aA}) Y^{l_a}(\hat{\mathbf{r}}_{ac})]^K [Y^{l_a}(\hat{\mathbf{k}}_a) Y'^{l_a}(\hat{\mathbf{k}}'_a)]_0^K \right\}_0 &= \sum_M \langle K \ K \ M \ - M | 0 \ 0 \rangle \\ &\times [Y^{l_a}(\hat{\mathbf{r}}_{aA}) Y^{l_a}(\hat{\mathbf{r}}_{ac})]_M^K [Y^{l_a}(\hat{\mathbf{k}}_a) Y'^{l_a}(\hat{\mathbf{k}}'_a)]_{-M}^K = \sum_M \frac{(-1)^{K+M}}{\sqrt{2K+1}} \\ &\times [Y^{l_a}(\hat{\mathbf{r}}_{aA}) Y^{l_a}(\hat{\mathbf{r}}_{ac})]_M^K [Y^{l_a}(\hat{\mathbf{k}}_a) Y'^{l_a}(\hat{\mathbf{k}}'_a)]_{-M}^K. \end{aligned} \quad (5.7.12)$$

Thus,

$$\begin{aligned} [Y^{l_a}(\hat{\mathbf{r}}_{aA}) Y^{l_a}(\hat{\mathbf{k}}_a)]_0^0 [Y'^{l_a}(\hat{\mathbf{r}}_{ac}) Y'^{l_a}(\hat{\mathbf{k}}'_a)]_0^0 &= \sqrt{\frac{1}{(2l'_a+1)(2l_a+1)}} \\ &\times \sum_{KM} (-1)^{K+M} [Y^{l_a}(\hat{\mathbf{r}}_{aA}) Y^{l_a}(\hat{\mathbf{r}}_{ac})]_M^K [Y^{l_a}(\hat{\mathbf{k}}_a) Y'^{l_a}(\hat{\mathbf{k}}'_a)]_{-M}^K. \end{aligned} \quad (5.7.13)$$

One can further simplify the above expression by choosing the direction of the initial momentum to be parallel to the z axis, so that $Y_m^{l_a}(\hat{\mathbf{k}}_a) = Y_m^{l_a}(\hat{\mathbf{z}}) = \sqrt{\frac{2l_a+1}{4\pi}} \delta_{m,0}$. Then,

$$\begin{aligned} [Y^{l_a}(\hat{\mathbf{r}}_{aA}) Y^{l_a}(\hat{\mathbf{k}}_a)]_0^0 [Y'^{l_a}(\hat{\mathbf{r}}_{ac}) Y'^{l_a}(\hat{\mathbf{k}}'_a)]_0^0 &= \sqrt{\frac{1}{4\pi(2l'_a+1)}} \sum_{KM} (-1)^{K+M} \\ &\times \langle l_a \ 0 \ l'_a \ - M | K \ - M \rangle [Y^{l_a}(\hat{\mathbf{r}}_{aA}) Y^{l_a}(\hat{\mathbf{r}}_{ac})]_M^K Y_{-M}^{l'_a}(\hat{\mathbf{k}}'_a). \end{aligned} \quad (5.7.14)$$

For particle b we have

$$Y_{m_b}^{l_b}(\hat{\mathbf{r}}_{bc}) \left[Y_{l_b}^{l_b}(\hat{\mathbf{r}}_{bc}) Y_{l_b}^{l_b}(\hat{\mathbf{k}}'_b) \right]_0^0 = Y_{m_b}^{l_b}(\hat{\mathbf{r}}_{cb}) \sum_m \frac{(-1)^{l_b' + m}}{\sqrt{2l_b' + 1}} Y_m^{l_b'}(\hat{\mathbf{r}}_{bc}) Y_{-m}^{l_b'}(\hat{\mathbf{k}}'_b), \quad (5.7.15)$$

and can write

$$Y_{m_b}^{l_b}(\hat{\mathbf{r}}_{bc}) Y_m^{l_b'}(\hat{\mathbf{r}}_{bc}) = \sum_{K'} \langle l_b \ m_b \ l'_b \ m | K' \ m_b + m \rangle \left[Y^{l_b}(\hat{\mathbf{r}}_{bc}) Y^{l_b'}(\hat{\mathbf{r}}_{bc}) \right]_{m_b+m}^{K'}. \quad (5.7.16)$$

In order to couple to 0 angular momentum with (5.7.14) we must only keep the term with $K' = K$, $m = -M - m_b$ so

$$\begin{aligned} Y_{m_b}^{l_b}(\hat{\mathbf{r}}_{bc}) \left[Y_{l_b}^{l_b}(\hat{\mathbf{r}}_{bc}) Y_{l_b}^{l_b}(\hat{\mathbf{k}}'_b) \right]_0^0 &= \frac{(-1)^{l_b' - M - m_b}}{\sqrt{2l_b' + 1}} \langle l_b \ m_b \ l'_b \ -M - m_b | K \ -M \rangle \\ &\times \left[Y^{l_b}(\hat{\mathbf{r}}_{bc}) Y^{l_b'}(\hat{\mathbf{r}}_{bc}) \right]_{-M}^K Y_{-M-m_b}^{l_b'}(\hat{\mathbf{k}}'_b), \end{aligned} \quad (5.7.17)$$

and (5.7.9) becomes

$$\begin{aligned} \frac{32\pi^2}{k_a k'_a k'_b} \sum_{KM} (-1)^{K+l'_b-m_b} \langle l_a \ 0 \ l'_a \ -M | K \ -M \rangle \langle l_b \ m_b \ l'_b \ -M - m_b | K \ -M \rangle \\ \times \sum_{l_a, l'_a, l'_b} \sqrt{(2l_a + 1)} e^{i(\sigma^{l_a} + \sigma^{l'_a} + \sigma^{l'_b})} Y_{-M-m_b}^{l_b'}(\hat{\mathbf{k}}'_b) Y_{-M}^{l'_a}(\hat{\mathbf{k}}'_a) \int d\mathbf{r}_{aA} d\mathbf{r}_{bc} u_{l_b}(r_{bc}) v(r_{ab}) \\ \times \frac{F_{l_a}(r_{aA}) F_{l'_a}(r_{ac}) F_{l'_b}(r_{bc})}{r_{ac} r_{aA} r_{bc}} \left[Y^{l_a}(\hat{\mathbf{r}}_{aA}) Y^{l'_a}(\hat{\mathbf{r}}_{ac}) \right]_M^K \left[Y^{l_b}(\hat{\mathbf{r}}_{bc}) Y^{l'_b}(\hat{\mathbf{r}}_{bc}) \right]_{-M}^K. \end{aligned} \quad (5.7.18)$$

Note that

$$\begin{aligned} \left[Y^{l_a}(\hat{\mathbf{r}}_{aA}) Y^{l'_a}(\hat{\mathbf{r}}_{ac}) \right]_M^K \left[Y^{l_b}(\hat{\mathbf{r}}_{bc}) Y^{l'_b}(\hat{\mathbf{r}}_{bc}) \right]_{-M}^K &= \sum_P \langle K \ M \ K \ -M | P \ 0 \rangle \\ &\times \left\{ \left[Y^{l_a}(\hat{\mathbf{r}}_{aA}) Y^{l'_a}(\hat{\mathbf{r}}_{ac}) \right]^K \left[Y^{l_b}(\hat{\mathbf{r}}_{bc}) Y^{l'_b}(\hat{\mathbf{r}}_{bc}) \right]_0^P \right\}_0^P, \end{aligned} \quad (5.7.19)$$

and that to survive the integration the rotational tensors must be coupled to $P = 0$. Keeping only this term in the sum over P , we get

$$\begin{aligned} \left[Y^{l_a}(\hat{\mathbf{r}}_{aA}) Y^{l'_a}(\hat{\mathbf{r}}_{ac}) \right]_M^K \left[Y^{l_b}(\hat{\mathbf{r}}_{bc}) Y^{l'_b}(\hat{\mathbf{r}}_{bc}) \right]_{-M}^K &= \\ \frac{(-1)^{K+M}}{\sqrt{2K+1}} \left\{ \left[Y^{l_a}(\hat{\mathbf{r}}_{aA}) Y^{l'_a}(\hat{\mathbf{r}}_{ac}) \right]^K \left[Y^{l_b}(\hat{\mathbf{r}}_{bc}) Y^{l'_b}(\hat{\mathbf{r}}_{bc}) \right]_0^K \right\}_0^0. \end{aligned} \quad (5.7.20)$$

The coordinate-dependent part of the latter expression is a rotationally invariant scalar, so it can be evaluated in any conventional “standard” configuration such as the one depicted in Fig. 5.7.2. It must then be multiplied by a factor resulting

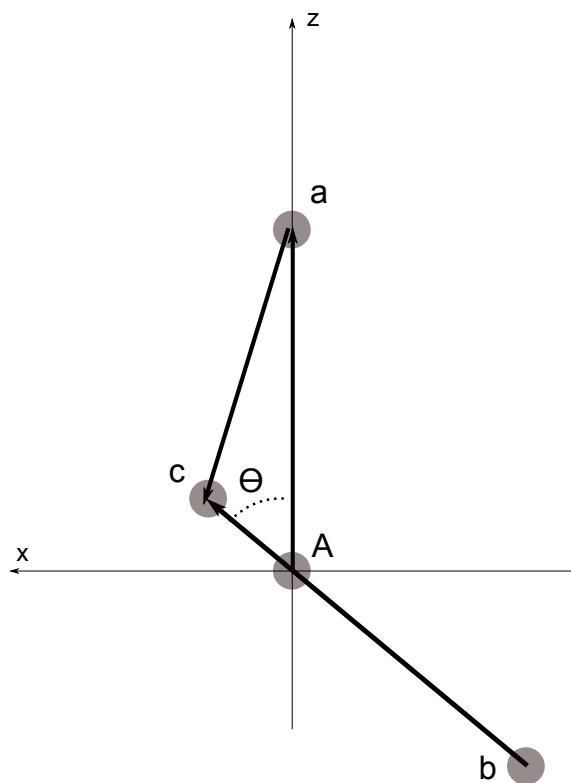


Figure 5.7.2: Coordinates in the “standard” configuration.

of the integration of the remaining angular variables, which accounts for the rigid rotations needed to connect any arbitrary configuration to one of this type. This factor turns out to be $8\pi^2$ (a 4π factor for all possible orientations of, say, \mathbf{r}_{aA} and a 2π factor for a complete rotation around its direction). According to Fig. 5.7.2,

$$\begin{aligned}\mathbf{r}_{bc} &= r_{bc} (\sin \theta \hat{x} + \cos \theta \hat{z}), \\ \mathbf{r}_{aA} &= -r_{aA} \hat{z}, \\ \mathbf{r}_{ac} &= \frac{b}{A} r_{bc} \sin \theta \hat{x} + \left(\frac{b}{A} r_{bc} \cos \theta - r_{aA} \right) \hat{z}.\end{aligned}\quad (5.7.21)$$

As \mathbf{r}_{aA} lies parallel to the z axis, $Y_{M_K}^{l_a}(\hat{\mathbf{r}}_{aA}) = \sqrt{\frac{2l_a+1}{4\pi}} \delta_{M_K,0}$ and

$$\begin{aligned}\left[Y^{l_a}(\hat{\mathbf{r}}_{aA}) Y^{l'_a}(\hat{\mathbf{r}}_{ac}) \right]_{M_K}^K &= \sum_m \langle l_a m | l'_a M_K - m | K M_K \rangle Y_m^{l_a}(\hat{\mathbf{r}}_{aA}) Y_{M_K-m}^{l'_a}(\hat{\mathbf{r}}_{ac}) = \\ &\sqrt{\frac{2l_a+1}{4\pi}} \langle l_a 0 | l'_a M_K | K M_K \rangle Y_{M_K}^{l'_a}(\hat{\mathbf{r}}_{ac}).\end{aligned}\quad (5.7.22)$$

Then

$$\begin{aligned}\left\{ \left[Y^{l_a}(\hat{\mathbf{r}}_{aA}) Y^{l'_a}(\hat{\mathbf{r}}_{ac}) \right]_{M_K}^K \left[Y^{l_b}(\hat{\mathbf{r}}_{bc}) Y^{l'_b}(\hat{\mathbf{r}}_{bc}) \right]_{-M_K}^K \right\}_0^0 &= \\ \sum_{M_K} \langle K M_K | K - M_K | 0 0 \rangle \left[Y^{l_a}(\hat{\mathbf{r}}_{aA}) Y^{l'_a}(\hat{\mathbf{r}}_{ac}) \right]_{M_K}^K \left[Y^{l_b}(\hat{\mathbf{r}}_{bc}) Y^{l'_b}(\hat{\mathbf{r}}_{bc}) \right]_{-M_K}^K &= \\ \sqrt{\frac{2l_a+1}{4\pi}} \sum_{M_K} \frac{(-1)^{K+M_K}}{\sqrt{2K+1}} \langle l_a 0 | l'_a M_K | K M_K \rangle &\\ \times \left[Y^{l_b}(\hat{\mathbf{r}}_{bc}) Y^{l'_b}(\hat{\mathbf{r}}_{bc}) \right]_{-M_K}^K Y_{M_K}^{l'_a}(\hat{\mathbf{r}}_{ac}). &\end{aligned}\quad (5.7.23)$$

Remembering the $8\pi^2$ factor, the term arising from (5.7.20) to be considered in the integral is

$$\begin{aligned}4\pi^{3/2} \frac{\sqrt{2l_a+1}}{2K+1} (-1)^K \sum_{M_K} (-1)^{M_K} \langle l_a 0 | l'_a M_K | K M_K \rangle &\\ \times \left[Y^{l_b}(\cos \theta, 0) Y^{l'_b}(\cos \theta, 0) \right]_{-M_K}^K Y_{M_K}^{l'_a}(\cos \theta_{ac}, 0), &\end{aligned}\quad (5.7.24)$$

with

$$\cos \theta_{ac} = \frac{\frac{b}{A} r_{bc} \cos \theta - r_{aA}}{\sqrt{\left(\frac{b}{A} r_{bc} \sin \theta \right)^2 + \left(\frac{b}{A} r_{bc} \cos \theta - r_{aA} \right)^2}}, \quad (5.7.25)$$

(see (5.7.21)). The final expression of the transition amplitude is

$$\begin{aligned} T_{m_b}(\mathbf{k}'_a, \mathbf{k}'_b) = & \frac{128\pi^{7/2}}{k_a k'_a k'_b} \sum_{KM} \frac{(-1)^{l'_b+m_b}}{2K+1} \langle l_a 0 | l'_a - M | K - M \rangle \langle l_b m_b | l'_b - M - m_b | K - M \rangle \\ & \times \sum_{l_a, l'_a, l'_b} (2l_a + 1) e^{i(\sigma^{l_a} + \sigma^{l'_a} + \sigma^{l'_b})} Y_{-M-m_b}^{l'_b}(\hat{\mathbf{k}}'_b) Y_{-M}^{l'_a}(\hat{\mathbf{k}}'_a) \mathcal{I}(l_a, l'_a, l'_b, K), \end{aligned} \quad (5.7.26)$$

where

$$\begin{aligned} \mathcal{I}(l_a, l'_a, l'_b, K) = & \int dr_{aA} dr_{bc} d\theta r_{aA} r_{bc} \frac{\sin \theta}{r_{ac}} u_{l_b}(r_{bc}) v(r_{ab}) F_{l_a}(r_{aA}) F_{l'_a}(r_{ac}) F_{l'_b}(r_{bc}) \\ & \times \sum_{M_K} (-1)^{M_K} \langle l_a 0 | l'_a | M_K | K | M_K \rangle [Y^{l_b}(\cos \theta, 0) Y^{l'_b}(\cos \theta, 0)]_{-M_K}^K Y_{M_K}^{l'_a}(\cos \theta_{ac}, 0) \end{aligned} \quad (5.7.27)$$

is a 3-dimensional integral that can be numerically evaluated with the help of, e.g., Gaussian integration.

5.7.2 Particles with spin

We now treat the case in which the clusters have a definite spin (see Fig. 5.7.3), and the complex optical potentials $U(r_{aA}), U(r_{cb}), U(r_{ac})$ contain now a spin-orbit term proportional to the product $\mathbf{l} \cdot \mathbf{s} = 1/2(j(j+1) - l(l+1) - 3/4)$ for particles with spin 1/2. In addition, the interaction $V(r_{ab}, \sigma_a, \sigma_b)$ between a and b is taken to be a separable function of the distance r_{ab} and of the spin orientations, $V(r_{ab}, \sigma_a, \sigma_b) = v(r_{ab})v_\sigma(\sigma_a, \sigma_b)$. Note that this ansatz rules out spin-orbit as well as tensor terms in the NN -interaction. For the time being we will assume that the spin-dependent interaction is rotationally invariant (scalar with respect to rotations), such as, e.g., $v_\sigma(\sigma_a, \sigma_b) \propto \sigma_a \cdot \sigma_b$. Again, this assumption excludes tensor terms in the interaction. The transition amplitude is then,

$$\begin{aligned} T_{m_a, m_b}^{m'_a, m'_b} = & \sum_{\sigma_a, \sigma_b} \int d\mathbf{r}_{aA} d\mathbf{r}_{bc} \chi_{m'_a}^{(-)*}(\mathbf{r}_{ac}, \sigma_a) \chi_{m'_b}^{(-)*}(\mathbf{r}_{bc}, \sigma_b) \\ & \times v(r_{ab}) v_\sigma(\sigma_a, \sigma_b) \chi_{m_a}^{(+)}(\mathbf{r}_{aA}, \sigma_a) \psi_{m_b}^{l_b, j_b}(\mathbf{r}_{bc}, \sigma_b). \end{aligned} \quad (5.7.28)$$

Distorted waves

The distorted waves in (5.7.28) $\chi_m(\mathbf{r}, \sigma) = \chi(\mathbf{r})\phi_m^{1/2}(\sigma)$ have a spin dependence contained in the spinor $\phi_m^{1/2}(\sigma)$, where σ is the spin degree of freedom and m the projection of the spin along the quantization axis. The superscript 1/2 reminds us that we are considering spin 1/2 particles, which have important consequences when dealing with the spin-orbit term of the optical potentials. As for the spin-dependent term $v_\sigma(\sigma_a, \sigma_b)$, the actual value of the spin of particles involved in the

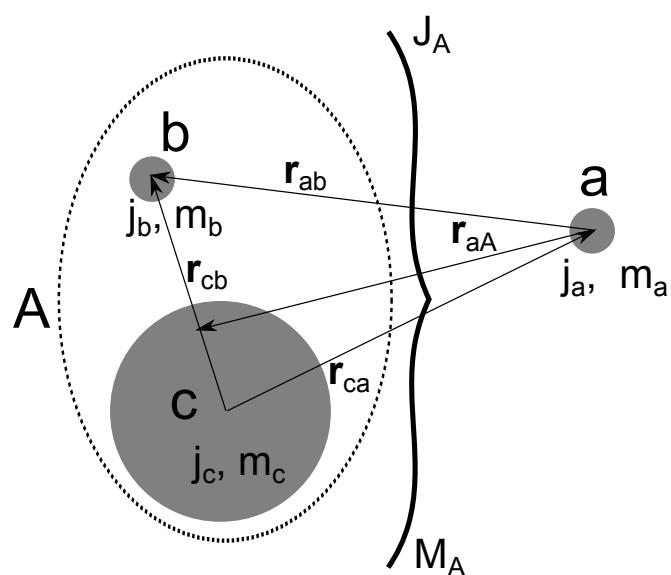


Figure 5.7.3: In the present case all three clusters a, b, c have definite spins and projections. The nucleus A is coupled to total spin J_A, M_A .

reaction process do not make much difference, *as long as this term is rotationally invariant*. Following (5.7.4),

$$\chi^{(+)}(\mathbf{k}, \mathbf{r})\phi_m(\sigma) = \sum_{l,j} \frac{4\pi}{kr} i^l \sqrt{2l+1} e^{i\sigma' l} F_{l,j}(r) [Y^l(\hat{\mathbf{r}}) Y^l(\hat{\mathbf{k}})]_0^0 \phi_m^{1/2}(\sigma). \quad (5.7.29)$$

Note that now one also sums over the total angular momentum j , as the radial functions $F_{l,j}(r)$ depend both on j as well as on l , in keeping with the fact that they are solutions of an optical potential containing a spin-orbit term proportional to $1/2(j(j+1) - l(l+1) - 3/4)$. One must then couple the radial and spin functions to total angular momentum j , noting that

$$\begin{aligned} [Y^l(\hat{\mathbf{r}}) Y^l(\hat{\mathbf{k}})]_0^0 \phi_m^{1/2}(\sigma) &= \sum_{m_l} \langle l m_l l - m_l | 0 0 \rangle Y_{m_l}^l(\hat{\mathbf{r}}) Y_{-m_l}^l(\hat{\mathbf{k}}) \phi_m^{1/2}(\sigma) = \\ &\sum_{m_l} \frac{(-1)^{l-m_l}}{\sqrt{2l+1}} Y_{m_l}^l(\hat{\mathbf{r}}) Y_{-m_l}^l(\hat{\mathbf{k}}) \phi_m^{1/2}(\sigma), \end{aligned} \quad (5.7.30)$$

and

$$Y_{m_l}^l(\hat{\mathbf{r}}) \phi_m^{1/2}(\sigma) = \sum_j \langle l m_l 1/2 m | j m_l + m \rangle [Y^l(\hat{\mathbf{r}}) \phi^{1/2}(\sigma)]_{m_l+m}^j, \quad (5.7.31)$$

we can write

$$\begin{aligned} [Y^l(\hat{\mathbf{r}}) Y^l(\hat{\mathbf{k}})]_0^0 \phi_m^{1/2}(\sigma) &= \sum_{m_l, j} \frac{(-1)^{l+m_l}}{\sqrt{2l+1}} \langle l m_l 1/2 m | j m_l + m \rangle \\ &\times [Y^l(\hat{\mathbf{r}}) \phi^{1/2}(\sigma)]_{m_l+m}^j Y_{-m_l}^l(\hat{\mathbf{k}}), \end{aligned} \quad (5.7.32)$$

and the distorted waves in (5.7.28) are

$$\begin{aligned} \chi_{m_a}^{(+)}(\mathbf{r}_{aA}, \mathbf{k}_a, \sigma_a) &= \sum_{l_a, m_{l_a}, j_a} \frac{4\pi}{k_a r_{aA}} i^{l_a} (-1)^{l_a+m_{l_a}} e^{i\sigma^{l_a}} F_{l_a, j_a}(r_{aA}) \\ &\times \langle l_a m_{l_a} 1/2 m_a | j_a m_{l_a} + m_a \rangle [Y^{l_a}(\hat{\mathbf{r}}_{aA}) \phi^{1/2}(\sigma_a)]_{m_{l_a}+m_a}^{j_a} Y_{-m_{l_a}}^{l_a}(\hat{\mathbf{k}}_a), \end{aligned} \quad (5.7.33)$$

$$\begin{aligned} \chi_{m'_b}^{(-)*}(\mathbf{r}_{bc}, \mathbf{k}'_b, \sigma_b) &= \sum_{l'_b, m'_{l'_b}, j'_b} \frac{4\pi}{k'_b r_{bc}} i^{-l'_b} (-1)^{l'_b+m'_{l'_b}} e^{i\sigma^{l'_b}} F_{l'_b, j'_b}(r_{bc}) \\ &\times \langle l'_b m'_{l'_b} 1/2 m'_b | j'_b m'_{l'_b} + m'_b \rangle [Y^{l'_b}(\hat{\mathbf{r}}_{bc}) \phi^{1/2}(\sigma_b)]_{m'_{l'_b}+m'_b}^{j'_b*} Y_{-m'_{l'_b}}^{l'_b*}(\hat{\mathbf{k}}'_b), \end{aligned} \quad (5.7.34)$$

$$\begin{aligned} \chi_{m'_a}^{(-)*}(\mathbf{r}_{ac}, \mathbf{k}'_a, \sigma_a) &= \sum_{l'_a, m'_{l'_a}, j'_a} \frac{4\pi}{k'_a r_{ac}} i^{-l'_a} (-1)^{l'_a+m'_{l'_a}} e^{i\sigma^{l'_a}} F_{l'_a, j'_a}(r_{ac}) \\ &\times \langle l'_a m'_{l'_a} 1/2 m'_a | j'_a m'_{l'_a} + m'_a \rangle [Y^{l'_a}(\hat{\mathbf{r}}_{ac}) \phi^{1/2}(\sigma_a)]_{m'_{l'_a}+m'_a}^{j'_a*} Y_{-m'_{l'_a}}^{l'_a*}(\hat{\mathbf{k}}'_a). \end{aligned} \quad (5.7.35)$$

The initial bound particle b wavefunction is

$$\psi_{m_b}^{l_b, j_b}(\mathbf{r}_{bc}, \sigma_b) = u_{l_b, j_b}(r_{bc}) \left[Y^{l_b}(\hat{\mathbf{r}}_{bc}) \phi^{1/2}(\sigma_b) \right]_{m_b}^{j_b}, \quad (5.7.36)$$

Substituting in (5.7.28), one obtains,

$$\begin{aligned} T_{m_a, m_b}^{m'_a, m'_b}(\mathbf{k}'_a, \mathbf{k}'_b) &= \frac{64\pi^3}{k_a k'_a k'_b} \sum_{\sigma_a, \sigma_b} \sum_{l_a, m_{l_a}, j_a} \sum_{l'_a, m'_{l'_a}, j'_a} \sum_{l'_b, m'_{l'_b}, j'_b} e^{i(\sigma_a l_a + \sigma'_a l'_a + \sigma'_b l'_b)} i^{l_a - l'_a - l'_b} (-1)^{l_a - m_{l_a} + l'_a - j'_a + l'_b - j'_b} \\ &\times \langle l'_a m'_{l'_a} | 1/2 m'_a | j'_a m'_{l'_a} + m'_a \rangle \langle l_a m_{l_a} | 1/2 m_a | j_a m_{l_a} + m_a \rangle \langle l'_b m'_{l'_b} | 1/2 m'_b | j'_b m'_{l'_b} + m'_b \rangle \\ &\times Y_{-m_{l_a}}^{l'_a}(\hat{\mathbf{k}}_a) Y_{-m'_{l'_a}}^{l'_a}(\hat{\mathbf{k}}'_a) Y_{-m'_{l'_b}}^{l'_b}(\hat{\mathbf{k}}'_b) \int d\mathbf{r}_{aA} d\mathbf{r}_{bc} \left[Y^{l'_a}(\hat{\mathbf{r}}_{ac}) \phi^{1/2}(\sigma_a) \right]_{-m'_{l'_a} - m'_a}^{j'_a} \left[Y^{l'_b}(\hat{\mathbf{r}}_{bc}) \phi^{1/2}(\sigma_b) \right]_{-m'_{l'_b} - m'_b}^{j'_b} \\ &\times \frac{F_{l_a, j_a}(r_{aA}) F_{l'_a, j'_a}(r_{ac}) F_{l'_b, j'_b}(r_{bc})}{r_{ac} r_{aA} r_{bc}} u_{l_b, j_b}(r_{bc}) v(r_{ab}) v_\sigma(\sigma_a, \sigma_b) \\ &\times \left[Y^{l_a}(\hat{\mathbf{r}}_{aa}) \phi^{1/2}(\sigma_a) \right]_{m_{l_a} + m_a}^{j_a} \left[Y^{l_b}(\hat{\mathbf{r}}_{bc}) \phi^{1/2}(\sigma_b) \right]_{m_b}^{j_b}, \quad (5.7.37) \end{aligned}$$

where use was made of the relation

$$\left[Y^l(\hat{\mathbf{r}}) \phi^{1/2}(\sigma) \right]_m^{j*} = (-1)^{j-m} \left[Y^l(\hat{\mathbf{r}}) \phi^{1/2}(\sigma) \right]_{-m}^j. \quad (5.7.38)$$

Recoupling of angular momenta

Let us now separate spatial and spin coordinates, noting that the spin functions must be coupled to $S = 0$, a consequence of the fact that the interaction $v_\sigma(\sigma_a, \sigma_b)$ is rotationally invariant. Starting with particle a ,

$$\begin{aligned} \left[Y^{l'_a}(\hat{\mathbf{r}}_{ac}) \phi^{1/2*}(\sigma_a) \right]_{-m'_{l'_a} - m'_a}^{j'_a} \left[Y^{l_a}(\hat{\mathbf{r}}_{aa}) \phi^{1/2}(\sigma_a) \right]_{m_{l_a} + m_a}^{j_a} &= \\ \sum_K ((l'_a \frac{1}{2})_{j'_a} (l_a \frac{1}{2})_{j_a} | (l_a l'_a)_K (\frac{1}{2} \frac{1}{2})_0)_K & \\ \times \left[Y^{l'_a}(\hat{\mathbf{r}}_{ac}) Y^{l_a}(\hat{\mathbf{r}}_{aa}) \right]_{-m'_{l'_a} - m'_a + m_{l_a} + m_a}^K \left[\phi^{1/2*}(\sigma_a) \phi^{1/2}(\sigma_a) \right]_0^0. \quad (5.7.39) \end{aligned}$$

For particle b ,

$$\begin{aligned} \left[Y^{l'_b}(\hat{\mathbf{r}}_{bc}) \phi^{1/2*}(\sigma_b) \right]_{-m'_{l'_b} - m'_b}^{j'_b} \left[Y^{l_b}(\hat{\mathbf{r}}_{bc}) \phi^{1/2}(\sigma_b) \right]_{m_b}^{j_b} &= \\ \sum_{K'} ((l'_b \frac{1}{2})_{j'_b} (l_b \frac{1}{2})_{j_b} | (l_b l'_b)_{K'} (\frac{1}{2} \frac{1}{2})_0)_{K'} & \\ \times \left[Y^{l'_b}(\hat{\mathbf{r}}_{bc}) Y^{l_b}(\hat{\mathbf{r}}_{bc}) \right]_{-m'_{l'_b} - m'_b + m_b}^{K'} \left[\phi^{1/2*}(\sigma_b) \phi^{1/2}(\sigma_b) \right]_0^0. \quad (5.7.40) \end{aligned}$$

The spin summation yields a constant factor,

$$\sum_{\sigma_a, \sigma_b} \left[\phi^{1/2*}(\sigma_a) \phi^{1/2}(\sigma_a) \right]_0^0 \left[\phi^{1/2*}(\sigma_b) \phi^{1/2}(\sigma_b) \right]_0^0 v_\sigma(\sigma_a, \sigma_b) \equiv T_\sigma, \quad (5.7.41)$$

and what we have yet to do is very similar to what we have done in the case of spinless particles. First of all note that the constrain of coupling all angular momenta to 0, imposes $K' = K$ and $m_{l_a} + m_a - m_{l'_a} - m'_a = m_{l'_b} + m'_b - m_b$ (see (5.7.39) and (5.7.40)). If we set $M = m_{l_a} + m_a - m_{l'_a} - m'_a$ and take, as before, $\hat{\mathbf{k}}_a \equiv \hat{z}$

$$\begin{aligned} T_{m_a, m_b}^{m'_a, m'_b}(\mathbf{k}'_a, \mathbf{k}'_b) &= \frac{32\pi^{5/2}}{k_a k'_a k'_b} T_\sigma \sum_{l_a, j_a} \sum_{l'_a, j'_a} \sum_{l'_b, j'_b} \sum_{K, M} e^{i(\sigma^{l_a} + \sigma^{l'_a} + \sigma^{l'_b})} i^{l_a - l'_a - l'_b} (-1)^{l_a + l'_a + l'_b - j'_a - j'_b} \\ &\times \sqrt{2l_a + 1} ((l'_a \frac{1}{2})_{j'_a} (l_a \frac{1}{2})_{j_a} |(l_a l'_a)_K (\frac{1}{2} \frac{1}{2})_0)_K ((l'_b \frac{1}{2})_{j'_b} (l_b \frac{1}{2})_{j_b} |(l_b l'_b)_K (\frac{1}{2} \frac{1}{2})_0)_K \\ &\times \langle l'_a m_a - m'_a - M \ 1/2 \ m'_a | j'_a \ m_a - M \rangle \langle l_a 0 \ 1/2 \ m_a | j_a \ m_a \rangle \langle l'_b m_b - m'_b + M \ 1/2 \ m'_b | j'_b \ M + m_b \rangle \\ &\times Y_{m'_b - m_b - M}^{l'_b}(\hat{\mathbf{k}}'_b) Y_{m'_a - m_a + M}^{l'_a}(\hat{\mathbf{k}}'_a) \int d\mathbf{r}_{aA} d\mathbf{r}_{bc} \frac{F_{l_a, j_a}(r_{aA}) F_{l'_a, j'_a}(r_{ac}) F_{l'_b, j'_b}(r_{bc})}{r_{ac} r_{aA} r_{bc}} \\ &\times u_{l_b, j_b}(r_{bc}) v(r_{ab}) \left[Y^{l_a}(\hat{\mathbf{r}}_{aA}) Y^{l'_a}(\hat{\mathbf{r}}_{ac}) \right]_M^K \left[Y^{l_b}(\hat{\mathbf{r}}_{bc}) Y^{l'_b}(\hat{\mathbf{r}}_{bc}) \right]_{-M}^K. \quad (5.7.42) \end{aligned}$$

The integral of the above expression is similar to the one in (5.7.18), so we obtain

$$\begin{aligned} T_{m_a, m_b}^{m'_a, m'_b}(\mathbf{k}'_a, \mathbf{k}'_b) &= \frac{128\pi^4}{k_a k'_a k'_b} T_\sigma \sum_{l_a, j_a} \sum_{l'_a, j'_a} \sum_{l'_b, j'_b} \sum_{K, M} e^{i(\sigma^{l_a} + \sigma^{l'_a} + \sigma^{l'_b})} i^{l_a - l'_a - l'_b} (-1)^{l_a + l'_a + l'_b - j'_a - j'_b} \\ &\times \frac{2l_a + 1}{2K + 1} ((l'_a \frac{1}{2})_{j'_a} (l_a \frac{1}{2})_{j_a} |(l_a l'_a)_K (\frac{1}{2} \frac{1}{2})_0)_K ((l'_b \frac{1}{2})_{j'_b} (l_b \frac{1}{2})_{j_b} |(l_b l'_b)_K (\frac{1}{2} \frac{1}{2})_0)_K \\ &\times \langle l'_a m_a - m'_a - M \ 1/2 \ m'_a | j'_a \ m_a - M \rangle \langle l'_b m_b - m'_b + M \ 1/2 \ m'_b | j'_b \ M + m_b \rangle \\ &\times \langle l_a 0 \ 1/2 \ m_a | j_a \ m_a \rangle Y_{m'_b - m_b - M}^{l'_b}(\hat{\mathbf{k}}'_b) Y_{m'_a - m_a + M}^{l'_a}(\hat{\mathbf{k}}'_a) \mathcal{I}(l_a, l'_a, l'_b, j_a, j'_a, j'_b, K), \quad (5.7.43) \end{aligned}$$

with

$$\begin{aligned} \mathcal{I}(l_a, l'_a, l'_b, j_a, j'_a, j'_b, K) &= \int dr_{aA} dr_{bc} d\theta r_{aA} r_{bc} \frac{\sin \theta}{r_{ac}} u_{l_b}(r_{bc}) v(r_{ab}) \\ &\times F_{l_a, j_a}(r_{aA}) F_{l'_a, j'_a}(r_{ac}) F_{l'_b, j'_b}(r_{bc}) \\ &\times \sum_{M_K} \langle l_a 0 \ l'_a \ M_K | K \ M_K \rangle \left[Y^{l_b}(\cos \theta, 0) Y^{l'_b}(\cos \theta, 0) \right]_{-M_K}^K Y_{M_K}^{l'_a}(\cos \theta_{ac}, 0). \quad (5.7.44) \end{aligned}$$

Again, this is a 3-dimensional integral that can be evaluated with the method of Gaussian quadratures. The transition amplitude $T_{m_a, m_b}^{m'_a, m'_b}(\mathbf{k}'_a, \mathbf{k}'_b)$ depends explicitly on the initial (m_a, m'_a) and final (m'_a, m'_b) polarizations of a, b . If the particle b is initially coupled to core c to total angular momentum J_A, M_A , the amplitude to be considered is rather

$$T_{m_a}^{m'_a, m'_b}(\mathbf{k}'_a, \mathbf{k}'_b) = \sum_{m_b} \langle j_b \ m_b \ j_c \ M_A - m_b | J_A \ M_A \rangle T_{m_a, m_b}^{m'_a, m'_b}(\mathbf{k}'_a, \mathbf{k}'_b), \quad (5.7.45)$$

and the multi-differential cross section for detecting particle c (or a) is

$$\frac{d\sigma}{d\mathbf{k}'_a d\mathbf{k}'_b} \Big|_{m_a}^{m'_a, m'_b} = \frac{k'_a \mu_{aA} \mu_{ac}}{k_a 4\pi^2 \hbar^4} \left| \sum_{m_b} \langle j_b m_b j_c M_A - m_b | J_A M_A \rangle T_{m_a, m_b}^{m'_a, m'_b}(\mathbf{k}'_a, \mathbf{k}'_b) \right|^2. \quad (5.7.46)$$

All spin-polarization observables (analysing powers, etc.,) can be derived from this expression. But let us now work out the expression of the cross section for an unpolarized beam (sum over initial spin orientations divided by the number of such orientations) and when we do not detect the final polarizations (sum over final spin orientations),

$$\begin{aligned} \frac{d\sigma}{d\mathbf{k}'_a d\mathbf{k}'_b} &= \frac{k'_a \mu_{aA} \mu_{ac}}{k_a 4\pi^2 \hbar^4} \frac{1}{(2J_A + 1)(2j_a + 1)} \\ &\times \sum_{\substack{m_a, m'_a \\ M_A, m'_b}} \left| \sum_{m_b} \langle j_b m_b j_c M_A - m_b | J_A M_A \rangle T_{m_a, m_b}^{m'_a, m'_b}(\mathbf{k}'_a, \mathbf{k}'_b) \right|^2. \end{aligned} \quad (5.7.47)$$

The sum above can be simplified a bit. Let us consider a single particular value of m_b in the sum over m_b ,

$$\begin{aligned} \sum_{m_a, m'_a, m'_b} \left| T_{m_a, m_b}^{m'_a, m'_b}(\mathbf{k}'_a, \mathbf{k}'_b) \right|^2 \sum_{M_A} \left| \langle j_b m_b j_c M_A - m_b | J_A M_A \rangle \right|^2 = \\ \frac{2J_A + 1}{2j_b + 1} \sum_{m_a, m_a, m'_b} \left| T_{m_a, m_b}^{m'_a, m'_b}(\mathbf{k}'_a, \mathbf{k}'_b) \right|^2 \\ \times \sum_{M_A} \left| \langle J_A - M_A j_c M_A - m_b | j_b m_b \rangle \right|^2, \end{aligned} \quad (5.7.48)$$

where we have used

$$\langle j_b m_b j_c M_A - m_b | J_A M_A \rangle = (-1)^{j_c - M_A + m_b} \sqrt{\frac{2J_A + 1}{2j_b + 1}} \langle J_A - M_A j_c M_A - m_b | j_b m_b \rangle. \quad (5.7.49)$$

As

$$\sum_{M_A} \left| \langle J_A - M_A j_c M_A - m_b | j_b m_b \rangle \right|^2 = 1, \quad (5.7.50)$$

we finally have

$$\frac{d\sigma}{d\mathbf{k}'_a d\mathbf{k}'_b} = \frac{k'_a \mu_{aA} \mu_{ac}}{k_a 4\pi^2 \hbar^4} \frac{1}{(2j_b + 1)(2j_a + 1)} \sum_{m_a, m'_a, m'_b} \left| \sum_{m_b} T_{m_a, m_b}^{m'_a, m'_b}(\mathbf{k}'_a, \mathbf{k}'_b) \right|^2. \quad (5.7.51)$$

Zero range approximation.

The zero range approximation (Eq. (5.6.15)) consists, in the present case to assume $v(r_{ab}) = D_0\delta(r_{ab})$. Then, (see (5.7.21))

$$\begin{aligned}\mathbf{r}_{aA} &= \frac{c}{A}\mathbf{r}_{bc}, \\ \mathbf{r}_{ac} &= \mathbf{r}_{bc}.\end{aligned}\quad (5.7.52)$$

The angular dependence of the integral can be readily evaluated. From (6.A.20), noting that $\hat{\mathbf{r}}_{aA} = \hat{\mathbf{r}}_{ac} = \hat{\mathbf{r}}_{bc} \equiv \hat{\mathbf{r}}$,

$$\begin{aligned}\left[Y^{l_a}(\hat{\mathbf{r}})Y^{l'_a}(\hat{\mathbf{r}})\right]_M^K \left[Y^{l_b}(\hat{\mathbf{r}})Y^{l'_b}(\hat{\mathbf{r}})\right]_{-M}^K = \\ \frac{(-1)^{K-M}}{\sqrt{2K+1}} \left\{ \left[Y^{l_a}(\hat{\mathbf{r}})Y^{l'_a}(\hat{\mathbf{r}})\right]^K \left[Y^{l_b}(\hat{\mathbf{r}})Y^{l'_b}(\hat{\mathbf{r}})\right]^K \right\}_0^0.\end{aligned}\quad (5.7.53)$$

We can as before evaluate this expression in the configuration shown in Fig. 5.7.2 ($\hat{\mathbf{r}} = \hat{z}$), but now the multiplicative factor is 4π . The corresponding contribution to the integral is

$$\frac{(-1)^K}{4\pi(2K+1)} \langle l_a 0 l'_a 0 | K 0 \rangle \sqrt{(2l_a + 1)(2l'_a + 1)(2l_b + 1)(2l'_b + 1)}, \quad (5.7.54)$$

and

$$\begin{aligned}T_{m_a, m_b}^{m'_a, m'_b}(\mathbf{k}'_a, \mathbf{k}'_b) &= \frac{16\pi^2}{k_a k'_a k'_b A} D_0 T_\sigma \sum_{l_a, j_a} \sum_{l'_a, j'_a} \sum_{l'_b, j'_b} \sum_{K, M} e^{i(\sigma^{l_a} + \sigma^{l'_a} + \sigma^{l'_b})} i^{l_a - l'_a - l'_b} (-1)^{l_a + l'_a + l'_b - j'_a - j'_b} \\ &\times \sqrt{(2l_a + 1)(2l'_a + 1)(2l_b + 1)(2l'_b + 1)} \langle l_a 0 l'_a 0 | K 0 \rangle \\ &\times \frac{2l_a + 1}{2K + 1} ((l'_a \frac{1}{2})_{j'_a} (l_a \frac{1}{2})_{j_a} | (l_a l'_a)_K (\frac{1}{2} \frac{1}{2})_0)_K ((l'_b \frac{1}{2})_{j'_b} (l_b \frac{1}{2})_{j_b} | (l_b l'_b)_K (\frac{1}{2} \frac{1}{2})_0)_K \\ &\times \langle l'_a m_a - M 1/2 m'_a | j'_a m_a - M \rangle \langle l'_b m_b - M'_b | j'_b M + m_b \rangle \\ &\times \langle l 0 1/2 m_a | j m_a \rangle Y_{M+m_b+m'_b}^{l'_b}(\hat{\mathbf{k}}'_b) Y_{m_a+m'_a-M}^{l'_a}(\hat{\mathbf{k}}'_a) I_{ZR}(l_a, l'_a, l'_b, j_a, j'_a, j'_b),\end{aligned}\quad (5.7.55)$$

where now the 1-dimensional integral to solve is

$$I_{ZR}(l_a, l'_a, l'_b, j_a, j'_a, j'_b) = \int dr u_l(r) F_{l_a, j_a}(\frac{c}{A}r) F_{l'_a, j'_a}(r) F_{l'_b, j'_b}(r)/r. \quad (5.7.56)$$

5.7.3 One-particle transfer

It may be interesting to state the expression for the one particle transfer reaction within the same context and using the same elements, in order to better compare these two type of experiments. In particle transfer, the final state of b is a bound

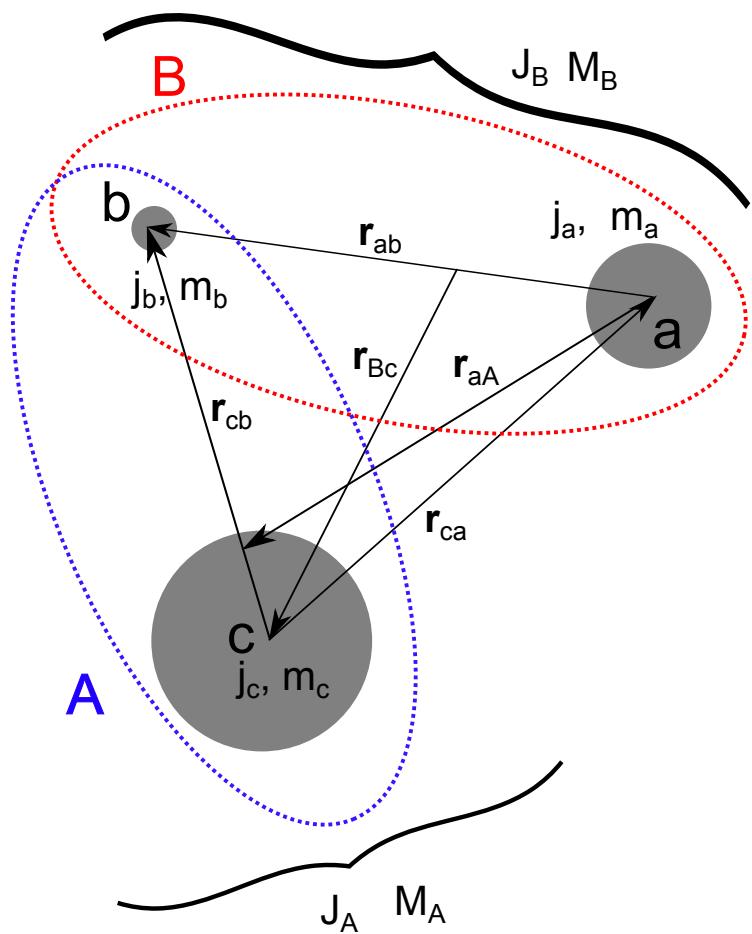


Figure 5.7.4: One particle transfer reaction $A(= c + b) + a \rightarrow B(= a + b) + c$.

state of the $B (= a + b)$ nucleus (cf. Fig. 5.7.4), and we can carry on in a similar way as done previously just by substituting the distorted wave (continuum) wave function (5.7.34) with

$$\psi_{m'_b}^{l'_b, j'_b *}(r_{ab}, \sigma_b) = u_{l'_b, j'_b}^*(r_{ab}) [Y_{-m'_b}^{l'_b}(\hat{\mathbf{r}}_{ab}) \phi^{1/2}(\sigma_b)]_{m'_b}^{j'_b *}, \quad (5.7.57)$$

so the transition amplitude is now

$$\begin{aligned} T_{m_a, m_b}^{m'_a, m'_b}(\mathbf{k}'_a) &= \frac{8\pi^{3/2}}{k_a k'_a} \sum_{\sigma_a, \sigma_b} \sum_{l_a, j_a} \sum_{l'_a, m'_{l'_a}, j'_a} e^{i(\sigma^{l_a} + \sigma'^{l'_a})} i^{l_a - l'_a} (-1)^{l_a + l'_a - j'_a - j'_b} \\ &\times \sqrt{2l_a + 1} \langle l'_a m_{l'_a} | 1/2 m'_a | j'_a m_{l'_a} + m'_a \rangle \langle l_a 0 1/2 m_a | j_a m_a \rangle \\ &\times Y_{-m'_{l'_a}}^{l'_a}(\hat{\mathbf{k}}'_a) \int d\mathbf{r}_{aA} d\mathbf{r}_{bc} [Y_{-m'_{l'_a}}^{l'_a}(\hat{\mathbf{r}}_{bc}) \phi^{1/2}(\sigma_a)]_{-m'_{l'_a} - m'_a}^{j'_a} [Y_{-m'_b}^{l'_b}(\hat{\mathbf{r}}_{ab}) \phi^{1/2}(\sigma_b)]_{-m'_b}^{j'_b} \\ &\times \frac{F_{l_a, j_a}(r_{aA}) F_{l'_a, j'_a}(r_{bc})}{r_{bc} r_{aA}} u_{l'_b, j'_b}^*(r_{ab}) u_{l_b, j_b}(r_{bc}) v(r_{ab}) v_\sigma(\sigma_a, \sigma_b) \\ &\times [Y^{l_a}(\hat{\mathbf{r}}_{aA}) \phi^{1/2}(\sigma_a)]_{m_a}^{j_a} [Y^{l_b}(\hat{\mathbf{r}}_{bc}) \phi^{1/2}(\sigma_b)]_{m_b}^{j_b}. \end{aligned} \quad (5.7.58)$$

Using (5.7.39), (5.7.40), (6.C.4), and setting $M = m_a - m'_a - m_{l'_a}$

$$\begin{aligned} T_{m_a, m_b}^{m'_a, m'_b}(\mathbf{k}'_a) &= \frac{8\pi^{3/2}}{k_a k'_a} T_\sigma \sum_{l_a, j_a} \sum_{l'_a, j'_a} \sum_{K, M} e^{i(\sigma^{l_a} + \sigma'^{l'_a})} i^{l_a - l'_a} (-1)^{l_a + l'_a - j'_a - j'_b} \\ &\times ((l'_a \frac{1}{2})_{j'_a} (l_a \frac{1}{2})_{j_a} | (l_a l'_a)_K (\frac{1}{2} \frac{1}{2})_0)_K ((l'_b \frac{1}{2})_{j'_b} (l_b \frac{1}{2})_{j_b} | (l_b l'_b)_K (\frac{1}{2} \frac{1}{2})_0)_K \\ &\times \sqrt{2l_a + 1} \langle l'_a m_a - m'_a - M | 1/2 m'_a | j'_a m_a - M \rangle \langle l_a 0 1/2 m_a | j_a m_a \rangle \\ &\times Y_{m_a - m'_a - M}^{l'_a}(\hat{\mathbf{k}}'_a) \int d\mathbf{r}_{aA} d\mathbf{r}_{bc} \frac{F_{l_a, j_a}(r_{aA}) F_{l'_a, j'_a}(r_{bc})}{r_{bc} r_{aA}} u_{l'_b, j'_b}^*(r_{ab}) u_{l_b, j_b}(r_{bc}) v(r_{ab}) \\ &\times [Y^{l_a}(\hat{\mathbf{r}}_{aA}) Y^{l'_a}(\hat{\mathbf{r}}_{bc})]_M^K [Y^{l_b}(\hat{\mathbf{r}}_{bc}) Y^{l'_b}(\hat{\mathbf{r}}_{ab})]_{-M}^K. \end{aligned} \quad (5.7.59)$$

Aside from (5.7.21), we also need

$$\mathbf{r}_{bc} = \frac{a + B}{B} \mathbf{r}_{aA} + \frac{b}{A} \mathbf{r}_{bc}. \quad (5.7.60)$$

From (5.7.20–5.7.25), one gets

$$\begin{aligned} T_{m_a, m_b}^{m'_a, m'_b}(\mathbf{k}'_a) &= \frac{32\pi^3}{k_a k'_a} T_\sigma \sum_{l_a, j_a} \sum_{l'_a, j'_a} \sum_{K, M} e^{i(\sigma^{l_a} + \sigma'^{l'_a})} i^{l_a - l'_a} (-1)^{l_a + l'_a - j'_a - j'_b} \\ &\times ((l'_a \frac{1}{2})_{j'_a} (l_a \frac{1}{2})_{j_a} | (l_a l'_a)_K (\frac{1}{2} \frac{1}{2})_0)_K ((l'_b \frac{1}{2})_{j'_b} (l_b \frac{1}{2})_{j_b} | (l_b l'_b)_K (\frac{1}{2} \frac{1}{2})_0)_K \\ &\times \frac{2l_a + 1}{2K + 1} \langle l'_a m_a - m'_a - M | 1/2 m'_a | j'_a m_a - M \rangle \\ &\times \langle l_a 0 1/2 m_a | j_a m_a \rangle Y_{m_a - m'_a - M}^{l'_a}(\hat{\mathbf{k}}'_a) \mathcal{I}(l_a, l'_a, j_a, j'_a, j'_b, K), \end{aligned} \quad (5.7.61)$$

with

$$\begin{aligned} I(l_a, l'_a, j_a, j'_a, K) = & \int dr_{aA} dr_{bc} d\theta r_{aA} r_{bc}^2 \frac{\sin \theta}{r_{Bc}} \\ & \times F_{l_a, j_a}(r_{aA}) F_{l'_a, j'_a}(r_{ac}) u_{l'_b, j'_b}^*(r_{ab}) u_{l_b, j_b}(r_{bc}) v(r_{ab}) \\ & \times \sum_{M_K} \langle l_a 0 | l'_a M_K | K M_K \rangle [Y^{l_b}(\cos \theta, 0) Y^{l'_b}(\cos \theta_{ab}, 0)]_{-M_K}^K Y_{M_K}^{l'_a}(\cos \theta_{Bc}, 0), \end{aligned} \quad (5.7.62)$$

where (see (5.7.21), (5.7.60) and Fig. 5.7.2)

$$\cos \theta_{ab} = \frac{-r_{aA} - \frac{c}{A} r_{bc} \cos \theta}{\sqrt{\left(\frac{c}{A} r_{bc} \sin \theta\right)^2 + \left(r_{aA} + \frac{c}{A} r_{bc} \cos \theta\right)^2}}, \quad (5.7.63)$$

$$\cos \theta_{Bc} = \frac{\frac{a+B}{B} r_{aA} + \frac{b}{A} r_{bc} \cos \theta}{\sqrt{\left(\frac{b}{A} r_{bc} \sin \theta\right)^2 + \left(\frac{a+B}{B} r_{aA} + \frac{b}{A} r_{bc} \cos \theta\right)^2}}, \quad (5.7.64)$$

and

$$r_{Bc} = \sqrt{\left(\frac{b}{A} r_{bc} \sin \theta\right)^2 + \left(\frac{a+B}{B} r_{aA} + \frac{b}{A} r_{bc} \cos \theta\right)^2}. \quad (5.7.65)$$

It is noted that Eq. (5.7.61) can also be used when particle b populates a resonant state of nucleus B which lies in the continuum.

5.8 Dynamical shell model in a nutshell

In the extreme shell model the nucleons move independently of each other, feeling the presence of the other nucleons when bouncing elastically off the nuclear surface of the average potential. The properly normalized probability for removing a nucleon from such orbitals is equal to 1 for ω equal to the (unperturbed) energy of the orbital in question and zero otherwise (see Fig. 5.8.1 (a)).

In the dynamical shell model⁶⁶, the nucleons can bounce inelastically off the nuclear surface setting the nucleus in a vibrational state, changing their state of motion as well as remaining in the same state. In this case the strength of the levels becomes, in general, distributed over a range of energies both below and above the Fermi energy. That is, the state k is found both in the system $(A - 1)$ populated in a pick-up process and in the system $(A + 1)$ populated through a stripping reaction, as indicated in Fig 5.8.1 (b).

The next question is whether this distribution is concentrated in discrete states or displays a continuous behavior. Both situations can in principle be found depending of whether the original single-particle state is close or far away from the

⁶⁶Mahaux et al. (1985) and references therein; in connection with Eq. (5.8.1) see Sect. 4.7 (in particular Eq. (4.7.20)) of this reference.

Fermi energy. The sum of the strength functions associated with all the states excited in the pick-up process of a nucleon with quantum numbers k gives the occupation number associated with the orbital, i.e. $\int_{-\infty}^{\epsilon_F} S_h(k; \omega) d\omega = n_k$, where S_h is the (hole) strength function (see Fig. 1.2.5). The full single-particle strength is found adding to this quantity that associated with the excitation of states in the $(A + 1)$ system where a particle with quantum numbers k is deposited in the target.

It is noted that the fact that in the dynamical shell model the single-particle strength is distributed not only over an energy range in the $(A - 1)$ system but also in the $(A + 1)$ system is intimately connected with the ground state correlations associated with the vibrational modes which produce particle-hole excitations in the ground state as well as pair addition and subtraction modes. In this way single-particle states which originally were filled become partially empty, and vice versa (see Fig. 1.8.3).

The transfer process shown in Fig. 5.8.1 (b) for both stripping and pick-up reactions are closely associated with the polarization and correlation contributions to the mass operator $\mathcal{M}(k; \omega) = \mathcal{V}(k; \omega) + i\mathcal{W}(k; \omega)$. The associated total strength distribution

$$\begin{aligned} S(k; \omega) &= S_h(k; \omega) + S_p(k; \omega), \\ &= \frac{1}{2\pi} \frac{\mathcal{W}(k; \omega) + \Delta}{(E_k + \mathcal{V}(k; \omega) - E)^2 + \frac{1}{4}(\mathcal{W}(k; \omega) + \Delta)^2}, \end{aligned} \quad (5.8.1)$$

can display a variety of shapes, as $\mathcal{V}(k; \omega)$ and $\mathcal{W}(k; \omega)$ are energy dependent.

The intermediate states of the polarization and correlation diagrams which, in the present discussion are the doorway states to the renormalization of the single-particle motion must be furthermore mixed with more complicated configurations. The extension to the complex plane of the associated expressions in terms of an imaginary parameter Δ is made to take these effects into account in some average way (see Sect. 2.3). In any case, most of the results are not dependent on the detailed value this quantity has. The picture shown in Fig. 5.8.1 (b), although accurate and controlled by only few parameters, is much too rich as compared to the originally extreme shell model picture.

Within this context, a major simplification can be achieved recognizing that the essential difference between the situation depicted in Figs 5.8.1 (a) and (b) is the value of the mean free path. In the first case it is infinite while in the second it is finite. That is, due to the coupling to the surface modes (and eventually pair modes) single-particle motion acquires a lifetime. In other terms, the coupling to doorway states leads to a breaking of the single-particle strength (see e.g. Fig. 5.2.3). The associated energy range (width Γ), over which this phenomenon takes place, provides insight into the time it takes for the different components to come out of phase ($\approx \hbar/\Gamma$).

Making the ansatz that the single-particle motion decays exponentially, a Breit-Wigner shape can be fitted to the different strength concentration resulting from the detailed calculation (Fig. 5.8.1 (b)) The centroid of the corresponding peaks can be

viewed as the energy of the dressed single-particle state. On the other hand, caution should be exercised in interpreting the area under the fitted curve as single-particle content (spectroscopic factor), if only because this quantity can become larger than one (Fig. 5.8.1 (c)). It is only for the levels not too far away from the Fermi energy that such interpretation can be reasonably accurate.

Another quantity which can be used to characterize the dressed single-particle states is the rate of change of the energy shift as a function of the energy, i.e. the effective mass. Again, for the orbitals close to the Fermi energy this quantity is inversely proportional to the area under the Breit-Wigner shapes. The energy range over which the effective mass deviates from one determine the region where the dynamical couplings change the density of single-particle levels (see Sect. 5.4.1, in particular Eq. (5.4.19)).

The trend of the effective masses shown in Fig. 5.8.1 (d) can be quantitatively understood as follows. For single-particle states close to the Fermi energy the intermediate states have all energies larger than the unperturbed single-particle energy, and the resulting shape is δ -like with a tail extending away from the Fermi energy. For single-particle orbitals 5–7 MeV away from ϵ_F there are a number of intermediate states which have the same energy of the initial state leading to vanishing energy denominators and thus to a marked structure in the strength function. Finally, for single-particle states far away from ϵ_F , the density of intermediate states is so large that the matrix elements of these couplings average out to a constant with a value much larger than typical distances between successive intermediate states. These are the conditions which lead to a Breit-Wigner shape.

It is then to be expected that the single-particle levels in the intermediate region will be associated with strength functions which deviate much from a smooth function and for which comparatively large errors can be made through a Breit-Wigner fitting. This is also seen from Fig. 5.8.1 (d) where the effective mass becomes smaller than one, indicating that the area of the fitted shapes are larger than that of the original strength function $S(k; \omega)$. This result emphasizes the difficulties found in trying to translate effective masses into spectroscopic factors in a general situation.

Starting from the extreme picture shown in Fig. 5.8.1 (a) where the relation $\omega = \hbar^2 k^2 / 2m$ holds, one arrives at the picture (c) where the relation $d\omega = \hbar^2 k \delta k / m^*$ again accounts for the main properties of the single-particle motion and where the effect of the couplings are contained in m^* (see Eq. (5.4.3)).

The question then arises of how to define an average quantity which depends only on the energy of the orbital and is state independent. Rather different prescriptions have been used to deal with the question. In any case the main result obtained was that the ω -dependent effective mass shows a well defined peak at the Fermi surface, its maximum value being considerably larger than one (≈ 1.4). The associated full width at half maximum is of the order of 10 MeV (i.e. ± 5 MeV around ϵ_F). This quantity is much smaller than the Fermi energy (~ 36 MeV) and is essentially determined by the energy of the low-lying collective modes and of the single-particle gap around closed shell system, e.g. ^{208}Pb .

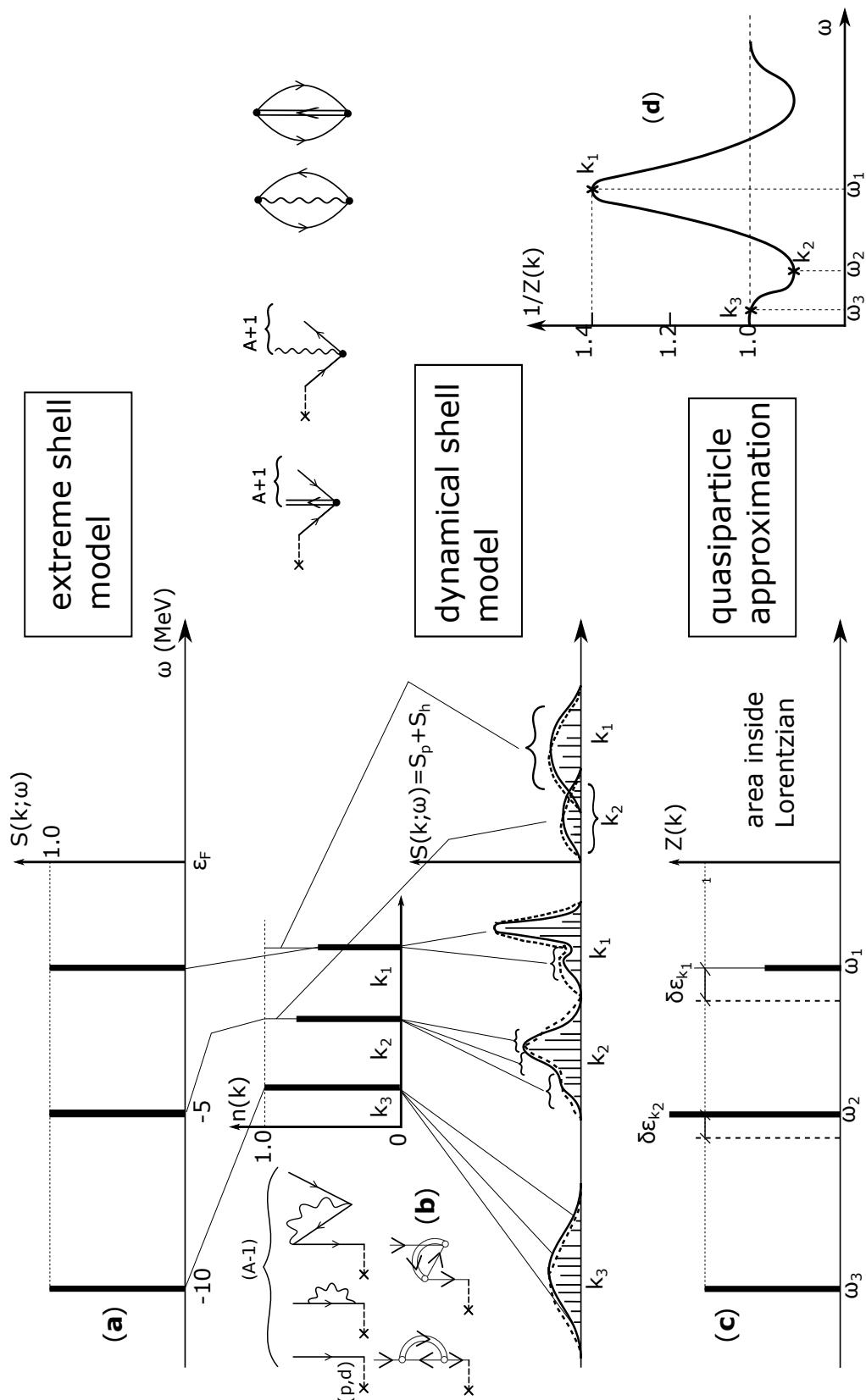


Figure 5.8.1

Caption Fig 5.8.1

Schematic representation of the main quantities characterizing the single-particle motion in the nuclear shell model taking into account the residual interaction among the particles at different levels of approximation. In **(a)** the bare NN -interaction is treated in the Hartree-Fock approximation and the particles feel the presence of the other particles through their own confinement in the average field. The strength function shows sharp peaks, each of them carrying the full strength of the states. In fact, the occupation number associated with each state k contains only one contribution. In **(b)** the particles still couple only with the average field. However, in this case, they can set the surface into vibrations by changing its state of motion (particle-vibration coupling mechanism). The strength associated with each orbital is distributed over a finite energy range. The corresponding occupation numbers arise from the sum of many contributions. Fitting a Breit-Wigner shape to each peak one can regain the simplicity of the extreme shell model by defining new levels with energy equal to that of the centroid and strength equal to that of the area covered by the Lorentzian shape (cf. **(c)**). In **(d)** the energy variation of the shift of the centroids is contained into an effective ω -mass according to the relation $\frac{m_\omega}{m} = \left(1 - \frac{\partial \Delta E}{\partial \hbar \omega}\right)$ (see Eq. (5.4.5)). The resulting curve resembles the shape obtained by calculating the inverse of the area below the different Lorentzians (quasiparticle approximation).

Chapter 6

Two-particle transfer

A unified picture of pairing and two-nucleon transfer, where the two subjects are blended together (which is what happens in nature)

B. F. Bayman

The present Chapter is structured in the following way. In Section 6.1 we present a summary of two-nucleon transfer reaction theory. It provides, together with Section 4.1 the elements needed to calculate the absolute two-nucleon transfer differential cross sections in second order DWBA, and thus to compare theory with experiment. For the practitioner in search of details and clarification we present in section 6.2 a derivation of the equations presented in section 6.1. These equations are implemented and made operative in the software COOPER used in the applications (cf. App. 7.A).

A number of Appendices are provided. Appendices 6.A–6.D contain relations used in Sect. 6.2 as well as in the derivation of some two-nucleon transfer spectroscopic amplitudes. Finally, Appendix 6.E provides a glimpse of original material due to Ben Bayman¹ which was instrumental to render quantitative the studies of two-nucleon transfer, studies which can now be carried out in terms of absolute differential cross sections and not only relative ones.

6.1 Summary of second order DWBA

Let us illustrate the theory of second order DWBA two-nucleon transfer reactions with the $A + t \rightarrow B(\equiv A + 2) + p$ reaction, in which $A + 2$ and A are even nuclei in their 0^+ ground state. The extension of the expressions to the transfer of pairs coupled to arbitrary angular momentum is discussed in subsection 6.2.10.

¹Bayman (1970), Bayman (1971), Bayman and Chen (1982).

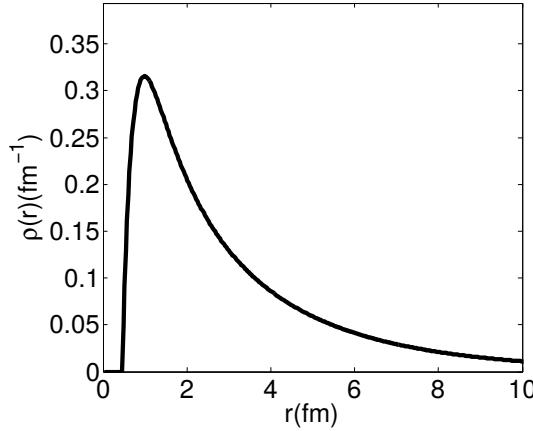


Figure 6.1.1: Radial function $\rho(r)$ (hard core 0.45 fm) entering the tritium wavefunction.

The wavefunction of the nucleus $A + 2$ can be written as

$$\Psi_{A+2}(\xi_A, \mathbf{r}_{A1}, \sigma_1, \mathbf{r}_{A2}, \sigma_2) = \psi_A(\xi_A) \sum_{l_i, j_i} [\phi_{l_i, j_i}^{A+2}(\mathbf{r}_{A1}, \sigma_1, \mathbf{r}_{A2}, \sigma_2)]_0^0, \quad (6.1.1)$$

where

$$[\phi_{l_i, j_i}^{A+2}(\mathbf{r}_{A1}, \sigma_1, \mathbf{r}_{A2}, \sigma_2)]_0^0 = \sum_{nm} a_{nm} [\varphi_{n, l_i, j_i}^{A+2}(\mathbf{r}_{A1}, \sigma_1) \varphi_{m, l_i, j_i}^{A+2}(\mathbf{r}_{A2}, \sigma_2)]_0^0, \quad (6.1.2)$$

while the wavefunctions $\varphi_{n, l_i, j_i}^{A+2}(\mathbf{r})$ are single-particle wavefunctions. The radial dependence of the wavefunction of the two neutrons in the triton is written as $\phi_t(\mathbf{r}_{p1}, \mathbf{r}_{p2}) = \rho(r_{p1})\rho(r_{p2})\rho(r_{12})$, where r_{p1}, r_{p2}, r_{12} are the distances between neutron 1 and the proton, neutron 2 and the proton and between neutrons 1 and 2 respectively, while $\rho(r)$ is the hard core ($r_{core} = 0.45$ fm) potential wavefunction depicted in² Fig 6.1.1.

The two-nucleon transfer differential cross section is written as

$$\frac{d\sigma}{d\Omega} = \frac{\mu_i \mu_f}{(4\pi\hbar^2)^2} \frac{k_f}{k_i} \left| T^{(1)}(\theta) + T_{succ}^{(2)}(\theta) - T_{NO}^{(2)}(\theta) \right|^2, \quad (6.1.3)$$

where³,

$$\begin{aligned} T^{(1)}(\theta) = 2 \sum_{l_i, j_i} \sum_{\sigma_1 \sigma_2} \int d\mathbf{r}_{tA} d\mathbf{r}_{p1} d\mathbf{r}_{A2} & [\phi_{l_i, j_i}^{A+2}(\mathbf{r}_{A1}, \sigma_1, \mathbf{r}_{A2}, \sigma_2)]_0^{0*} \chi_{pB}^{(-)*}(\mathbf{r}_{pB}) \\ & \times v(\mathbf{r}_{p1}) \phi_t(\mathbf{r}_{p1}, \sigma_1, \mathbf{r}_{p2}, \sigma_2) \chi_{tA}^{(+)}(\mathbf{r}_{tA}), \end{aligned} \quad (6.1.4a)$$

²Tang and Herndon (1965).

³See Bayman and Chen (1982) and App. 6.E.

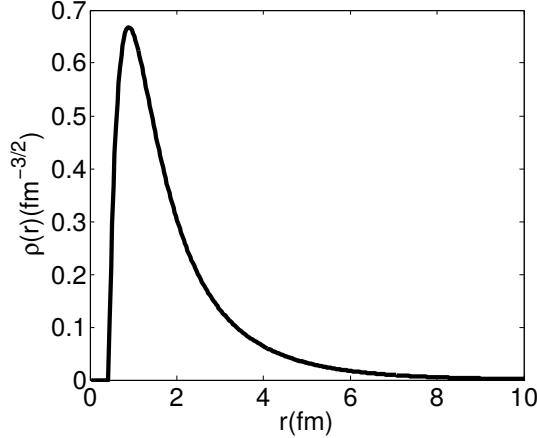


Figure 6.1.2: Radial wavefunction $\rho_d(r)$ (hard core 0.45 fm) entering the deuteron wavefunction.

$$\begin{aligned}
 T_{succ}^{(2)}(\theta) = & 2 \sum_{l_i, j_i} \sum_{l_f, j_f, m_f} \sum_{\sigma_1' \sigma_2'} \int d\mathbf{r}_{dF} d\mathbf{r}_{p1} d\mathbf{r}_{A2} [\phi_{l_i, j_i}^{A+2}(\mathbf{r}_{A1}, \sigma_1, \mathbf{r}_{A2}, \sigma_2)]_0^{0*} \chi_{pB}^{(-)*}(\mathbf{r}_{pB}) v(\mathbf{r}_{p1}) \\
 & \times \phi_d(\mathbf{r}_{p1}, \sigma_1) \varphi_{l_f, j_f, m_f}^{A+1}(\mathbf{r}_{A2}, \sigma_2) \int d\mathbf{r}'_{dF} d\mathbf{r}'_{p1} d\mathbf{r}'_{A2} G(\mathbf{r}_{dF}, \mathbf{r}'_{dF}) \\
 & \times \phi_d(\mathbf{r}'_{p1}, \sigma'_1)^* \varphi_{l_f, j_f, m_f}^{A+1*}(\mathbf{r}'_{A2}, \sigma'_2) \frac{2\mu_{dF}}{\hbar^2} v(\mathbf{r}'_{p2}) \phi_d(\mathbf{r}'_{p1}, \sigma'_1) \phi_d(\mathbf{r}'_{p2}, \sigma'_2) \chi_{tA}^{(+)}(\mathbf{r}'_{tA}),
 \end{aligned} \tag{6.1.4b}$$

$$\begin{aligned}
 T_{NO}^{(2)}(\theta) = & 2 \sum_{l_i, j_i} \sum_{l_f, j_f, m_f} \sum_{\sigma_1' \sigma_2'} \int d\mathbf{r}_{dF} d\mathbf{r}_{p1} d\mathbf{r}_{A2} [\phi_{l_i, j_i}^{A+2}(\mathbf{r}_{A1}, \sigma_1, \mathbf{r}_{A2}, \sigma_2)]_0^{0*} \chi_{pB}^{(-)*}(\mathbf{r}_{pB}) v(\mathbf{r}_{p1}) \\
 & \times \phi_d(\mathbf{r}_{p1}, \sigma_1) \varphi_{l_f, j_f, m_f}^{A+1}(\mathbf{r}_{A2}, \sigma_2) \int d\mathbf{r}'_{p1} d\mathbf{r}'_{A2} d\mathbf{r}'_{dF} \\
 & \times \phi_d(\mathbf{r}'_{p1}, \sigma'_1)^* \varphi_{l_f, j_f, m_f}^{A+1*}(\mathbf{r}'_{A2}, \sigma'_2) \phi_d(\mathbf{r}'_{p1}, \sigma'_1) \phi_d(\mathbf{r}'_{p2}, \sigma'_2) \chi_{tA}^{(+)}(\mathbf{r}'_{tA}).
 \end{aligned} \tag{6.1.4c}$$

The quantities $\mu_i, \mu_f (k_i, k_f)$ are the reduced masses (relative linear momenta) in both entrance (initial, i) and exit (final, f) channels, respectively. In the above expressions, $\varphi_{l_f, j_f, m_f}^{A+1}(\mathbf{r}_{A1})$ are the wavefunctions describing the intermediate states of the nucleus $F (\equiv (A + 1))$, generated as solutions of a Woods-Saxon potential, $\phi_d(\mathbf{r}_{p2})$ being the deuteron bound wavefunction (see Fig. 6.1.2)⁴. Note that some or all of the single-particle states described by the wavefunctions $\varphi_{l_f, j_f, m_f}^{A+1}(\mathbf{r}_{A1})$ may lie in the continuum (case in which the nucleus F is loosely bound or unbound). Although there are a number of ways to exactly treat such states, discretization processes may be sufficiently accurate. They can be implemented by,

⁴Tang and Herndon (1965).

for example, embedding the Woods–Saxon potential in a spherical box of sufficiently large radius. In actual calculations involving e.g. the halo nucleus ^{11}Li , and where $|F\rangle = |^{10}\text{Li}\rangle$, one achieved convergence making use of approximately 20 continuum states and a box of 30 fm of radius. Concerning the components of the triton wavefunction describing the relative motion of the dineutron, it was generated with the $p - n$ interaction⁵

$$v(r) = -v_0 \exp(-k(r - r_c)) \quad r > r_c \quad (6.1.5)$$

$$v(r) = \infty \quad r < r_c, \quad (6.1.6)$$

where $k = 2.5 \text{ fm}^{-1}$ and $r_c = 0.45 \text{ fm}$, the depth v_0 being adjusted to reproduce the experimental separation energies. The positive-energy wavefunctions $\chi_{IA}^{(+)}(\mathbf{r}_{IA})$ and $\chi_{pB}^{(-)}(\mathbf{r}_{pB})$ are the ingoing distorted wave in the initial channel and the outgoing distorted wave in the final channel respectively. They are continuum solutions of the Schrödinger equation associated with the corresponding optical potentials.

The transition potential responsible for the transfer of the pair is, in the *post* representation (cf. Fig. 6.5.1),

$$V_\beta = v_{pB} - U_\beta, \quad (6.1.7)$$

where v_{pB} is the interaction between the proton and nucleus B , and U_β is the optical potential in the final channel. We make the assumption that v_{pB} can be decomposed into a term containing the interaction between A and p and the potential describing the interaction between the proton and each of the transferred nucleons, namely

$$v_{pB} = v_{pA} + v_{p1} + v_{p2}, \quad (6.1.8)$$

where v_{p1} and v_{p2} is the hard–core potential (6.1.5). The transition potential is

$$V_\beta = v_{pA} + v_{p1} + v_{p2} - U_\beta. \quad (6.1.9)$$

Assuming that $\langle\beta|v_{pA}|\alpha\rangle \approx \langle\beta|U_\beta|\alpha\rangle$ (i.e, assuming that the matrix element of the core–core interaction between the initial and final states is very similar to the matrix element of the real part of the optical potential), one obtains the final expression of the transfer potential in the *post* representation, namely,

$$V_\beta \simeq v_{p1} + v_{p2} = v(\mathbf{r}_{p1}) + v(\mathbf{r}_{p2}). \quad (6.1.10)$$

We make the further approximation of using the same interaction potential in all the (i.e. initial, intermediate and final) channels.

The extension to a heavy-ion reaction $A + a (\equiv b + 2) \longrightarrow B (\equiv A + 2) + b$ imply no essential modifications in the formalism. The deuteron and triton wavefunctions appearing in Eqs. (6.1.4a), (6.1.4b) and (6.1.4c) are to be substituted with the corresponding wavefunctions $\Psi_{b+2}(\xi_b, \mathbf{r}_{b1}, \sigma_1, \mathbf{r}_{b2}, \sigma_2)$, constructed in a similar

⁵Tang and Herndon (1965).

way as those appearing in (6.1.1 and 6.1.2). The interaction potential used in Eqs. (6.1.4a), (6.1.4b) and (6.1.4c) will now be the Saxon–Woods used to define the initial (final) state in the post (prior) representation, instead of the proton–neutron interaction (6.1.5).

The Green's function $G(\mathbf{r}_{dF}, \mathbf{r}'_{dF})$ appearing in (6.1.4b) propagates the intermediate channel d, F . It can be expanded in partial waves as,

$$G(\mathbf{r}_{dF}, \mathbf{r}'_{dF}) = i \sum_l \sqrt{2l+1} \frac{f_l(k_{dF}, r_<)g_l(k_{dF}, r_>)}{k_{dF}r_{dF}r'_{dF}} \left[Y^l(\hat{r}_{dF})Y^l(\hat{r}'_{dF}) \right]_0^0. \quad (6.1.11)$$

The $f_l(k_{dF}, r)$ and $g_l(k_{dF}, r)$ are the regular and the irregular solutions of a Schrödinger equation for a suitable optical potential and an energy equal to the kinetic energy of the intermediate state. In most cases of interest, the result is hardly altered if we use the same energy of relative motion for all the intermediate states. This representative energy is calculated when both intermediate nuclei are in their corresponding ground states. It is of notice that the validity of this approximation can break down in some particular cases. If, for example, some relevant intermediate state become off shell, its contribution is significantly quenched. An interesting situation can arise when this happens to all possible intermediate states, so they can only be virtually populated.

6.2 Detailed derivation of second order DWBA

6.2.1 Simultaneous transfer: distorted waves

For a (t, p) reaction, the triton is represented by an incoming distorted wave. We make the assumption that the two neutrons are in an $S = L = 0$ state, and that the relative motion of the proton with respect to the dineutron is also $l = 0$. Consequently, the total spin of the triton is entirely due to the spin of the proton. We will explicitly treat it, as we will consider a spin–orbit term in the optical potential acting between the triton and the target⁶.

Following (6.B.1), we can write the triton distorted wave as

$$\psi_{m_t}^{(+)}(\mathbf{R}, \mathbf{k}_i, \sigma_p) = \sum_{l_t} \exp(i\sigma_{l_t}^t) g_{l_t j_t} Y_0^{l_t}(\hat{\mathbf{R}}) \frac{\sqrt{4\pi(2l_t + 1)}}{k_i R} \chi_{m_t}(\sigma_p), \quad (6.2.1)$$

where use was made of $Y_0^{l_t}(\hat{\mathbf{k}}_i) = i^{l_t} \sqrt{\frac{2l_t + 1}{4\pi}} \delta_{m_t, 0}$, in keeping with the fact that \mathbf{k}_i is oriented along the z -axis. Note the phase difference with eq. (7) of Bayman (1971), due to the use of time-reversal rather than Condon–Shortley phase convention. Making use of the relation

$$Y_0^{l_t}(\hat{\mathbf{R}}) \chi_{m_t}(\sigma_p) = \sum_{j_t} \langle l_t 0 1/2 m_t | j_t m_t \rangle \left[Y^l(\hat{\mathbf{R}}) \chi(\sigma_p) \right]_{m_t}^{j_t}, \quad (6.2.2)$$

⁶In what follows we will use the notation of Bayman (1971), see also App. 6.E.

we have

$$\begin{aligned} \psi_{m_t}^{(+)}(\mathbf{R}, \mathbf{k}_t, \sigma_p) &= \sum_{l_t, j_t} \exp(i\sigma_{l_t}^t) \frac{\sqrt{4\pi(2l_t + 1)}}{k_t R} g_{l_t j_t}(R) \\ &\quad \times \langle l_t 0 1/2 m_t | j_t m_t \rangle [Y^{l_t}(\hat{\mathbf{R}})\chi(\sigma_p)]_{m_t}^{j_t}. \end{aligned} \quad (6.2.3)$$

We now turn our attention to the outgoing proton distorted wave, which, following (6.B.3) can be written as

$$\psi_{m_p}^{(-)}(\zeta, \mathbf{k}_f, \sigma_p) = \sum_{l_p, j_p} \frac{4\pi}{k_f \zeta} i^{l_p} \exp(-i\sigma_{l_p}^p) f_{l_p j_p}^*(\zeta) \sum_m Y_m^{l_p}(\hat{\zeta}) Y_m^{l_p*}(\hat{\mathbf{k}}_f) \chi_{m_p}(\sigma_p). \quad (6.2.4)$$

Making use of the relation

$$\begin{aligned} \sum_m Y_m^{l_p}(\hat{\zeta}) Y_m^{l_p*}(\hat{\mathbf{k}}_f) \chi_{m_p}(\sigma_p) &= \sum_{m, j_p} Y_m^{l_p*}(\hat{\mathbf{k}}_f) \langle l_p m 1/2 m_p | j_p m + m_p \rangle \\ &\quad \times [Y^{l_p}(\hat{\zeta}) \chi_{m_p}(\sigma_p)]_{m+m_p}^{j_p} \\ &= \sum_{m, j_p} Y_{m-m_p}^{l_p*}(\hat{\mathbf{k}}_f) \langle l_p m - m_p 1/2 m_p | j_p m \rangle [Y^{l_p}(\hat{\zeta}) \chi_{m_p}(\sigma_p)]_m^{j_p}, \end{aligned} \quad (6.2.5)$$

one obtains

$$\begin{aligned} \psi_{m_p}^{(-)}(\zeta, \mathbf{k}_f, \sigma_p) &= \frac{4\pi}{k_f \zeta} \sum_{l_p, j_p, m} i^{l_p} \exp(-i\sigma_{l_p}^p) f_{l_p j_p}^*(\zeta) Y_{m-m_p}^{l_p*}(\hat{\mathbf{k}}_f) \\ &\quad \times \langle l_p m - m_p 1/2 m_p | j_p m \rangle [Y^{l_p}(\hat{\zeta}) \chi(\sigma_p)]_m^{j_p}. \end{aligned} \quad (6.2.6)$$

6.2.2 matrix element for the transition amplitude

We now turn our attention to the evaluation of

$$\begin{aligned} \langle \Psi_f^{(-)}(\mathbf{k}_f) | V(r_{1p}) | \Psi_i^{(+)}(k_i, \hat{\mathbf{z}}) \rangle &= \frac{(4\pi)^{3/2}}{k_i k_f} \sum_{l_p, l_t, j_p, j_t, m} ((\lambda \frac{1}{2})_k (\lambda \frac{1}{2})_k | (\lambda \lambda)_0 (\frac{1}{2} \frac{1}{2})_0)_0 \sqrt{2l_t + 1} \\ &\quad \times \langle l_p m - m_p 1/2 m_p | j_p m \rangle \langle l_t 0 1/2 m_t | j_t m_t \rangle i^{-l_p} \exp[i(\sigma_{l_p}^p + \sigma_{l_t}^t)] \\ &\quad \times 2Y_{m-m_p}^{l_p}(\hat{\mathbf{k}}_f) \sum_{\sigma_1 \sigma_2 \sigma_p} \int \frac{d\zeta d\mathbf{r} d\boldsymbol{\eta}}{\zeta R} u_{\lambda k}(r_1) u_{\lambda k}(r_2) [Y^\lambda(\hat{\mathbf{r}}_1) Y^\lambda(\hat{\mathbf{r}}_2)]_0^{0*} \\ &\quad \times f_{l_p j_p}(\zeta) g_{l_t j_t}(R) [\chi(\sigma_1) \chi(\sigma_2)]_0^{0*} [Y^{l_p}(\hat{\zeta}) \chi(\sigma_p)]_m^{j_p*} V(r_{1p}) \\ &\quad \times \theta_0^0(\mathbf{r}, \mathbf{s}) [\chi(\sigma_1) \chi(\sigma_2)]_0^0 [Y^{l_t}(\hat{\mathbf{R}}) \chi(\sigma_p)]_{m_t}^{j_t}, \end{aligned} \quad (6.2.7)$$

where

$$\begin{aligned}\mathbf{r} &= \mathbf{r}_2 - \mathbf{r}_1, \\ \mathbf{s} &= \frac{1}{2}(\mathbf{r}_1 + \mathbf{r}_2) - \mathbf{r}_p, \\ \boldsymbol{\eta} &= \frac{1}{2}(\mathbf{r}_1 + \mathbf{r}_2), \\ \zeta &= \mathbf{r}_p - \frac{\mathbf{r}_1 + \mathbf{r}_2}{A+2}.\end{aligned}\tag{6.2.8}$$

The sum over σ_1, σ_2 in (6.2.7) is found to be equal to 1. We will now simplify the term $[Y^{l_p}(\hat{\zeta})\chi(\sigma_p)]_m^{j_p*} [Y^{l_t}(\hat{\mathbf{R}})\chi(\sigma_p)]_{m_t}^{j_t}$, noting that, (6.A.13)

$$[Y^{l_p}(\hat{\zeta})\chi(\sigma_p)]_m^{j_p*} = (-1)^{1/2-\sigma_p+j_p-m} [Y^{l_p}(\hat{\zeta})\chi(-\sigma_p)]_{-m}^{j_p}. \tag{6.2.9}$$

and that

$$\begin{aligned}[Y^{l_p}(\hat{\zeta})\chi(-\sigma_p)]_{-m}^{j_p} [Y^{l_t}(\hat{\mathbf{R}})\chi(\sigma_p)]_{m_t}^{j_t} &= \sum_{JM} \langle j_p \ -m \ j_t \ m_t | J \ M \rangle \\ &\times \left\{ [Y^{l_p}(\hat{\zeta})\chi(-\sigma_p)]_{-m}^{j_p} [Y^{l_t}(\hat{\mathbf{R}})\chi(\sigma_p)]_{m_t}^{j_t} \right\}_M^J\end{aligned}\tag{6.2.10}$$

The only term which does not vanish after the integration is performed is the one in which the angular and spin functions are coupled to $L = 0, S = 0, J = 0$. Thus,

$$\begin{aligned}\langle j_p \ -m \ j_t \ m_t | 0 \ 0 \rangle \left\{ [Y^{l_p}(\hat{\zeta})\chi(-\sigma_p)]_{-m}^{j_p} [Y^{l_t}(\hat{\mathbf{R}})\chi(\sigma_p)]_{m_t}^{j_t} \right\}_0^0 \delta_{l_p l_t} \delta_{j_p j_t} \delta_{m m_t} \\ = \frac{(-1)^{j_p+m_t}}{\sqrt{2j_p+1}} \left\{ [Y^{l_p}(\hat{\zeta})\chi(-\sigma_p)]_{-m}^{j_p} [Y^{l_t}(\hat{\mathbf{R}})\chi(\sigma_p)]_{m_t}^{j_t} \right\}_0^0 \delta_{l_p l_t} \delta_{j_p j_t} \delta_{m m_t}.\end{aligned}\tag{6.2.11}$$

Coupling separately the spin and angular functions, one obtains

$$\begin{aligned}\left\{ [Y^{l_p}(\hat{\zeta})\chi(-\sigma_p)]_{-m}^{j_p} [Y^{l_t}(\hat{\mathbf{R}})\chi(\sigma_p)]_{m_t}^{j_t} \right\}_0^0 \\ = ((l\frac{1}{2})_j (l\frac{1}{2})_j | (ll)_0 (\frac{1}{2}\frac{1}{2})_0)_0 \left[\chi(-\sigma_p)\chi(\sigma_p) \right]_0^0 \left[Y^{l_p}(\hat{\zeta})Y^{l_t}(\hat{\mathbf{R}}) \right]_0^0.\end{aligned}\tag{6.2.12}$$

We substitute (6.2.9), (6.2.30),(6.2.31) in (6.2.7) to obtain

$$\begin{aligned}
 \langle \Psi_f^{(-)}(\mathbf{k}_f) | V(r_{1p}) | \Psi_i^{(+)}(k_i, \hat{\mathbf{z}}) \rangle &= -\frac{(4\pi)^{3/2}}{k_i k_f} \sum_{lj} ((\lambda \frac{1}{2})_k (\lambda \frac{1}{2})_l | (\lambda \lambda)_0 (\frac{1}{2} \frac{1}{2})_0)_0 \sqrt{\frac{2l+1}{2j+1}} \\
 &\times \langle l m_t - m_p 1/2 m_p | j m_t \rangle \langle l 0 1/2 m_t | j m_t \rangle i^{-l} \exp[i(\sigma_l^p + \sigma_l^t)] \\
 &\times 2Y_{m_t - m_p}^l(\hat{\mathbf{k}}_f) \int \frac{d\zeta d\mathbf{r} d\mathbf{n}}{\zeta R} u_{\lambda k}(r_1) u_{\lambda l}(r_2) [Y^\lambda(\hat{\mathbf{r}}_1) Y^\lambda(\hat{\mathbf{r}}_2)]_0^{0*} \\
 &\times f_{lj}(\zeta) g_{lj}(R) [Y^l(\hat{\zeta}) Y^l(\hat{\mathbf{R}})]_0^0 V(r_{1p}) \theta_0^0(\mathbf{r}, \mathbf{s}) \\
 &\times ((l \frac{1}{2})_j (l \frac{1}{2})_j | (ll)_0 (\frac{1}{2} \frac{1}{2})_0)_0 \sum_{\sigma_p} (-1)^{1/2 - \sigma_p} [\chi(-\sigma_p) \chi(\sigma_p)]_0^0.
 \end{aligned} \tag{6.2.13}$$

The last sum over σ_p leads to

$$\begin{aligned}
 \sum_{\sigma_p} (-1)^{1/2 - \sigma_p} [\chi(-\sigma_p) \chi(\sigma_p)]_0^0 &= \sum_{\sigma_p m} (-1)^{1/2 - \sigma_p} \langle 1/2 m 1/2 - m | 0 0 \rangle \\
 &\times \chi_m(-\sigma_p) \chi_{-m}(\sigma_p) \\
 &= \frac{1}{\sqrt{2}} \sum_{\sigma_p m} (-1)^{1/2 - \sigma_p} (-1)^{1/2 - m} \delta_{m, -\sigma_p} \delta_{-m, \sigma_p} = -\sqrt{2}.
 \end{aligned} \tag{6.2.14}$$

The 9j-symbols can be evaluated to find

$$\begin{aligned}
 ((\lambda \frac{1}{2})_k (\lambda \frac{1}{2})_l | (\lambda \lambda)_0 (\frac{1}{2} \frac{1}{2})_0)_0 &= \sqrt{\frac{2k+1}{2(2\lambda+1)}} \\
 ((l \frac{1}{2})_j (l \frac{1}{2})_j | (ll)_0 (\frac{1}{2} \frac{1}{2})_0)_0 &= \sqrt{\frac{2j+1}{2(2l+1)}},
 \end{aligned} \tag{6.2.15}$$

and consequently,

$$\begin{aligned}
 \langle \Psi_f^{(-)}(\mathbf{k}_f) | V(r_{1p}) | \Psi_i^{(+)}(k_i, \hat{\mathbf{z}}) \rangle &= \frac{(4\pi)^{3/2}}{k_i k_f} \sum_{lj} \sqrt{\frac{2k+1}{2\lambda+1}} \\
 &\times \langle l m_t - m_p 1/2 m_p | j m_t \rangle \langle l 0 1/2 m_t | j m_t \rangle i^{-l} \exp[i(\sigma_l^p + \sigma_l^t)] \\
 &\times \sqrt{2} Y_{m_t - m_p}^l(\hat{\mathbf{k}}_f) \int \frac{d\zeta d\mathbf{r} d\mathbf{n}}{\zeta R} u_{\lambda k}(r_1) u_{\lambda l}(r_2) [Y^\lambda(\hat{\mathbf{r}}_1) Y^\lambda(\hat{\mathbf{r}}_2)]_0^{0*} \\
 &\times f_{lj}(\zeta) g_{lj}(R) [Y^l(\hat{\zeta}) Y^l(\hat{\mathbf{R}})]_0^0 V(r_{1p}) \theta_0^0(\mathbf{r}, \mathbf{s}).
 \end{aligned} \tag{6.2.16}$$

The values of the Clebsh–Gordan coefficients are, for $j = l - 1/2$,

$$\begin{aligned}
 \langle l m_t - m_p 1/2 m_p | l - 1/2 m_t \rangle \langle l 0 1/2 m_t | l - 1/2 m_t \rangle \\
 = \begin{cases} \frac{l}{2l+1} & \text{if } m_t = m_p \\ -\frac{\sqrt{l(l+1)}}{2l+1} & \text{if } m_t = -m_p \end{cases}
 \end{aligned} \tag{6.2.17}$$

and, for $j = l + 1/2$:

$$\begin{aligned} & \langle l m_t - m_p 1/2 m_p | l + 1/2 m_t \rangle \langle l 0 1/2 m_t | l + 1/2 m_t \rangle \\ &= \begin{cases} \frac{l+1}{2l+1} & \text{if } m_t = m_p \\ \frac{\sqrt{l(l+1)}}{2l+1} & \text{if } m_t = -m_p \end{cases} \quad (6.2.18) \end{aligned}$$

One thus can write,

$$\begin{aligned} \langle \Psi_f^{(-)}(\mathbf{k}_f) | V(r_{1p}) | \Psi_i^{(+)}(k_i, \hat{\mathbf{z}}) \rangle &= \frac{(4\pi)^{3/2}}{k_i k_f} \sum_l \frac{1}{(2l+1)} \sqrt{\frac{(2k+1)}{(2\lambda+1)}} \exp[i(\sigma_l^p + \sigma_l^t)] t^{-l} \\ &\times \sqrt{2} Y_{m_t - m_p}^l(\hat{\mathbf{k}}_f) \int \frac{d\zeta d\mathbf{r} d\boldsymbol{\eta}}{\zeta R} u_{\lambda k}(r_1) u_{\lambda k}(r_2) [Y^\lambda(\hat{\mathbf{r}}_1) Y^\lambda(\hat{\mathbf{r}}_2)]_0^{0*} \\ &\times V(r_{1p}) \theta_0^0(\mathbf{r}, \mathbf{s}) [Y^l(\hat{\zeta}) Y^l(\hat{\mathbf{R}})]_0^0 \\ &\times \left[\left(f_{ll+1/2}(\zeta) g_{ll+1/2}(R)(l+1) + f_{ll-1/2}(\zeta) g_{ll-1/2}(R)l \right) \delta_{m_p, m_t} \right. \\ &\quad \left. + \left(f_{ll+1/2}(\zeta) g_{ll+1/2}(R) \sqrt{l(l+1)} - f_{ll-1/2}(\zeta) g_{ll-1/2}(R) \sqrt{l(l+1)} \right) \delta_{m_p, -m_t} \right]. \quad (6.2.19) \end{aligned}$$

We can further simplify this expression using

$$\begin{aligned} [Y^\lambda(\hat{\mathbf{r}}_1) Y^\lambda(\hat{\mathbf{r}}_2)]_0^{0*} &= [Y^\lambda(\hat{\mathbf{r}}_1) Y^\lambda(\hat{\mathbf{r}}_2)]_0^0 = \sum_m \langle \lambda m \lambda - m | 0 0 \rangle Y_m^\lambda(\hat{\mathbf{r}}_1) Y_{-m}^\lambda(\hat{\mathbf{r}}_2) \\ &= \sum_m (-1)^{\lambda-m} \langle \lambda m \lambda - m | 0 0 \rangle Y_m^\lambda(\hat{\mathbf{r}}_1) Y_m^{\lambda*}(\hat{\mathbf{r}}_2) \\ &= \frac{1}{\sqrt{2\lambda+1}} \sum_m Y_m^\lambda(\hat{\mathbf{r}}_1) Y_m^{\lambda*}(\hat{\mathbf{r}}_2) \\ &= \frac{\sqrt{(2\lambda+1)}}{4\pi} P_\lambda(\cos \theta_{12}). \quad (6.2.20) \end{aligned}$$

Note that when using Condon–Shortley phases this last expression is to be multiplied by $(-1)^\lambda$, and that

$$\begin{aligned} [Y^l(\hat{\zeta}) Y^l(\hat{\mathbf{R}})]_0^0 &= \sum_m \langle l m l - m | 0 0 \rangle Y_m^l(\hat{\zeta}) Y_{-m}^l(\hat{\mathbf{R}}) \\ &= \frac{1}{\sqrt{(2l+1)}} \sum_m (-1)^{l+m} Y_m^l(\hat{\zeta}) Y_{-m}^l(\hat{\mathbf{R}}). \quad (6.2.21) \end{aligned}$$

Because the integral of the above expression is independent of m , one can eliminate the m -sum and multiply by $2l+1$ the $m=0$ term, leading to

$$\begin{aligned} [Y^l(\hat{\zeta}) Y^l(\hat{\mathbf{R}})]_0^0 &\Rightarrow (-1)^l \sqrt{(2l+1)} Y_0^l(\hat{\zeta}) Y^l(\hat{\mathbf{R}}) \\ &= \sqrt{(2l+1)} Y_0^l(\hat{\zeta}) Y_0^{l*}(\hat{\mathbf{R}}). \quad (6.2.22) \end{aligned}$$

We now change the integration variables from $(\zeta, \mathbf{r}, \boldsymbol{\eta})$ to $(\mathbf{R}, \alpha, \beta, \gamma, r_{12}, r_{1p}, r_{2p})$, the quantity

$$\left| \frac{\partial(\mathbf{r}, \boldsymbol{\eta}, \zeta)}{\partial(\mathbf{R}, \alpha, \beta, \gamma, r_{12}, r_{1p}, r_{2p})} \right| = r_{12} r_{1p} r_{2p} \sin \beta, \quad (6.2.23)$$

being the Jacobian of the transformation. Finally,

$$\begin{aligned} \langle \Psi_f^{(-)}(\mathbf{k}_f) | V(r_{1p}) | \Psi_i^{(+)}(k_i, \hat{\mathbf{z}}) \rangle &= \frac{\sqrt{8\pi}}{k_i k_f} \sum_l \sqrt{\frac{2k+1}{2l+1}} \exp[i(\sigma_l^p + \sigma_l^t)] t^{-l} \\ &\times Y_{m_t - m_p}^l(\hat{\mathbf{k}}_f) \int d\mathbf{R} Y_0^{l*}(\hat{\mathbf{R}}) \int \frac{d\alpha d\beta d\gamma dr_{12} dr_{1p} dr_{2p} \sin \beta}{\zeta R} Y_0^l(\hat{\zeta}) \\ &\times u_{\lambda k}(r_1) u_{\lambda k}(r_2) V(r_{1p}) \theta_0^0(\mathbf{r}, \mathbf{s}) P_\lambda(\cos \theta_{12}) r_{12} r_{1p} r_{2p} \\ &\times \left[(f_{ll+1/2}(\zeta) g_{ll+1/2}(R)(l+1) + f_{ll-1/2}(\zeta) g_{ll-1/2}(R)l) \delta_{m_p, m_t} \right. \\ &\quad \left. + (f_{ll+1/2}(\zeta) g_{ll+1/2}(R) \sqrt{l(l+1)} - f_{ll-1/2}(\zeta) g_{ll-1/2}(R) \sqrt{l(l+1)}) \delta_{m_p, -m_t} \right]. \end{aligned} \quad (6.2.24)$$

It is noted that the second integral is a function of solely \mathbf{R} transforming under rotations as $Y_0^l(\hat{\mathbf{R}})$, in keeping with the fact that the full dependence on the orientation of \mathbf{R} is contained in the spherical harmonic $Y_0^l(\hat{\zeta})$. The second integral can thus be cast into the form

$$\begin{aligned} A(R) Y_0^l(\hat{\mathbf{R}}) &= \int d\alpha d\beta d\gamma dr_{12} dr_{1p} dr_{2p} \sin \beta \\ &\times F(\alpha, \beta, \gamma, r_{12}, r_{1p}, r_{2p}, R_x, R_y, R_z). \end{aligned} \quad (6.2.25)$$

To evaluate $A(R)$, we set \mathbf{R} along the z -axis

$$\begin{aligned} A(R) &= 2\pi i^{-l} \sqrt{\frac{4\pi}{2l+1}} \int d\beta d\gamma dr_{12} dr_{1p} dr_{2p} \sin \beta \\ &\times F(\alpha, \beta, \gamma, r_{12}, r_{1p}, r_{2p}, 0, 0, R), \end{aligned} \quad (6.2.26)$$

where a factor 2π results from the integration over α , the integrand not depending on α . Substituting (6.2.25) and (6.2.26) in (6.2.24) and, after integration over the angular variables of \mathbf{R} , we obtain

$$\begin{aligned} \langle \Psi_f^{(-)}(\mathbf{k}_f) | V(r_{1p}) | \Psi_i^{(+)}(k_i, \hat{\mathbf{z}}) \rangle &= 2 \frac{(2\pi)^{3/2}}{k_i k_f} \sum_l \sqrt{\frac{2k+1}{2l+1}} \exp[i(\sigma_l^p + \sigma_l^t)] t^{-l} \\ &\times Y_{m_t - m_p}^l(\hat{\mathbf{k}}_f) \int dR d\beta d\gamma dr_{12} dr_{1p} dr_{2p} R \sin \beta r_{12} r_{1p} r_{2p} \\ &\times u_{\lambda k}(r_1) u_{\lambda k}(r_2) V(r_{1p}) \theta_0^0(\mathbf{r}, \mathbf{s}) P_\lambda(\cos \theta_{12}) P_l(\cos \theta_\zeta) \\ &\times \left[(f_{ll+1/2}(\zeta) g_{ll+1/2}(R)(l+1) + f_{ll-1/2}(\zeta) g_{ll-1/2}(R)l) \delta_{m_p, m_t} \right. \\ &\quad \left. + (f_{ll+1/2}(\zeta) g_{ll+1/2}(R) \sqrt{l(l+1)} - f_{ll-1/2}(\zeta) g_{ll-1/2}(R) \sqrt{l(l+1)}) \delta_{m_p, -m_t} \right] / \zeta, \end{aligned} \quad (6.2.27)$$

where use was made of the relation

$$Y_0^l(\hat{\zeta}) = i^l \sqrt{\frac{2l+1}{4\pi}} P_l(\cos \theta_\zeta). \quad (6.2.28)$$

The final expression of the differential cross section involves a sum over the spin orientations:

$$\frac{d\sigma}{d\Omega}(\hat{\mathbf{k}}_f) = \frac{k_f}{k_i} \frac{\mu_i \mu_f}{(2\pi\hbar^2)^2} \frac{1}{2} \sum_{m_t m_p} |\langle \Psi_f^{(-)}(\mathbf{k}_f) | V(r_{1p}) | \Psi_i^{(+)}(k_i, \hat{\mathbf{z}}) \rangle|^2. \quad (6.2.29)$$

When $m_p = 1/2, m_t = 1/2$ or $m_p = -1/2, m_t = -1/2$, the terms proportional to δ_{m_p, m_t} including the factor

$$|Y_{m_t - m_p}^l(\hat{\mathbf{k}}_f) \delta_{m_p, m_t}| = |Y_0^l(\hat{\mathbf{k}}_f)| = \left| i^l \sqrt{\frac{2l+1}{4\pi}} P_l^0(\cos \theta) \right|, \quad (6.2.30)$$

in the case in which $m_p = -1/2, m_t = 1/2$

$$|Y_{m_t - m_p}^l(\hat{\mathbf{k}}_f) \delta_{m_p, -m_t}| = |Y_1^l(\hat{\mathbf{k}}_f)| = \left| i^l \sqrt{\frac{2l+1}{4\pi}} \frac{1}{l(l+1)} P_l^1(\cos \theta) \right|, \quad (6.2.31)$$

and

$$|Y_{m_t - m_p}^l(\hat{\mathbf{k}}_f) \delta_{m_p, -m_t}| = |Y_{-1}^l(\hat{\mathbf{k}}_f)| = |Y_1^l(\hat{\mathbf{k}}_f)| = \left| i^l \sqrt{\frac{2l+1}{4\pi}} \frac{1}{l(l+1)} P_l^1(\cos \theta) \right|, \quad (6.2.32)$$

when $m_p = 1/2, m_t = -1/2$. Taking the squared modulus of (6.2.27), the sum over m_t and m_p yields a factor 2 multiplying each one of the 2 different terms of the sum ($m_t = m_p$ and $m_t = -m_p$). This is equivalent to multiply each amplitude by $\sqrt{2}$, so the final constant that multiply the amplitudes is

$$\frac{8\pi^{3/2}}{k_i k_f}. \quad (6.2.33)$$

Now, for the triton wavefunction we use

$$\theta_0^0(\mathbf{r}, \mathbf{s}) = \rho(r_{1p}) \rho(r_{2p}) \rho(r_{12}), \quad (6.2.34)$$

$\rho(r)$ being a Tang–Herndon (1965) wave function also used by Bayman (1971). We obtain

$$\frac{d\sigma}{d\Omega}(\hat{\mathbf{k}}_f) = \frac{1}{2E_i^{3/2} E_f^{1/2}} \sqrt{\frac{\mu_f}{\mu_i}} (|I_{\lambda k}^{(0)}(\theta)|^2 + |I_{\lambda k}^{(1)}(\theta)|^2), \quad (6.2.35)$$

where

$$\begin{aligned} I_{\lambda k}^{(0)}(\theta) &= \sum_l P_l^0(\cos \theta) \sqrt{2k+1} \exp[i(\sigma_l^p + \sigma_l^t)] \\ &\times \int dR d\beta d\gamma dr_{12} dr_{1p} dr_{2p} R \sin \beta \rho(r_{1p}) \rho(r_{2p}) \rho(r_{12}) \\ &\times u_{\lambda k}(r_1) u_{\lambda k}(r_2) V(r_{1p}) P_\lambda(\cos \theta_{12}) P_l(\cos \theta_\zeta) r_{12} r_{1p} r_{2p} \\ &\times (f_{ll+1/2}(\zeta) g_{ll+1/2}(R) (l+1) + f_{ll-1/2}(\zeta) g_{ll-1/2}(R) l) / \zeta, \end{aligned} \quad (6.2.36)$$

and

$$\begin{aligned} I_{\lambda k}^{(1)}(\theta) &= \sum_l P_l^1(\cos \theta) \sqrt{2k+1} \exp[i(\sigma_l^p + \sigma_l^t)] \\ &\times \int dR d\beta d\gamma dr_{12} dr_{1p} dr_{2p} R \sin \beta \rho(r_{1p}) \rho(r_{2p}) \rho(r_{12}) \\ &\times u_{\lambda k}(r_1) u_{\lambda k}(r_2) V(r_{1p}) P_\lambda(\cos \theta_{12}) P_l(\cos \theta_\zeta) r_{12} r_{1p} r_{2p} \\ &\times (f_{ll+1/2}(\zeta) g_{ll+1/2}(R) - f_{ll-1/2}(\zeta) g_{ll-1/2}(R)) / \zeta. \end{aligned} \quad (6.2.37)$$

Note that the absence of the $(-1)^l$ factor with respect to what is found in the work of Bayman⁷, is due to the use of time-reversed phases instead of Condon–Shortley phasing. This is compensated in the total result by a similar difference in the expression of the spectroscopic amplitudes. This ensures that, in either case, the contribution of all the single particle transitions tend to have the same phase for superfluid nuclei, adding coherently to enhance the transfer cross section.

Heavy-ion Reactions

In dealing with a heavy ion reaction, $\theta_0^0(\mathbf{r}, \mathbf{s})$ are be the spatial part of the wavefunction

$$\begin{aligned} \Psi(\mathbf{r}_{b1}, \mathbf{r}_{b2}, \sigma_1, \sigma_2) &= [\psi^{j_i}(\mathbf{r}_{b1}, \sigma_1) \psi^{j_i}(\mathbf{r}_{b2}, \sigma_2)]_0^0 \\ &= \theta_0^0(\mathbf{r}, \mathbf{s}) [\chi(\sigma_1) \chi(\sigma_2)]_0^0, \end{aligned} \quad (6.2.38)$$

where $\mathbf{r}_{b1}, \mathbf{r}_{b2}$ are the positions of the two neutrons with respect to the b core. It can be shown to be

$$\theta_0^0(\mathbf{r}, \mathbf{s}) = \frac{u_{l_i j_i}(r_{b1}) u_{l_i j_i}(r_{b2})}{4\pi} \sqrt{\frac{2j_i + 1}{2}} P_{l_i}(\cos \theta_i), \quad (6.2.39)$$

where θ_i is the angle between \mathbf{r}_{b1} and \mathbf{r}_{b2} . Neglecting the spin-orbit term in the optical potential, as is usually done for heavy ion reactions, one obtains

$$\frac{d\sigma}{d\Omega}(\hat{\mathbf{k}}_f) = \frac{\mu_f \mu_i}{16\pi^2 \hbar^4 k_i^3 k_f^3} |T^{(1)}(\theta)|^2, \quad (6.2.40)$$

⁷Bayman (1971).

where

$$\begin{aligned}
T^{(1)}(\theta) = & \sum_l (2l+1) P_l(\cos \theta) \sqrt{(2j_i+1)(2j_f+1)} \exp[i(\sigma_l^p + \sigma_l^t)] \\
& \times \int dR d\beta d\gamma dr_{12} dr_{b1} dr_{b2} R \sin \beta u_{l_i j_i}(r_{b1}) u_{l_i j_i}(r_{b2}) \\
& \times u_{l_f j_f}(r_{A1}) u_{l_f j_f}(r_{A2}) V(r_{b1}) P_\lambda(\cos \theta_{12}) P_l(\cos \theta_\zeta) \\
& \times r_{12} r_{b1} r_{b2} P_{l_i}(\cos \theta_i) \frac{f_l(\zeta) g_l(R)}{\zeta},
\end{aligned} \tag{6.2.41}$$

obtained by using Eq. (6.2.39) in Eq. (6.2.7) instead of (6.2.34), $\mathbf{r}_{A1}, \mathbf{r}_{A2}$ being the coordinates of the two transferred neutrons with respect to the A core.

For control, in what follows we work out the same transition amplitude but starting from the distorted waves for a reaction taking place between spinless nuclei, namely

$$\psi^{(+)}(\mathbf{r}_{Aa}, \mathbf{k}_{Aa}) = \sum_l \exp(i\sigma_l^i) g_l Y_0^l(\hat{\mathbf{r}}_{aA}) \frac{\sqrt{4\pi(2l+1)}}{k_{aA} r_{aA}}, \tag{6.2.42}$$

and

$$\psi^{(-)}(\mathbf{r}_{bB}, \mathbf{k}_{bB}) = \frac{4\pi}{k_{bB} r_{bB}} \sum_{\tilde{l}} i^{\tilde{l}} \exp(-i\sigma_{\tilde{l}}^f) f_{\tilde{l}}^*(r_{bB}) \sum_m Y_m^{\tilde{l}*}(\hat{\mathbf{k}}_{bB}) Y_m^{\tilde{l}}(\hat{\mathbf{r}}_{bB}). \tag{6.2.43}$$

One can then write,

$$\begin{aligned}
T_{2N}^{1step} = & \langle \Psi_f^{(-)}(\mathbf{k}_{bB}) | V(r_{1p}) | \Psi_i^{(+)}(k_{aA}, \hat{\mathbf{z}}) \rangle = \frac{(4\pi)^{3/2}}{k_{aA} k_{bB}} \sum_{\tilde{l}lm} ((l_f \frac{1}{2})_{j_f} (l_f \frac{1}{2})_{j_f} | (l_f l_f)_0 (\frac{1}{2} \frac{1}{2})_0)_0 \\
& \times ((l_i \frac{1}{2})_{j_i} (l_i \frac{1}{2})_{j_i} | (l_i l_i)_0 (\frac{1}{2} \frac{1}{2})_0)_0 \sqrt{2l+1} i^{-l_p} \exp[i(\sigma_{\tilde{l}}^f + \sigma_l^i)] \\
& \times 2Y_m^{\tilde{l}}(\hat{\mathbf{k}}_{bB}) \sum_{\sigma_1 \sigma_2} \int \frac{d\mathbf{r}_{bB} d\mathbf{r} d\boldsymbol{\eta}}{r_{bB} r_{aA}} u_{l_f j_f}(r_{A1}) u_{l_f j_f}(r_{A2}) u_{l_i j_i}(r_{b1}) u_{l_i j_i}(r_{b2}) \\
& \times [Y^{l_f}(\hat{\mathbf{r}}_{A1}) Y^{l_f}(\hat{\mathbf{r}}_{A2})]_0^{0*} [Y^{l_i}(\hat{\mathbf{r}}_{b1}) Y^{l_i}(\hat{\mathbf{r}}_{b2})]_0^0 \\
& \times f_{\tilde{l}}(r_{bB}) g_l(r_{aA}) [\chi(\sigma_1) \chi(\sigma_2)]_0^{0*} Y_m^{\tilde{l}*}(\hat{\mathbf{r}}_{bB}) V(r_{1p}) \\
& \times [\chi(\sigma_1) \chi(\sigma_2)]_0^0 Y_0^l(\hat{\mathbf{r}}_{aA}),
\end{aligned} \tag{6.2.44}$$

which, after a number of simplifications becomes

$$\begin{aligned} \langle \Psi_f^{(-)}(\mathbf{k}_{bB}) | V(r_{1p}) | \Psi_i^{(+)}(k_{aA}, \hat{\mathbf{z}}) \rangle &= \frac{(4\pi)^{3/2}}{k_{aA} k_{bB}} \sum_{\tilde{l}\tilde{m}} \sqrt{\frac{(2j_f + 1)(2j_i + 1)}{(2l_f + 1)(2l_i + 1)}} \\ &\times \sqrt{2l + 1} i^{-\tilde{l}} \exp[i(\sigma_l^f + \sigma_l^i)] \\ &\times Y_m^{\tilde{l}}(\hat{\mathbf{k}}_{bB}) \int \frac{d\mathbf{r}_{bB} d\mathbf{r} d\boldsymbol{\eta}}{r_{bB} r_{aA}} u_{l_f j_f}(r_{A1}) u_{l_f j_f}(r_{A2}) u_{l_i j_i}(r_{b1}) u_{l_i j_i}(r_{b2}) \\ &\times [Y^{l_f}(\hat{\mathbf{r}}_{A1}) Y^{l_f}(\hat{\mathbf{r}}_{A2})]_0^{0*} [Y^{l_i}(\hat{\mathbf{r}}_{b1}) Y^{l_i}(\hat{\mathbf{r}}_{b2})]_0^0 \\ &\times f_l(r_{bB}) g_l(r_{aA}) Y_m^{\tilde{l}*}(\hat{\mathbf{r}}_{bB}) V(r_{1p}) Y_0^l(\hat{\mathbf{r}}_{aA}), \end{aligned} \quad (6.2.45)$$

where $l = \tilde{l}$ and $m = 0$. Making use of Legendre polynomials leads to,

$$\begin{aligned} \langle \Psi_f^{(-)}(\mathbf{k}_{bB}) | V(r_{1p}) | \Psi_i^{(+)}(k_{aA}, \hat{\mathbf{z}}) \rangle &= \frac{(4\pi)^{-1/2}}{k_{aA} k_{bB}} \sum_l \sqrt{(2j_f + 1)(2j_i + 1)} \\ &\times \sqrt{2l + 1} i^{-l} \exp[i(\sigma_l^f + \sigma_l^i)] Y_0^l(\hat{\mathbf{k}}_{bB}) \\ &\times \int \frac{d\mathbf{r}_{bB} d\mathbf{r} d\boldsymbol{\eta}}{r_{bB} r_{aA}} u_{l_f j_f}(r_{A1}) u_{l_f j_f}(r_{A2}) u_{l_i j_i}(r_{b1}) u_{l_i j_i}(r_{b2}) \\ &\times P_{l_f}(\cos \theta_A) P_{l_i}(\cos \theta_b) \\ &\times f_l(r_{bB}) g_l(r_{aA}) Y_0^{l*}(\hat{\mathbf{r}}_{bB}) V(r_{1p}) Y_0^l(\hat{\mathbf{r}}_{aA}). \end{aligned} \quad (6.2.46)$$

Changing the integration variables and proceeding as in last section, (implying the multiplicative factor $2\pi \sqrt{\frac{4\pi}{2l+1}}$), the above expression becomes

$$\begin{aligned} \langle \Psi_f^{(-)}(\mathbf{k}_{bB}) | V(r_{1p}) | \Psi_i^{(+)}(k_{aA}, \hat{\mathbf{z}}) \rangle &= \frac{2\pi}{k_{aA} k_{bB}} \sum_l \sqrt{(2j_f + 1)(2j_i + 1)} \\ &\times i^{-l} \exp[i(\sigma_l^f + \sigma_l^i)] Y_0^l(\hat{\mathbf{k}}_{bB}) \\ &\times \int dr_{aA} d\beta d\gamma dr_{12} dr_{b1} dr_{b2} r_{aA} \sin \beta r_{12} r_{b1} r_{b2} \\ &\times P_{l_f}(\cos \theta_A) P_{l_i}(\cos \theta_b) u_{l_f j_f}(r_{A1}) u_{l_f j_f}(r_{A2}) u_{l_i j_i}(r_{b1}) u_{l_i j_i}(r_{b2}) \\ &\times f_l(r_{bB}) g_l(r_{aA}) Y_0^{l*}(\hat{\mathbf{r}}_{bB}) V(r_{1p}) / r_{bB}, \end{aligned} \quad (6.2.47)$$

which eventually can be recasted, through the use of Legendre polynomials, in the expression,

$$\begin{aligned} T_{2N}^{1step} = \langle \Psi_f^{(-)}(\mathbf{k}_{bB}) | V(r_{1p}) | \Psi_i^{(+)}(k_{aA}, \hat{\mathbf{z}}) \rangle &= \frac{1}{2k_{aA} k_{bB}} \sum_l \sqrt{(2j_f + 1)(2j_i + 1)} \\ &\times i^{-l} \exp[i(\sigma_l^f + \sigma_l^i)] P_l(\cos \theta) (2l + 1) \\ &\times \int dr_{aA} d\beta d\gamma dr_{12} dr_{b1} dr_{b2} r_{aA} \sin \beta r_{12} r_{b1} r_{b2} \\ &\times P_{l_f}(\cos \theta_A) P_{l_i}(\cos \theta_b) u_{l_f j_f}(r_{A1}) u_{l_f j_f}(r_{A2}) V(r_{1p}) \\ &\times u_{l_i j_i}(r_{b1}) u_{l_i j_i}(r_{b2}) f_l(r_{bB}) g_l(r_{aA}) P_l(\cos \theta_{if}) / r_{bB}, \end{aligned} \quad (6.2.48)$$

expression which gives the same results as (6.2.41)

6.2.3 Coordinates for the calculation of simultaneous transfer

In what follows we explicit the coordinates used in the calculation of the above equations. Making use of the notation of Bayman (1971), we find the expression of the variables appearing in the integral as functions of the integration variables $r_{1p}, r_{2p}, r_{12}, R, \beta, \gamma$ (remember that $\mathbf{R} = R \hat{\mathbf{z}}$, see last section). \mathbf{R} being the center of mass coordinate. Thus, one can write

$$\mathbf{R} = \frac{1}{3}(\mathbf{r}_1 + \mathbf{r}_2 + \mathbf{r}_p) = \frac{1}{3}(\mathbf{R} + \mathbf{d}_1 + \mathbf{R} + \mathbf{d}_2 + \mathbf{R} + \mathbf{d}_p), \quad (6.2.49)$$

so

$$\mathbf{d}_1 + \mathbf{d}_2 + \mathbf{d}_p = 0. \quad (6.2.50)$$

Together with

$$\mathbf{d}_1 + \mathbf{r}_{12} = \mathbf{d}_2 \quad \mathbf{d}_2 + \mathbf{r}_{2p} = \mathbf{d}_p, \quad (6.2.51)$$

we find

$$\mathbf{d}_1 = \frac{1}{3}(2\mathbf{r}_{12} + \mathbf{r}_{2p}), \quad (6.2.52)$$

and

$$d_1^2 = \frac{1}{9}(4r_{12}^2 + r_{2p}^2 + 4\mathbf{r}_{12}\mathbf{r}_{2p}). \quad (6.2.53)$$

Making use of

$$\begin{aligned} \mathbf{r}_{12} + \mathbf{r}_{2p} &= \mathbf{r}_{1p} \\ r_{1p}^2 &= r_{12}^2 + r_{2p}^2 + 2\mathbf{r}_{12}\mathbf{r}_{2p} \\ 2\mathbf{r}_{12}\mathbf{r}_{2p} &= r_{1p}^2 - r_{12}^2 - r_{2p}^2. \end{aligned} \quad (6.2.54)$$

one obtains

$$d_1 = \frac{1}{3} \sqrt{2r_{12}^2 + 2r_{1p}^2 - r_{2p}^2}. \quad (6.2.55)$$

Similarly,

$$d_2 = \frac{1}{3} \sqrt{2r_{12}^2 + 2r_{2p}^2 - r_{1p}^2} \quad d_p = \frac{1}{3} \sqrt{2r_{2p}^2 + 2r_{1p}^2 - r_{12}^2}. \quad (6.2.56)$$

We now express the angle α between \mathbf{d}_1 and \mathbf{r}_{12} . We have

$$-\mathbf{d}_1 \mathbf{r}_{12} = r_{12} d_1 \cos(\alpha), \quad (6.2.57)$$

and

$$\begin{aligned} \mathbf{d}_1 + \mathbf{r}_{12} &= \mathbf{d}_2 \\ d_1^2 + r_{12}^2 + 2\mathbf{d}_1 \mathbf{r}_{12} &= d_2^2. \end{aligned} \quad (6.2.58)$$

Consequently,

$$\cos(\alpha) = \frac{d_1^2 + r_{12}^2 - d_2^2}{2r_{12}d_1}. \quad (6.2.59)$$

The complete determination of $\mathbf{r}_1, \mathbf{r}_2, \mathbf{r}_{12}$ can be made by writing their expression in a simple configuration, in which the triangle lies in the xz -plane with \mathbf{d}_1 pointing along the positive z -direction, and $\mathbf{R} = 0$. Then, a first rotation $\mathcal{R}_z(\gamma)$ of an angle γ around the z -axis, a second rotation $\mathcal{R}_y(\beta)$ of an angle β around the y -axis, and a translation along \mathbf{R} will bring the vectors to the most general configuration. In other words,

$$\begin{aligned} \mathbf{r}_1 &= \mathbf{R} + \mathcal{R}_y(\beta)\mathcal{R}_z(\gamma)\mathbf{r}'_1, \\ \mathbf{r}_{12} &= \mathcal{R}_y(\beta)\mathcal{R}_z(\gamma)\mathbf{r}'_{12}, \\ \mathbf{r}_2 &= \mathbf{r}_1 + \mathbf{r}_{12}, \end{aligned} \quad (6.2.60)$$

with

$$\mathbf{r}'_1 = \begin{bmatrix} 0 \\ 0 \\ d_1 \end{bmatrix}, \quad (6.2.61)$$

$$\mathbf{r}'_{12} = r_{12} \begin{bmatrix} \sin(\alpha) \\ 0 \\ -\cos(\alpha) \end{bmatrix}, \quad (6.2.62)$$

and the rotation matrixes are

$$\mathcal{R}_y(\beta) = \begin{bmatrix} \cos(\beta) & 0 & \sin(\beta) \\ 0 & 1 & 0 \\ -\sin(\beta) & 0 & \cos(\beta) \end{bmatrix}, \quad (6.2.63)$$

and

$$\mathcal{R}_z(\gamma) = \begin{bmatrix} \cos(\gamma) & -\sin(\gamma) & 0 \\ \sin(\gamma) & \cos(\gamma) & 0 \\ 0 & 0 & 1 \end{bmatrix}. \quad (6.2.64)$$

then

$$\mathbf{r}_1 = \begin{bmatrix} d_1 \sin(\beta) \\ 0 \\ R + d_1 \cos(\beta) \end{bmatrix}, \quad (6.2.65)$$

$$\mathbf{r}_{12} = \begin{bmatrix} r_{12} \cos(\beta) \cos(\gamma) \sin(\alpha) - r_{12} \sin(\beta) \cos(\alpha) \\ r_{12} \sin(\gamma) \sin(\alpha) \\ -r_{12} \sin(\beta) \cos(\gamma) \sin(\alpha) - r_{12} \cos(\alpha) \cos(\beta) \end{bmatrix}, \quad (6.2.66)$$

$$\mathbf{r}_2 = \begin{bmatrix} d_1 \sin(\beta) + r_{12} \cos(\beta) \cos(\gamma) \sin(\alpha) - r_{12} \sin(\beta) \cos(\alpha) \\ r_{12} \sin(\gamma) \sin(\alpha) \\ R + d_1 \cos(\beta) - r_{12} \sin(\beta) \cos(\gamma) \sin(\alpha) - r_{12} \cos(\alpha) \cos(\beta) \end{bmatrix}. \quad (6.2.67)$$

We also need $\cos(\theta_{12})$, ζ and $\cos(\theta_\zeta)$, θ_{12} being the angle between \mathbf{r}_1 and \mathbf{r}_2 , $\zeta = \mathbf{r}_p - \frac{\mathbf{r}_1 + \mathbf{r}_2}{A+2}$ the position of the proton with respect to the final nucleus, and θ_ζ the angle between ζ and the z -axis:

$$\cos(\theta_{12}) = \frac{\mathbf{r}_1 \cdot \mathbf{r}_2}{r_1 r_2}, \quad (6.2.68)$$

and

$$\zeta = 3\mathbf{R} - \frac{A+3}{A+2}(\mathbf{r}_1 + \mathbf{r}_2), \quad (6.2.69)$$

where we have used (6.2.49).

For heavy ions, we find instead

$$\mathbf{R} = \frac{1}{m_a} (\mathbf{r}_{A1} + \mathbf{r}_{A2} + m_b \mathbf{r}_{Ab}), \quad (6.2.70)$$

$$\mathbf{d}_1 = \frac{1}{m_a} (m_b \mathbf{r}_{b2} - (m_b + 1) \mathbf{r}_{12}), \quad (6.2.71)$$

$$d_1 = \frac{1}{m_a} \sqrt{(m_b + 1)r_{12}^2 + m_b(m_b + 1)r_{b1}^2 - m_b r_{b2}^2}, \quad (6.2.72)$$

$$d_2 = \frac{1}{m_a} \sqrt{(m_b + 1)r_{12}^2 + m_b(m_b + 1)r_{b2}^2 - m_b r_{b1}^2}, \quad (6.2.73)$$

and

$$\zeta = \frac{m_a}{m_b} \mathbf{R} - \frac{m_B + m_b}{m_b m_B} (\mathbf{r}_{A1} + \mathbf{r}_{A2}). \quad (6.2.74)$$

The rest of the formulae are identical to the (t, p) ones. We list them for convenience,

$$\mathbf{r}_{A1} = \begin{bmatrix} d_1 \sin(\beta) \\ 0 \\ R + d_1 \cos(\beta) \end{bmatrix}, \quad (6.2.75)$$

$$\mathbf{r}_{A2} = \begin{bmatrix} d_1 \sin(\beta) + r_{12} \cos(\beta) \cos(\gamma) \sin(\alpha) - r_{12} \sin(\beta) \cos(\alpha) \\ r_{12} \sin(\gamma) \sin(\alpha) \\ R + d_1 \cos(\beta) - r_{12} \sin(\beta) \cos(\gamma) \sin(\alpha) - r_{12} \cos(\alpha) \cos(\beta) \end{bmatrix}. \quad (6.2.76)$$

We we also find

$$\mathbf{r}_{b1} = \frac{1}{m_b} (\mathbf{r}_{A2} + (m_b + 1) \mathbf{r}_{A1} - m_a \mathbf{R}), \quad (6.2.77)$$

and

$$\mathbf{r}_{b2} = \frac{1}{m_b} (\mathbf{r}_{A1} + (m_b + 1) \mathbf{r}_{A2} - m_a \mathbf{R}). \quad (6.2.78)$$

One can readily obtain

$$\cos \theta_{12} = \frac{r_{A1}^2 + r_{A2}^2 - r_{12}^2}{2r_{A1}r_{A2}}, \quad (6.2.79)$$

and

$$\cos \theta_i = \frac{r_{b1}^2 + r_{b2}^2 - r_{12}^2}{2r_{b1}r_{b2}}. \quad (6.2.80)$$

6.2.4 Matrix element for the transition amplitude (alternative derivation)

In what follows we work out an alternative derivation of T_{2N}^{1step} , more closely related to heavy ion reactions⁸, that is

$$\begin{aligned}
 T^{(1)}(\theta) = & 2 \frac{(4\pi)^{3/2}}{k_{Aa} k_{Bb}} \sum_{l_p j_p m_l j_p} i^{-l_p} \exp[i(\sigma_{l_p}^p + \sigma_{l_t}^t)] \sqrt{2l_t + 1} \\
 & \times \langle l_p m - m_p | 1/2 m_p | j_p m \rangle \langle l_t 0 1/2 m_t | j_t m_t \rangle Y_{m-m_p}^{l_p}(\hat{\mathbf{k}}_{Bb}) \\
 & \times \sum_{\sigma_1 \sigma_2 \sigma_p} \int d\mathbf{r}_{Cc} d\mathbf{r}_{b1} d\mathbf{r}_{A2} [\psi^{j_f}(\mathbf{r}_{A1}, \sigma_1) \psi^{j_f}(\mathbf{r}_{A2}, \sigma_2)]_0^{0*} \\
 & \times v(r_{b1}) [\psi^{j_i}(\mathbf{r}_{b1}, \sigma_1) \psi^{j_i}(\mathbf{r}_{b2}, \sigma_2)]_0^0 \frac{g_{l_t j_t}(r_{Aa}) f_{l_p j_p}(r_{Bb})}{r_{Aa} r_{Bb}} \\
 & \times [Y^{l_t}(\hat{\mathbf{r}}_{Aa}) \chi(\sigma_p)]_{m_t}^{j_t} [Y^{l_p}(\hat{\mathbf{r}}_{Bb}) \chi(\sigma_p)]_m^{j_p*}. \tag{6.2.81}
 \end{aligned}$$

As shown above one can write,

$$\begin{aligned}
 & \sum_{\sigma_p} \langle l_p m - m_p | 1/2 m_p | j_p m \rangle \langle l_t 0 1/2 m_t | j_t m_t \rangle [Y^{l_t}(\hat{\mathbf{r}}_{Aa}) \chi(\sigma_p)]_{m_t}^{j_t} [Y^{l_p}(\hat{\mathbf{r}}_{Bb}) \chi(\sigma_p)]_m^{j_p*} \\
 & = -\frac{\delta_{l_p, l_t} \delta_{j_p, j_t} \delta_{m, m_t}}{\sqrt{2l+1}} [Y^l(\hat{\mathbf{r}}_{Aa}) Y^l(\hat{\mathbf{r}}_{Bb})]_0^0 \begin{cases} \frac{l}{2l+1} & \text{if } m_t = m_p \\ -\frac{\sqrt{l(l+1)}}{2l+1} & \text{if } m_t = -m_p \end{cases} \tag{6.2.82}
 \end{aligned}$$

when $j = l - 1/2$ and

$$\begin{aligned}
 & \sum_{\sigma_p} \langle l_p m - m_p | 1/2 m_p | j_p m \rangle \langle l_t 0 1/2 m_t | j_t m_t \rangle [Y^{l_t}(\hat{\mathbf{r}}_{Aa}) \chi(\sigma_p)]_{m_t}^{j_t} [Y^{l_p}(\hat{\mathbf{r}}_{Bb}) \chi(\sigma_p)]_m^{j_p*} \\
 & = -\frac{\delta_{l_p, l_t} \delta_{j_p, j_t} \delta_{m, m_t}}{\sqrt{2l+1}} [Y^l(\hat{\mathbf{r}}_{Aa}) Y^l(\hat{\mathbf{r}}_{Bb})]_0^0 \begin{cases} \frac{l+1}{2l+1} & \text{if } m_t = m_p \\ \frac{\sqrt{l(l+1)}}{2l+1} & \text{if } m_t = -m_p \end{cases} \tag{6.2.83}
 \end{aligned}$$

⁸Bayman and Chen (1982).

if $j = l + 1/2$. One then gets

$$\begin{aligned}
 T^{(1)}(\mu = 0; \theta) = & 2 \frac{(4\pi)^{3/2}}{k_{Aa} k_{Bb}} \sum_l i^{-l} \frac{\exp[i(\sigma_l^p + \sigma_l^t)]}{2l+1} Y_{m_t - m_p}^l(\hat{\mathbf{k}}_{Bb}) \\
 & \times \sum_{\sigma_1 \sigma_2} \int \frac{d\mathbf{r}_{Cc} d\mathbf{r}_{b1} d\mathbf{r}_{A2}}{r_{Aa} r_{Bb}} [\psi^{j_f}(\mathbf{r}_{A1}, \sigma_1) \psi^{j_f}(\mathbf{r}_{A2}, \sigma_2)]_0^{0*} \\
 & \times v(r_{b1}) [\psi^{j_i}(\mathbf{r}_{b1}, \sigma_1) \psi^{j_i}(\mathbf{r}_{b2}, \sigma_2)]_0^0 [Y^l(\hat{\mathbf{r}}_{Aa}) Y^l(\hat{\mathbf{r}}_{Bb})]_0^0 \\
 & \times \left[(f_{ll+1/2}(r_{Bb}) g_{ll+1/2}(r_{Aa})(l+1) + f_{ll-1/2}(r_{Bb}) g_{ll-1/2}(r_{Aa})l) \delta_{m_p, m_t} \right. \\
 & \left. + (f_{ll+1/2}(r_{Bb}) g_{ll+1/2}(r_{Aa}) \sqrt{l(l+1)} - f_{ll-1/2}(r_{Bb}) g_{ll-1/2}(r_{Aa}) \sqrt{l(l+1)}) \delta_{m_p, -m_t} \right]. \tag{6.2.84}
 \end{aligned}$$

Making use of the relations,

$$\begin{aligned}
 & [\psi^{j_f}(\mathbf{r}_{A1}, \sigma_1) \psi^{j_f}(\mathbf{r}_{A2}, \sigma_2)]_0^{0*} \\
 & = ((l_f \frac{1}{2})_{j_f} (l_f \frac{1}{2})_{j_f} | (l_f l_f)_0 (\frac{1}{2} \frac{1}{2})_0)_0 u_{l_f}(r_{A1}) u_{l_f}(r_{A2}) \\
 & \times [Y^{l_f}(\hat{\mathbf{r}}_{A1}) Y^{l_f}(\hat{\mathbf{r}}_{A2})]_0^{0*} [\chi(\sigma_1) \chi(\sigma_2)]_0^{0*} \\
 & = \sqrt{\frac{2j_f + 1}{2(2l_f + 1)}} u_{l_f}(r_{A1}) u_{l_f}(r_{A2}) \\
 & \times [Y^{l_f}(\hat{\mathbf{r}}_{A1}) Y^{l_f}(\hat{\mathbf{r}}_{A2})]_0^{0*} [\chi(\sigma_1) \chi(\sigma_2)]_0^{0*} \\
 & = \sqrt{\frac{2j_f + 1}{2}} \frac{u_{l_f}(r_{A1}) u_{l_f}(r_{A2})}{4\pi} P_{l_f}(\cos \omega_A) [\chi(\sigma_1) \chi(\sigma_2)]_0^{0*}, \tag{6.2.85}
 \end{aligned}$$

and

$$\begin{aligned}
 & [\psi^{j_i}(\mathbf{r}_{b1}, \sigma_1) \psi^{j_i}(\mathbf{r}_{b2}, \sigma_2)]_0^0 \\
 & = ((l_i \frac{1}{2})_{j_i} (l_i \frac{1}{2})_{j_i} | (l_i l_i)_0 (\frac{1}{2} \frac{1}{2})_0)_0 u_{l_i}(r_{b1}) u_{l_i}(r_{b2}) \\
 & \times [Y^{l_i}(\hat{\mathbf{r}}_{b1}) Y^{l_i}(\hat{\mathbf{r}}_{b2})]_0^0 [\chi(\sigma_1) \chi(\sigma_2)]_0^0 \\
 & = \sqrt{\frac{2j_i + 1}{2(2l_i + 1)}} u_{l_i}(r_{b1}) u_{l_i}(r_{b2}) \\
 & \times [Y^{l_i}(\hat{\mathbf{r}}_{b1}) Y^{l_i}(\hat{\mathbf{r}}_{b2})]_0^0 [\chi(\sigma_1) \chi(\sigma_2)]_0^0 \\
 & = \sqrt{\frac{2j_i + 1}{2}} \frac{u_{l_i}(r_{b1}) u_{l_i}(r_{b2})}{4\pi} P_{l_i}(\cos \omega_b) [\chi(\sigma_1) \chi(\sigma_2)]_0^0, \tag{6.2.86}
 \end{aligned}$$

where ω_A is the angle between \mathbf{r}_{A1} and \mathbf{r}_{A2} , and ω_b is the angle between \mathbf{r}_{b1} and \mathbf{r}_{b2} . Consequently

$$\begin{aligned} T^{(1)}(\theta) = & (4\pi)^{-3/2} \frac{\sqrt{(2j_i+1)(2j_f+1)}}{k_{Aa}k_{Bb}} \sum_l i^{-l} \frac{\exp[i(\sigma_l^p + \sigma_l^t)]}{\sqrt{2l+1}} Y_m^l(\hat{\mathbf{k}}_{Bb}) \\ & \times \int \frac{d\mathbf{r}_{Cc} d\mathbf{r}_{b1} d\mathbf{r}_{A2}}{r_{Aa} r_{Bb}} P_{l_f}(\cos \omega_A) P_{l_i}(\cos \omega_b) P_l(\cos \omega_{if}) \\ & \times v(r_{b1}) u_{l_i}(r_{b1}) u_{l_i}(r_{b2}) u_{l_f}(r_{A1}) u_{l_f}(r_{A2}) \\ & \times \left[(f_{ll+1/2}(r_{Bb}) g_{ll+1/2}(r_{Aa})(l+1) + f_{ll-1/2}(r_{Bb}) g_{ll-1/2}(r_{Aa})l) \delta_{m_p, m_t} \right. \\ & \left. + (f_{ll+1/2}(r_{Bb}) g_{ll+1/2}(r_{Aa}) \sqrt{l(l+1)} - f_{ll-1/2}(r_{Bb}) g_{ll-1/2}(r_{Aa}) \sqrt{l(l+1)}) \delta_{m_p, -m_t} \right], \end{aligned} \quad (6.2.87)$$

where ω_{if} is the angle between \mathbf{r}_{Aa} and \mathbf{r}_{Bb} . For heavy ions, we can consider that the optical potential does not have a spin-orbit term, and the distorted waves are independent of j . We thus have

$$\begin{aligned} T^{(1)}(\theta) = & (4\pi)^{-3/2} \frac{\sqrt{(2j_i+1)(2j_f+1)}}{k_{Aa}k_{Bb}} \sum_l i^{-l} \exp[i(\sigma_l^p + \sigma_l^t)] Y_0^l(\hat{\mathbf{k}}_{Bb}) \sqrt{2l+1} \\ & \times \int \frac{d\mathbf{r}_{Cc} d\mathbf{r}_{b1} d\mathbf{r}_{A2}}{r_{Aa} r_{Bb}} P_{l_f}(\cos \omega_A) P_{l_i}(\cos \omega_b) P_l(\cos \omega_{if}) \\ & \times v(r_{b1}) u_{l_i}(r_{b1}) u_{l_i}(r_{b2}) u_{l_f}(r_{A1}) u_{l_f}(r_{A2}) f_l(r_{Bb}) g_l(r_{Aa}). \end{aligned} \quad (6.2.88)$$

Changing variables one obtains,

$$\begin{aligned} T^{(1)}(\theta) = & (4\pi)^{-1} \frac{\sqrt{(2j_i+1)(2j_f+1)}}{k_{Aa}k_{Bb}} \sum_l \exp[i(\sigma_l^p + \sigma_l^t)] P_l(\cos \theta) (2l+1) \\ & \times \int dr_{1A} dr_{2A} dr_{Aa} d(\cos \beta) d(\cos \omega_A) d\gamma r_{1A}^2 r_{2A}^2 r_{Aa}^2 \\ & \times P_{l_f}(\cos \omega_A) P_{l_i}(\cos \omega_b) P_l(\cos \omega_{if}) v(r_{b1}) \\ & \times u_{l_i}(r_{b1}) u_{l_i}(r_{b2}) u_{l_f}(r_{A1}) u_{l_f}(r_{A2}) f_l(r_{Bb}) g_l(r_{Aa}). \end{aligned} \quad (6.2.89)$$

6.2.5 Coordinates used to derive Eq. (6.2.89)

We determine the relation between the integration variables in (6.2.87) and the coordinates needed to evaluate the quantities in the integrand. Noting that

$$\mathbf{r}_{Aa} = \frac{\mathbf{r}_{A1} + \mathbf{r}_{A2} + m_b \mathbf{r}_{Ab}}{m_b + 2}, \quad (6.2.90)$$

one has

$$\mathbf{r}_{b1} = \mathbf{r}_{bA} + \mathbf{r}_{A1} = \frac{(m_b + 1)\mathbf{r}_{A1} + \mathbf{r}_{A2} - (m_b + 2)\mathbf{r}_{Aa}}{m_b}, \quad (6.2.91)$$

$$\mathbf{r}_{b2} = \mathbf{r}_{bA} + \mathbf{r}_{A2} = \frac{(m_b + 1)\mathbf{r}_{A2} + \mathbf{r}_{A1} - (m_b + 2)\mathbf{r}_{Aa}}{m_b}, \quad (6.2.92)$$

and

$$\begin{aligned} \mathbf{r}_{Cc} &= \mathbf{r}_{CA} + \mathbf{r}_{A1} + \mathbf{r}_{1c} = -\frac{1}{m_A + 1}\mathbf{r}_{A2} + \mathbf{r}_{A1} - \frac{m_b}{m_b + 1}\mathbf{r}_{b1} \\ &= \frac{m_b + 2}{m_b + 1}\mathbf{r}_{Aa} - \frac{m_b + 2 + m_A}{(m_b + 1)(m_A + 1)}\mathbf{r}_{A2} \end{aligned} \quad (6.2.93)$$

Since,

$$\mathbf{r}_{AB} = \frac{\mathbf{r}_{A1} + \mathbf{r}_{A2}}{m_A + 2}, \quad (6.2.94)$$

one obtains

$$\mathbf{r}_{Bb} = \mathbf{r}_{BA} + \mathbf{r}_{Ab} = \frac{m_b + 2}{m_b}\mathbf{r}_{Aa} - \frac{m_A + m_b + 2}{(m_A + 2)m_b}(\mathbf{r}_{A1} + \mathbf{r}_{A2}). \quad (6.2.95)$$

Using the same rotations as those used in Section 6.2.3 one gets,

$$\mathbf{r}_{A1} = r_{A1} \begin{bmatrix} \sin \alpha \\ 0 \\ \cos \alpha \end{bmatrix}, \quad (6.2.96)$$

and

$$\mathbf{r}_{A2} = r_{A2} \begin{bmatrix} -\cos \alpha \cos \gamma \sin \omega_A + \sin \alpha \cos \omega_A \\ -\sin \gamma \sin \omega_A \\ \sin \alpha \cos \gamma \sin \omega_A + \cos \alpha \cos \omega_A \end{bmatrix}, \quad (6.2.97)$$

with

$$\cos \alpha = \frac{r_{A1}^2 - d_1^2 + r_{Aa}^2}{2r_{A1}r_{Aa}}, \quad (6.2.98)$$

and

$$d_1 = \sqrt{r_{A1}^2 - r_{Aa}^2 \sin^2 \beta} - r_{Aa} \cos \beta. \quad (6.2.99)$$

Note that though β, r_{1A}, r_{Aa} are independent integration variables, they have to fulfill the condition

$$r_{Aa} \sin \beta \leq r_{A1}, \quad \text{for } 0 \leq \beta \leq \pi. \quad (6.2.100)$$

The expression of the remaining quantities appearing in the integral are now straightforward,

$$\begin{aligned} r_{b1} &= m_b^{-1} |(m_b + 1)\mathbf{r}_{A1} + \mathbf{r}_{A2} - (m_b + 2)\mathbf{r}_{Aa}| \\ &= m_b^{-1} \left((m_b + 2)^2 r_{Aa}^2 + (m_b + 1)^2 r_{A1}^2 + r_{A2}^2 \right. \\ &\quad \left. - 2(m_b + 2)(m_b + 1)\mathbf{r}_{Aa} \mathbf{r}_{A1} - 2(m_b + 2)\mathbf{r}_{Aa} \mathbf{r}_{A2} + 2(m_b + 1)\mathbf{r}_{A1} \mathbf{r}_{A2} \right)^{1/2}, \end{aligned} \quad (6.2.101)$$

$$\begin{aligned}
r_{b2} &= m_b^{-1} |(m_b + 1)\mathbf{r}_{A2} + \mathbf{r}_{A1} - (m_b + 2)\mathbf{r}_{Aa}| \\
&= m_b^{-1} \left((m_b + 2)^2 r_{Aa}^2 + (m_b + 1)^2 r_{A2}^2 + r_{A1}^2 \right. \\
&\quad \left. - 2(m_b + 2)(m_b + 1)\mathbf{r}_{Aa} \mathbf{r}_{A2} - 2(m_b + 2)\mathbf{r}_{Aa} \mathbf{r}_{A1} + 2(m_b + 1)\mathbf{r}_{A2} \mathbf{r}_{A1} \right)^{1/2}, \\
\end{aligned} \tag{6.2.102}$$

$$\begin{aligned}
r_{Bb} &= \left| \frac{m_b + 2}{m_b} \mathbf{r}_{Aa} - \frac{m_A + m_b + 2}{(m_A + 2)m_b} (\mathbf{r}_{A1} + \mathbf{r}_{A2}) \right| \\
&= \left[\left(\frac{m_b + 2}{m_b} \right)^2 r_{Aa}^2 + \left(\frac{m_A + m_b + 2}{(m_A + 2)m_b} \right)^2 (r_{A1}^2 + r_{A2}^2 + 2\mathbf{r}_{A1} \mathbf{r}_{A2}) \right. \\
&\quad \left. - 2 \frac{(m_b + 2)(m_A + m_b + 2)}{(m_A + 2)m_b^2} \mathbf{r}_{Aa} (\mathbf{r}_{A1} + \mathbf{r}_{A2}) \right]^{1/2}, \\
\end{aligned} \tag{6.2.103}$$

$$\begin{aligned}
r_{Cc} &= \left| \frac{m_b + 2}{m_b + 1} \mathbf{r}_{Aa} - \frac{m_b + 2 + m_A}{(m_b + 1)(m_A + 1)} \mathbf{r}_{A2} \right| \\
&= \left[\left(\frac{m_a}{(m_a - 1)} \right)^2 r_{Aa}^2 + \left(\frac{m_A + m_a}{(m_A + 1)(m_a - 1)} \right)^2 r_{A2}^2 \right. \\
&\quad \left. - 2 \frac{m_A m_a + m_a^2}{(m_A + 1)(m_a - 1)^2} \mathbf{r}_{Aa} \mathbf{r}_{A2} \right]^{1/2}, \\
\end{aligned} \tag{6.2.104}$$

$$\cos \omega_b = \frac{\mathbf{r}_{b1} \mathbf{r}_{b2}}{r_{b1} r_{b2}}, \tag{6.2.105}$$

$$\cos \omega_{if} = \frac{\mathbf{r}_{Aa} \mathbf{r}_{Bb}}{r_{Aa} r_{Bb}}, \tag{6.2.106}$$

with

$$\mathbf{r}_{Aa} \mathbf{r}_{A1} = r_{Aa} r_{A1} \cos \alpha, \tag{6.2.107}$$

$$\mathbf{r}_{Aa} \mathbf{r}_{A2} = r_{Aa} r_{A2} (\sin \alpha \cos \gamma \sin \omega_A + \cos \alpha \cos \omega_A), \tag{6.2.108}$$

$$\mathbf{r}_{A1} \mathbf{r}_{A2} = r_{A1} r_{A2} \cos \omega_A. \tag{6.2.109}$$

6.2.6 Successive transfer

The successive two-neutron transfer amplitudes can be written as (Bayman and Chen (1982)):

$$\begin{aligned}
T_{succ}^{(2)}(\theta) = & \frac{4\mu_{Cc}}{\hbar^2} \sum_{\substack{\sigma_1 \sigma_2 \\ \sigma'_1 \sigma'_2 \\ KM}} \int d^3 r_{Cc} d^3 r_{b1} d^3 r_{A2} d^3 r'_{Cc} d^3 r'_{b1} d^3 r'_{A2} \chi^{(-)*}(\mathbf{k}_{Bb}, \mathbf{r}_{Bb}) \\
& \times [\psi^{j_f}(\mathbf{r}_{A1}, \sigma_1) \psi^{j_f}(\mathbf{r}_{A2}, \sigma_2)]_0^{0*} v(r_{b1}) [\psi^{j_f}(\mathbf{r}_{A2}, \sigma_2) \psi^{j_i}(\mathbf{r}_{b1}, \sigma_1)]_M^K \\
& \times G(\mathbf{r}_{Cc}, \mathbf{r}'_{Cc}) [\psi^{j_f}(\mathbf{r}'_{A2}, \sigma'_2) \psi^{j_i}(\mathbf{r}'_{b1}, \sigma'_1)]_M^{K*} v(r'_{c2}) \\
& \times [\psi^{j_i}(\mathbf{r}'_{b1}, \sigma'_1) \psi^{j_i}(\mathbf{r}'_{b2}, \sigma'_2)]_0^0 \chi^{(+)}(\mathbf{r}'_{Aa}). \tag{6.2.110}
\end{aligned}$$

It is of notice that the time-reversal phase convention is used throughout. Expanding the Green function and the distorted waves in a basis of angular momentum eigenstate one can write,

$$\chi^{(-)*}(\mathbf{k}_{Bb}, \mathbf{r}_{Bb}) = \sum_{\tilde{l}} \frac{4\pi}{k_{Bb} r_{Bb}} i^{-\tilde{l}} e^{i\sigma_f^{\tilde{l}}} F_{\tilde{l}} \sum_m Y_m^{\tilde{l}}(\hat{r}_{Bb}) Y_m^{\tilde{l}*}(\hat{k}_{Bb}), \tag{6.2.111}$$

the sum over m being

$$\sum_m (-1)^{\tilde{l}-m} Y_m^{\tilde{l}}(\hat{r}_{Bb}) Y_{-m}^{\tilde{l}}(\hat{k}_{Bb}) = \sqrt{2\tilde{l}+1} [Y^{\tilde{l}}(\hat{r}_{Bb}) Y^{\tilde{l}}(\hat{k}_{Bb})]_0^0, \tag{6.2.112}$$

where we have used (6.A.2) and (6.A.18), so

$$\chi^{(-)*}(\mathbf{k}_{Bb}, \mathbf{r}_{Bb}) = \sum_{\tilde{l}} \sqrt{2\tilde{l}+1} \frac{4\pi}{k_{Bb} r_{Bb}} i^{-\tilde{l}} e^{i\sigma_f^{\tilde{l}}} F_{\tilde{l}}(r_{Bb}) [Y^{\tilde{l}}(\hat{r}_{Bb}) Y^{\tilde{l}}(\hat{k}_{Bb})]_0^0. \tag{6.2.113}$$

Similarly,

$$\chi^{(+)}(\mathbf{r}'_{Aa}) = \sum_l i^l \sqrt{2l+1} \frac{4\pi}{k_{Aa} r'_{Aa}} e^{i\sigma_i^l} F_l(r'_{Aa}) [Y^l(\hat{r}'_{Aa}) Y^l(\hat{k}_{Aa})]_0^0 \tag{6.2.114}$$

where we have taken into account the choice $\hat{k}_{Aa} \equiv \hat{z}$. The Green function can be written as

$$G(\mathbf{r}_{Cc}, \mathbf{r}'_{Cc}) = i \sum_{l_c} \sqrt{2l_c+1} \frac{f_{l_c}(k_{Cc}, r_{<}) P_{l_c}(k_{Cc}, r_{>})}{k_{Cc} r_{Cc} r'_{Cc}} [Y^{l_c}(\hat{r}_{Cc}) Y^{l_c}(\hat{r}'_{Cc})]_0^0. \tag{6.2.115}$$

Finally

$$\begin{aligned}
T_{succ}^{(2)}(\theta) = & \frac{4\mu_{Cc}(4\pi)^2 i}{\hbar^2 k_{Aa} k_{Bb} k_{Cc}} \sum_{l, l_c, \tilde{l}} e^{i(\sigma'_l + \sigma'_f)l - \tilde{l}} \sqrt{(2l+1)(2l_c+1)(2\tilde{l}+1)} \\
& \times \sum_{\sigma'_1 \sigma'_2} \int d^3 r_{Cc} d^3 r_{b1} d^3 r_{A2} d^3 r'_{Cc} d^3 r'_{b1} d^3 r'_{A2} v(r_{b1}) v(r'_{c2}) \left[Y^{\tilde{l}}(\hat{r}_{Bb}) Y^{\tilde{l}}(\hat{k}_{Bb}) \right]_0^0 \\
& \times \left[Y^l(\hat{r}'_{Aa}) Y^l(\hat{k}'_{Aa}) \right]_0^0 \left[Y^{l_c}(\hat{r}_{Cc}) Y^{l_c}(\hat{r}'_{Cc}) \right]_0^0 \frac{F_l(r_{Bb})}{r_{Bb}} \frac{F_l(r'_{Aa})}{r'_{Aa}} \\
& \times \frac{f_{l_c}(k_{Cc}, r_<) P_{l_c}(k_{Cc}, r_>)}{r_{Cc} r'_{Cc}} \left[\psi^{j_f}(\mathbf{r}_{A1}, \sigma_1) \psi^{j_f}(\mathbf{r}_{A2}, \sigma_2) \right]_0^{0*} \\
& \times \left[\psi^{j_i}(\mathbf{r}'_{b1}, \sigma'_1) \psi^{j_i}(\mathbf{r}'_{b2}, \sigma'_2) \right]_0^0 \sum_{KM} \left[\psi^{j_f}(\mathbf{r}_{A2}, \sigma_2) \psi^{j_i}(\mathbf{r}_{b1}, \sigma_1) \right]_M^K \\
& \times \left[\psi^{j_f}(\mathbf{r}'_{A2}, \sigma'_2) \psi^{j_i}(\mathbf{r}'_{b1}, \sigma'_1) \right]_M^{K*}.
\end{aligned} \tag{6.2.116}$$

Let us now perform the integration over \mathbf{r}_{A2} ,

$$\begin{aligned}
& \sum_{\sigma_1, \sigma_2} \int d\mathbf{r}_{A2} \left[\psi^{j_f}(\mathbf{r}_{A1}, \sigma_1) \psi^{j_f}(\mathbf{r}_{A2}, \sigma_2) \right]_0^{0*} \left[\psi^{j_f}(\mathbf{r}_{A2}, \sigma_2) \psi^{j_i}(\mathbf{r}_{b1}, \sigma_1) \right]_M^K \\
& = \sum_{\sigma_1, \sigma_2} (-1)^{1/2-\sigma_1+1/2-\sigma_2} \int d\mathbf{r}_{A2} \left[\psi^{j_f}(\mathbf{r}_{A1}, -\sigma_1) \psi^{j_f}(\mathbf{r}_{A2}, -\sigma_2) \right]_0^0 \left[\psi^{j_f}(\mathbf{r}_{A2}, \sigma_2) \psi^{j_i}(\mathbf{r}_{b1}, \sigma_1) \right]_M^K \\
& = - \sum_{\sigma_1, \sigma_2} (-1)^{1/2-\sigma_1+1/2-\sigma_2} \int d\mathbf{r}_{A2} \left[\psi^{j_f}(\mathbf{r}_{A2}, -\sigma_2) \psi^{j_f}(\mathbf{r}_{A1}, -\sigma_1) \right]_0^0 \left[\psi^{j_f}(\mathbf{r}_{A2}, \sigma_2) \psi^{j_i}(\mathbf{r}_{b1}, \sigma_1) \right]_M^K \\
& = - ((j_f j_f)_0 (j_f j_i)_K | (j_f j_f)_0 (j_f j_i)_K)_K \sum_{\sigma_1, \sigma_2} (-1)^{1/2-\sigma_1+1/2-\sigma_2} \\
& \quad \times \int d\mathbf{r}_{A2} \left[\psi^{j_f}(\mathbf{r}_{A2}, -\sigma_2) \psi^{j_f}(\mathbf{r}_{A2}, \sigma_2) \right]_0^0 \left[\psi^{j_f}(\mathbf{r}_{A1}, -\sigma_1) \psi^{j_i}(\mathbf{r}_{b1}, \sigma_1) \right]_M^K \\
& = \frac{1}{2j_f + 1} \sqrt{2j_f + 1} ((l_f \frac{1}{2})_{j_f} (l_i \frac{1}{2})_{j_i} | (l_f l_i)_K (\frac{1}{2} \frac{1}{2})_0)_K \\
& \quad \times u_{l_f}(r_{A1}) u_{l_i}(r_{b1}) \left[Y^{l_f}(\hat{r}_{A1}) Y^{l_i}(\hat{r}_{b1}) \right]_M^K \sum_{\sigma_1} (-1)^{1/2-\sigma_1} \left[\chi^{1/2}(-\sigma_1) \chi^{1/2}(\sigma_1) \right]_0^0 \\
& = - \sqrt{\frac{2}{2j_f + 1}} ((l_f \frac{1}{2})_{j_f} (l_i \frac{1}{2})_{j_i} | (l_f l_i)_K (\frac{1}{2} \frac{1}{2})_0)_K \left[Y^{l_f}(\hat{r}_{A1}) Y^{l_i}(\hat{r}_{b1}) \right]_M^K u_{l_f}(r_{A1}) u_{l_i}(r_{b1}),
\end{aligned} \tag{6.2.117}$$

where we have evaluated the 9 j -symbol

$$((j_f j_f)_0 (j_f j_i)_K | (j_f j_f)_0 (j_f j_i)_K)_K = \frac{1}{2j_f + 1}, \tag{6.2.118}$$

as well as (6.A.19). We proceed in a similar way to evaluate the integral over \mathbf{r}'_{b1} ,

$$\begin{aligned}
& \sum_{\sigma'_1, \sigma'_2} \int d\mathbf{r}'_{b1} [\psi^{j_i}(\mathbf{r}'_{b1}, \sigma'_1) \psi^{j_i}(\mathbf{r}'_{b2}, \sigma'_2)]_0^0 [\psi^{j_f}(\mathbf{r}'_{A2}, \sigma'_2) \psi^{j_i}(\mathbf{r}'_{b1}, \sigma'_1)]_M^{K*} \\
&= -(-1)^{K-M} \sum_{\sigma'_1, \sigma'_2} \int d\mathbf{r}'_{b1} [\psi^{j_f}(\mathbf{r}'_{A2}, -\sigma'_2) \psi^{j_i}(\mathbf{r}'_{b1}, -\sigma'_1)]_{-M}^K \\
&\quad \times [\psi^{j_i}(\mathbf{r}'_{b2}, \sigma'_2) \psi^{j_i}(\mathbf{r}'_{b1}, \sigma'_1)]_0^0 (-1)^{1/2-\sigma'_1+1/2-\sigma'_2} \\
&= -(-1)^{K-M} ((j_f j_i)_K (j_i j_i)_0 | (j_f j_i)_K (j_i j_i)_0)_K (-\sqrt{2 j_i + 1}) \\
&\quad \times ((l_f \frac{1}{2})_{j_f} (l_i \frac{1}{2})_{j_i} | (l_f l_i)_K (\frac{1}{2} \frac{1}{2})_0)_K (-\sqrt{2}) u_{l_f}(r'_{A2}) u_{l_i}(r'_{b2}) [Y^{l_f}(\hat{r}'_{A2}) Y^{l_i}(\hat{r}'_{b2})]_{-M}^K \\
&= -\sqrt{\frac{2}{2 j_i + 1}} ((l_f \frac{1}{2})_{j_f} (l_i \frac{1}{2})_{j_i} | (l_f l_i)_K (\frac{1}{2} \frac{1}{2})_0)_K [Y^{l_f}(\hat{r}'_{A2}) Y^{l_i}(\hat{r}'_{b2})]_M^{K*} u_{l_f}(r'_{A2}) u_{l_i}(r'_{b2}). \tag{6.2.119}
\end{aligned}$$

Setting the different elements together one obtains

$$\begin{aligned}
T_{succ}^{(2)}(\theta) &= \frac{4\mu_{Cc}(4\pi)^2 i}{\hbar^2 k_{Aa} k_{Bb} k_{Cc}} \frac{2}{\sqrt{(2j_i + 1)(2j_f + 1)}} \sum_{K,M} ((l_f \frac{1}{2})_{j_f} (l_i \frac{1}{2})_{j_i} | (l_f l_i)_K (\frac{1}{2} \frac{1}{2})_0)_K^2 \\
&\quad \times \sum_{l_c, l, \tilde{l}} e^{i(\sigma'_i + \sigma_f^l)} \sqrt{(2l_c + 1)(2l + 1)(2\tilde{l} + 1)} i^{l-\tilde{l}} \\
&\quad \times \int d^3 r_{Cc} d^3 r_{b1} d^3 r'_{Cc} d^3 r'_{A2} v(r_{b1}) v(r'_{c2}) u_{l_f}(r_{A1}) u_{l_i}(r_{b1}) u_{l_f}(r'_{A2}) u_{l_i}(r'_{b2}) \\
&\quad \times [Y^{l_f}(\hat{r}'_{A2}) Y^{l_i}(\hat{r}'_{b2})]_M^{K*} [Y^{l_f}(\hat{r}_{A1}) Y^{l_i}(\hat{r}_{b1})]_M^K \frac{F_l(r'_{Aa}) F_{\tilde{l}}(r'_{Bb}) f_{l_c}(k_{Cc}, r_<) P_{l_c}(k_{Cc}, r_>)}{r'_{Aa} r_{Bb} r_{Cc} r'_{Cc}} \\
&\quad \times [Y^{\tilde{l}}(\hat{r}_{Bb}) Y^{\tilde{l}}(\hat{k}_{Bb})]_0^0 [Y^l(\hat{r}'_{Aa}) Y^l(\hat{k}_{Aa})]_0^0 [Y^{l_c}(\hat{r}_{Cc}) Y^{l_c}(\hat{r}'_{Cc})]_0^0. \tag{6.2.120}
\end{aligned}$$

We now proceed to write this expression in a more compact way. For this purpose one writes

$$\begin{aligned}
& [Y^{\tilde{l}}(\hat{r}_{Bb}) Y^{\tilde{l}}(\hat{k}_{Bb})]_0^0 [Y^l(\hat{r}'_{Aa}) Y^l(\hat{k}_{Aa})]_0^0 = \\
& ((l l)_0 (\tilde{l} \tilde{l})_0 | (l \tilde{l})_0 (l \tilde{l})_0)_0 [Y^{\tilde{l}}(\hat{r}_{Bb}) Y^l(\hat{r}'_{Aa})]_0^0 [Y^{\tilde{l}}(\hat{k}_{Bb}) Y^l(\hat{k}_{Aa})]_0^0 \\
& = \frac{\delta_{\tilde{l}l}}{2l + 1} [Y^l(\hat{r}_{Bb}) Y^l(\hat{r}'_{Aa})]_0^0 [Y^l(\hat{k}_{Bb}) Y^l(\hat{k}_{Aa})]_0^0. \tag{6.2.121}
\end{aligned}$$

Taking into account the relations

$$[Y^l(\hat{k}_{Bb}) Y^l(\hat{k}_{Aa})]_0^0 = \frac{(-1)^l}{\sqrt{4\pi}} Y_0^l(\hat{k}_{Bb}) i^l, \tag{6.2.122}$$

and

$$\begin{aligned}
& \left[Y^l(\hat{r}_{Bb}) Y^l(\hat{r}'_{Aa}) \right]_0^0 \left[Y^{l_c}(\hat{r}_{Cc}) Y^{l_c}(\hat{r}'_{Cc}) \right]_0^0 = \\
& ((ll)_0(l_c l_c)_0 | (ll_c)_K (ll_c)_K)_0 \left\{ \left[Y^l(\hat{r}_{Bb}) Y^{l_c}(\hat{r}_{Cc}) \right]^K \left[Y^l(\hat{r}'_{Aa}) Y^{l_c}(\hat{r}'_{Cc}) \right]^K \right\}_0^0 \\
& = \sqrt{\frac{2K+1}{(2l+1)(2l_c+1)}} \\
& \times \sum_{M'} \frac{(-1)^{K+M'}}{\sqrt{2K+1}} \left[Y^l(\hat{r}_{Bb}) Y^{l_c}(\hat{r}_{Cc}) \right]_{-M'}^K \left[Y^l(\hat{r}'_{Aa}) Y^{l_c}(\hat{r}'_{Cc}) \right]_{M'}^K \\
& = \sqrt{\frac{1}{(2l+1)(2l_c+1)}} \\
& \times \sum_{M'} \left[Y^l(\hat{r}_{Bb}) Y^{l_c}(\hat{r}_{Cc}) \right]_{M'}^{K*} \left[Y^l(\hat{r}'_{Aa}) Y^{l_c}(\hat{r}'_{Cc}) \right]_{M'}^K.
\end{aligned} \tag{6.2.123}$$

It is of notice that the integrals

$$\int d\hat{r}_{Cc} d\hat{r}_{b1} \left[Y^l(\hat{r}_{Bb}) Y^{l_c}(\hat{r}_{Cc}) \right]_M^{K*} \left[Y^{l_f}(\hat{r}_{A1}) Y^{l_i}(\hat{r}_{b1}) \right]_M^K, \tag{6.2.124}$$

and

$$\int d\hat{r}'_{Cc} d\hat{r}'_{A2} \left[Y^l(\hat{r}'_{Aa}) Y^{l_c}(\hat{r}'_{Cc}) \right]_M^K \left[Y^{l_f}(\hat{r}'_{A2}) Y^{l_i}(\hat{r}'_{b2}) \right]_M^{K*}, \tag{6.2.125}$$

over the angular variables do not depend on M . Let us see why this is so with the help of (6.2.124),

$$\begin{aligned}
& \left[Y^l(\hat{r}_{Bb}) Y^{l_c}(\hat{r}_{Cc}) \right]_M^{K*} \left[Y^{l_f}(\hat{r}_{A1}) Y^{l_i}(\hat{r}_{b1}) \right]_M^K = (-1)^{K-M} \left[Y^l(\hat{r}_{Bb}) Y^{l_c}(\hat{r}_{Cc}) \right]_{-M}^K \\
& \times \left[Y^{l_f}(\hat{r}_{A1}) Y^{l_i}(\hat{r}_{b1}) \right]_M^K = (-1)^{K-M} \sum_J \langle K \ K \ M \ -M | J \ 0 \rangle \\
& \times \left\{ \left[Y^l(\hat{r}_{Bb}) Y^{l_c}(\hat{r}_{Cc}) \right]^K \left[Y^{l_f}(\hat{r}_{A1}) Y^{l_i}(\hat{r}_{b1}) \right]^K \right\}_0^J.
\end{aligned} \tag{6.2.126}$$

After integration, only the term

$$\begin{aligned}
& (-1)^{K-M} \langle K \ K \ M \ -M | 0 \ 0 \rangle \left\{ \left[Y^l(\hat{r}_{Bb}) Y^{l_c}(\hat{r}_{Cc}) \right]^K \left[Y^{l_f}(\hat{r}_{A1}) Y^{l_i}(\hat{r}_{b1}) \right]^K \right\}_0^0 = . \\
& \frac{1}{\sqrt{2K+1}} \left\{ \left[Y^l(\hat{r}_{Bb}) Y^{l_c}(\hat{r}_{Cc}) \right]^K \left[Y^{l_f}(\hat{r}_{A1}) Y^{l_i}(\hat{r}_{b1}) \right]^K \right\}_0^0
\end{aligned} \tag{6.2.127}$$

corresponding to $J = 0$ survives, which is indeed independent of M . We can thus omit the sum over M in (6.2.120) and multiply by $(2K + 1)$, obtaining

$$\begin{aligned} T_{succ}^{(2)}(\theta) &= \frac{64\mu_{Cc}(\pi)^{3/2}i}{\hbar^2 k_{Aa} k_{Bb} k_{Cc}} \frac{i^{-l}}{\sqrt{(2j_i + 1)(2j_f + 1)}} \\ &\times \sum_K (2K + 1)((l_f \frac{1}{2})_{j_f} (l_i \frac{1}{2})_{j_i} | (l_f l_i)_K (\frac{1}{2} \frac{1}{2})_0)_K^2 \\ &\times \sum_{l_c, l} \frac{e^{i(\sigma_i^l + \sigma_f^l)}}{\sqrt{(2l + 1)}} Y_0^l(\hat{k}_{Bb}) S_{K, l, l_c}, \end{aligned} \quad (6.2.128)$$

where

$$\begin{aligned} S_{K, l, l_c} &= \int d^3 r_{Cc} d^3 r_{b1} v(r_{b1}) u_{l_f}(r_{A1}) u_{l_i}(r_{b1}) \frac{s_{K, l, l_c}(r_{Cc})}{r_{Cc}} \frac{F_l(r_{Bb})}{r_{Bb}} \\ &\times [Y^{l_f}(\hat{r}_{A1}) Y^{l_i}(\hat{r}_{b1})]_M^K [Y^{l_c}(\hat{r}_{Cc}) Y^l(\hat{r}_{Bb})]_M^{K*}, \end{aligned} \quad (6.2.129)$$

and

$$\begin{aligned} s_{K, l, l_c}(r_{Cc}) &= \int_{r_{Cc, fixed}} d^3 r'_{Cc} d^3 r'_{A2} v(r'_{c2}) u_{l_f}(r'_{A2}) u_{l_i}(r'_{b2}) \frac{F_l(r'_{Aa})}{r'_{Aa}} \frac{f_l(k_{Cc}, r_{<}) P_{l_c}(k_{Cc}, r_{>})}{r'_{Cc}} \\ &\times [Y^{l_f}(\hat{r}'_{A2}) Y^{l_i}(\hat{r}'_{b2})]_M^{K*} [Y^{l_c}(\hat{r}'_{Cc}) Y^l(\hat{r}'_{Aa})]_M^K. \end{aligned} \quad (6.2.130)$$

It can be shown that the integrand in (6.2.129) is independent of M . Consequently, one can sum over M and divide by $(2K + 1)$, to get

$$\begin{aligned} \frac{1}{2K + 1} v(r_{b1}) u_{l_f}(r_{A1}) u_{l_i}(r_{b1}) \frac{s_{K, l, l_c}(r_{Cc})}{r_{Cc}} \frac{F_l(r_{Bb})}{r_{Bb}} \\ \times \sum_M [Y^{l_f}(\hat{r}_{A1}) Y^{l_i}(\hat{r}_{b1})]_M^K [Y^{l_c}(\hat{r}_{Cc}) Y^l(\hat{r}_{Bb})]_M^{K*}. \end{aligned} \quad (6.2.131)$$

This integrand is rotationally invariant (it is proportional to a T_M^L spherical tensor with $L = 0, M = 0$), so one can evaluate it in the “standard” configuration in which \mathbf{r}_{Cc} is directed along the z -axis and multiply by $8\pi^2$ (see Bayman and Chen (1982)), obtaining the final expression for S_{K, l, l_c} :

$$\begin{aligned} S_{K, l, l_c} &= \frac{4\pi^{3/2} \sqrt{2l_c + 1}}{2K + 1} i^{-l_c} \\ &\times \int r_{Cc}^2 dr_{Cc} r_{b1}^2 dr_{b1} \sin \theta d\theta v(r_{b1}) u_{l_f}(r_{A1}) u_{l_i}(r_{b1}) \\ &\times \frac{s_{K, l, l_c}(r_{Cc})}{r_{Cc}} \frac{F_l(r_{Bb})}{r_{Bb}} \\ &\times \sum_M \langle l_c 0 l M | K M \rangle [Y^{l_f}(\hat{r}_{A1}) Y^{l_i}(\theta + \pi, 0)]_M^K Y_M^{l*}(\hat{r}_{Bb}). \end{aligned} \quad (6.2.132)$$

Similarly, one has

$$\begin{aligned}
 s_{K,l,l_c}(r_{Cc}) &= \frac{4\pi^{3/2} \sqrt{2l_c + 1}}{2K + 1} i^{l_c} \\
 &\times \int r'_{Cc}^2 dr'_{Cc} r'^2_{A2} dr'_{A2} \sin \theta' d\theta' v(r'_{c2}) u_{l_f}(r'_{A2}) u_{l_i}(r'_{b2}) \\
 &\times \frac{F_l(r'_{Aa})}{r'_{Aa}} \frac{f_{l_c}(k_{Cc}, r_<) P_{l_c}(k_{Cc}, r_>)}{r'_{Cc}} \\
 &\times \sum_M \langle l_c 0 l M | K M \rangle [Y^{l_f}(\hat{r}'_{A2}) Y^{l_i}(\hat{r}'_{b2})]_M^{K*} Y_M^l(\hat{r}'_{Aa}).
 \end{aligned} \tag{6.2.133}$$

Introducing the further approximations $\mathbf{r}_{A1} \approx \mathbf{r}_{C1}$ and $\mathbf{r}_{b2} \approx \mathbf{r}_{c2}$, one obtains the final expression

$$\begin{aligned}
 T_{succ}^{(2)}(\theta) &= \frac{1024\mu_{Cc}\pi^{9/2}i}{\hbar^2 k_{Aa} k_{Bb} k_{Cc}} \frac{1}{\sqrt{(2j_i+1)(2j_f+1)}} \\
 &\times \sum_K \frac{1}{2K+1} ((l_f \frac{1}{2})_{j_f} (l_i \frac{1}{2})_{j_i} | (l_f l_i)_K (\frac{1}{2} \frac{1}{2})_0)_K^2 \\
 &\times \sum_{l_c, l} e^{i(\sigma_i^l + \sigma_f^l)} \frac{(2l_c + 1)}{\sqrt{2l + 1}} Y_0^l(\hat{k}_{Bb}) S_{K,l,l_c},
 \end{aligned} \tag{6.2.134}$$

with

$$\begin{aligned}
 S_{K,l,l_c} &= \int r_{Cc}^2 dr_{Cc} r_{b1}^2 dr_{b1} \sin \theta d\theta v(r_{b1}) u_{l_f}(r_{C1}) u_{l_i}(r_{b1}) \\
 &\times \frac{s_{K,l,l_c}(r_{Cc})}{r_{Cc}} \frac{F_l(r_{Bb})}{r_{Bb}} \\
 &\times \sum_M \langle l_c 0 l M | K M \rangle [Y^{l_f}(\hat{r}_{C1}) Y^{l_i}(\theta + \pi, 0)]_M^K Y_M^{l*}(\hat{r}_{Bb}),
 \end{aligned} \tag{6.2.135}$$

and

$$\begin{aligned}
 s_{K,l,l_c}(r_{Cc}) &= \int r'_{Cc}^2 dr'_{Cc} r'^2_{A2} dr'_{A2} \sin \theta' d\theta' v(r'_{c2}) u_{l_f}(r'_{A2}) u_{l_i}(r'_{c2}) \\
 &\times \frac{F_l(r'_{Aa})}{r'_{Aa}} \frac{f_{l_c}(k_{Cc}, r_<) P_{l_c}(k_{Cc}, r_>)}{r'_{Cc}} \\
 &\times \sum_M \langle l_c 0 l M | K M \rangle [Y^{l_f}(\hat{r}'_{A2}) Y^{l_i}(\hat{r}'_{c2})]_M^{K*} Y_M^l(\hat{r}'_{Aa}).
 \end{aligned} \tag{6.2.136}$$

6.2.7 Coordinates for the successive transfer

In the standard configuration in which the integrals (6.2.135) and (6.2.136) are to be evaluated, we have

$$\mathbf{r}_{Cc} = r_{Cc} \hat{\mathbf{z}}, \quad \mathbf{r}_{b1} = r_{b1}(-\cos \theta \hat{\mathbf{z}} - \sin \theta \hat{\mathbf{x}}). \tag{6.2.137}$$

Now,

$$\begin{aligned}\mathbf{r}_{C1} &= \mathbf{r}_{Cc} + \mathbf{r}_{c1} = \mathbf{r}_{Cc} + \frac{m_b}{m_b + 1} \mathbf{r}_{b1} \\ &= \left(r_{Cc} - \frac{m_b}{m_b + 1} r_{b1} \cos \theta \right) \hat{\mathbf{z}} - \frac{m_b}{m_b + 1} r_{b1} \sin \theta \hat{\mathbf{x}},\end{aligned}\quad (6.2.138)$$

and

$$\mathbf{r}_{Bb} = \mathbf{r}_{BC} + \mathbf{r}_{Cb} = -\frac{1}{m_B} \mathbf{r}_{C1} + \mathbf{r}_{Cb}. \quad (6.2.139)$$

Substituting the relation

$$\mathbf{r}_{Cb} = \mathbf{r}_{Cc} + \mathbf{r}_{cb} = \mathbf{r}_{Cc} - \frac{1}{m_b + 1} \mathbf{r}_{b1}, \quad (6.2.140)$$

in (6.2.139) one gets

$$\mathbf{r}_{Bb} = \left(\frac{m_B - 1}{m_B} r_{Cc} + \frac{m_b + m_B}{m_B(m_b + 1)} r_{b1} \cos \theta \right) \hat{\mathbf{z}} + \frac{m_b + m_B}{m_B(m_b + 1)} r_{b1} \sin \theta \hat{\mathbf{x}}. \quad (6.2.141)$$

The primed variables are arranged in a similar fashion,

$$\mathbf{r}'_{Cc} = r'_{Cc} \hat{\mathbf{z}}, \quad \mathbf{r}'_{A2} = r'_{A2} (-\cos \theta' \hat{\mathbf{z}} - \sin \theta' \hat{\mathbf{x}}). \quad (6.2.142)$$

Thus,

$$\mathbf{r}'_{c2} = \left(-r'_{Cc} - \frac{m_A}{m_A + 1} r'_{A2} \cos \theta' \right) \hat{\mathbf{z}} - \frac{m_A}{m_A + 1} r'_{A2} \sin \theta' \hat{\mathbf{x}}, \quad (6.2.143)$$

and

$$\mathbf{r}'_{Aa} = \left(\frac{m_a - 1}{m_a} r'_{Cc} - \frac{m_A + m_a}{m_a(m_A + 1)} r'_{A2} \cos \theta' \right) \hat{\mathbf{z}} - \frac{m_A + m_a}{m_a(m_A + 1)} r'_{A2} \sin \theta' \hat{\mathbf{x}}. \quad (6.2.144)$$

6.2.8 Simplifying the vector coupling

We will now turn our attention to the vector-coupled quantities in (6.2.135) and (6.2.136),

$$\sum_M \langle l_c 0 l M | K M \rangle \left[Y^{l_f}(\hat{r}_{C1}) Y^{l_i}(\theta + \pi, 0) \right]_M^K Y_M^{l*}(\hat{r}_{Bb}), \quad (6.2.145)$$

and

$$\sum_M \langle l_c 0 l M | K M \rangle \left[Y^{l_f}(\hat{r}'_{A2}) Y^{l_i}(\hat{r}'_{c2}) \right]_M^{K*} Y_M^l(\hat{r}'_{Aa}). \quad (6.2.146)$$

We can express them both as

$$\sum_M f(M), \quad (6.2.147)$$

where e.g. in the case of (6.2.145), one has

$$f(M) = \langle l_c 0 l M | K M \rangle \left[Y^{l_f}(\hat{r}_{C1}) Y^{l_i}(\theta + \pi, 0) \right]_M^K Y_M^{l^*}(\hat{r}_{Bb}). \quad (6.2.148)$$

Note that all the vectors that come into play in the above expressions are in the (x, z) -plane. Consequently, the azimuthal angle ϕ is always equal to zero. Under these circumstances and for time-reversed phases, $(Y_M^{L^*}(\theta, 0) = (-1)^L Y_M^L(\theta, 0))$ one has

$$f(-M) = (-1)^{l_c + l_f + l_i + l} f(M). \quad (6.2.149)$$

Consequently,

$$\begin{aligned} \sum_M \langle l_c 0 l M | K M \rangle f(M) &= \langle l_c 0 l 0 | K 0 \rangle f(0) \\ &+ \sum_{M>0} \langle l_c 0 l M | K M \rangle f(M) \left(1 + (-1)^{l_c + l + l_i + l_f} \right). \end{aligned} \quad (6.2.150)$$

Consequently, in the case in which $l_c + l + l_i + l_f$ is odd, we have only to evaluate the $M = 0$ contribution. This consideration is useful to restrict the number of numerical operations needed to calculate the transition amplitude.

6.2.9 non-orthogonality term

We write the non-orthogonality contribution to the transition amplitude (see Bayman and Chen (1982)):

$$\begin{aligned} T_{NO}^{(2)}(\theta) = 2 \sum_{\substack{\sigma_1 \sigma_2 \\ \sigma'_1 \sigma'_2 \\ KM}} \int d^3 r_{Cc} d^3 r_{b1} d^3 r_{A2} d^3 r'_{b1} d^3 r'_{A2} \chi^{(-)*}(\mathbf{k}_{Bb}, \mathbf{r}_{Bb}) \\ \times \left[\psi^{j_f}(\mathbf{r}_{A1}, \sigma_1) \psi^{j_f}(\mathbf{r}_{A2}, \sigma_2) \right]_0^{0*} v(r_{b1}) \left[\psi^{j_f}(\mathbf{r}_{A2}, \sigma_2) \psi^{j_i}(\mathbf{r}_{b1}, \sigma_1) \right]_M^K \\ \times \left[\psi^{j_f}(\mathbf{r}'_{A2}, \sigma'_2) \psi^{j_i}(\mathbf{r}'_{b1}, \sigma'_1) \right]_M^{K*} \left[\psi^{j_i}(\mathbf{r}'_{b1}, \sigma'_1) \psi^{j_i}(\mathbf{r}'_{b2}, \sigma'_2) \right]_0^0 \chi^{(+)}(\mathbf{r}'_{Aa}). \end{aligned} \quad (6.2.151)$$

This expression is equivalent to (6.2.110) if we make the replacement

$$\frac{2\mu_{Cc}}{\hbar^2} G(\mathbf{r}_{Cc}, \mathbf{r}'_{Cc}) v(r'_{A2}) \rightarrow \delta(\mathbf{r}_{Cc} - \mathbf{r}'_{Cc}). \quad (6.2.152)$$

Looking at the partial-wave expansions of $G(\mathbf{r}_{Cc}, \mathbf{r}'_{Cc})$ and $\delta(\mathbf{r}_{Cc} - \mathbf{r}'_{Cc})$ (see App. 6.A), we find that we can use the above expressions for the successive transfer with the replacement

$$i \frac{2\mu_{Cc}}{\hbar^2} \frac{f_{l_c}(k_{Cc}, r_<) P_{l_c}(k_{Cc}, r_>)}{k_{Cc}} \rightarrow \delta(r_{Cc} - r'_{Cc}). \quad (6.2.153)$$

We thus have

$$\begin{aligned} T_{2NT}^{NO} = & \frac{512\pi^{9/2}}{k_{Aa}k_{Bb}} \frac{1}{\sqrt{(2j_i+1)(2j_f+1)}} \\ & \times \sum_K ((l_f \frac{1}{2})_{j_f} (l_i \frac{1}{2})_{j_i} | (l_f l_i)_K (\frac{1}{2} \frac{1}{2})_0)_K^2 \\ & \times \sum_{l_c, l} e^{i(\sigma_i^l + \sigma_f^l)} \frac{(2l_c + 1)}{\sqrt{2l + 1}} Y_0^l(\hat{k}_{Bb}) S_{K, l, l_c}, \end{aligned} \quad (6.2.154)$$

with

$$\begin{aligned} S_{K, l, l_c} = & \int r_{Cc}^2 dr_{Cc} r_{b1}^2 dr_{b1} \sin \theta d\theta v(r_{b1}) u_{l_f}(r_{C1}) u_{l_i}(r_{b1}) \\ & \times \frac{s_{K, l, l_c}(r_{Cc})}{r_{Cc}} \frac{F_l(r_{Bb})}{r_{Bb}} \\ & \times \sum_M \langle l_c 0 l M | K M \rangle [Y^{l_f}(\hat{r}_{C1}) Y^{l_i}(\theta + \pi, 0)]_M^K Y_M^{l*}(\hat{r}_{Bb}), \end{aligned} \quad (6.2.155)$$

and

$$\begin{aligned} s_{K, l, l_c}(r_{Cc}) = & r_{Cc} \int dr'_{A2} r'^2_{A2} \sin \theta' d\theta' u_{l_f}(r'_{A2}) u_{l_i}(r'_{c2}) \frac{F_l(r'_{Aa})}{r'_{Aa}} \\ & \times \sum_M \langle l_c 0 l M | K M \rangle [Y^{l_f}(\hat{r}'_{A2}) Y^{l_i}(\hat{r}'_{c2})]_M^{K*} Y_M^l(\hat{r}'_{Aa}). \end{aligned} \quad (6.2.156)$$

6.2.10 Arbitrary orbital momentum transfer

We will now examine the case in which the two transferred nucleons carry an angular momentum Λ different from 0. Let us assume that two nucleons coupled to angular momentum Λ in the initial nucleus A are transferred into a final state of zero angular momentum in nucleus B . The transition amplitude is given by the integral

$$\begin{aligned} 2 \sum_{\sigma_1 \sigma_2} \int d\mathbf{r}_{cC} d\mathbf{r}_{A2} d\mathbf{r}_{b1} \chi^{(-)*}(\mathbf{r}_{bB}) [\psi^{j_f}(\mathbf{r}_{A1}, \sigma_1) \psi^{j_f}(\mathbf{r}_{A2}, \sigma_2)]_0^{0*} \\ \times v(r_{b1}) \Psi^{(+)}(\mathbf{r}_{aA}, \mathbf{r}_{b1}, \mathbf{r}_{b2}, \sigma_1, \sigma_2). \end{aligned} \quad (6.2.157)$$

If we neglect core excitations, the above expression is exact as long as $\Psi^{(+)}(\mathbf{r}_{aA}, \mathbf{r}_{b1}, \mathbf{r}_{b2}, \sigma_1, \sigma_2)$ is the exact wavefunction. We can instead obtain an approximation for the transfer amplitude using

$$\begin{aligned} \Psi^{(+)}(\mathbf{r}_{aA}, \mathbf{r}_{b1}, \mathbf{r}_{b2}, \sigma_1, \sigma_2) \approx & \chi^{(+)}(\mathbf{r}_{aA}) [\psi^{j_{i1}}(\mathbf{r}_{b1}, \sigma_1) \psi^{j_{i2}}(\mathbf{r}_{b2}, \sigma_2)]_\mu^\Lambda \\ & + \sum_{K, M} \mathcal{U}_{K, M}(\mathbf{r}_{cC}) [\psi^{j_f}(\mathbf{r}_{A2}, \sigma_2) \psi^{j_{i1}}(\mathbf{r}_{b1}, \sigma_1)]_M^K \end{aligned} \quad (6.2.158)$$

as an approximation for the incoming state. The first term of (6.2.158) gives rise to the simultaneous amplitude, while from second one leads to both the successive and the non-orthogonality contributions. To extract the amplitude $\mathcal{U}_{K,M}(\mathbf{r}_{cC})$, we define $f_{KM}(\mathbf{r}_{cC})$ as the scalar product

$$f_{KM}(\mathbf{r}_{cC}) = \left\langle \left[\psi^{j_f}(\mathbf{r}_{A2}, \sigma_2) \psi^{j_{i1}}(\mathbf{r}_{b1}, \sigma_1) \right]_M^K \middle| \Psi^{(+)}(\mathbf{r}_{aA}, \mathbf{r}_{b1}, \mathbf{r}_{b2}, \sigma_1, \sigma_2) \right\rangle \quad (6.2.159)$$

for fixed \mathbf{r}_{cC} , which can be seen to obey the equation

$$\begin{aligned} & \left(\frac{\hbar^2}{2\mu_{cC}} k_{cC}^2 + \frac{\hbar^2}{2\mu_{cC}} \nabla_{r_{cC}}^2 - U(r_{cC}) \right) f_{KM}(\mathbf{r}_{cC}) \\ &= \left\langle \left[\psi^{j_f}(\mathbf{r}_{A2}, \sigma_2) \psi^{j_{i1}}(\mathbf{r}_{b1}, \sigma_1) \right]_M^K \middle| v(r_{c2}) \middle| \Psi^{(+)}(\mathbf{r}_{aA}, \mathbf{r}_{b1}, \mathbf{r}_{b2}, \sigma_1, \sigma_2) \right\rangle. \end{aligned} \quad (6.2.160)$$

The solution can be written in terms of the Green function $G(\mathbf{r}_{cC}, \mathbf{r}'_{cC})$ defined by

$$\left(\frac{\hbar^2}{2\mu_{cC}} k_{cC}^2 + \frac{\hbar^2}{2\mu_{cC}} \nabla_{r_{cC}}^2 - U(r_{cC}) \right) G(\mathbf{r}_{cC}, \mathbf{r}'_{cC}) = \frac{\hbar^2}{2\mu_{cC}} \delta(\mathbf{r}_{cC} - \mathbf{r}'_{cC}). \quad (6.2.161)$$

Thus,

$$\begin{aligned} f_{KM}(\mathbf{r}_{cC}) &= \frac{2\mu_{cC}}{\hbar^2} \int d\mathbf{r}'_{cC} G(\mathbf{r}_{cC}, \mathbf{r}'_{cC}) \left\langle \left[\psi^{j_f}(\mathbf{r}'_{A2}, \sigma'_2) \psi^{j_{i1}}(\mathbf{r}'_{b1}, \sigma'_1) \right]_M^K \middle| v(r_{c2}) \middle| \Psi^{(+)}(\mathbf{r}'_{aA}, \mathbf{r}'_{b1}, \mathbf{r}'_{b2}, \sigma'_1, \sigma'_2) \right\rangle \\ &\approx \frac{2\mu_{cC}}{\hbar^2} \sum_{\sigma'_1 \sigma'_2} \int d\mathbf{r}'_{cC} d\mathbf{r}'_{A2} d\mathbf{r}'_{b1} G(\mathbf{r}_{cC}, \mathbf{r}'_{cC}) \left[\psi^{j_f}(\mathbf{r}'_{A2}, \sigma'_2) \psi^{j_{i1}}(\mathbf{r}'_{b1}, \sigma'_1) \right]_M^{K*} \\ &\quad \times v(r'_{c2}) \chi^{(+)}(\mathbf{r}'_{aA}) \left[\psi^{j_{i1}}(\mathbf{r}'_{b1}, \sigma'_1) \psi^{j_{i2}}(\mathbf{r}'_{b2}, \sigma'_2) \right]_\mu^\Lambda = \mathcal{U}_{K,M}(\mathbf{r}_{cC}) \\ &\quad + \left\langle \left[\psi^{j_f}(\mathbf{r}'_{A2}, \sigma_2) \psi^{j_{i1}}(\mathbf{r}'_{b1}, \sigma_1) \right]_M^K \middle| \chi^{(+)}(\mathbf{r}'_{aA}) \left[\psi^{j_{i1}}(\mathbf{r}'_{b1}, \sigma'_1) \psi^{j_{i2}}(\mathbf{r}'_{b2}, \sigma'_2) \right]_\mu^\Lambda \right\rangle. \end{aligned} \quad (6.2.162)$$

Therefore

$$\begin{aligned} \mathcal{U}_{K,M}(\mathbf{r}_{cC}) &= \frac{2\mu_{cC}}{\hbar^2} \sum_{\sigma'_1 \sigma'_2} \int d\mathbf{r}'_{cC} d\mathbf{r}'_{A2} d\mathbf{r}'_{b1} G(\mathbf{r}_{cC}, \mathbf{r}'_{cC}) \left[\psi^{j_f}(\mathbf{r}'_{A2}, \sigma'_2) \psi^{j_{i1}}(\mathbf{r}'_{b1}, \sigma'_1) \right]_M^{K*} \\ &\quad \times v(r'_{c2}) \chi^{(+)}(\mathbf{r}'_{aA}) \left[\psi^{j_{i1}}(\mathbf{r}'_{b1}, \sigma'_1) \psi^{j_{i2}}(\mathbf{r}'_{b2}, \sigma'_2) \right]_\mu^\Lambda \\ &\quad - \left\langle \left[\psi^{j_f}(\mathbf{r}'_{A2}, \sigma_2) \psi^{j_{i1}}(\mathbf{r}'_{b1}, \sigma_1) \right]_M^K \middle| \chi^{(+)}(\mathbf{r}'_{aA}) \left[\psi^{j_{i1}}(\mathbf{r}'_{b1}, \sigma'_1) \psi^{j_{i2}}(\mathbf{r}'_{b2}, \sigma'_2) \right]_\mu^\Lambda \right\rangle. \end{aligned} \quad (6.2.163)$$

When we substitute $\mathcal{U}_{K,M}(\mathbf{r}_{cC})$ into (6.2.158) and (6.2.157), the first term gives rise to the successive amplitude for the two-particle transfer, while the second term is responsible for the non-orthogonal contribution.

6.2.11 Successive transfer contribution

We need to evaluate the integral

$$\begin{aligned} T_{succ}^{(2)}(\theta; \mu) = & \frac{4\mu_{cC}}{\hbar^2} \sum_{\sigma_1 \sigma_2} \sum_{KM} \int d\mathbf{r}_{cC} d\mathbf{r}_{A2} d\mathbf{r}_{b1} d\mathbf{r}'_{cC} d\mathbf{r}'_{A2} d\mathbf{r}'_{b1} [\psi^{j_f}(\mathbf{r}_{A1}, \sigma_1) \psi^{j_f}(\mathbf{r}_{A2}, \sigma_2)]_0^{0*} \\ & \times \chi^{(-)*}(\mathbf{r}_{bB}) G(\mathbf{r}_{cC}, \mathbf{r}'_{cC}) [\psi^{j_f}(\mathbf{r}'_{A2}, \sigma'_2) \psi^{j_{i1}}(\mathbf{r}'_{b1}, \sigma'_1)]_M^{K*} \chi^{(+)}(\mathbf{r}'_{aA}) v(r'_{c2}) v(r_{b1}) \\ & \times [\psi^{j_{i1}}(\mathbf{r}'_{b1}, \sigma'_1) \psi^{j_{i2}}(\mathbf{r}'_{b2}, \sigma'_2)]_\mu^K [\psi^{j_f}(\mathbf{r}_{A2}, \sigma_2) \psi^{j_{i1}}(\mathbf{r}_{b1}, \sigma_1)]_M^K, \end{aligned} \quad (6.2.164)$$

where we must substitute the Green function and the distorted waves by their partial wave expansions (see App. 6.B). The integral over \mathbf{r}'_{b1} is:

$$\begin{aligned} & \sum_{\sigma'_1} \int d\mathbf{r}'_{b1} [\psi^{j_f}(\mathbf{r}'_{A2}, \sigma'_2) \psi^{j_{i1}}(\mathbf{r}'_{b1}, \sigma'_1)]_M^{K*} [\psi^{j_{i1}}(\mathbf{r}'_{b1}, \sigma'_1) \psi^{j_{i2}}(\mathbf{r}'_{b2}, \sigma'_2)]_\mu^K \\ &= \sum_{\sigma'_1} \int d\mathbf{r}'_{b1} (-1)^{-M+j_f+j_{i1}-\sigma_1-\sigma_2} [\psi^{j_{i1}}(\mathbf{r}'_{b1}, -\sigma'_1) \psi^{j_f}(\mathbf{r}'_{A2}, -\sigma'_2)]_{-M}^K [\psi^{j_{i1}}(\mathbf{r}'_{b1}, \sigma'_1) \psi^{j_{i2}}(\mathbf{r}'_{b2}, \sigma'_2)]_\mu^K \\ &= \sum_{\sigma'_1} \int d\mathbf{r}'_{b1} (-1)^{-M+j_f+j_{i1}-\sigma_1-\sigma_2} \sum_P \langle K \Lambda - M \mu | P \mu - M \rangle ((j_{i1} j_f)_K (j_{i1} j_{i2})_\Lambda | (j_{i1} j_{i1})_0 (j_f j_{i2})_P)_P \\ & \quad \times [\psi^{j_{i1}}(\mathbf{r}'_{b1}, -\sigma'_1) \psi^{j_{i1}}(\mathbf{r}'_{b1}, \sigma'_1)]_0^0 [\psi^{j_f}(\mathbf{r}'_{A2}, -\sigma'_2) \psi^{j_{i2}}(\mathbf{r}'_{b2}, \sigma'_2)]_{\mu-M}^P \\ &= (-1)^{-M+j_f+j_{i1}} \sqrt{2j_{i1}+1} u_{l_f}(r_{A2}) u_{l_{i2}}(r'_{b2}) \sum_P \langle K \Lambda - M \mu | P \mu - M \rangle \\ & \quad \times ((j_{i1} j_f)_K (j_{i1} j_{i2})_\Lambda | (j_{i1} j_{i1})_0 (j_f j_{i2})_P)_P ((l_f \frac{1}{2})_{j_f} (l_{i2} \frac{1}{2})_{j_{i2}} | (l_f l_{i2})_P (\frac{1}{2} \frac{1}{2})_0)_P \\ & \quad \times [Y^{l_f}(\hat{\mathbf{r}}'_{A2}) Y^{l_{i2}}(\hat{\mathbf{r}}'_{b2})]_{\mu-M}^P u_{l_f}(r_{A2}) u_{l_{i2}}(r_{b2}). \end{aligned} \quad (6.2.165)$$

Integrating over \mathbf{r}_{A2} (see (6.2.117)) leads to,

$$\begin{aligned} & \sum_{\sigma_2} \int d\mathbf{r}_{A2} [\psi^{j_f}(\mathbf{r}_{A1}, \sigma_1) \psi^{j_f}(\mathbf{r}_{A2}, \sigma_2)]_0^{0*} [\psi^{j_f}(\mathbf{r}_{A2}, \sigma_2) \psi^{j_{i1}}(\mathbf{r}_{b1}, \sigma_1)]_M^K \\ &= - \sqrt{\frac{2}{2j_f+1}} ((l_f \frac{1}{2})_{j_f} (l_{i1} \frac{1}{2})_{j_{i1}} | (l_f l_{i1})_K (\frac{1}{2} \frac{1}{2})_0)_K [Y^{l_f}(\hat{\mathbf{r}}_{A1}) Y^{l_{i1}}(\hat{\mathbf{r}}_{b1})]_M^K u_{l_f}(r_{A1}) u_{l_{i1}}(r_{b1}). \end{aligned} \quad (6.2.166)$$

Let us examine the term

$$\sum_M (-1)^M \langle K \Lambda - M \mu | P \mu - M \rangle [Y^{l_f}(\hat{\mathbf{r}}_{A1}) Y^{l_{i1}}(\hat{\mathbf{r}}_{b1})]_M^K [Y^{l_f}(\hat{\mathbf{r}}'_{A2}) Y^{l_{i2}}(\hat{\mathbf{r}}'_{b2})]_{\mu-M}^P. \quad (6.2.167)$$

Making use of the relation

$$\langle l_1 l_2 m_1 m_2 | L M_L \rangle = (-1)^{l_2-m_2} \sqrt{\frac{2L+1}{2l_1+1}} \langle L l_2 - M_L m_2 | l_1 - m_1 \rangle, \quad (6.2.168)$$

the expression (6.2.168) is equivalent to,

$$(-1)^K \sqrt{\frac{2P+1}{2\Lambda+1}} \left\{ \left[Y^{l_f}(\hat{\mathbf{r}}'_{A2}) Y^{l_{i2}}(\hat{\mathbf{r}}'_{b2}) \right]^P \left[Y^{l_f}(\hat{\mathbf{r}}_{A1}) Y^{l_{i1}}(\hat{\mathbf{r}}_{b1}) \right]^K \right\}_\mu^\Lambda. \quad (6.2.169)$$

We now recouple the term

$$\left[Y^{l_a}(\hat{\mathbf{r}}'_{aA}) Y^{l_a}(\hat{\mathbf{k}}_{aA}) \right]_0^0 \left[Y^{l_b}(\hat{\mathbf{r}}_{bB}) Y^{l_b}(\hat{\mathbf{k}}_{bB}) \right]_0^0, \quad (6.2.170)$$

arising from the partial wave expansion of the incoming and outgoing distorted waves to have,

$$((l_a l_a)_0 (l_b l_b)_0 | (l_a l_b)_\Lambda (l_a l_b)_\Lambda)_0 \left\{ \left[Y^{l_a}(\hat{\mathbf{r}}'_{aA}) Y^{l_b}(\hat{\mathbf{r}}_{bB}) \right]^\Lambda \left[Y^{l_a}(\hat{\mathbf{k}}_{aA}) Y^{l_b}(\hat{\mathbf{k}}_{bB}) \right]_0^\Lambda \right\}_0^0. \quad (6.2.171)$$

The only term which does not vanish upon integration is

$$\frac{(-1)^{\Lambda-\mu}}{\sqrt{(2l_a+1)(2l_b+1)}} \left[Y^{l_a}(\hat{\mathbf{r}}'_{aA}) Y^{l_b}(\hat{\mathbf{r}}_{bB}) \right]_{-\mu}^\Lambda \left[Y^{l_a}(\hat{\mathbf{k}}_{aA}) Y^{l_b}(\hat{\mathbf{k}}_{bB}) \right]_\mu^\Lambda. \quad (6.2.172)$$

Again, the only term surviving

$$\left\{ \left[Y^{l_f}(\hat{\mathbf{r}}'_{A2}) Y^{l_{i2}}(\hat{\mathbf{r}}'_{b2}) \right]^P \left[Y^{l_f}(\hat{\mathbf{r}}_{A1}) Y^{l_{i1}}(\hat{\mathbf{r}}_{b1}) \right]^K \right\}_\mu^\Lambda \left[Y^{l_a}(\hat{\mathbf{r}}'_{aA}) Y^{l_b}(\hat{\mathbf{r}}_{bB}) \right]_{-\mu}^\Lambda \quad (6.2.173)$$

is

$$\begin{aligned} & \frac{(-1)^{\Lambda+\mu}}{\sqrt{2\Lambda+1}} \left[\left\{ \left[Y^{l_f}(\hat{\mathbf{r}}'_{A2}) Y^{l_{i2}}(\hat{\mathbf{r}}'_{b2}) \right]^P \right. \right. \\ & \left. \left. \left[Y^{l_f}(\hat{\mathbf{r}}_{A1}) Y^{l_{i1}}(\hat{\mathbf{r}}_{b1}) \right]^K \right\}^\Lambda \left[Y^{l_a}(\hat{\mathbf{r}}'_{aA}) Y^{l_b}(\hat{\mathbf{r}}_{bB}) \right]_0^\Lambda \right]. \end{aligned} \quad (6.2.174)$$

We now couple this last term with the term $\left[Y^{l_c}(\hat{\mathbf{r}}'_{cC}) Y^{l_c}(\hat{\mathbf{r}}_{cC}) \right]_0^0$, arising from the partial wave expansion of the Green function. That is,

$$\begin{aligned}
& \left\{ \left[Y^{l_f}(\hat{\mathbf{r}}'_{A2}) Y^{l_{i2}}(\hat{\mathbf{r}}'_{b2}) \right]^P \left[Y^{l_f}(\hat{\mathbf{r}}_{A1}) Y^{l_{i1}}(\hat{\mathbf{r}}_{b1}) \right]^K \right\}^\Lambda \left[Y^{l_a}(\hat{\mathbf{r}}'_{aA}) Y^{l_b}(\hat{\mathbf{r}}_{bB}) \right]^\Lambda \Big|_0^0 \left[Y^{l_c}(\hat{\mathbf{r}}'_{cC}) Y^{l_c}(\hat{\mathbf{r}}_{cC}) \right]_0^0 \\
& = ((l_a l_b)_\Lambda (l_c l_c)_0 | (l_a l_c)_P (l_b l_c)_K)_\Lambda \left\{ \left[Y^{l_f}(\hat{\mathbf{r}}'_{A2}) Y^{l_{i2}}(\hat{\mathbf{r}}'_{b2}) \right]^P \left[Y^{l_f}(\hat{\mathbf{r}}_{A1}) Y^{l_{i1}}(\hat{\mathbf{r}}_{b1}) \right]^K \right\}^\Lambda \\
& \left\{ \left[Y^{l_a}(\hat{\mathbf{r}}'_{aA}) Y^{l_c}(\hat{\mathbf{r}}'_{cC}) \right]^P \left[Y^{l_b}(\hat{\mathbf{r}}_{bB}) Y^{l_c}(\hat{\mathbf{r}}_{cC}) \right]^K \right\}^\Lambda \Big|_0^0 = ((l_a l_b)_\Lambda (l_c l_c)_0 | (l_a l_c)_P (l_b l_c)_K)_\Lambda \\
& \times ((PK)_\Lambda (PK)_\Lambda | (PP)_0 (KK)_0)_0 \left\{ \left[Y^{l_f}(\hat{\mathbf{r}}'_{A2}) Y^{l_{i2}}(\hat{\mathbf{r}}'_{b2}) \right]^P \left[Y^{l_a}(\hat{\mathbf{r}}'_{aA}) Y^{l_c}(\hat{\mathbf{r}}'_{cC}) \right]^P \right\}_0^0 \\
& \times \left\{ \left[Y^{l_f}(\hat{\mathbf{r}}_{A1}) Y^{l_{i1}}(\hat{\mathbf{r}}_{b1}) \right]^K \left[Y^{l_b}(\hat{\mathbf{r}}_{bB}) Y^{l_c}(\hat{\mathbf{r}}_{cC}) \right]^K \right\}_0^0 = ((l_a l_b)_\Lambda (l_c l_c)_0 | (l_a l_c)_P (l_b l_c)_K)_\Lambda \\
& \times \sqrt{\frac{2\Lambda + 1}{(2K + 1)(2P + 1)}} \left\{ \left[Y^{l_f}(\hat{\mathbf{r}}'_{A2}) Y^{l_{i2}}(\hat{\mathbf{r}}'_{b2}) \right]^P \left[Y^{l_a}(\hat{\mathbf{r}}'_{aA}) Y^{l_c}(\hat{\mathbf{r}}'_{cC}) \right]^P \right\}_0^0 \\
& \times \left\{ \left[Y^{l_f}(\hat{\mathbf{r}}_{A1}) Y^{l_{i1}}(\hat{\mathbf{r}}_{b1}) \right]^K \left[Y^{l_b}(\hat{\mathbf{r}}_{bB}) Y^{l_c}(\hat{\mathbf{r}}_{cC}) \right]^K \right\}_0^0. \tag{6.2.175}
\end{aligned}$$

Collecting all the contributions (including the constants and phases arising from the partial wave expansion of the distorted waves and the Green function), we get

$$\begin{aligned}
T_{succ}^{(2)}(\theta; \mu) &= (-1)^{j_f + j_{i1}} \frac{2048\pi^5 \mu_{Cc}}{\hbar^2 k_{Aa} k_{Bb} k_{Cc}} \sqrt{\frac{(2j_{i1} + 1)}{(2\Lambda + 1)(2j_f + 1)}} \sum_{K,P} ((l_f \frac{1}{2})_{j_f} (l_i \frac{1}{2})_{j_{i1}} | (l_f l_i)_P (\frac{1}{2} \frac{1}{2})_0)_P \\
&\quad \times ((l_f \frac{1}{2})_{j_f} (l_i \frac{1}{2})_{j_{i1}} | (l_f l_i)_K (\frac{1}{2} \frac{1}{2})_0)_K ((j_{i1} j_f)_K (j_{i1} j_{i2})_0 | (j_{i1} j_{i1})_0 (j_f j_{i2})_P)_P \\
&\quad \times \frac{(-1)^K}{(2K + 1) \sqrt{2P + 1}} \sum_{l_c, l_a, l_b} ((l_a l_b)_\Lambda (l_c l_c)_0 | (l_a l_c)_P (l_b l_c)_K)_\Lambda e^{i(\sigma_i^{l_a} + \sigma_f^{l_b})} i^{l_a - l_b} \\
&\quad \times (2l_c + 1)^{3/2} \left[Y^{l_a}(\hat{\mathbf{k}}_{aA}) Y^{l_b}(\hat{\mathbf{k}}_{bB}) \right]_\mu^K S_{K,P,l_a,l_b,l_c}, \tag{6.2.176}
\end{aligned}$$

with (note that we have reduced the dimensionality of the integrals in the same fashion as for the $L=0$ -angular momentum transfer calculation, see (6.2.132))

$$\begin{aligned}
S_{K,P,l_a,l_b,l_c} &= \int r_{Cc}^2 dr_{Cc} r_{b1}^2 dr_{b1} \sin \theta d\theta v(r_{b1}) u_{l_f}(r_{C1}) u_{l_i}(r_{b1}) \\
&\quad \times \frac{s_{P,l_a,l_c}(r_{Cc})}{r_{Cc}} \frac{F_{l_b}(r_{Bb})}{r_{Bb}} \\
&\quad \times \sum_M \langle l_c 0 l_b M | K M \rangle \left[Y^{l_f}(\hat{r}_{C1}) Y^{l_{i1}}(\theta + \pi, 0) \right]_M^K Y^{l_b}_{-M}(\hat{r}_{Bb}), \tag{6.2.177}
\end{aligned}$$

and

$$\begin{aligned} s_{P,l_a,l_c}(r_{Cc}) &= \int r'_{Cc} dr'_{Cc} r'_{A2} dr'_{A2} \sin \theta' d\theta' v(r'_{c2}) u_{l_f}(r'_{A2}) u_{l_i}(r'_{c2}) \\ &\times \frac{F_{l_a}(r'_{Aa})}{r'_{Aa}} \frac{f_{l_c}(k_{Cc}, r_<) P_{l_c}(k_{Cc}, r_>)}{r'_{Cc}} \\ &\times \sum_M \langle l_c 0 l_a M | P M \rangle \left[Y^{l_f}(\hat{r}'_{A2}) Y^{l_i}(r'_{c2}) \right]_M^P Y_{-M}^{l_a}(\hat{r}'_{Aa}). \end{aligned} \quad (6.2.178)$$

We have evaluated the transition matrix element for a particular projection μ of the initial angular momentum of the two transferred nucleons. If they are coupled to a core of angular momentum J_f to total angular momentum J_i, M_i , the fraction of the initial wavefunction with projection μ is $\langle \Lambda \mu J_f M_i - \mu | J_i M_i \rangle$, and the cross section will be

$$\frac{d\sigma}{d\Omega}(\hat{\mathbf{k}}_{bB}) = \frac{k_{bB}}{k_{aA}} \frac{\mu_{aA}\mu_{bB}}{(2\pi\hbar^2)^2} \left| \sum_\mu \langle \Lambda \mu J_f M_i - \mu | J_i M_i \rangle T_{succ}^{(2)}(\theta; \mu) \right|^2. \quad (6.2.179)$$

For a non polarized incident beam,

$$\frac{d\sigma}{d\Omega}(\hat{\mathbf{k}}_{bB}) = \frac{k_{bB}}{k_{aA}} \frac{\mu_{aA}\mu_{bB}}{(2\pi\hbar^2)^2} \frac{1}{2J_i + 1} \sum_{M_i} \left| \sum_\mu \langle \Lambda \mu J_f M_i - \mu | J_i M_i \rangle T_{succ}^{(2)}(\theta; \mu) \right|^2. \quad (6.2.180)$$

This would be the differential cross section for a transition to a definite final state M_f . If we do not measure M_f we have to sum for all M_f ,

$$\frac{d\sigma}{d\Omega}(\hat{\mathbf{k}}_{bB}) = \frac{k_{bB}}{k_{aA}} \frac{\mu_{aA}\mu_{bB}}{(2\pi\hbar^2)^2} \frac{1}{2J_i + 1} \sum_\mu |T_{succ}^{(2)}(\theta; \mu)|^2 \sum_{M_i, M_f} |\langle \Lambda \mu J_f M_f | J_i M_i \rangle|^2. \quad (6.2.181)$$

The sum over M_i, M_f of the Clebsh–Gordan coefficients gives $(2J_i + 1)/(2\Lambda + 1)$ (see Eq. (6.A.26)). One then gets,

$$\frac{d\sigma}{d\Omega}(\hat{\mathbf{k}}_{bB}) = \frac{k_{bB}}{k_{aA}} \frac{\mu_{aA}\mu_{bB}}{(2\pi\hbar^2)^2} \frac{1}{(2\Lambda + 1)} \sum_\mu |T_{succ}^{(2)}(\theta; \mu)|^2, \quad (6.2.182)$$

where one can write

$$\begin{aligned} T_{succ}^{(2)}(\theta; \mu) &= \sum_{l_a, l_b} C_{l_a, l_b} \left[Y^{l_a}(\hat{\mathbf{k}}_{aA}) Y^{l_b}(\hat{\mathbf{k}}_{bB}) \right]_\mu^\Lambda \\ &= \sum_{l_a, l_b} C_{l_a, l_b} i^{l_a} \sqrt{\frac{2l_a + 1}{4\pi}} \langle l_a l_b 0 \mu | \Lambda \mu \rangle Y_\mu^{l_b}(\hat{\mathbf{k}}_{bB}). \end{aligned} \quad (6.2.183)$$

Note that (6.2.182) takes into account only the spins of the heavy nucleus. In a (t, p) or (p, t) reaction, we have to sum over the spins of the proton and of the triton

and divide by 2. If a spin-orbit term is present in the optical potential, the sum yields the combination of terms shown in Section (6.2.2),

$$\frac{d\sigma}{d\Omega}(\hat{\mathbf{k}}_{bB}) = \frac{k_{bB}}{k_{aA}} \frac{\mu_{aA}\mu_{bB}}{(2\pi\hbar^2)^2} \frac{1}{2(2\Lambda+1)} \sum_{\mu} |A_{\mu}|^2 + |B_{\mu}|^2. \quad (6.2.184)$$

6.3 ZPF and Pauli principle at the basis of medium polarization effects: self-energy, vertex corrections and induced interaction

Of all quantal phenomena, zero point fluctuations (ZPF), closely connected with virtual states, are likely the most representative of the essential difference existing between quantum and classical mechanics. In fact, ZPF are intimately connected with the complementary principle⁹, and thus with indeterminacy¹⁰ and non-commutative¹¹ relations, and with the probabilistic interpretation¹² of the (modulus squared) of the wavefunctions, solution of Schrödinger's or Dirac's equations¹³. Pauli principle¹⁴ brings about essential modifications of the virtual fluctuations of the many-body system, modifications which are instrumental in the dressing and interweaving of the elementary modes of excitation¹⁵.

In Fig. 6.3.1, NFT diagrams are given which correspond to the lowest order medium polarization effects renormalizing the properties of a particle-hole collective mode (wavy line), correlated particle-hole excitation ((up-going)-(down-going) arrowed lines) calculated within the random phase approximation (RPA,QRPA), and leading to the particle-vibration coupling vertex (formfactor and strength, i.e. transition density (solid dot), see inset (I), bottom). The action of an external field on the zero point fluctuations (ZPF) of the vacuum (inset (II)), forces a virtual process to become real, leading to a collective vibration by annihilating a (virtual, spontaneous) particle-hole excitation (backwards RPA Y -amplitude) or, in the time ordered process, by creating a particle-hole excitation which eventually, through the particle-vibration coupling vertex, materializes into the collective (coherent) state (forwardsgoing X -amplitudes). Now, oyster-like diagrams associated with the vacuum ZPF can occur at any time (see inset (III) of Fig. 6.3.1). It is of notice that while virtual states can violate energy conservation, they have to respect conservation rules, as well as Pauli principle. For example, a virtual state cannot violate angular momentum neither allow for the presence of two fermions in the same quantal state. The process shown in the inset III (α) leads, through Pauli principle correcting diagrams (exchange of fermionic arrowed lines) to self-energy

⁹Bohr (1928).

¹⁰Heisenberg (1927).

¹¹Born and Jordan (1925), Born et al. (1926).

¹²Born (1926a).

¹³Schrödinger (1926a), Dirac (1930).

¹⁴Pauli (1925).

¹⁵Within the present context, see also Schrieffer (1964).

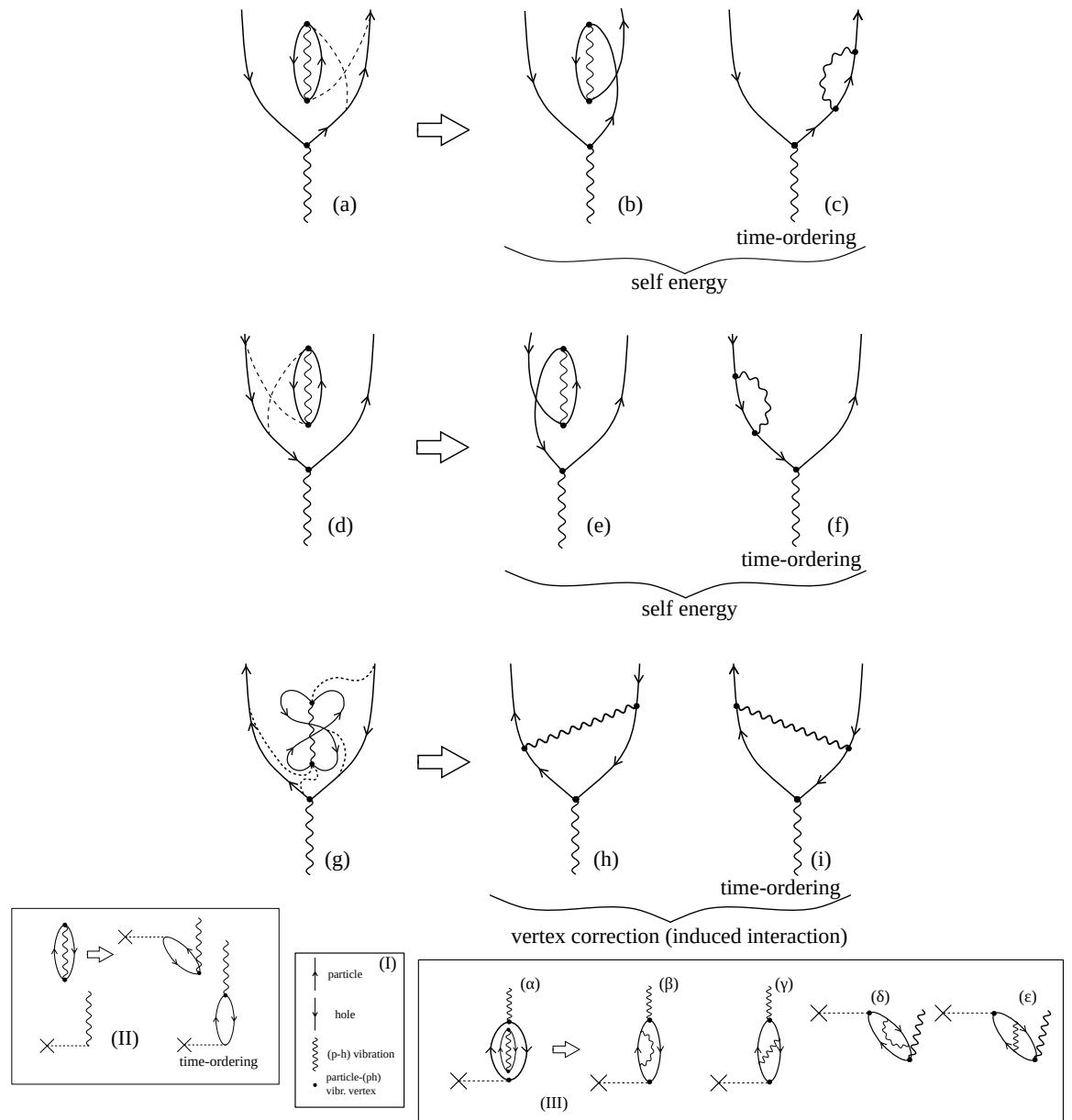


Figure 6.3.1: Nuclear field theory (NFT) diagrams describing renormalization processes associated with ZPF. For details see caption to Fig. 6.3.2.

(inset III (β), (δ)) and vertex corrections (induced $p - h$ interaction; inset III (γ), (ε), see also graphs (c), (f), (h) and (i) of Fig. 6.3.1) processes. Similar processes are found in the interplay between ZPF and pair addition modes as shown in Fig. 6.3.2. Note the parallel between diagrams of Figs. 6.3.1 (g)–(i) and of Figs. 6.3.2 (d)–(f).

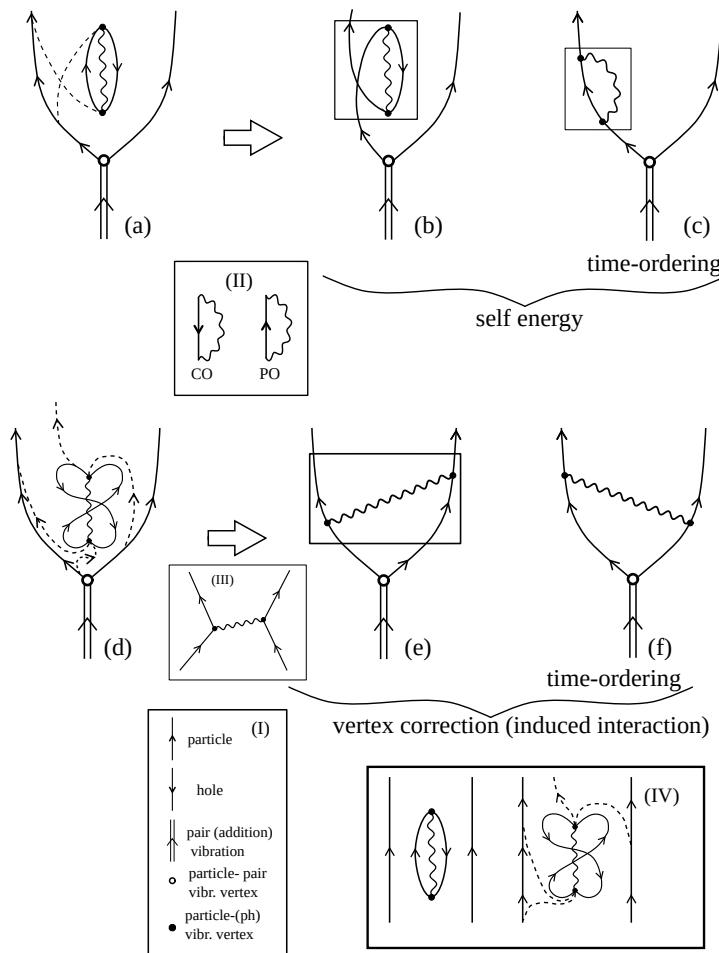


Figure 6.3.2: Pauli effects associated (p-h) ZPF dressing a pairing vibrational (pair addition) mode (see inset I) in terms of self-energy (graphs (a)–(c); correlation (CO) and polarization (PO) diagrams, inset II) and vertex correction (graphs (d)–(f); induced particle-particle (pairing) interaction processes (inset (III)), associated with phonon exchange between nucleons (inset (IV)).

The collective vibrational modes can be viewed as coherent states¹⁶ exhausting a consistent fraction of the EWSR.

6.4 Coherence and effective formfactors

In what follows we shall work out a simplified derivation of the simultaneous two-nucleon transfer amplitude, within the framework of first order DWBA specially

¹⁶See e.g. Glauber (1969, 2007).

suites to discuss correlation aspects of pair transfer in general, and of the associated effective formfactors in particular¹⁷.

We will concentrate on (t, p) reaction, namely reactions of the type $A(\alpha, \beta)B$ where $\alpha = \beta + 2$ and $B = A + 2$. The intrinsic wave functions are in this case

$$\begin{aligned}\psi_\alpha &= \psi_{M_i}^{J_i}(\xi_A) \sum_{ss'_f} \left[\chi^s(\sigma_\alpha) \chi^{s'_f}(\sigma_\beta) \right]_{M_{s_i}}^{s_i} \phi_t^{L=0}(r_{12}, r_{1p}, r_{2p}) \\ &= \psi_{M_i}^{J_i}(\xi_A) \sum_{M_s M'_{s_f}} (s M'_s s'_f M'_{s_f} | s_i M_{s_i}) \chi_{M'_s}^s(\sigma_\alpha) \chi_{M'_{s_f}}^{s'_f}(\sigma_\beta) \\ &\quad \times \phi_t^{L=0}(r_{12}, r_{1p}, r_{2p})\end{aligned}\quad (6.4.1)$$

while

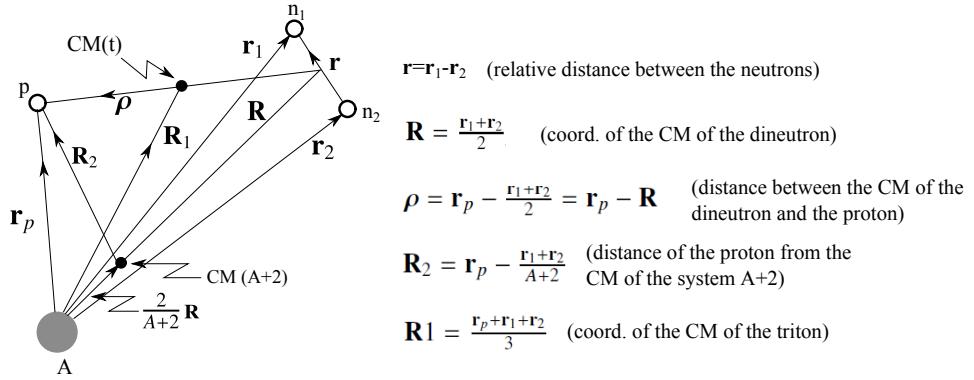


Figure 6.4.1: Coordinate system used in the calculation of the two–nucleon transfer amplitude.

$$\begin{aligned}\psi_\beta &= \psi_{M_f}^{J_f}(\xi_{A+2}) \chi_{M_{s_f}}^{s_f}(\sigma_\beta) \\ &= \sum_{\substack{n_1 l_1 j_1 \\ n_2 l_2 j_2}} B(n_1 l_1 j_1, n_2 l_2 j_2; J J'_i J_f) \left[\phi^J(j_1 j_2) \phi^{J'_i}(\xi_A) \right]_{M_f}^{J_f} \\ &\quad \times \chi_{M_{s_f}}^{s_f}(\sigma_\beta).\end{aligned}\quad (6.4.2)$$

Making use of the above equation one can define the spectroscopic amplitude B as

$$\begin{aligned}B(n_1 l_1 j_1, n_2 l_2 j_2; J J'_i J_f) \\ = \left\langle \psi^{J_f}(\xi_{A+2}) \left[\phi^J(j_1 j_2) \phi^{J_i}(\xi_A) \right]^{J_f} \right\rangle,\end{aligned}\quad (6.4.3)$$

¹⁷Glendenning (1965); Bayman and Kallio (1967).

where

$$\phi^J(j_1 j_2) = \frac{[\phi_{j_1}(\mathbf{r}_1)\phi_{j_2}(\mathbf{r}_2)]^J - [\phi_{j_1}(\mathbf{r}_2)\phi_{j_2}(\mathbf{r}_1)]^J}{\sqrt{1 + \delta(j_1, j_2)}}, \quad (6.4.4)$$

is an antisymmetrized, normalized wave function of the two transferred particles. The function $\chi_{M_s}^s(\sigma_\beta)$ appearing both in eq. (6.4.1) and (6.4.2) is the spin wave function of the proton while

$$\chi^s(\sigma_\alpha) = [\chi^{s_1}(\sigma_{n_1})\chi^{s_2}(\sigma_{n_2})]^s, \quad (6.4.5)$$

is the spin function of the two-neutron system.

A simple description of the intrinsic degrees of freedom of the triton is obtained by using a wavefunction symmetric in the coordinates of all particles, i.e.

$$\begin{aligned} \phi_t^{L=0}(r_{12}, r_{1p}, r_{2p}) &= N_t e^{[(r_1-r_2)^2 + (r_1-r_p)^2 + (r_2-r_p)^2]} \\ &= \phi_{000}(\mathbf{r})\phi_{000}(\boldsymbol{\rho}), \end{aligned} \quad (6.4.6)$$

where

$$\phi_{000}(\mathbf{r}) = R_{nl}(\nu^{1/2}r)Y_{lm}(\hat{\mathbf{r}}). \quad (6.4.7)$$

The coordinate $\boldsymbol{\rho}$ is the radius vector which measures the distance between the center of mass of the dineutron and the proton, while the vector \mathbf{r} is the dineutron relative coordinate (cf. Fig. 6.4.1). To obtain the DWBA cross section we have to calculate the integral

$$T(\theta) = \int d\xi_A d\mathbf{r}_1 d\mathbf{r}_2 d\mathbf{r}_p \chi_p^{(-)}(\mathbf{R}_2) \psi_\beta^*(\xi_{A+2}, \sigma_\beta) V_\beta \psi_\alpha(\xi_A, \sigma_\alpha, \sigma_\beta) \psi_t^{(+)}(\mathbf{R}_1), \quad (6.4.8)$$

where the final state effective interaction $V_\beta(\rho)$ is assumed to depend only on the distance ρ between the center of mass of the di-neutron and of the proton.

To carry out the integral (6.4.8) we transform the wave function (6.4.4) into center of mass and relative coordinates. If we assume that both $\phi_{j_1}(\mathbf{r}_1)$ and $\phi_{j_2}(\mathbf{r}_2)$ are harmonic oscillator wave functions (used as a basis to expand the Saxon-Woods single-particle wavefunctions), this transformation can be carried with the aid of the Moshinsky brackets. If $|n_1 l_1, n_2 l_2; \lambda \mu\rangle$ is a complete system of wave functions in the harmonic oscillator basis, depending on \mathbf{r}_1 and \mathbf{r}_2 and $|nl, NL; \lambda \mu\rangle$ is the corresponding one depending on \mathbf{r} and \mathbf{R} , we can write

$$\begin{aligned} |n_1 l_1, n_2 l_2; \lambda \mu\rangle &= \sum_{nlNL} (|nl, NL; \lambda \mu\rangle \langle nl, NL; \lambda \mu|) |n_1 l_1, n_2 l_2; \lambda \mu\rangle \\ &= \sum_{nlNL} |nl, NL; \lambda \mu\rangle \langle nl, NL; \lambda \mu| n_1 l_1, n_2 l_2; \lambda \rangle. \end{aligned} \quad (6.4.9)$$

The labels n, l are the principal and angular momentum quantum numbers of the relative motion, while N, L are the corresponding ones corresponding to the center

of mass motion of the two-neutron system. Because of energy and parity conservation we have

$$\begin{aligned} 2n_1 + l_1 + 2n_2 + l_2 &= 2n + l + 2N + L, \\ (-1)^{l_1+l_2} &= (-1)^{l+L}. \end{aligned} \quad (6.4.10)$$

The coefficients $\langle nl, NL, L | n_1 l_1, n_2 l_2, L \rangle$ were first discussed by Moshinsky¹⁸.

With the help of eq. (6.4.9) we can write the wave function $\psi_{M_f}^{J_f}(\xi_{A+2})$ as

$$\begin{aligned} \psi_{M_f}^{J_f}(\xi_{A+2}) &= \sum_{\substack{n_1 l_1 j_1 \\ n_2 l_2 j_2 \\ JJ_i}} B(n_1 l_1 j_1, n_2 l_2 j_2; JJ'_f J_f) [\phi^J(j_1 j_2) \phi^{J'_i}(\xi_A)]_{M_f}^{J_f} \\ &= \sum_{\substack{n_1 l_1 j_1 \\ n_2 l_2 j_2 \\ JJ_i}} \sum_{JJ'_i} B(n_1 l_1 j_1, n_2 l_2 j_2; JJ'_i J_f) \\ &\quad \times \sum_{M_J M'_{J_i}} \langle JM_J J'_i M_{J_i} | J_f M_{J_f} \rangle \psi_{M'_{J_i}}^{J'_i}(\xi_A) \\ &\quad \times \sum_{LS'} \langle S' LJ | j_1 j_2 J \rangle \sum_{M_L M'_S} \langle LM_L S' M'_S | JM_J \rangle \chi_{M'_S}^{S'}(\sigma_\alpha) \\ &\quad \times \sum_{nLN\Lambda} \langle nl, N\Lambda, L | n_1 l_1, n_2 l_2, L \rangle \\ &\quad \times \sum_{m_l M_\Lambda} \langle lm_l \Lambda M_\Lambda | LM_L \rangle \phi_{nlm_l}(\mathbf{r}) \phi_{N\Lambda M_\Lambda}(\mathbf{R}), \end{aligned} \quad (6.4.11)$$

Integration over \vec{r} gives

$$\langle \phi_{nlm_l}(\mathbf{r}) | \phi_{000}(\mathbf{r}) \rangle = \delta(l, 0) \delta(m_l, 0) \Omega_n, \quad (6.4.12)$$

where

$$\Omega_n = \int R_{nl}(\nu_1^{1/2} r) R_{00}(\nu_2^{1/2} r) r^2 dr. \quad (6.4.13)$$

Note that there is no selection rule in the principal quantum number n , as the potential in which the two neutrons move in the triton has a frequency ν_2 which is different from the one that the two neutrons are subjected to, when moving in the system A (non-orthogonality effect).

Integration over ξ_A and multiplication of the spin functions gives

$$\begin{aligned} (\psi_{M_{J_i}}^{J_i}, V'_\beta(\rho) \psi_{M'_{J_i}}^{J'_i}) &= \delta(J_i, J'_i) \delta(M_{J_i}, M_{J'_i}) V(\rho), \\ (\chi_{M_S}^S(\sigma_\alpha), \chi_{M'_S}^{S'}(\sigma_\alpha)) &= \delta(S, S') \delta(M_S, M_{S'}), \\ (\chi_{M_{S_f}}^{S_f}(\sigma_\beta), \chi_{M'_{S_f}}^{S'_f}(\sigma_\beta)) &= \delta(S_f, S'_f) \delta(M_{S_f}, M_{S'_f}). \end{aligned} \quad (6.4.14)$$

¹⁸Moshinsky (1959).

The integral (6.4.8) can then be written as

$$\begin{aligned} T(\theta) = & \sum_{\substack{n_1 l_1 j_1 \\ n_2 l_2 j_2}} \sum_{JM_J} \sum_{nN} \sum_S B(n_1 l_1 j_1, n_2 l_2 j_2; JJ_i J_f) \\ & \times \langle JM_J J_i M_{J_i} | J_f M_{J_f} \rangle \langle S L J | j_1 j_2 J \rangle \\ & \times \langle LM_L S M_S | JM_J \rangle \langle n0, NL, L | n_1 l_1, n_2 l_2, L \rangle \\ & \times \langle S M_S S_f M_{S_f} | S_i M_{S_i} \rangle \Omega_n \\ & \times \int d\mathbf{R} d\mathbf{r}_p \chi_t^{(+)*}(\mathbf{R}_1) \phi_{NLM_L}^*(\mathbf{R}) V(\rho) \phi_{000}(\boldsymbol{\rho}) \chi_t^{(+)}(\mathbf{R}_1), \end{aligned} \quad (6.4.15)$$

where we have approximated V'_β by an effective interaction depending on $\rho = |\boldsymbol{\rho}|$. We now define the effective two-nucleon transfer form factor as

$$\begin{aligned} u_{LSJ}^{J_i J_f}(R) = & \sum_{\substack{n_1 l_1 j_1 \\ n_2 l_2 j_2}} \sum_{nN} B(n_1 l_1 j_1, n_2 l_2 j_2; JJ_i J_f) \langle S L J | j_1 j_2 J \rangle \\ & \langle n0, NL, L | n_1 l_1, n_2 l_2; L \rangle \Omega_n R_{NL}(R). \end{aligned} \quad (6.4.16)$$

One can then rewrite Eq. (6.4.15) as

$$\begin{aligned} T(\theta) = & \sum_J \sum_L \sum_S (JM_J J_i M_{J_i} | J_f M_{J_f} \rangle) (S M_S S_f M_{S_f} | S_i M_{S_i} \rangle) (LM_L S M_S | JM_J \rangle) \\ & \times \int d\mathbf{R} d\mathbf{r}_p \chi_p^{*(-)}(\mathbf{R}_2) u_{LSJ}^{J_i J_f}(R) Y_{LM_L}^* V(\rho) \phi_{000}(\boldsymbol{\rho}) \chi_t^{(+)}(\mathbf{R}_1). \end{aligned} \quad (6.4.17)$$

Because the di-neutron has $S = 0$, we have that

$$(LM_L 00 | JM_J) = \delta(J, L) \delta(M_L, M_J), \quad (6.4.18)$$

and the summations over S and L disappear from Eq. (6.4.17). Let us now make also here, as done in Sect. 5.6, Eq. (5.6.15) for one-particle transfer reactions, the zero range approximation, that is,

$$V(\rho) \phi_{000}(\boldsymbol{\rho}) = D_0 \delta(\boldsymbol{\rho}), \quad (6.4.19)$$

where D_0 is an empirical parameter ($D_0^2 = (31.6 \pm 9.3) \times 10^4$ MeV 2 fm 2) determined to reproduce, in average, the observed absolute cross sections¹⁹ with only the simultaneous transfer. This means that the proton interacts with the center of mass of the di-neutron, only when they are at the same point in space. Within this approximation (cf. Fig. 6.4.1)

$$\begin{aligned} \mathbf{R} = & \mathbf{R}_1 = \mathbf{r}, \\ \mathbf{R}_2 = & \frac{A}{A+2} \mathbf{R}. \end{aligned} \quad (6.4.20)$$

¹⁹Broglio et al. (1973).

Then Eq. (6.4.15) can be written as

$$T = D_0 \sum_L (LM_L J_i M_{J_i} | J_f M_{J_f}) \\ \times \int d\mathbf{R} \chi_p^{*(-)} \left(\frac{A}{A+2} \mathbf{R} \right) u_L^{J_i J_f}(R) Y_{LM_L}^*(\hat{\mathbf{R}}) \chi_t^{(+)}(\mathbf{R}) \quad (6.4.21)$$

From Eq. (6.4.21) it is seen that the change in parity implied by the reaction is given by $\Delta\pi = (-1)^L$. Consequently, the selection rules for (t, p) and (p, t) reactions in zero-range approximation are,

$$\begin{aligned} \Delta S &= 0 \\ \Delta J = \Delta L &= L \\ \Delta\pi &= (-1)^L \end{aligned} \quad (6.4.22)$$

i.e. only normal parity states are excited. The integral appearing in Eq. (6.4.21) has the same structure as the DWBA integral appearing in Eq. (5.6.16) which was derived for the case of one-nucleon transfer reactions.

The difference between the two processes manifests itself through the different structure of the two form factors. While $u_l(r)$ is a single-particle bound state wave function (cf. Eq. (5.6.1a)), $u_L^{J_i J_f}$ is a coherent summation *over the center of mass states of motion of the two transferred neutrons* (see Eq. (6.4.16)). In other words, an effective quantity (function). It is of notice that this difference between single-particle transfer formfactors and simultaneous two-nucleon transfer effective formfactors, essentially vanishes when one considers dressed particles resulting from the coupling to collective vibrations and leading to renormalized energies, single-particle content, and radial wavefunction. In view of the relation (4.3.1), closely connected with Fig. 4.3.1 (a) found at the basis of the fact that successive transfer of entangled nucleons over distances of the order of the correlation length, the above “historic” result²⁰ is certainly dated. It is also found at the basis of a series of misunderstandings in working out the variety of potentialities concerning the parallel between BCS²¹ and nuclear BCS²², let alone regarding Cooper pair tunneling across a Josephson junction, and two-nucleon transfer in heavy ion collisions between superfluid nuclei at energies below the Coulomb barrier. The reason to bring it here is to set in evidence subtle aspects of the correlations existing between partner nucleons of a Cooper pair to be also found in the formfactors associated with successive transfer although, arguably, in a less evident fashion.

Examples of two-nucleon transfer form factors are given in²³ Figs 6.4.2, 6.4.3 and 6.4.4.

²⁰See Potel et al. (2013a), in particular App. C and Fig. 10 of this reference.

²¹Cooper and Feldman (2011).

²²Broglia and Zelevinsky (2013).

²³Broglia and Riedel (1967).

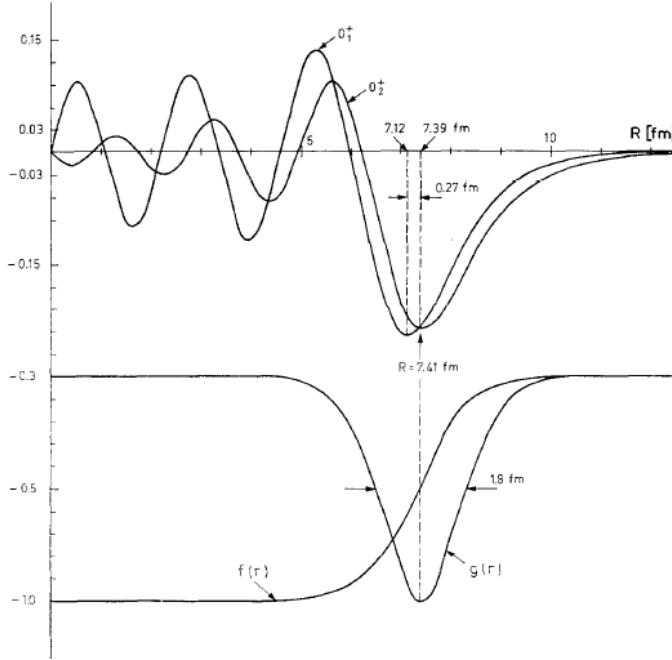


Figure 6.4.2: The upper part of the figure shows the modified formfactor for the $^{206}\text{Pb}(\text{t},\text{p})^{208}\text{Pb}$ transition to the ground state (0_1^+) and the pairing vibrational state (0_2^+) at 4.87 MeV. Both curves are matched with appropriate Hankel functions. In the lower part the form factors of the real ($f(r)$) and the imaginary ($g(r)$) part of the optical potential used to calculate the differential cross sections (see Fig. 4.5.4), are given.

6.5 Relative importance of successive and simultaneous transfer and non-orthogonality corrections

In what follows we discuss the relative importance of successive and simultaneous two-neutron transfer and of non-orthogonality corrections associated with the reaction

$$\alpha \equiv a (= b + 2) + A \rightarrow b + B (= A + 2) \equiv \beta, \quad (6.5.1)$$

in the limits of independent particles and of strongly correlated Cooper pairs, making use of the semiclassical approximation²⁴, in which case the two-particle transfer differential cross section can be written as

$$\frac{d\sigma_{\alpha \rightarrow \beta}}{d\Omega} = P_{\alpha \rightarrow \beta}(t = +\infty) \sqrt{\left(\frac{d\sigma_\alpha}{d\Omega} \right)_{el}} \sqrt{\left(\frac{d\sigma_\beta}{d\Omega} \right)_{el}}, \quad (6.5.2)$$

²⁴For details cf. Broglia and Winther (2004).

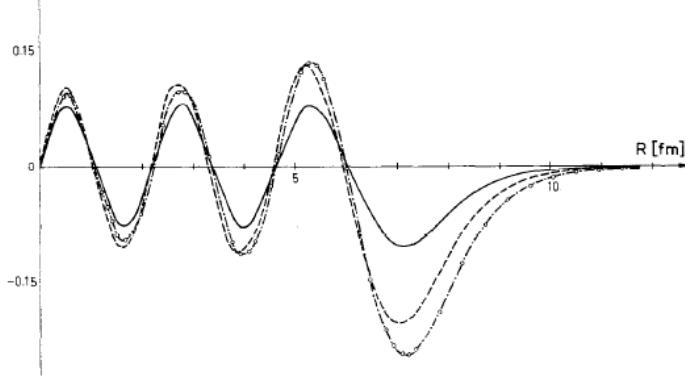


Figure 6.4.3: Modified formfactor for the transition to the ground state ($^{206}\text{Pb}(\text{t},\text{p})^{208}\text{Pb(gs)}$; see Fig. 4.5.4) calculated in different spectroscopic models (pure shell-model configuration ——, shell model plus pairing residual interaction ——, including ground state correlations -o-o-).

where P is the absolute value squared of a quantum mechanical transition amplitude. It gives the probability that the system at $t = +\infty$ is found in the final channel. The quantities $(d\sigma/d\Omega)_{el}$ are the classical elastic cross sections in the center of mass system, calculated in terms of the deflection function, namely the functional relating the impact parameter and the scattering angle.

The transfer amplitude can be written as

$$a(t = +\infty) = a^{(1)}(\infty) - a^{(NO)}(\infty) + \tilde{a}^{(2)}(\infty), \quad (6.5.3)$$

where $\tilde{a}^{(2)}(\infty)$ at $t = +\infty$ labels the successive transfer amplitude expressed in the post-prior representation (see below). The simultaneous transfer amplitude is given by (see Fig. 6.5.1 (I))

$$\begin{aligned} a^{(1)}(\infty) &= \frac{1}{i\hbar} \int_{-\infty}^{\infty} dt (\psi^b \psi^B, (V_{bB} - < V_{bB} >) \psi^a \psi^A) \times \exp\left[\frac{i}{\hbar}(E^{bB} - E^{aA})t\right] \\ &\approx \frac{2}{i\hbar} \int_{-\infty}^{\infty} dt \left(\phi^{B(A)}(S_{(2n)}^B; \mathbf{r}_{1A}, \mathbf{r}_{2A}), U(r_{1b}) e^{i(\sigma_1 + \sigma_2)} \phi^{a(b)}(S_{(2n)}^a; \mathbf{r}_{1b}, \mathbf{r}_{2b}) \right) \\ &\quad \times \exp\left[\frac{i}{\hbar}(E^{bB} - E^{aA})t + \gamma(t)\right] \end{aligned} \quad (6.5.4)$$

where

$$\sigma_1 + \sigma_2 = \frac{1}{\hbar} \frac{2m_n}{m_A} (m_{aA} \mathbf{v}_{aA}(t) + m_{bB} \mathbf{v}_{bB}(t)) \cdot (\mathbf{r}_{1\alpha} - \mathbf{r}_{2\alpha}), \quad (6.5.5)$$

$\exp(i(\sigma_1 + \sigma_2))$ taking care of recoil effects (Galilean transformation associated with the mismatch between entrance and exit channels). The phase $\gamma(t)$ is related

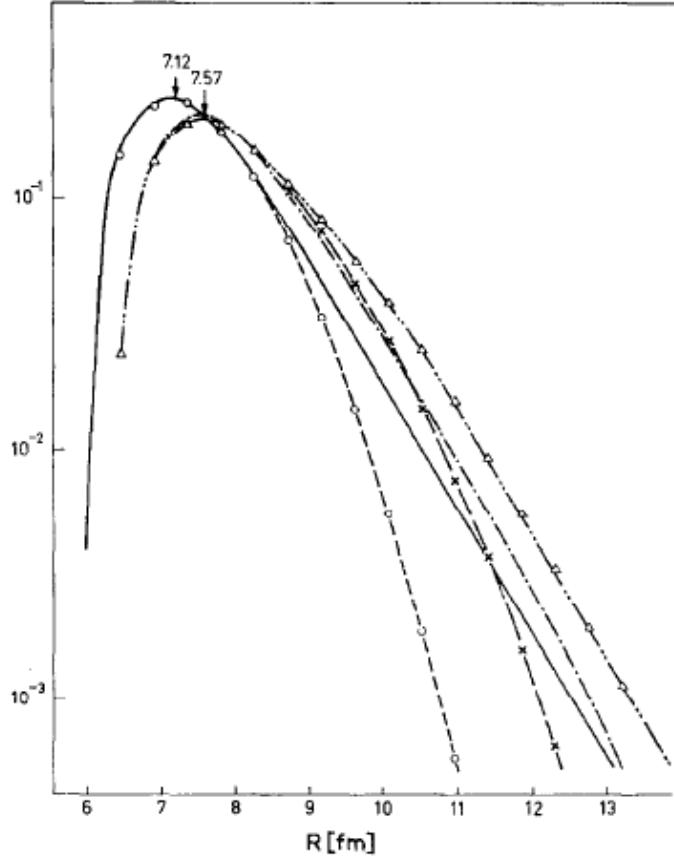


Figure 6.4.4: Asymptotic behavior of the modified formfactor for the $^{206}(\text{t},\text{p})^{208}\text{Pb}(\text{gs})$ ground state transition for oscillator plus Hankel wave functions (continuous solid curve), oscillator wave functions alone (dash point dashed curve), and Saxon-Woods wave functions with a variety of asymptotic matchings.

with the effective Q -value of the reaction. In the above expression, ϕ indicates an antisymmetrized, correlated two-particle (Cooper pair) wavefunction, $S(2n)$ being the two-neutron separation energy (see Fig. 6.5.3), $U(r_{1b})$ being the single particle potential generated by nucleus b . The contribution arising from non-orthogonality effects can be written as (see Fig. 6.5.1 (II))

$$\begin{aligned}
 a^{(NO)}(\infty) &= \frac{1}{i\hbar} \int_{-\infty}^{\infty} dt (\psi^b \psi^B, (V_{bB} - \langle V_{bB} \rangle) \psi^f \psi^F) (\psi^f \psi^F, \psi^a \psi^A) \exp\left[\frac{i}{\hbar}(E^{bB} - E^{aA})t\right] \\
 &\approx \frac{2}{i\hbar} \int_{-\infty}^{\infty} \phi^{B(F)}(S_{(n)}^B, \mathbf{r}_{1A}), U(r_{1b}) e^{i\sigma_1} (\phi^{f(b)}(S^f(n), \mathbf{r}_{1b}) \\
 &\quad \times \phi^{F(A)}(S^F(n), \mathbf{r}_{2A}) e^{i\sigma_2} \phi^{a(f)}(S^a(n), \mathbf{r}_{2b})) \exp\left[\frac{i}{\hbar}(E^{bB} - E^{aA})t + \gamma(t)\right],
 \end{aligned} \tag{6.5.6}$$

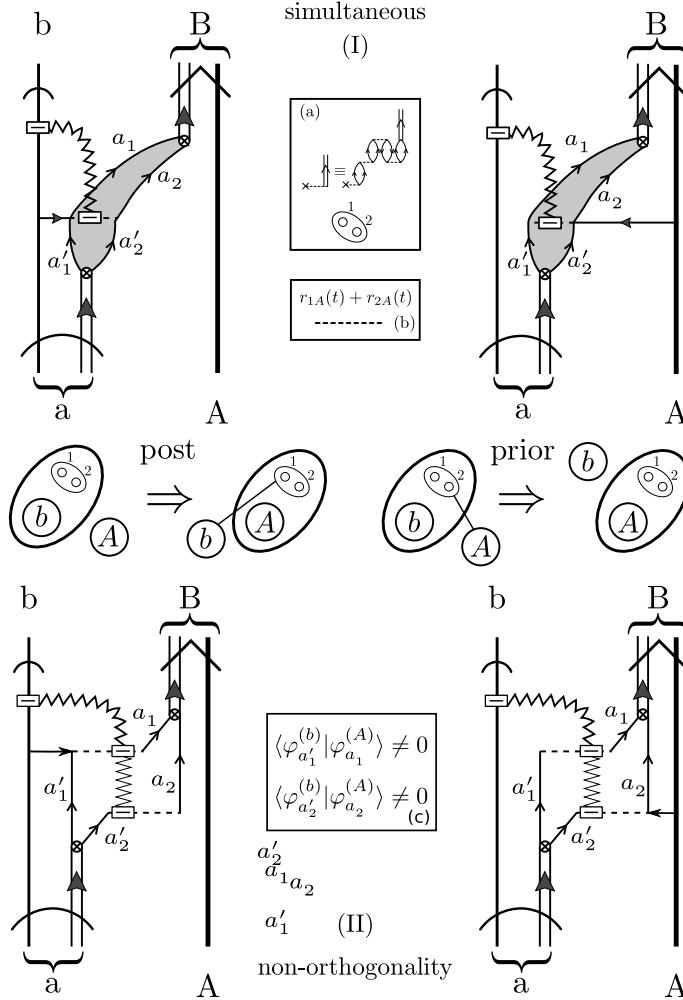


Figure 6.5.1: Graphical representation of simultaneous (I) and non-orthogonality (II) transfer processes. For details see text and caption to Fig. 6.5.2. In this figure, as well as in Fig. 6.5.2, and at variance with, e.g. Fig. 2.7.2, the particle-pair vibration coupling vertex is represented by a crossed circle. Concerning the jagged line representing recoil see App. 2.A.

the reaction channel $f = (b + 1) + F (= A + 1)$ having been introduced, the quantity $S(n)$ being the one-neutron separation energy (see Fig. 6.5.3). The summation over $f (\equiv a'_1, a'_2)$ and $F (\equiv a_1, a_2)$ involves a restricted number of states, namely the valence shells in nuclei B and a . It is of notice that the NFT_(s+r) diagrams appearing in Fig. 6.5.1 (II) are the only ones we have encountered in which to a particle-recoil coupling vertex (dashed open rectangle) and thus the starting or/and ending of a recoil mode (jagged line) does not correspond the action of the nuclear

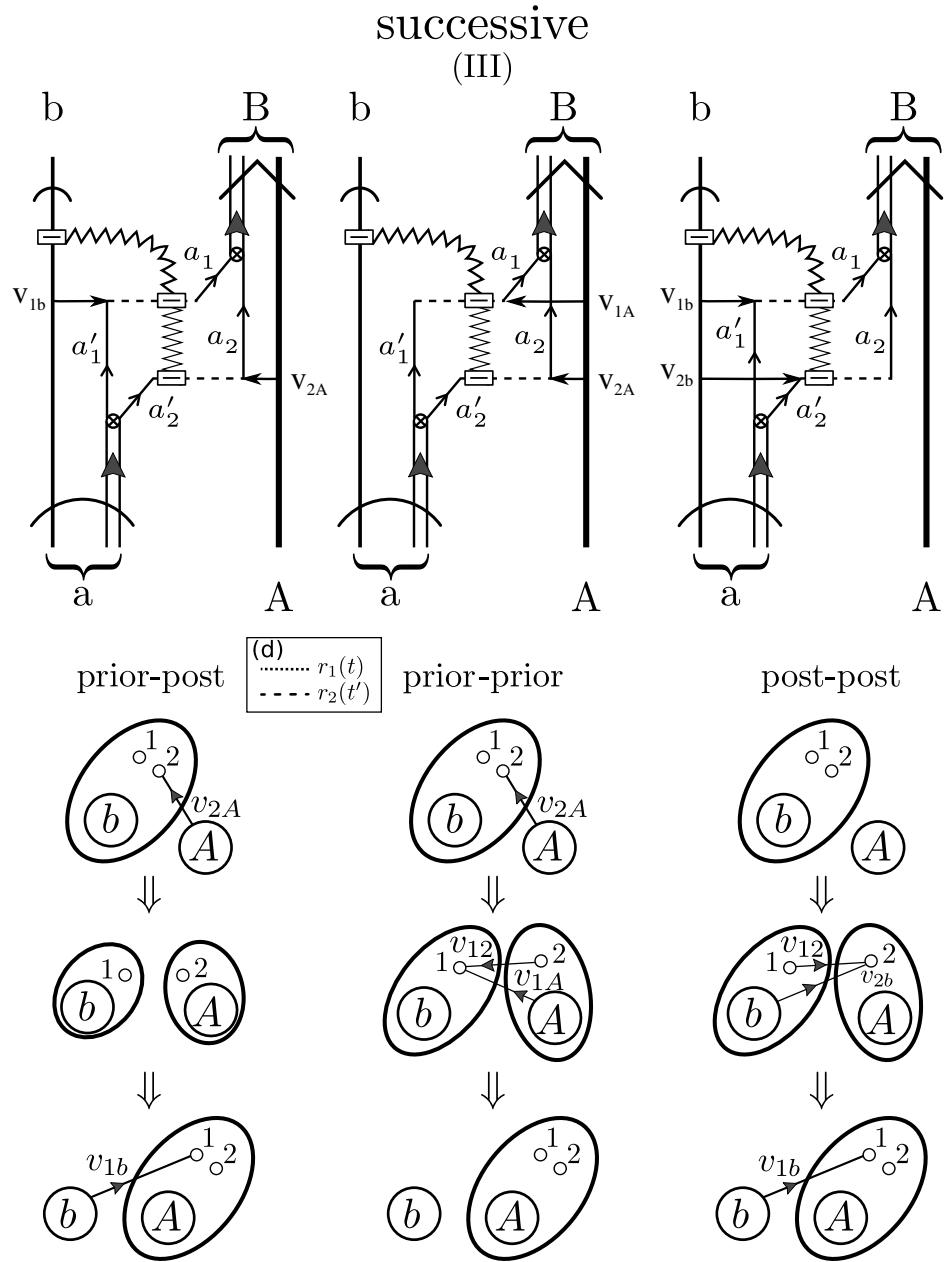


Figure 6.5.2: Graphical representation of the successive transfer of two nucleons. For details see text. Because of energy conservation, the different choices concerning the interactions inducing transfer, lead to identical results.

interaction or mean field (horizontal short arrow) inducing a transfer process.

The successive transfer amplitude $\tilde{a}_\infty^{(2)}$ written making use of the post-prior representation is equal to (see Fig. 6.5.2)

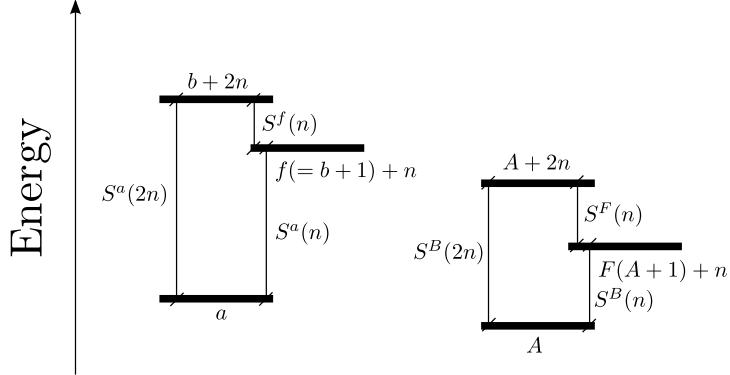


Figure 6.5.3: One- and two-neutron separation energies $S(n)$ and $S(2n)$ associated with the channels $\alpha \equiv a (= b + 2) + A \rightarrow \gamma \equiv f (= b + 1) + F (= A + 1) \rightarrow \beta \equiv b + B (= A + 2)$.

$$\begin{aligned} \tilde{a}^{(2)}(\infty) = & \frac{1}{i\hbar} \int_{-\infty}^{\infty} dt (\psi^b \psi^B, (V_{bB} - < V_{bB} >) e^{i\sigma_1} \psi^f \psi^F) \\ & \times \exp\left[\frac{i}{\hbar}(E^{bB} - E^{fF})t + \gamma_1(t)\right] \\ & \times \frac{1}{i\hbar} \int_{-\infty}^t dt' (\psi^f \psi^F, (V_{fF} - < V_{fF} >) e^{i\sigma_2} \psi^a \psi^A) \\ & \times \exp\left[\frac{i}{\hbar}(E^{fF} - E^{aA})t' + \gamma_2(t')\right]. \end{aligned} \quad (6.5.7)$$

To gain insight into the relative importance of the three terms contributing to Eq. (6.5.3) we discuss two situations, namely, the independent-particle model and the strong-correlation limits.

Before doing so, let us comment on the graphical description of the transfer amplitudes (6.5.4) (6.5.6) and (6.5.7) displayed in Figs. 6.5.1 and 6.5.2. The time arrow is assumed to point upwards: (I) Simultaneous transfer, in which one particle is transferred by the nucleon-nucleon interaction acting either in the entrance $\alpha \equiv a + A$ channel (prior) or in the final $\beta \equiv b + B$ channel (post), while the other particle follows suit making use of the particle-particle correlation (grey area) which binds the Cooper pair (see upper inset labelled (a)), represented by a solid arrow on a double line, to the projectile (curved arrowed lines) or to the target (opened arrowed lines). The above argument provides the explanation why when e.g. v_{1b} acts on one nucleon, the other nucleon also reacts instantaneously. In fact a Cooper pair displays generalized rigidity (emergent property in gauge space). A crossed open circle represents the particle-pair vibration coupling. The associated strength, together with an energy denominator, determines the amplitude $X_{a'_1 a'_2}$ with which the pair mode (Cooper pair) is in the (time reversed generalized to include different number of nodes) two particle configuration $a'_1 a'_2$. In the transfer

process, the relative motion orbit changes, the readjustement of the corresponding trajectory mismatch being operated by a Galilean transformation induced by the operator $(\exp\{\mathbf{k} \cdot (\mathbf{r}_{1A}(t) + \mathbf{r}_{2A}(t))\})$. This phenomenon, known as recoil process, is represented by a jagged line which provides information on the two transferred nucleons (single time appearing as argument of both single-particle coordinates r_1 and r_2 ; see Fig. 6.5.1 inset labeled (b)). In other words, information on the coupling of structure and reaction modes. **(II)** Non-orthogonality contribution. While one of the nucleons of the Cooper pairs is transferred under the action of v , the other goes, uncorrelatedly over, profiting of the non-orthogonality of the associated single-particle wavefunctions (see inset (c)). **(III)** Successive transfer. In this case, there are two time dependences associated with the acting of the nucleon-nucleon interaction twice (see Fig. 6.5.2 inset (d)).

6.5.1 Independent particle limit

In the independent particle limit, the two transferred particles do not interact among themselves but for antisymmetrization. Thus, the separation energies fulfill the relations (see Fig. 6.5.3)

$$S^B(2n) = 2S^B(n) = 2S^F(n), \quad (6.5.8)$$

and

$$S^a(2n) = 2S^a(n) = 2S^f(n). \quad (6.5.9)$$

In this case

$$\phi^{B(A)}(S^B(2n), \mathbf{r}_{1A}, \mathbf{r}_{2A}) = \sum_{a_1 a_2} \phi_{a_1}^{B(F)}(S^B(n), \mathbf{r}_{1A}) \phi_{a_2}^{F(A)}(S^F(n), \mathbf{r}_{2A}), \quad (6.5.10)$$

and

$$\phi^{a(b)}(S^a(2n), \mathbf{r}_{1b}, \mathbf{r}_{2b}) = \sum_{a'_1 a'_2} \phi_{a'_1}^{a(f)}(S^a(n), \mathbf{r}_{1b}) \phi_{a'_2}^{f(b)}(S^f(n), \mathbf{r}_{2b}), \quad (6.5.11)$$

where $(a_1, a_2) \equiv F$ and $(a'_1, a'_2) \equiv f$ span, as mentioned above, shells in nuclei B and a respectively.

Inserting Eqs. (6.5.8–6.5.11) in Eq. (6.5.4) one can show that

$$a^{(1)}(\infty) = a^{(NO)}(\infty). \quad (6.5.12)$$

It can be further demonstrated that within the present approximation, $\text{Im } \tilde{a}^{(2)} = 0$, and that

$$\begin{aligned} \tilde{a}^{(2)}(\infty) &= \frac{1}{i\hbar} \int_{-\infty}^{\infty} dt (\psi^b \psi^B, (V_{bB} - \langle V_{bB} \rangle) e^{i\sigma_1} \psi^f \psi^F) \\ &\quad \times \exp\left[\frac{i}{\hbar} (E^{bB} - E^{fF}) t + \gamma_1(t)\right] \\ &\quad \times \frac{1}{i\hbar} \int_{-\infty}^{\infty} dt' (\psi^f \psi^F, (V_{fF} - \langle V_{fF} \rangle) e^{i\sigma_2} \psi^a \psi^A) \\ &\quad \times \exp\left[\frac{i}{\hbar} (E^{fF} - E^{aA}) t' + \gamma_2(t)\right]. \end{aligned} \quad (6.5.13)$$

The total absolute differential cross section (6.5.2), where $P = |a(\infty)|^2 = |\tilde{a}^{(2)}|^2$, is then equal to the product of two one-particle transfer cross sections (see Fig. 5.1.1), associated with the (virtual) reaction channels

$$\alpha \equiv a + A \rightarrow f + F \equiv \gamma, \quad (6.5.14)$$

and

$$\gamma \equiv f + F \rightarrow b + B \equiv \beta. \quad (6.5.15)$$

In fact, Eq.(6.5.13) involves no time ordering and consequently the two processes above are completely independent of each other. This result was expected because being $v_{12} = 0$, the transfer of one nucleon cannot influence, aside from selecting the initial state for the second step, the behaviour of the other nucleon.

6.5.2 Strong correlation (cluster) limit

The second limit to be considered is the one in which the correlation between the two nucleons is so strong that (see Fig. 6.5.3)

$$S^a(2n) \approx S^a(n) \gg S^f(n), \quad (6.5.16)$$

and

$$S^B(2n) \approx S^B(n) \gg S^F(n). \quad (6.5.17)$$

That is, the magnitude of the one-nucleon separation energy is strongly modified by the pair breaking.

There is a different, although equivalent way to express (6.5.3) which is more convenient to discuss the strong coupling limit. In fact, making use of the posterior representation one can write

$$\begin{aligned} a^{(2)}(t) &= \tilde{a}^{(2)}(t) - a^{(NO)}(t) = \frac{1}{i\hbar} \int_{-\infty}^{\infty} dt (\psi^b \psi^B, (V_{bB} - < V_{bB} >) e^{i\sigma_1} \psi^f \psi^F) \\ &\quad \times \exp[\frac{i}{\hbar} (E^{bB} - E^{fF}) t + \gamma_1(t)] \\ &\quad \times \frac{1}{i\hbar} \int_{-\infty}^t dt' (\psi^f \psi^F, (V_{aA} - < V_{aA} >) \psi^a \psi^A) \\ &\quad \times \exp[\frac{i}{\hbar} (E^{fF} - E^{aA}) t' + \gamma_2(t')]. \end{aligned} \quad (6.5.18)$$

The relations (6.5.16), (6.5.17) imply

$$E^{fF} - E^{aA} = S^a(n) - S^F(n) \gg \frac{\hbar}{\tau}, \quad (6.5.19)$$

where τ is the collision time. Consequently the real part of $a^{(2)}(\infty)$ vanishes exponentially with the Q -value of the intermediate transition, while the imaginary part vanishes inversely proportional to this energy. One can thus write,

$$\operatorname{Re} a^{(2)}(\infty) \approx 0, \quad (6.5.20)$$

and

$$\begin{aligned} a^{(2)}(\infty) \approx & \frac{1}{i\hbar} \frac{\tau}{< E^{fF} > - E^{bB}} \sum_{fF} (\psi^b \psi^B, (V_{bB} - < V_{bB} >) \psi^f \psi^F)_{t=0} \\ & \times (\psi^f \psi^F, (V_{aA} - < V_{aA} >) \psi^a \psi^A)_{t=0}, \end{aligned} \quad (6.5.21)$$

where one has utilized the fact that $E^{bB} \approx E^{aA}$. For $v_{12} \rightarrow \infty$, ($< E^{fF} > - E^{bB} \rightarrow \infty$) and, consequently,

$$\lim_{v_{12} \rightarrow \infty} a^{(2)}(\infty) = 0. \quad (6.5.22)$$

Thus the total two-nucleon transfer amplitude is equal, in the strong coupling limit, to the amplitude $a^{(1)}(\infty)$.

Summing up, only when successive transfer and non-orthogonal corrections are included in the description of the two-nucleon transfer process, does one obtain a consistent description of the process, which correctly converges to the weak and strong correlation limiting values.

As already stated (Sect. 4.5), actual nuclei are, as a rule, closer to the independent particle limit than to the cluster one. Thus, it is not surprising that successive constitutes the major contribution to two-nucleon transfer processes. Nonetheless, the slight deviation from independent particle motion arising from pairing correlations, entangles the fermions (nucleons) partners of a Cooper pair, in a subtle fashion. This can be seen by expressing the pair transfer operator in terms of quasiparticles. Making use of the relation

$$a_\nu^\dagger = U_\nu \alpha_\nu^\dagger + V_\nu \alpha_{\tilde{\nu}}, \quad (6.5.23)$$

one obtains

$$P_\nu^\dagger = a_\nu^\dagger a_{\tilde{\nu}}^\dagger = U_\nu^2 \alpha_\nu^\dagger \alpha_{\tilde{\nu}}^\dagger - U_\nu V_\nu (\alpha_\nu^\dagger \alpha_\nu + \alpha_{\tilde{\nu}}^\dagger \alpha_{\tilde{\nu}}) - V_\nu^2 \alpha_{\tilde{\nu}} \alpha_\nu + U_\nu V_\nu. \quad (6.5.24)$$

Because one-quasiparticle transfer amplitude is $\sim U_\nu$ and the associated cross section is $\sim U_\nu^2$ (i.e. $\sim T^2$, T being the associated transition amplitude), it is natural that the two-quasiparticle transfer amplitude is $\sim U_\nu^2$ and that the associated cross section is $\sim U_\nu^4$ (i.e. $\sim T^4$). As apparent from (6.5.24), there exists a different possibility to transfer a pair of nucleons, leaving the quasiparticle distribution unchanged²⁵. Namely, through the last term of (6.5.24). Thus, a transfer amplitude $\sim U_\nu V_\nu$ and associated two-nucleon transfer cross section $\sim (U_\nu V_\nu)^2$ ($\sim T^2$). A result which one finds at the basis of (4.3.1).

²⁵This was the main message of Josephson (Josephson (1962)), and was the possibility which was missed by Giaever (Giaever (1973)), who interpreted supercurrents with zero voltage drop through the oxide barrier separating two superconductors as a metallic short. As he (Giaever) states: “to make an experimental discovery is not enough to observe something, one must also realize the significance of the observation”. It was also missed by Cohen, Falicov and Phillips (Cohen et al. (1962)), who developed the many-body tunneling Hamiltonian (eventually extended by Josephson) and who found four contributions to the current which correspond to normal (carriers of charge e) current of type S-Q. That is, implying quasiparticle excitation.

Appendix 6.A Spherical harmonics and angular momenta

With Condon–Shortley phases

$$Y_m^l(\hat{z}) = \delta_{m,0} \sqrt{\frac{2l+1}{4\pi}}, \quad Y_m^{l*}(\hat{r}) = (-1)^m Y_{-m}^l(\hat{r}). \quad (6.A.1)$$

Time-reversed phases consist in multiplying Condon–Shortley phases with a factor i^l , so

$$Y_m^l(\hat{z}) = \delta_{m,0} i^l \sqrt{\frac{2l+1}{4\pi}}, \quad Y_m^{l*}(\hat{r}) = (-1)^{l-m} Y_{-m}^l(\hat{r}). \quad (6.A.2)$$

With this phase convention, the relation with the associated Legendre polynomials includes an extra i^l factor with respect to the Condon–Shortley phase,

$$Y_m^l(\theta, \phi) = i^l \sqrt{\frac{2l+1}{4\pi}} \frac{(l-m)!}{(l+m)!} P_l^m(\cos \theta) e^{im\phi}. \quad (6.A.3)$$

6.A.1 Addition theorem

The addition theorem for the spherical harmonics states that

$$P_l(\cos \theta_{12}) = \frac{4\pi}{2l+1} \sum_m Y_m^l(\mathbf{r}_1) Y_m^{l*}(\mathbf{r}_2), \quad (6.A.4)$$

where θ_{12} is the angle between the vectors \mathbf{r}_1 and \mathbf{r}_2 . This result is independent of the phase convention. With *time-reversed phases*,

$$P_l(\cos \theta_{12}) = \frac{4\pi}{\sqrt{2l+1}} \left[Y^l(\hat{\mathbf{r}}_1) Y^l(\hat{\mathbf{r}}_2) \right]_0^0. \quad (6.A.5)$$

With *Condon–Shortley phases*,

$$P_l(\cos \theta_{12}) = (-1)^l \frac{4\pi}{\sqrt{2l+1}} \left[Y^l(\hat{\mathbf{r}}_1) Y^l(\hat{\mathbf{r}}_2) \right]_0^0. \quad (6.A.6)$$

6.A.2 Expansion of the delta function

The Dirac delta function can be expanded in multipoles, yielding

$$\begin{aligned} \delta(\mathbf{r}_2 - \mathbf{r}_1) &= \sum_l \delta(r_1 - r_2) \frac{2l+1}{4\pi r_1^2} P_l(\cos \theta_{12}) \\ &= \sum_l \delta(r_1 - r_2) \frac{1}{r_1^2} \sum_m Y_m^l(\mathbf{r}_1) Y_m^{l*}(\mathbf{r}_2). \end{aligned} \quad (6.A.7)$$

This result is independent of the phase convention. With *time-reversed phases*,

$$\delta(\mathbf{r}_2 - \mathbf{r}_1) = \sum_l \delta(r_1 - r_2) \frac{\sqrt{2l+1}}{r_1^2} \left[Y^l(\hat{\mathbf{r}}_1) Y^l(\hat{\mathbf{r}}_2) \right]_0^0. \quad (6.A.8)$$

6.A.3 Coupling and complex conjugation

If $\Psi_{M_1}^{I_1*} = (-1)^{I_1-M_1} \Psi_{-M_1}^{I_1}$ and $\Phi_{M_2}^{I_2*} = (-1)^{I_2-M_2} \Phi_{-M_2}^{I_2}$, as it happens to be the case for spherical harmonics with time-reversed phases, then

$$\begin{aligned} [\Psi^{I_1} \Phi^{I_2}]_M^{I*} &= \sum_{\substack{M_1 M_2 \\ (M_1 + M_2 = M)}} \langle I_1 I_2 M_1 M_2 | IM \rangle \Psi_{M_1}^{I_1*} \Phi_{M_2}^{I_2*} \\ &= \sum_{\substack{M_1 M_2 \\ (M_1 + M_2 = M)}} (-1)^{I-M_1-M_2} \langle I_1 I_2 - M_1 - M_2 | I - M \rangle \Psi_{-M_1}^{I_1} \Phi_{-M_2}^{I_2} \\ &= (-1)^{I-M} \sum_{\substack{M_1 M_2 \\ (M_1 + M_2 = M)}} \langle I_1 I_2 - M_1 - M_2 | I - M \rangle \Psi_{-M_1}^{I_1} \Phi_{-M_2}^{I_2} \\ &= (-1)^{I-M} [\Psi^{I_1} \Phi^{I_2}]_{-M}^I, \end{aligned} \quad (6.A.9)$$

where we have used (6.A.23).

Let us care now about the spinor functions $\chi_m^{1/2}(\sigma)$, which have the form

$$\chi^{1/2}(\sigma = 1/2) = \begin{bmatrix} 1 \\ 0 \end{bmatrix} \quad \chi^{1/2}(\sigma = -1/2) = \begin{bmatrix} 0 \\ 1 \end{bmatrix}, \quad (6.A.10)$$

or

$$\chi_m^{1/2}(\sigma) = \delta_{m,\sigma}. \quad (6.A.11)$$

Thus, $\chi_m^{1/2*}(\sigma) = \chi_m^{1/2}(\sigma) = \delta_{m,\sigma}$, but we can also write

$$\chi_m^{1/2*}(\sigma) = (-1)^{1/2-m+1/2-\sigma} \chi_{-m}^{1/2}(-\sigma). \quad (6.A.12)$$

This trick enable us to write

$$[Y^l(\hat{r}) \chi^{1/2}(\sigma)]_M^{J*} = (-1)^{1/2-\sigma+J-M} [Y^l(\hat{r}) \chi^{1/2}(-\sigma)]_{-M}^J, \quad (6.A.13)$$

which can be derived in a similar way as (6.A.9).

6.A.4 Angular momenta coupling

Relation between Clebsh–Gordan and $3j$ coefficients:

$$\langle j_1 j_2 m_1 m_2 | JM \rangle = (-1)^{j_1-j_2+M} \sqrt{2J+1} \begin{pmatrix} j_1 & j_2 & J \\ m_1 & m_2 & -M \end{pmatrix}. \quad (6.A.14)$$

Relation between Wigner and $9j$ coefficients:

$$\begin{aligned} ((j_1 j_2)_{j_{12}} (j_3 j_4)_{j_{34}} | (j_1 j_3)_{j_{13}} (j_2 j_4)_{j_{24}})_j &= \\ \sqrt{(2j_{12}+1)(2j_{13}+1)(2j_{24}+1)(2j_{34}+1)} &\left\{ \begin{array}{ccc} j_1 & j_2 & j_{12} \\ j_3 & j_4 & j_{34} \\ j_{13} & j_{24} & j \end{array} \right\}. \end{aligned} \quad (6.A.15)$$

6.A.5 Integrals

Let us now prove

$$\int d\Omega \left[Y^l(\hat{r}) Y^l(\hat{r}) \right]_M^I = \delta_{M,0} \delta_{I,0} \sqrt{2l+1}. \quad (6.A.16)$$

$$\begin{aligned} \int d\Omega \left[Y^l(\hat{r}) Y^l(\hat{r}) \right]_M^I &= \sum_{\substack{m_1, m_2 \\ (m_1+m_2=M)}} \langle l \ l \ m_1 \ m_2 | IM \rangle \int d\Omega Y_{m_1}^l(\hat{r}) Y_{m_2}^l(\hat{r}) \\ &= \sum_{\substack{m_1, m_2 \\ (m_1+m_2=M)}} (-1)^{l+m_1} \langle l \ l - m_1 \ m_2 | IM \rangle \int d\Omega Y_{m_1}^{l*}(\hat{r}) Y_{m_2}^l(\hat{r}) \\ &= \delta_{M,0} \sum_m (-1)^{l+m} \langle l \ l - m \ m | IO \rangle \\ &= \delta_{M,0} \sqrt{2l+1} \sum_m \langle l \ l - m \ m | IO \rangle \langle l \ l - m \ m | 00 \rangle \\ &= \delta_{M,0} \delta_{I,0} \sqrt{2l+1}, \end{aligned} \quad (6.A.17)$$

where we have used

$$\langle l \ l - m \ m | 0 \ 0 \rangle = \frac{(-1)^{l+m}}{\sqrt{2l+1}} \quad (6.A.18)$$

Let us now prove

$$\sum_{\sigma} \int d\Omega (-1)^{1/2-\sigma} \left[\Psi^j(\hat{r}, -\sigma) \Psi^j(\hat{r}, \sigma) \right]_M^I = -\delta_{M,0} \delta_{I,0} \sqrt{2j+1}. \quad (6.A.19)$$

$$\begin{aligned} \sum_{\sigma} \int d\Omega (-1)^{1/2-\sigma} \left[\Psi^j(\hat{r}, -\sigma) \Psi^j(\hat{r}, \sigma) \right]_M^I &= \sum_{\substack{m_1, m_2 \\ (m_1+m_2=M)}} \langle j \ j \ m_1 \ m_2 | IM \rangle \sum_{\sigma} \int d\Omega \Psi_{m_1}^j(\hat{r}, -\sigma) \Psi_{m_2}^j(\hat{r}, \sigma) \\ &= \sum_{\substack{m_1, m_2 \\ (m_1+m_2=M)}} \langle j \ j \ m_1 \ m_2 | IM \rangle \sum_{\sigma} (-1)^{j+m_1} \int d\Omega \Psi_{-m_1}^{j*}(\hat{r}, \sigma) \Psi_{m_2}^j(\hat{r}, \sigma) \\ &= \sum_{\substack{m_1, m_2 \\ (m_1+m_2=M)}} \langle j \ j \ m_1 \ m_2 | IM \rangle (-1)^{j+m_1} \delta_{-m_1, m_2} \\ &= \delta_{M,0} \sum_m (-1)^{j+m} \langle j \ j \ m - m | IO \rangle \\ &= -\delta_{M,0} \sqrt{2j+1} \sum_m (-1)^{j+m} \langle j \ j \ m - m | IO \rangle \langle j \ j \ m - m | 00 \rangle \\ &= -\delta_{M,0} \delta_{I,0} \sqrt{2j+1}. \end{aligned} \quad (6.A.20)$$

6.A.6 Symmetry properties

Note also another useful property

$$[\Psi^{I_1} \Psi^{I_2}]_M^I = (-1)^{I_1+I_2-I} [\Psi^{I_2} \Psi^{I_1}]_M^I, \quad (6.A.21)$$

by virtue of the symmetry property of the Clebsh-Gordan coefficients

$$\langle I_1 I_2 m_1 m_2 | IM \rangle = (-1)^{I_1+I_2-I} \langle I_2 I_1 m_2 m_1 | IM \rangle. \quad (6.A.22)$$

Here's another symmetry property of the Clebsh-Gordan coefficients

$$\langle I_1 I_2 m_1 m_2 | IM \rangle = (-1)^{I_1+I_2-I} \langle I_1 I_2 -m_2 -m_1 | I - M \rangle. \quad (6.A.23)$$

Another one, which can be derived from the simpler properties of $3j$ -symbols

$$\langle I_1 I_2 m_1 m_2 | IM \rangle = (-1)^{I_1-m_1} \sqrt{\frac{2I+1}{2I_2+1}} \langle I_1 I m_1 -M | I_2 m_2 \rangle. \quad (6.A.24)$$

Let us use this last property to calculate sums of the type

$$\sum_{m_1, m_3} |\langle I_1 I_2 m_1 m_2 | I_3 m_3 \rangle|^2. \quad (6.A.25)$$

Using (6.A.24), we have

$$\begin{aligned} \sum_{m_1, m_3} |\langle I_1 I_2 m_1 m_2 | I_3 m_3 \rangle|^2 &= \\ \frac{2I_3 + 1}{2I_2 + 1} \sum_{m_1, m_3} |\langle I_1 I_3 m_1 -m_3 | I_2 m_2 \rangle|^2 &= \frac{2I_3 + 1}{2I_2 + 1}, \end{aligned} \quad (6.A.26)$$

since

$$\sum_{m_1, m_3} |\langle I_1 I_3 m_1 -m_3 | I_2 m_2 \rangle|^2 = \sum_{m_1, m_3} |\langle I_1 I_3 m_1 m_3 | I_2 m_2 \rangle|^2 = 1. \quad (6.A.27)$$

Appendix 6.B distorted waves

Let us have a closer look at the partial wave expansion of the distorted waves

$$\chi^{(+)}(\mathbf{k}, \mathbf{r}) = \sum_l \frac{4\pi}{kr} i^l e^{i\sigma l} F_l \sum_m Y_m^l(\hat{r}) Y_m^{l*}(\hat{k}). \quad (6.B.1)$$

Notice the very important fact that *this definition is independent of the phase convention*, since the l -dependent phase is multiplied by its complex conjugate.

$$\chi^{(-)}(\mathbf{k}, \mathbf{r}) = \chi^{(+)*}(-\mathbf{k}, \mathbf{r}) = \sum_l \frac{4\pi}{kr} i^{-l} e^{-i\sigma l} F_l^* \sum_m Y_m^{l*}(\hat{r}) Y_m^l(-\hat{k}). \quad (6.B.2)$$

As $Y_m^l(-\hat{k}) = (-1)^l Y_m^l(\hat{k})$, we have

$$\chi^{(-)}(\mathbf{k}, \mathbf{r}) = \sum_l \frac{4\pi}{kr} i^l e^{-i\sigma^l} F_l^* \sum_m Y_m^{l*}(\hat{r}) Y_m^l(\hat{k}), \quad (6.B.3)$$

which is also independent of the phase convention. With time-reversed phase convention

$$\chi^{(+)}(\mathbf{k}, \mathbf{r}) = \sum_l \frac{4\pi}{kr} i^l \sqrt{2l+1} e^{i\sigma^l} F_l [Y^l(\hat{r}) Y^l(\hat{k})]_0^0, \quad (6.B.4)$$

while with Condon–Shortley phase convention we get an extra $(-1)^l$ factor:

$$\chi^{(+)}(\mathbf{k}, \mathbf{r}) = \sum_l \frac{4\pi}{kr} i^{-l} \sqrt{2l+1} e^{i\sigma^l} F_l [Y^l(\hat{r}) Y^l(\hat{k})]_0^0. \quad (6.B.5)$$

The partial-wave expansion of the Green function $G(\mathbf{r}, \mathbf{r}')$ is

$$G(\mathbf{r}, \mathbf{r}') = i \sum_l \frac{f_l(k, r_<) P_l(k, r_>)}{krr'} \sum_m Y_m^l(\hat{r}) Y_m^{l*}(\hat{r}'), \quad (6.B.6)$$

where $f_l(k, r_<)$ and $P_l(k, r_>)$ are the regular and the irregular solutions of the homogeneous problem respectively. With *time-reversed* phase convention

$$G(\mathbf{r}, \mathbf{r}') = i \sum_l \sqrt{2l+1} \frac{f_l(k, r_<) P_l(k, r_>)}{krr'} [Y^l(\hat{r}) Y^l(\hat{r}')]_0^0. \quad (6.B.7)$$

Appendix 6.C hole states and time reversal

Let us consider the state $|(jm)^{-1}\rangle$ obtained by removing a ψ_{jm} single-particle state from a $J = 0$ closed shell $|0\rangle$. The antisymmetrized product state

$$\sum_m \mathcal{A}\{\psi_{jm}|(jm)^{-1}\} \propto |0\rangle \quad (6.C.1)$$

is clearly proportional to $|0\rangle$. This gives us the transformation rules of $|(jm)^{-1}\rangle$ under rotations, which must be such that, when multiplied by a j, m spherical tensor and summed over m , yields a $j = 0$ tensor. It can be seen that these properties imply that $|(jm)^{-1}\rangle$ transforms like $(-1)^{j+m} T_{j-m}$, T_{j-m} being a spherical tensor. It also follows that the hole state $|(j\bar{m})^{-1}\rangle$ transforms like a j, m spherical tensor if $\psi_{j\bar{m}}$ is defined as the \mathcal{R} -conjugate to ψ_{jm} by the relation

$$\psi_{j\bar{m}} \equiv (-1)^{j+m} \psi_{j-m}. \quad (6.C.2)$$

In other words, with the latter definition a *hole state* transforms under rotations with the right phase. We will now show that \mathcal{R} -conjugation is equivalent to a rotation of spin and spatial coordinates through an angle $-\pi$ about the y -axis:

$$e^{i\pi J_y} \psi_{jm} = (-1)^{j+m} \psi_{j-m} \equiv \psi_{j\bar{m}}. \quad (6.C.3)$$

Let us begin by calculating $e^{i\pi L_y} Y_l^m$. The rotation matrix about the y -axis is

$$R_y(\theta) = \begin{pmatrix} \cos(\theta) & 0 & \sin(\theta) \\ 0 & 1 & 0 \\ -\sin(\theta) & 0 & \cos(\theta) \end{pmatrix}, \quad (6.C.4)$$

so for $R_y(-\pi)$ we get

$$R_y(-\pi) = \begin{pmatrix} -1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & -1 \end{pmatrix}. \quad (6.C.5)$$

When applied to the generic direction $(\sin(\theta)\cos(\phi), \sin(\theta)\sin(\phi), \cos(\theta))$, we obtain $(-\sin(\theta)\cos(\phi), \sin(\theta)\sin(\phi), -\cos(\theta))$, which corresponds to making the substitutions

$$\theta \rightarrow \pi - \theta, \quad \phi \rightarrow \pi - \phi. \quad (6.C.6)$$

When we substitute these angular transformations in the spherical harmonic $Y_l^m(\theta, \phi)$, we obtain the rotated $Y_l^m(\theta, \phi)$:

$$e^{i\pi L_y} Y_l^m = (-1)^{l+m} Y_l^{-m}. \quad (6.C.7)$$

Let us now turn our attention to the spin coordinates rotation $e^{i\pi s_y} \chi_m$. The rotation matrix in spin space is

$$\begin{pmatrix} \cos(\theta/2) & -\sin(\theta/2) \\ \sin(\theta/2) & \cos(\theta/2) \end{pmatrix}, \quad (6.C.8)$$

which, for $\theta = -\pi$ is

$$\begin{pmatrix} 0 & 1 \\ -1 & 0 \end{pmatrix}. \quad (6.C.9)$$

Applying it to the spinors, we find the rule

$$e^{i\pi s_y} \chi_m = (-1)^{1/2+m} \chi_{-m}, \quad (6.C.10)$$

so

$$\begin{aligned} e^{i\pi J_y} \psi_{jm} &= \sum_{m_l m_s} \langle l \ m_l \ 1/2 \ m_s | j \ m \rangle \ e^{i\pi L_y} Y_l^{m_l} \ e^{i\pi s_y} \chi_{m_s} \\ &= \sum_{m_l m_s} (-1)^{1/2+m_s+l+m_l} \langle l \ m_l \ 1/2 \ m_s | j \ m \rangle \ Y_l^{-m_l} \ \chi_{-m_s} \\ &= \sum_{m_l m_s} (-1)^{1+m-j+2l} \langle l \ -m_l \ 1/2 \ -m_s | j \ -m \rangle \ Y_l^{-m_l} \ \chi_{-m_s} \\ &= (-1)^{m+j} \psi_{j-m} \equiv \psi_{j\bar{m}}, \end{aligned} \quad (6.C.11)$$

where we have used $(-1)^{1+m-j+2l} = -(-1)^{m-j} = (-1)^{m+j}$, as j, m are always half-integers and l is always an integer.

We now turn our attention to the time reversal operation, which amounts to the transformations

$$\mathbf{r} \rightarrow \mathbf{r}, \quad \mathbf{p} \rightarrow -\mathbf{p}. \quad (6.C.12)$$

This is enough to define the operator of time reversal of a spinless particle (see Messiah). In the position representation, in which \mathbf{r} is real and \mathbf{p} pure imaginary, this (unitary antilinear) operator is the complex conjugation operator.

As angular momentum $\mathbf{l} = \mathbf{r} \times \mathbf{p}$ changes sign under time reversal, so does spin:

$$\mathbf{s} \rightarrow -\mathbf{s}, \quad (6.C.13)$$

which, along with (6.C.12), completes the set of rules that define the time reversal operation on a particle with spin. In the representation of eigenstates of \mathbf{s}^2 and s_z , complex conjugation alone changes only the sign of s_y , so an additional rotation of $-\pi$ around the y -axis is necessary to change the sign of s_x, s_z and implement the transformation (6.C.13). If we call K the time-reversal operator, we have

$$K\psi_{jm} = e^{i\pi s_y} \psi_{jm}^*. \quad (6.C.14)$$

This is completely general and independent of the phase convention. It only depends on the fact that we have used the \mathbf{r} representation for the spatial wave function and the representation of the eigenstates of \mathbf{s}^2 and s_z for the spin part. *If we use time-reversal phases for the spherical harmonics* (see(6.A.2)),

$$Y_m^{l*} = (-1)^{l+m} Y_{-m}^l = e^{i\pi L_y} Y_m^l. \quad (6.C.15)$$

So we can write

$$K\psi_{jm} = e^{i\pi J_y} \psi_{jm} = \psi_{j\bar{m}}. \quad (6.C.16)$$

Note again that this last expression is valid only if we use time-reversal phases for the spherical harmonics. Only in this case time-reversal coincides with \mathcal{R} -conjugation and hole states.

In BCS theory, the quasiparticles are defined in terms of linear combinations of particles and holes. With time-reversal phases, holes are equivalent to time-reversed states, and we get the definitions

$$\begin{aligned} \alpha_\nu^\dagger &= u_\nu a_\nu^\dagger - v_\nu a_{\bar{\nu}}^\dagger & a_\nu^\dagger &= u_\nu \alpha_\nu^\dagger + v_\nu \alpha_{\bar{\nu}} \\ a_{\bar{\nu}}^\dagger &= u_\nu a_\nu^\dagger + v_\nu a_{\bar{\nu}}^\dagger & a_{\bar{\nu}}^\dagger &= u_\nu \alpha_{\bar{\nu}}^\dagger - v_\nu \alpha_\nu \\ \alpha_\nu &= u_\nu a_\nu - v_\nu a_{\bar{\nu}}^\dagger & a_\nu &= u_\nu \alpha_\nu + v_\nu \alpha_{\bar{\nu}}^\dagger \\ a_{\bar{\nu}} &= u_\nu a_{\bar{\nu}} + v_\nu a_\nu^\dagger & a_{\bar{\nu}} &= u_\nu \alpha_{\bar{\nu}} - v_\nu \alpha_\nu^\dagger \end{aligned} \quad (6.C.17)$$

Appendix 6.D Spectroscopic amplitudes in the BCS approximation

The creation operator of a pair of fermions coupled to J, M can be expressed in second quantization as

$$T^\dagger(j_1, j_2, JM) = N \sum_m \langle j_1 m \ j_2 M - m | J M \rangle a_{j_1 m}^\dagger a_{j_2 M-m}^\dagger, \quad (6.D.1)$$

where N is a normalization constant. To determine it, we write the wave function resulting from the action of (6.D.1) on the vacuum

$$\begin{aligned}\Psi = T^\dagger(j_1, j_2, JM)|0\rangle &= \frac{N}{\sqrt{2}} \sum_m \langle j_1 m j_2 M - m | J M \rangle \\ &\times (\phi_{j_1 m}(\mathbf{r}_1) \phi_{j_2 M-m}(\mathbf{r}_2) - \phi_{j_2 M-m}(\mathbf{r}_1) \phi_{j_1 m}(\mathbf{r}_2)).\end{aligned}\quad (6.D.2)$$

The norm is

$$\begin{aligned}|\Psi|^2 &= \frac{N^2}{2} \sum_{mm'} \langle j_1 m j_2 M - m | J M \rangle \langle j_1 m' j_2 M - m' | J M \rangle \\ &\times (\phi_{j_1 m}(\mathbf{r}_1) \phi_{j_2 M-m}(\mathbf{r}_2) - \phi_{j_2 M-m}(\mathbf{r}_1) \phi_{j_1 m}(\mathbf{r}_2)) \\ &\times (\phi_{j_1 m'}(\mathbf{r}_1) \phi_{j_2 M-m'}(\mathbf{r}_2) - \phi_{j_2 M-m'}(\mathbf{r}_1) \phi_{j_1 m'}(\mathbf{r}_2)).\end{aligned}\quad (6.D.3)$$

Integrating we get

$$\begin{aligned}1 &= \frac{N^2}{2} \sum_{mm'} \langle j_1 m j_2 M - m | J M \rangle \langle j_1 m' j_2 M - m' | J M \rangle \\ &\times (2\delta_{m,m'} - 2\delta_{j_1,j_2}\delta_{m,M-m'}) \\ &= N^2 \left(\sum_m \langle j_1 m j_2 M - m | J M \rangle^2 \right. \\ &\quad \left. - \delta_{j_1,j_2} \sum_m \langle j_1 m j_2 M - m | J M \rangle \langle j_1 M - m j_2 m | J M \rangle \right) \\ &= N^2 (1 - \delta_{j_1,j_2}(-1)^{2j-J}),\end{aligned}\quad (6.D.4)$$

where we have used the closure condition for Clebsh–Gordan coefficients and (6.A.22), and δ_{j_1,j_2} must be interpreted as a δ function regarding all the quantum numbers but the magnetic one. We see that two fermions with identical quantum numbers (but the magnetic one) *cannot couple to J odd*. If J is even, the normalization constant is

$$N = \frac{1}{\sqrt{1 + \delta_{j_1,j_2}}}.\quad (6.D.5)$$

To sum up,

$$P^\dagger(j_1, j_2, JM) = \frac{1}{\sqrt{1 + \delta_{j_1,j_2}}} \sum_m \langle j_1 m j_2 M - m | J M \rangle a_{j_1 m}^\dagger a_{j_2 M-m}^\dagger.\quad (6.D.6)$$

The spectroscopic amplitude for finding in a $A + 2, J_f, M_f$ nucleus a couple of nucleons with quantum numbers j_1, j_2 coupled to J on top of a A, J_i nucleus is

$$B(j_1, j_2(J)) = \sum_{M, M_i} \langle J_i M_i JM | J_f M_f \rangle \langle \Psi_{J_f M_f} | T^\dagger(j_1, j_2, JM) | \Psi_{J_i M_i} \rangle.\quad (6.D.7)$$

This is completely general. It depends on the structure model only through the way the $A + 2$ and A nuclei are treated. We now want to turn our attention to the expression of $B(J, j_1, j_2)$ in the BCS approximation when both the $A + 2$ and the A are 0^+ , zero-quasiparticle ground states. In order to do this, we write (6.D.6) in terms of quasiparticle operators using (6.C.17)²⁶:

$$\begin{aligned} T^\dagger(j_1, j_2, JM) = & \frac{1}{\sqrt{1 + \delta_{j_1, j_2}}} \sum_{m_1, m_2} \langle j_1 \ m_1 \ j_2 \ m_2 | J \ M \rangle \left(U_{j_1} U_{j_2} \alpha_{j_1 m_1}^\dagger \alpha_{j_2 m_2}^\dagger \right. \\ & + (-1)^{j_1+j_2-M} V_{j_1} V_{j_2} \alpha_{j_1-m_1} \alpha_{j_2-m_2} \\ & + (-1)^{j_2-m_2} U_{j_1} V_{j_2} \alpha_{j_1 m_1}^\dagger \alpha_{j_2-m_2} \\ & - (-1)^{j_1-m_1} V_{j_1} U_{j_2} \alpha_{j_2 m_2}^\dagger \alpha_{j_1-m_1} \\ & \left. + (-1)^{j_1-m_1} V_{j_1} U_{j_2} \delta_{j_1 j_2} \delta_{-m_1 m_2} \right). \end{aligned} \quad (6.D.8)$$

If both nuclei are in the $|BCS\rangle$ ground state (zero-quasiparticles state), the only term that survives is the last one in the above expression, and (6.D.7) becomes

$$\begin{aligned} B_j = B(j^2(0)) &= \frac{1}{\sqrt{2}} \sum_m \langle j \ m \ j \ -m | 0 \ 0 \rangle (-1)^{j-m} V_j U_j \\ &= \frac{1}{\sqrt{2}} \sum_m \frac{(-1)^{j-m}}{\sqrt{(2j+1)}} (-1)^{j-m} V_j U_j \\ &= \frac{1}{\sqrt{2}} \sum_m \frac{1}{\sqrt{(2j+1)}} V_j U_j. \end{aligned} \quad (6.D.9)$$

After carrying out the summation one finds,

$$B_j = B(j^2(0)) = \sqrt{j+1/2} V_j U_j. \quad (6.D.10)$$

Note that in this final expression V_j refers to the $A + 2$ nucleus, while U_j is related to the A nucleus. In practice, it does not make a big difference to calculate both for the same nucleus.

Appendix 6.E Derivation of two-nucleon transfer transition amplitudes including recoil, non-orthogonality and successive transfer.

In the present Appendix we reproduce with the permission of the author the first (manuscript) page (Fig. 6.E.1) of what, arguably, was the first complete deriva-

²⁶In what follows, we use the phase convention $\alpha_{j\bar{m}} = (-1)^{j-m} \alpha_{j-m}$ instead of $\alpha_{j\bar{m}} = (-1)^{j+m} \alpha_{j-m}$, consistent with (6.C.2). Had we stick to the definition (6.C.2), the amplitude $B(0, j, j)$ calculated below would have a minus sign, which would not have any physical consequence.

↑ →

$$A + a \xrightarrow{(b+n_1+n_2)} C + c \xrightarrow{(A+n_2)} b + B \xrightarrow{(A+n_1+n_2)}$$

We need

$$2 \sum_{\sigma_1 \sigma_2} \int d^3 \vec{r}_{C_c} d^3 \vec{r}_{b_1} d^3 \vec{r}_{A_2} [\Psi^{j_{f_1}}(\vec{r}_{A_1}, \sigma_1) \Psi^{j_{f_2}}(\vec{r}_{A_2}, \sigma_2)]_M^* \chi^{(+)}_i(\vec{r}_{Bb}) U(r_{b_1}) \Psi^{(+)}_i$$

$\downarrow \chi_i$

$H = H_C + H_c - \frac{\hbar^2}{2\mu_{C_c}} \nabla_{C_c}^2 + U(r_{C_c}) + V(r_{c_2}), \quad (E - H) \Psi^{(+)} = O = [E - H]$

Let $f_{K_M}(\vec{r}_{C_c}) = \langle [\Psi^{j_{f_2}}(\vec{r}_{A_2}, \sigma_2) \Psi^{j_i}(\vec{r}_{b_1}, \sigma_1)]_M^K | \Psi^{(+)}(\vec{r}_{A_2}, \sigma_2, \vec{r}_{b_1}, \vec{r}_i; \vec{r}_{C_c}) \rangle_{\text{fixed}}$

$$\begin{aligned} f_{K_M}(\vec{r}_{C_c}) &= \frac{2\mu_{C_c}}{\hbar^2} \int d^3 \vec{r}'_{C_c} G(\vec{r}_{C_c}, \vec{r}'_{C_c}) \langle [\Psi^{j_{f_2}}(\vec{r}'_{A_2}, \sigma'_2) \Psi^{j_i}(\vec{r}'_{b_1}, \sigma'_1)]_M^K | V(r_{C'_2}) | \Psi \rangle_{\text{fixed}} \\ &\approx \frac{2\mu_c}{\hbar^2} \sum_{\sigma'_1 \sigma'_2} \int d^3 \vec{r}'_{C_c} d^3 \vec{r}'_{A_2} d^3 \vec{r}'_{b_1} G(\vec{r}_{C_c}, \vec{r}'_{C_c}) [\Psi^{j_{f_2}}(\vec{r}'_{A_2}, \sigma'_2) \Psi^{j_i}(\vec{r}'_{b_1}, \sigma'_1)]_M^K V(r_{C'_2}) \chi^{(+)}_i(\vec{r}'_{A_2}) \\ &\quad \times \langle [\Psi^{j_i}(\vec{r}'_{b_1}, \sigma'_1) \Psi^{j'_i}(\vec{r}'_{b_2})]_O^K \rangle + u_{K_M} \\ \text{Thus } u_{K_M}(\vec{r}_{C_c}) &= \frac{2\mu_{C_c}}{\hbar^2} \sum_{\sigma'_1 \sigma'_2} \int d^3 \vec{r}'_{C_c} d^3 \vec{r}'_{A_2} d^3 \vec{r}'_{b_1} G(\vec{r}_{C_c}, \vec{r}'_{C_c}) [\Psi^{j_{f_2}}(\vec{r}'_{A_2}, \sigma'_2) \Psi^{j_i}(\vec{r}'_{b_1}, \sigma'_1)]_M^K V(r_{C'_2}) \end{aligned}$$

↓

Figure 6.E.1: First manuscript page of Ben Bayman's derivation of the two-nucleon transfer reaction amplitude, in second order DWBA approximation.

tion²⁷ of the different contributions needed to calculate absolute two-nucleon transfer cross sections in a systematic way²⁸.

²⁷Bayman (1970) (unpublished).

²⁸Bayman (1971) and Bayman and Chen (1982)). Within this context we refer to Broglia et al. (1973) and Potel et al. (2013a) in particular Fig. 10 of this reference.

Chapter 7

Probing Cooper pairs with pair transfer

... I advocated the application of the word *phenomenon* exclusively to refer to the observation obtained under specified circumstances, including the account of the whole experimental arrangement.

Niels Bohr

Pairing density, both static ($n_{st} = e^{-2i\phi}\alpha'_0/V$) and dynamic¹ (α_{dyn}/V), can exist only to the extent that normal density exists. At the basis of *normal density* one finds independent nucleon (fermion) motion in a non local mean field, created by all nucleons. At the basis of pair (*abnormal*) density one finds the conversion of this motion into long range correlated motion of extended, strongly overlapping, weakly bound pairs of nucleons (quasi-bosons), moving in time reversal states close to the Fermi energy displaying *off diagonal long range order*. Because the mean square radius of a Cooper pair is considerably larger than nuclear dimensions, the average potential acts on them as a strong external field and affects their structure and dynamics. A fact whose implications have become physically clearer in the study of the structure of the two-neutron halo nucleus ^{11}Li and, quantitatively more accurate, by probing the system through pair transfer. *Said it differently, Cooper pairs are at the basis of both static and dynamic spontaneous breaking of gauge invariance, namely of (pairs) of particle number. To study the specific consequence of a broken symmetry one needs a probe which itself breaks the symmetry. In the present case a probe which changes in two particle number, like a heavy or light ion reaction inducing pair transfer.*

Within the above scenario we apply, in what follows, the formalism worked out in the previous chapters, to analyze Cooper pair transfer reactions. In the calcula-

¹See Figs. 3.5.7 and 4.4.2.

tion of the corresponding absolute differential cross sections, use has been made of the computer code COOPER (App. 7.A). A number of examples are considered covering nuclei throughout the mass table. Namely, two-particle transfer in ^{10}Be , ^7Li , ^{11}Li , ^{48}Ca , and ^{206}Pb (systems around closed shells), and on single open shell superfluid medium heavy nuclei (^{60}Ni and Sn-isotopes), induced both by light and heavy ions.

7.1 The $^1\text{H}(^{11}\text{Li}, ^9\text{Li})^3\text{H}$ reaction: evidence for phonon mediated pairing

We start by discussing the analysis of the two-neutron pickup² reaction $^1\text{H}(^{11}\text{Li}, ^9\text{Li})^3\text{H}$. The results provide evidence for a new mechanism of pairing correlations in nuclei: pygmy resonance (low-energy $E1$ -strength) mediated pairing interaction³, which strongly renormalizes the bare, $NN^{-1}S_0$ interaction. This is but a particular embodiment of phonon mediated pairing interaction found in nuclei, throughout the mass table⁴. Attention is also paid to the population of the $1/2^-$ first excited state of ^9Li lying⁵ at 2.69 MeV (see Figs. 2.7.2–2.7.5 and 7.1.1).

The main difference between light halo exotic nuclei and medium heavy nuclei lying along the valley of stability is the role fluctuations play in dressing particles (quasiparticles) and in renormalizing their properties (mass, charge, etc.) and their interactions, like the pairing interaction. In fact, in the case of e.g. Sn isotopes, mean field effects are dominant, while in the case of halo exotic nuclei renormalization effects can be as large as mean field ones. Concerning the pairing interaction, bare and induced contributions are about equal in the case of i.e. Sn-isotopes, while the second one is the overwhelming contribution in the case of ^{11}Li . The collective modes acting as glue of the Cooper pairs are mainly of quadrupole type in Sn and of dipole (pygmy resonance)⁶ type in the case of Li. In what follows we summarize aspects of structure and reaction which, interweaved shed insight on this single Cooper pair halo nucleus, and allows to compare theory with experiment in terms of absolute differential cross sections.

7.1.1 Structure

Within the scenario presented in Chapter 3 (Sect. 3.6) the wavefunction describing the structure of the halo neutrons in the ground state of ^{11}Li (the $p_{3/2}$ proton being assumed to act only as a spectator) can be written as

$$|0\rangle_v = |0\rangle + \alpha|(p_{1/2}, s_{1/2})_{1^-} \otimes 1^-; 0\rangle + \beta|(s_{1/2}, d_{5/2})_{2^+} \otimes 2^+; 0\rangle, \quad (7.1.1)$$

²Tanihata et al. (2008).

³Barranco et al. (2001); Potel et al. (2010).

⁴See footnote 51 Ch. 3 and refs. therein.

⁵To assess the direct character of the $1/2^-$ excitation process, the importance of (two-step) inelastic and of knockout channels were considered and found to be small (see Sect. 7.1.2).

⁶Within this context see Broglia et al. (2019a).

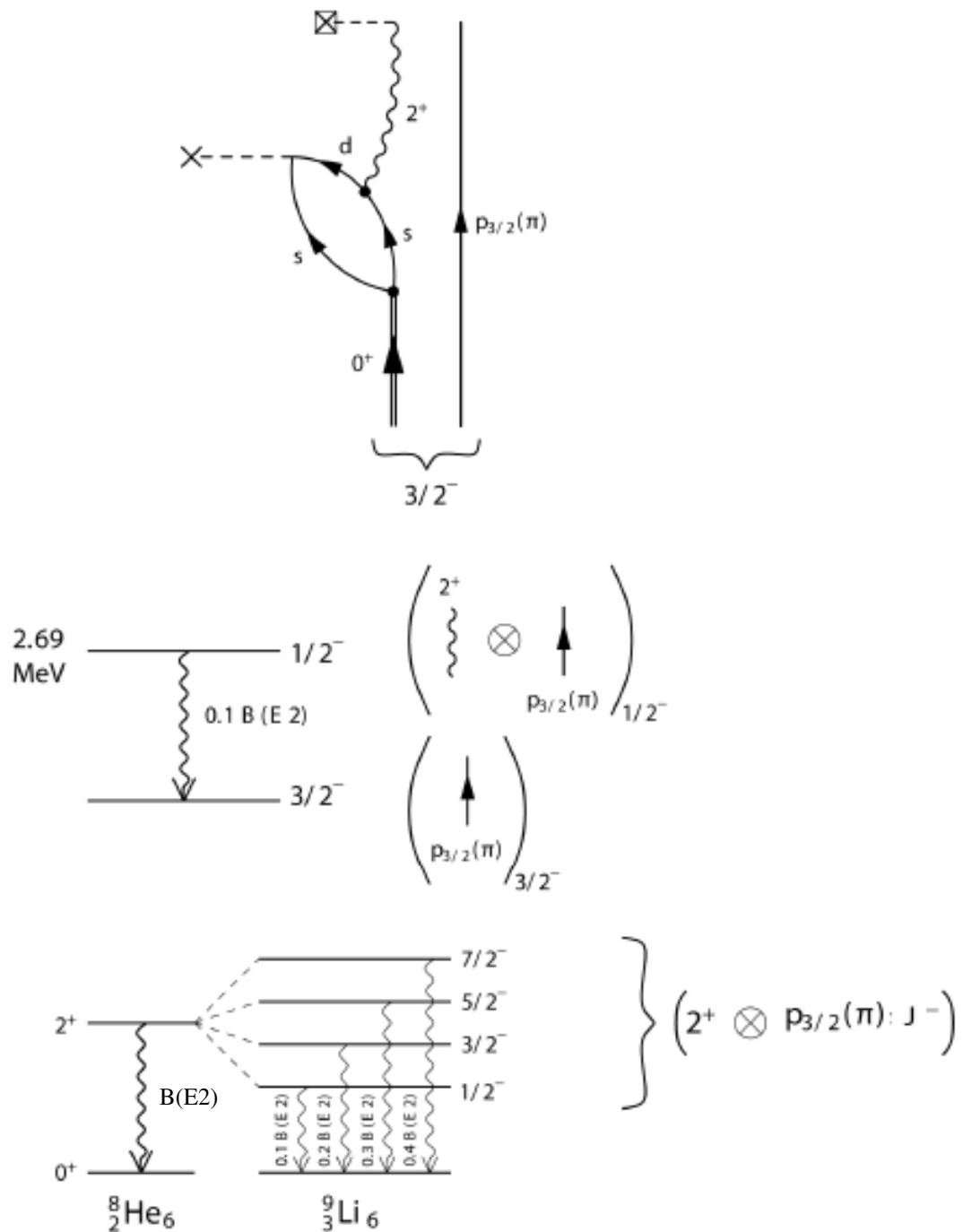


Figure 7.1.1: *Gedankenexperiments* (two-particle transfer γ -decay coincidence) aimed at better individuating the couplings involved in the neutron halo Cooper pair correlations in ^{11}Li and of the $1/2^-$ first excited state of ^9Li populated in the $^1\text{H}(^{11}\text{Li}, ^9\text{Li})^3\text{H}$ reaction (Tanihata et al. (2008); Barranco et al. (2001); Potel et al. (2010)).

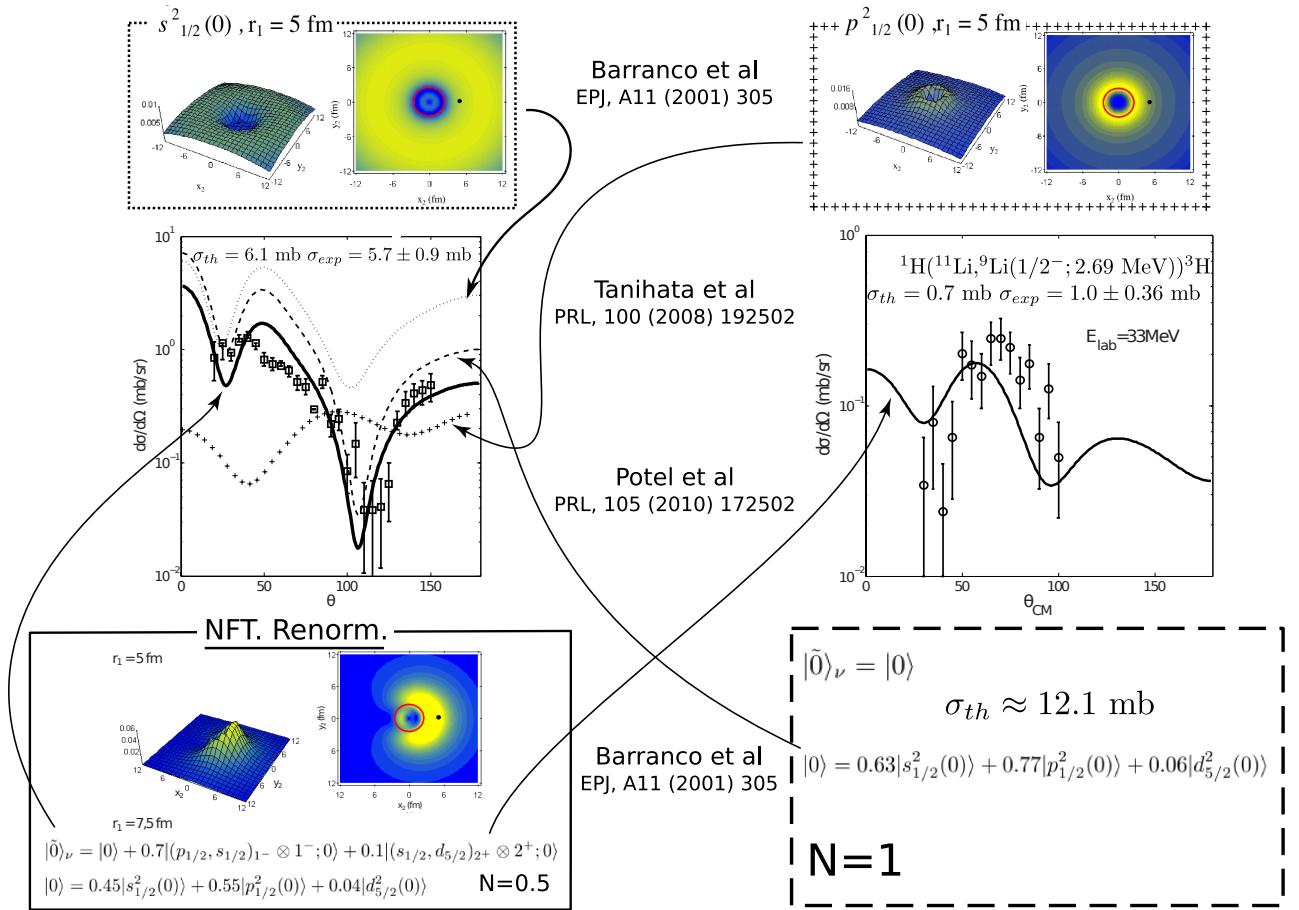


Figure 7.1.2: Absolute, two-nucleon transfer differential cross section associated with the ground state and the first excited state of ${}^9\text{Li}$, excited in the reaction ${}^1\text{H}({}^{11}\text{Li}, {}^9\text{Li})^3\text{H}$ (Tanihata et al., 2008) in comparison with the predicted differential cross sections (Potel et al., 2010) worked out making use of spectroscopic amplitudes and Cooper pair wavefunctions calculated with NFT, and of the optical potential collected in Table 7.1.1.

with

$$\alpha = 0.7, \quad \text{and} \quad \beta = 0.1, \quad (7.1.2)$$

and

$$|0\rangle = 0.45|s_{1/2}^2(0)\rangle + 0.55|p_{1/2}^2(0)\rangle + 0.04|d_{5/2}^2(0)\rangle, \quad (7.1.3)$$

$|1^-\rangle$ and $|2^+\rangle$ being the (QRPA) states describing the pygmy dipole resonance of ^{11}Li and the low-lying quadrupole vibration of the core. The corresponding NFT diagrams are shown in (a)–(e) of Fig. 2.7.4. The wavy curve represents both the quadrupole and dipole modes exchanged between the two halo neutrons, the horizontal dashed line the bare nuclear pairing interaction.

$^{11}\text{Li}(p,t)^9\text{Li}$												
	V	W	V_{so}	W_d	r_1	a_1	r_2	a_2	r_3	a_3	r_4	a_4
$p, {}^{11}\text{Li}^d)$	63.62	0.33	5.69	8.9	1.12	0.68	1.12	0.52	0.89	0.59	1.31	0.52
$d, {}^{10}\text{Li}^b)$	90.76	1.6	3.56	10.58	1.15	0.75	1.35	0.64	0.97	1.01	1.4	0.66
$t, {}^9\text{Li}^c)$	152.47	12.59	1.9	12.08	1.04	0.72	1.23	0.72	0.53	0.24	1.03	0.83

Table 7.1.1: Optical potentials (cf. Tanihata et al. (2008)) used in the calculation of the absolute differential cross sections displayed in Fig. 7.1.2.

We are then in presence of a paradigmatic nuclear embodiment of Cooper's model which is at the basis of BCS theory: a single weakly bound neutron pair on top of the Fermi surface, namely the ${}^9\text{Li}$ core. But the analogy goes beyond these aspects, and covers also the very nature of the interaction acting between Cooper pair partners nucleons (pairing interaction). Due to the the high polarizability of the system under study and of the small overlap of halo and core single particle wavefunctions, most of the Cooper pair correlation energy stems from the exchange of collective vibrations, the role of the strongly screened bare pairing interaction being, in this case, minor (see Sect. 3.6). In other words, we are in the presence of a new realization of the Cooper model in which a totally novel Bardeen-Pines-Frölich-like phonon induced pairing interaction is generated by a self induced collective vibration of the nuclear medium.

In connection with (7.1.1), it is revealing that the states populated in the inverse kinematics, two-neutron pick up reaction ${}^1\text{H}({}^{11}\text{Li}, {}^9\text{Li}){}^3\text{H}$ are⁷, aside from the ground $|3/2^-\text{gs}({}^9\text{Li})\rangle$ state, the first excited $|1/2^-, 2.69 \text{ MeV}({}^9\text{Li})\rangle$ level. The associated absolute differential cross sections calculated making use of the spectroscopic amplitudes displayed in Eqs. (7.1.1)–(7.1.3) and the optical potentials collected in Table 7.1.1, provide a quantitative account of the experimental data within the NFT scenario (Fig. 7.1.2). It is of notice that: (a) assuming the soft $E1$ -mode to be non-operative, that is, $\alpha = 0$ and normalizing to 1 (7.1.3) (see Fig. 7.1.2 lower right panel labeled $N = 1$), theory overpredicts experiment by a factor of 2; (b) the absolute direct two-particle transfer cross section associated with the $1/2^-$ state, reflects the value of the amplitude α .

⁷Tanihata et al. (2008).

7.1.2 Reaction

In the reaction ${}^1\text{H}({}^{11}\text{Li}, {}^9\text{Li}){}^3\text{H}$, the $1/2^-$ (2.69 MeV) first excited state of ${}^9\text{Li}$ can, in principle, not only be populated through the direct two-particle transfer reaction channel (sum of successive and simultaneous contributions plus non-orthogonality corrections, see Fig. 2.7.3 and Table 7.1.2 column **2**), but also, in principle, through other channels involving breakup and inelastic processes. In particular, that in which one of the two neutrons moving in the $1/2^-$ orbital (component $|p_{1/2}^2(0)\rangle$ of (7.1.3), see also Potel et al. (2010), Fig 1 (f) of this reference) is picked up (forming a virtual deuteron, the other $1/2^-$ going into the continuum, i.e. breaking up from the ${}^9\text{Li}$ core) together with a neutron from the $p_{3/2}$ orbital, which eventually is filled by the continuum neutron falling into it, and leaving the core in the quadrupole excited state. This is in keeping with the fact that the main amplitude of this vibration is⁸ $X(1p_{3/2}^{-1}, 1p_{1/2})$ (≈ 1 ; see column **3** of Table 7.1.2). In the second process involving breakup, the proton field (${}^1\text{H}$) acting once breaks the halo Cooper pair (assumed again to be in the $|p_{1/2}^2(0)\rangle$ component), forcing one of the $1/2^-$ halo neutrons to populate a $p_{1/2}$ continuum state, the other $1/2^-$ neutron assumed to do the same while, acting for a second time, it picks up one of the neutrons moving in the continuum and another one from those occupying the $p_{3/2}$ orbital of ${}^9\text{Li}$, which is eventually filled by the second continuum $p_{1/2}$ neutron, leaving again the core in the quadrupole mode of excitation (see column **4** Table 7.1.2).

In the remaining channel considered, two-particles are transferred to the ground state of ${}^9\text{Li}$, and final state interaction (FSI) between the outgoing triton and ${}^9\text{Li}$ results in the inelastic and Coulomb excitation of the $1/2^-$ state (Table 7.1.2, column **5**).

Making use of the NFT spectroscopic amplitudes, and of computer codes developed to microscopically take into account the different processes mentioned above, that is the different reaction channels and continuum states up to 50 MeV of excitation energy, the corresponding transfer amplitudes and associated probabilities p_l were calculated. In Table 7.1.2 the probabilities $p_l = |S_l^{(c)}|^2$ associated with each of the processes mentioned above are displayed, where the amplitude $S_l^{(c)}$ is related to the total cross section associated with each of the channels c by the expression⁹

$$\sigma_c = \frac{\pi}{k^2} \sum_l (2l+1) |S_l^{(c)}|^2, \quad (7.1.4)$$

k being the wave number of the relative motion between the reacting nuclei.

In keeping with the small values of p_l , the contributions associated with the different reaction channels were calculated making use of second order perturbation theory. In the case of the $1/2^-$ (2.69 MeV) first excited state of ${}^9\text{Li}$ the absolute

⁸Barranco et al. (2001).

⁹Satchler (1980); Landau and Lifshitz (1965).

$l \backslash c$	1	2	3	4	5
0	4.35×10^{-3}	1.79×10^{-4}	4.81×10^{-6}	2.90×10^{-11}	3.79×10^{-8}
1	3.50×10^{-3}	9.31×10^{-4}	1.47×10^{-5}	1.87×10^{-9}	1.09×10^{-6}
2	7.50×10^{-4}	8.00×10^{-5}	2.45×10^{-5}	1.25×10^{-8}	1.21×10^{-6}
3	6.12×10^{-4}	9.81×10^{-5}	1.51×10^{-6}	6.50×10^{-10}	2.20×10^{-7}
4	1.10×10^{-4}	1.18×10^{-5}	2.21×10^{-7}	4.80×10^{-11}	1.46×10^{-8}
5	3.65×10^{-5}	2.16×10^{-7}	7.42×10^{-9}	6.69×10^{-13}	9.63×10^{-10}
6	1.35×10^{-5}	6.05×10^{-8}	2.88×10^{-10}	8.04×10^{-15}	1.08×10^{-11}
7	4.93×10^{-6}	7.78×10^{-8}	6.01×10^{-11}	4.05×10^{-16}	5.26×10^{-13}
8	2.43×10^{-6}	2.62×10^{-8}	7.4×10^{-12}	1.26×10^{-17}	9.70×10^{-11}

Table 7.1.2: Probabilities p_l associated with the processes described in the text for each partial wave l . The different channels are labeled by a channel number c . It can be seen that aside from channels **1** (successive) and **2** (simultaneous+non-orthogonality), leading to the direct population of ${}^9\text{Li}(gs)$ and ${}^9\text{Li}(1/2^-; 2.6 \text{ MeV})$ states (Fig. 7.1.2), all other channels provide a negligible contribution to the population of the ${}^9\text{Li}$ first excited state (see also Figs. 2.7.3 (b), (d), 2.7.4 (f) and 2.7.5).

differential cross section is thus,

$$\frac{d\sigma}{d\Omega}(\theta) = \frac{\mu^2}{16\pi^3\hbar^4} \left| \sum_l (2l+1)P_l(\theta) \sum_{c=2}^5 T_l^{(c)} \right|^2, \quad (7.1.5)$$

where μ is the reduced mass and $T_l^{(c)}$ are the transition matrix elements associated with the different channels and for each partial wave of DWBA¹⁰.

Making use of all the elements discussed above, multistep transfer¹¹, breakup and inelastic channels were calculated, and the results displayed in Figs. 7.1.3 and 7.1.4 and in Table 7.1.2. Theory provides an overall account of the experimental findings. In particular, in connection with the $1/2^-$ state, this result essentially emerges from cancellations and coherence effects taking place between the three terms contributing to the multistep two-particle transfer cross section (Fig. 7.1.4), weighted by the corresponding nuclear structure amplitudes (see Eqs. (7.1.1)–(7.1.3)). In fact, and as shown in Figs. 7.1.3 and 7.1.4, the contributions of breakup and inelastic processes to the population of the $1/2^-$ (2.69 MeV) first excited state of ${}^9\text{Li}$ are negligible as compared with the process depicted in Fig. 7.1.1. In the case of the breakup channels, this is a consequence of the low bombarding energy of the ${}^{11}\text{Li}$ beam (inverse kinematics), combined with the small overlap between

¹⁰Satchler (1980).

¹¹Bayman and Chen (1982), Igarashi et al. (1991), Bayman and Feng (1973) as well as Broglia and Winther (2004).

continuum (resonant) neutron $p_{1/2}$ wavefunctions and bound state wavefunctions. In the case of the final state inelastic process, it is again a consequence of the relative low bombarding energy. In fact, the adiabaticity parameters ξ_C, ξ_N ¹² associated with Coulomb excitation and inelastic excitation in the $t+{}^9\text{Li}$ channel are larger than 1, implying an adiabatic cutoff. In other words, the quadrupole mode is essentially only polarized during the reaction but not excited.

Second order calculations of inelastic, break up and final state interaction channels, which in principle can provide alternative routes for the population of the first excited state of ${}^9\text{Li}$ to the direct one predicted by the wavefunction (7.1.1) (β component), lead to absolute cross sections which are smaller by few orders of magnitude than that shown in Fig. 7.1.3. One then can posit that quadrupole core polarization effects in $|\text{gs}({}^{11}\text{Li})\rangle$ is essential to account for the observation of the $|1/2^-, 2.69 \text{ MeV}\rangle$ state, and provides evidence for phonon mediated pairing in nuclei.

7.1.3 Further examples of pairing vibrational states

Before concluding this section we provide in Fig. 7.1.5 further examples of pairing vibrational states around ${}^9\text{Li}$, ${}^{10}\text{Be}$, ${}^{48}\text{Ca}_{28}$ and ${}^{208}\text{Pb}_{126}$, $N = 6$, $N = 28$ and $N = 126$ neutron closed shell systems. The fact that among the (p,t) and (t,p) absolute differential cross sections one also finds the ${}^{208}\text{Pb}({}^{16}\text{O}, {}^{18}\text{O}){}^{206}\text{Pb}(\text{gs})$ absolute differential cross section is in keeping with the fact that the formalism to treat both light and heavy ions two–nucleon transfer reactions and their connection is well known¹³ and rather homogeneous¹⁴. Thus, it has been implemented in the software COOPER as a standard option (cf. App. 7.A). As observed, theory provides a quantitative account of experimental findings. Within this context see also Sect. 7.3.

7.2 Pairing rotational bands

The $|\text{BCS}\rangle$ state, which can be viewed as the intrinsic state with deformation α_0 , of a pairing rotational band (see Eq. (4.7.25)) is a coherent state (see Eq. (1.4.39)). One then expects the associated observable ($\sim |\alpha_0|^2$), namely the absolute two-nucleon transfer differential cross section, to be stable with respect to the models used to calculate the corresponding spectroscopic amplitudes B_ν . That is, $B_\nu U_\nu V_\nu$ and, as a consequence, the order parameter $\alpha_0 = \sum_{\nu>0} U_\nu V_\nu$.

In Table 7.2.1 we collect the two–nucleon spectroscopic amplitudes associated with the reactions ${}^{A+2}\text{Sn}(p,t){}^A\text{Sn}$, for A in the interval 112–126. Making use of these results and of global optical parameters (see Table 7.2.2), the absolute dif-

¹²See eqs. (IV.12) and (IV.14) of ref. Broglia and Winther (2004).

¹³Broglia and Winther (2004), Bayman and Chen (1982) and Thompson (1988) and references therein.

¹⁴Potel et al. (2013a,b).

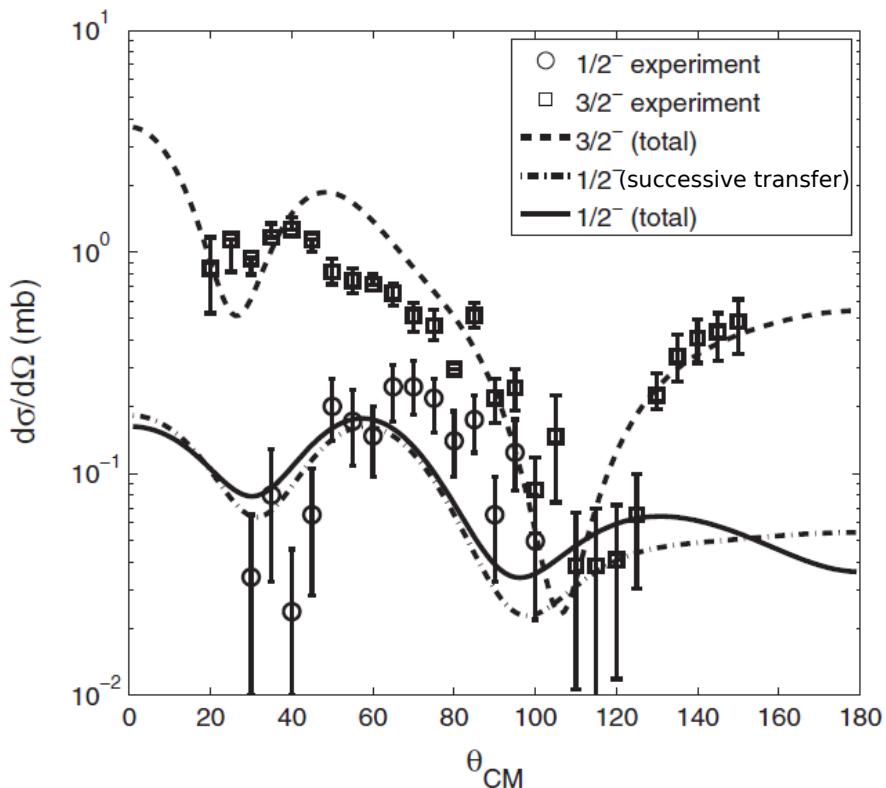


Figure 7.1.3: Experimental (Tanihata et al. (2008)) and theoretical differential cross sections (including successive transfer as well as breakup and inelastic channels, Potel et al. (2010)) of the ${}^1\text{H}({}^{11}\text{Li}, {}^9\text{Li}){}^3\text{H}$ reaction populating the ground state ($3/2^-$) and the first excited state ($1/2^-$; 2.69 MeV) of ${}^9\text{Li}$. Also shown (dash-dotted curve) is the differential cross section associated with this last state but taking into account only successive transfer. The optical potentials used are from (Tanihata et al., 2008; An and Cai, 2006), see Table 7.1.1. The absolute cross sections associated with the ground state ($3/2^-$) is predicted to be 6.1 mb (exp: 5.7 ± 0.9 mb) while that corresponding to the first excited state ($1/2^-$; 2.69 MeV) being 0.7 mb (exp: 1.0 ± 0.36 mb).

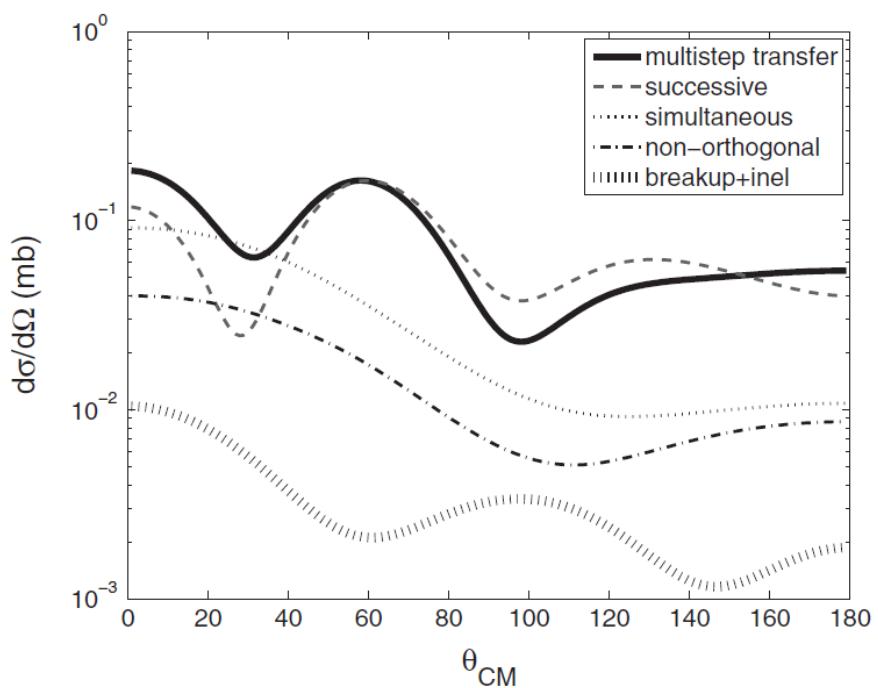


Figure 7.1.4: Successive, simultaneous and non-orthogonality contributions (prior representation) to the ${}^1H({}^{11}Li, {}^9Li){}^3H$ differential cross section associated with the population of the $1/2^-$ state of 9Li , displayed in Fig. 7.1.3. Also shown is the (incoherent) sum of the breakup ($c = 3$ and 4) and inelastic ($c = 5$) channel contributions.

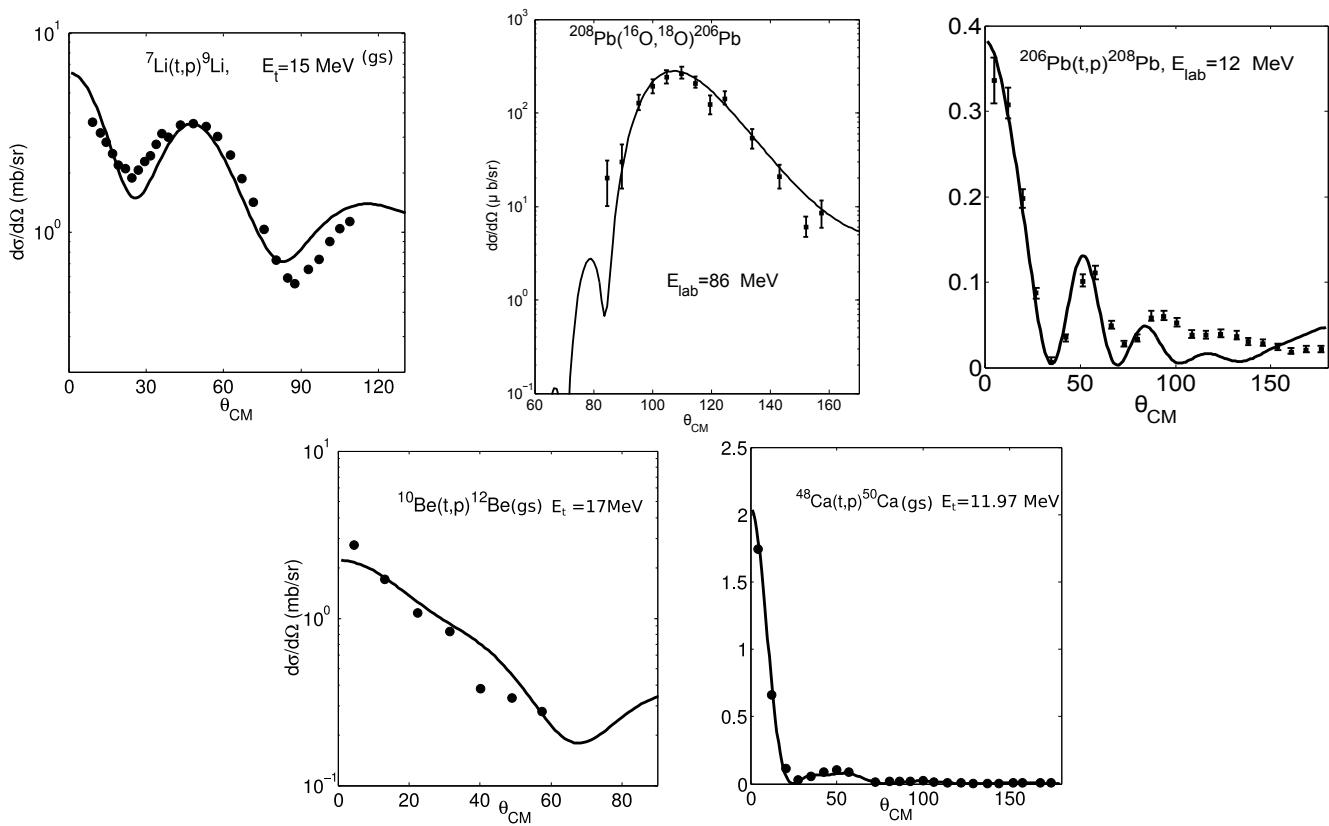


Figure 7.1.5: Absolute two-particle transfer differential cross sections for a number of reactions around closed shell associated with monopole pair addition and removal modes, and populating the ground state of the final system. Making use of spectroscopic amplitudes calculated as described in Sect. 3.5 in the particular case of $N = 6$ (Li,Be), $N = 48$ (Ca) and $N = 126$ (Pb), of global optical parameters and of the software COOPER, the absolute differential cross sections were calculated and are displayed in comparison with the experimental data (after Potel et al. (2013a)).

ferential cross section ${}^{A+2}\text{Sn}(p,t){}^A\text{Sn(gs)}$ were calculated. They are shown in Fig. 7.2.1 in comparison with the data.

	${}^{112}\text{Sn}$	${}^{114}\text{Sn}$	${}^{116}\text{Sn}$	${}^{118}\text{Sn}$	${}^{120}\text{Sn}$	${}^{122}\text{Sn}$	${}^{124}\text{Sn}$
$1d_{5/2}$	0.664	0.594	0.393	0.471	0.439	0.394	0.352
$0g_{7/2}$	0.958	0.852	0.542	0.255	0.591	0.504	0.439
$2s_{1/2}$	0.446	0.477	0.442	0.487	0.451	0.413	0.364
$1d_{3/2}$	0.542	0.590	0.695	0.706	0.696	0.651	0.582
$0h_{11/2}$	0.686	0.720	1.062	0.969	1.095	1.175	1.222

Table 7.2.1: Two-nucleon transfer spectroscopic amplitudes $\langle BCS(N)|T_\nu^{r\dagger}(j_\nu^2(0))|BCS(N+2)\rangle = \sqrt{(2j_\nu + 1)/2} U'_\nu(N)V'_\nu(N+2)$, associated with the reactions connecting the ground states (members of a pairing rotational band) of two superfluid Sn-nuclei ${}^{A+2}\text{Sn}(p,t){}^A\text{Sn(gs)}$.

	${}^A\text{Sn}(p,t){}^{A-2}\text{Sn}$											
	V	W	V_{so}	W_d	r_1	a_1	r_2	a_2	r_3	a_3	r_4	a_4
$p, {}^A\text{Sn}^a)$	50	5	3	6	1.35	0.65	1.2	0.5	1.25	0.7	1.3	0.6
$d, {}^{A-1}\text{Sn}^b)$	78.53	12	3.62	10.5	1.1	0.6	1.3	0.5	0.97	0.9	1.3	0.61
$t, {}^{A-2}\text{Sn}^a)$	176	20	8	8	1.14	0.6	1.3	0.5	1.1	0.8	1.3	0.6

Table 7.2.2: Optical potentials used in the calculations of the absolute differential cross sections displayed in Fig. 7.2.1.

7.2.1 Structure-reaction: stability of the order parameter α_0

The order parameter associated with distortion in gauge space can be written as

$$a'_0 = \sum_{j_a} \sqrt{\frac{2j_a + 1}{2}} B(j_a^2(0), N \rightarrow N+2), \quad (7.2.1)$$

where

$$B(j_a^2(0), N \rightarrow N+2) = \sqrt{\frac{2j_a + 1}{2}} U'_{j_a}(N)V'_{j_a}(N+2) \quad (7.2.2)$$

is the two-nucleon spectroscopic amplitude associated with the ($j^2(0)$) pair transfer between members of a pairing rotational band. Thus

$$\alpha'_0 = \sum_{j_a, m_a > 0} U'_{j_a} V'_{j_a} = \sum_{j_a} \frac{2j_a + 1}{2} U'_{j_a} V'_{j_a} = e^{2i\phi} \sum_{j_a} \frac{2j_a + 1}{2} U_{j_a} V_{j_a} = e^{2i\phi} \alpha_0, \quad (7.2.3)$$

defines a privileged orientation in gauge space. *Within the unified description of structure and reactions of the present monograph, the quantities (7.2.2) are the*

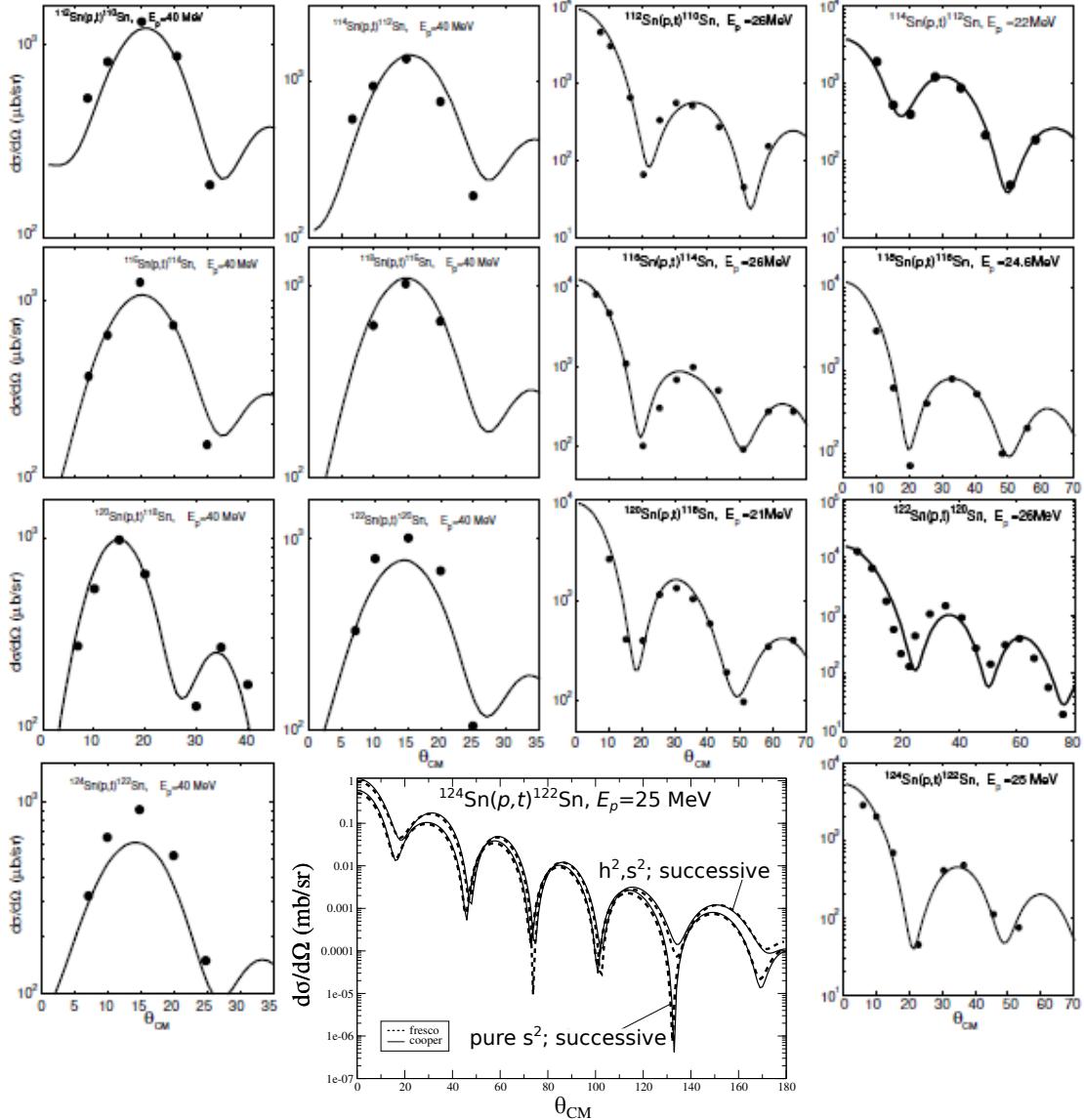


Figure 7.2.1: Predicted (continuous curve, Potel et al. (2013a,b)) absolute differential $A+2^{\text{Sn}}(p,t)^A\text{Sn}(\text{gs})$ cross sections for bombarding energies $E_p=40$ MeV (in the two left columns) and $21 \text{ MeV} \leq E_p \leq 26 \text{ MeV}$ (in the two right columns) in comparison with the experimental data (solid dots, Bassani et al. (1965), Guazzoni et al. (1999), Guazzoni et al. (2004), Guazzoni et al. (2006), Guazzoni et al. (2008), Guazzoni et al. (2011), Guazzoni et al. (2012)). In the center of the lowest row, the absolute differential cross section (successive transfer displayed in the angular interval $0^\circ \leq \theta_{CM} \leq 180^\circ$) associated with the process $^{124}\text{Sn}(p,t)^{122}\text{Sn}(f)$ and calculated with both the software COOPER (dashed bold) and FRESCO (thin continuous line) (Thompson (1988)) are displayed. This last calculation has been performed both for the $B_{1/2}s_{1/2}^2(0)$ contribution alone, and for the contribution corresponding to the $(B_{1/2}s_{1/2}^2(0) + B_{11/2}h_{11/2}^2(0))$ configuration, where the B coefficients are the spectroscopic amplitudes reported in the last column of Table 7.2.1.

weighting factors of the successive, simultaneous and non-orthogonality formfactors (Eq. 7.4.1) involved in the calculation of the corresponding transfer amplitudes and associated absolute transfer differential cross sections, making use of global optical parameters. Consequently, in discussing the properties of the order parameter α_0 , in particular in this section, one has in mind the two-nucleon transfer formfactors. The B -coefficients weight the corresponding radial functions, as well as the contributions of energy denominators (Green functions) associated with the different contributions and intermediate channels.

Empirical renormalization is also operative in connection with these channels, in terms of selecting the associated *effective Q-value*. Also of the distorted waves, again by selecting among possible embodiments of the *optical potential*. To be noted that, within this connection, one is talking about modest changes in the absolute differential cross sections to be used to predict new experiments, or to compare with the experimental data, as compared with the results obtained with standard choices (see App. 7.A, Fig. 7.A.1).

7.2.2 A two-nucleon transfer physical sum rule

In what follows we analyze the stability of a'_0 within the framework of three schemes to calculate the B -amplitude¹⁵ associated with the reaction $^{120}\text{Sn}(p, t)^{118}\text{Sn}(\text{gs})$. The first one corresponds to BCS approximation making use of a pairing interaction of constant matrix elements. Starting from the HF solution of a Skyrme interaction, namely Sly4, the gap and number equations are solved in the pairing approximation with $G = 0.26$ MeV leading to the empirical value of the three-point expression of the pairing gap $\Delta^{\exp} \approx 1.4$ MeV. The B -coefficients for the valence orbitals are reported in the fourth column of Table 7.2.3.

In the second calculational scheme, and making use of the same Skyrme interaction and of the v_{14} Argonne, 1S_0 NN -potential and neglecting the influence of the bare pairing force in the mean field, the HFB equation was solved. As a result, this step corresponds to an extended BCS calculation over the HF basis, allowing for the interference between states of equal quantum numbers $a (\equiv l j)$, but different number of nodes (k, k'). One includes (N_a) states (for each a) up to ≈ 600 MeV, to properly take into account the repulsive core of v_{14} and be able to accurately calculate Δ^{HFB} ($= 1.08$ MeV). The resulting B -coefficients are displayed in the third column of Table 7.2.3.

Going beyond mean field and including the particle-vibration coupling leading to retardation phenomena, both in the quasiparticle self-energy and in the induced pairing interaction, within the framework of renormalized NFT ((NFT)_{ren}) propagating the dressing of the quasiparticle states with the help of the Nambu-Gorkov (NG) equation leads, in the canonical basis, to the renormalized spectroscopic amplitudes shown in the second column of Table 7.2.3.

Making use of these B -coefficients, together with the global optical parameters

¹⁵Potel et al. (2017).

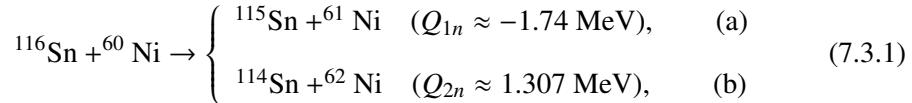
$a \equiv \{l_j\}$	NFT(NG)	HFB v_{14}	BCS(G)
$d_{5/2}$	0.22	0.29	0.41
$g_{7/2}$	0.46	0.47	0.57
$s_{1/2}$	0.37	0.34	0.41
$d_{3/2}$	0.59	0.60	0.66
$h_{11/2}$	0.95	1.0	1.03

Table 7.2.3: Two-nucleon spectroscopic amplitudes associated with the reaction $^{120}\text{Sn}(p, t)^{118}\text{Sn}(\text{gs})$. Note the small difference between the values of the fourth column with those reported in the column labeled ^{120}Sn of Table 7.2.1. They are due to a small difference in the value of the pairing coupling constant used here (Potel et al. (2017)) and that employed in Potel et al. (2017) and Potel et al. (2013b).

reported in Table 7.2.2, the corresponding absolute differential cross sections associated with the reaction $^{120}\text{Sn}(p, t)^{118}\text{Sn}(\text{gs})$ at 21 MeV of bombarding energy were calculated. They are displayed in Fig. 7.2.2 in comparison with the experimental findings. In keeping with the unified structure-reaction physical interpretation of α_0 mentioned following Eq. (7.2.3) (see also first paragraph of Sect. 7.2), and the results shown in Figs 7.2.2 and 7.2.3, namely the fact that the absolute cross section ratio $|\sigma_i - \sigma_{exp}|/\sigma_{exp}$ is equal to 0.09, 0.13 and 0.07 ($i = \text{BCS}, \text{HFB}, \text{NFT(NG)}$), one can posit that the relative errors of the associated two-nucleon transfer amplitudes $\alpha_0(\sim \sqrt{\sigma})$ are 4.5%, 6.5% and 3.5%. In other words, spontaneous breaking of gauge symmetry, a feature which is embodied in the three descriptions used (BCS, HFB, NFT(NG)) albeit, at very different levels of many-body refinement, seems to give rise to a new emergent property: a physical sum rule resulting from the intertwining of structure both in configuration- and in 3D-space (spectroscopic amplitudes and radial wavefunction) and reaction aspects of pairing in nuclei, and closely connected with the fact that $|\text{BCS}\rangle$ is a coherent state.

7.3 Heavy ion reactions between superfluid nuclei and correlation length

One- and two- neutron transfer reactions between superfluid nuclei have been studied with magnetic and gamma spectrometers heavy ion reactions between superfluid nuclei¹⁶,



for eight bombarding energies within the energy range $140.60 \text{ MeV} \leq E_{cm} \leq 167.95 \text{ MeV}$. That is, from energies above the Coulomb barrier ($E_B = 157.60 \text{ MeV}$), to

¹⁶Montanari et al. (2014, 2016).

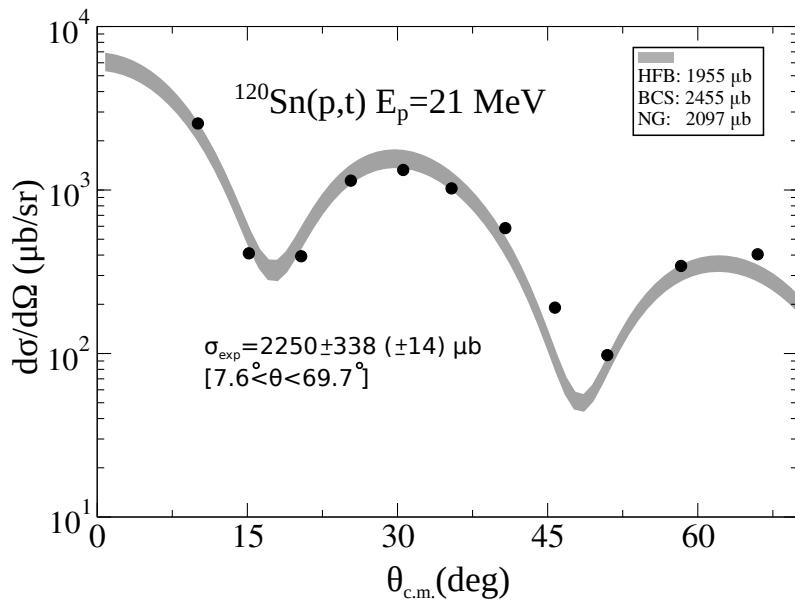


Figure 7.2.2: Absolute differential cross sections associated with the reaction $^{120}\text{Sn}(p,t)^{118}\text{Sn}(\text{gs})$ calculated making use of the BCS, HFB and renormalised NFT(NG) spectroscopic amplitudes and global optical parameters (grey thin area), in comparison with the experimental findings (solid dots) Guazzoni et al. (2008).

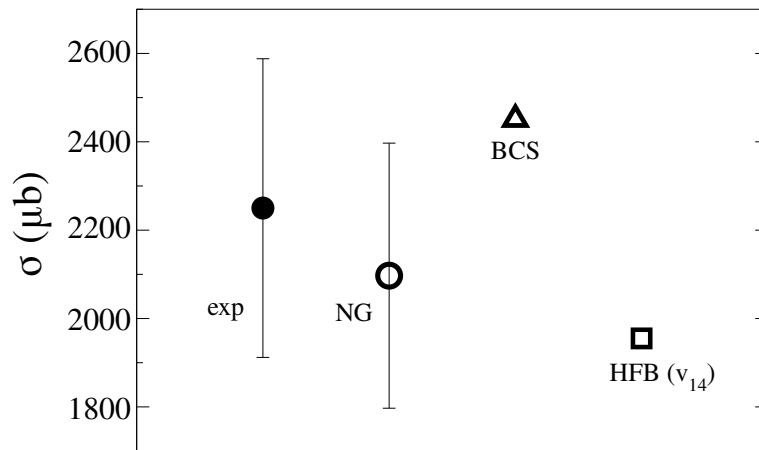


Figure 7.2.3: Integrated absolute cross sections associated with the reaction $^{120}\text{Sn}(p,t)^{118}\text{Sn}(\text{gs})$. The error ascribed to the NFT(NG) theoretical result stems from the variation the spin contribution of the different Skyrme interactions induce in the B -coefficients.

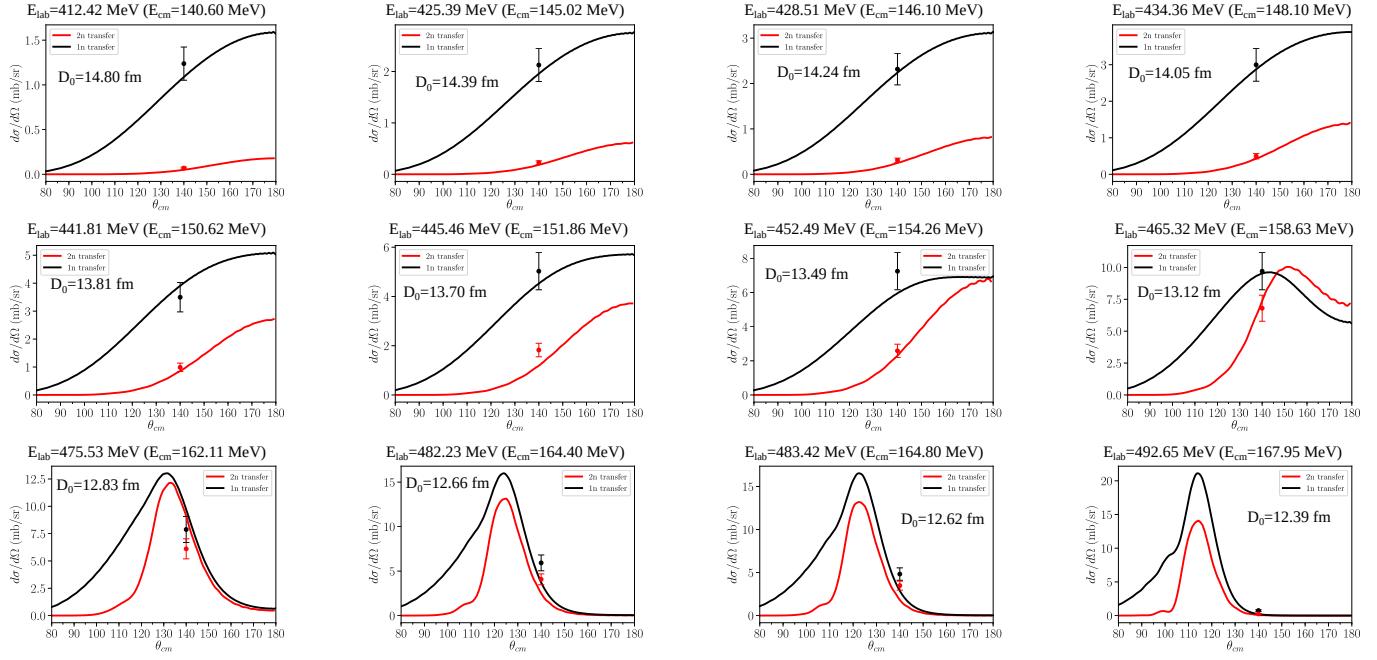


Figure 7.3.1: Angular distribution associated with the reactions $^{60}\text{Ni} + ^{116}\text{Sn} \rightarrow ^{62}\text{Ni} + ^{114}\text{Sn}$ (continuous line) and $^{60}\text{Ni} + ^{116}\text{Sn} \rightarrow ^{61}\text{Ni} + ^{115}\text{Sn}$, (dashed line) in comparison with the experimental data (Montanari et al., 2014).

well below it. The Cooper pair transfer channel (7.3.1 (b)) is dominated by the ground-ground state transition, while the theoretical calculation of the differential cross section associated with channel (7.3.1 (a)) indicates the incoherent contribution of about ten quasiparticle states of ^{61}Ni lying at energies $\lesssim 2.640$ MeV. A finding which is consistent with the fact that the value of the pairing gap of $^{60,62}\text{Ni}$, is of the order of 1.3–1.5 MeV. That is, in the case of the reaction 7.3.1 (a), we are in presence of a S-Q like transfer (see Fig. 4.6.1).

The analysis of the data associated with the reactions (7.3.1) carried out in Montanari et al. (2014, 2016) makes use of a powerful semiclassical approximation in which also the optical potential was calculated. The short wavelength of relative motion (de Broglie reduced wavelength $\lambda = 0.36/2\pi \text{ fm} \approx 0.06 \text{ fm}$), allows to accurately determine the distance of closest approach. Making use of the U, V occupation amplitudes for both Sn and Ni, as well as the optical potential given in Montanari et al. (2014), we have calculated, within the framework of first and second order DWBA, the absolute one- and two- nucleon transfer differential cross sections. In the second case, including both successive (dominant channel) and simultaneous transfer, properly corrected by non-orthogonality. Theory is compared with experiment in Fig. 7.3.1. As expected, it provides an overall account of the experimental findings.

As it emerges from direct inspection of Fig. 7.3.1, the distance of closest ap-

proach lying within the interval $13.12 \text{ fm} \leq D_0 \leq 13.49 \text{ fm}$, is the largest one for which $d\sigma/d\Omega|_{2n}$ is, within a factor of about 0.6 ($\approx (\pi/4)^2$) of the same order of $d\sigma/d\Omega|_{1n}$. In keeping with Eq. (4.5.12) ($v_F/c \approx 0.27$, $\Delta \approx 1.3 - 1.5 \text{ MeV}$) and Eq. (4.3.1) one can posit that the above interval provides a sensible bound to the size of the nuclear Cooper pair correlation length. Already increasing D_0 by 0.6 fm ($D_0 = 14.05$) σ_{1n} becomes a factor 6 larger than σ_{2n} . A consequence of forcing Cooper pair partners to tunnel across a barrier of “width” (distance of closest approach) $D > \xi$, leading to a strain which plays a role similar to that played by the critical bias $V_{eq} \approx 2\Delta/e$ (that is giving to the center of mass of the Cooper pairs a momentum $q \approx \hbar/\xi$) and resulting in the rupture of the transferred Cooper pair and in the unfreezing of the quasiparticle degrees of freedom. In other words, resulting in the transition from the S-S to the S-Q transfer regimes (Fig. 4.6.1). Within this context we refer to Eq. (4.9.6) and Fig. (4.9.2).

7.4 The structure of “observable” Cooper pairs

In his Waynflete lectures on Cause and Chance, Max Born¹⁷, to whom we owe the statistical interpretation of quantum mechanics¹⁸, states “... quantum mechanics does not describe an objective state in an independent external world, but the aspect of this world gained by considering it from a certain subjective standpoint, or with certain experimental means and arrangements”. It is within this context that we tried in previous sections to get insight concerning the structure of nuclear Cooper pairs. Specifically, in terms of two-nucleon transfer reactions. Being even more *subjective* (concrete), we were interested in shedding light on the structure of one of the 5-6 Cooper pairs participating in the BCS condensate (intrinsic state in gauge space) of the Sn-isotopes (ground state rotational band¹⁹) through pair transfer processes, both in light and in heavy ion reactions. That is $A+2\text{Sn}(p,t)^A\text{Sn(gs)}$ processes in general, and $^{120}\text{Sn}(p,t)^{118}\text{Sn}$ in particular, and also $^{116}\text{Sn} + ^{60}\text{Ni} \rightarrow ^{114}\text{Sn} + ^{62}\text{Ni}$. From a strict observational perspective, concerning Cooper pairs, one can only refer to the information two-nucleon transfer absolute differential cross sections carry on these entities. On the other hand, leaving out the discussion regarding the microscopic calculation of the optical potential, the carriers mediating information between structure and absolute differential cross sections, e.g. between target and outgoing particle in a standard laboratory setup, are the distorted waves. These functions can be studied independently of the transfer processes under consideration, in elastic scattering experiments. Consequently, the non-local, correlated formfactors (see Figs. 4.1.1 and 4.1.2),

$$F(\mathbf{r}_1, \mathbf{r}_2, \mathbf{r}_{Ap}) = F_{succ}(\mathbf{r}_1, \mathbf{r}_2, \mathbf{r}_{Ap}) + F_{sim}(\mathbf{r}_1, \mathbf{r}_2, \mathbf{r}_{Ap}) + F_{NO}(\mathbf{r}_1, \mathbf{r}_2, \mathbf{r}_{Ap}), \quad (7.4.1)$$

¹⁷Born (1948).

¹⁸Born (1964); Pais (1986).

¹⁹Potel et al. (2013b, 2017)

sum of the corresponding functions associated with successive and simultaneous transfer and with the non-orthogonality correction, and calculated with a variety of different sets of two-nucleon spectroscopic amplitudes (the same in each case for the three F -functions appearing in the rhs of Eq. (7.4.1)), can be compared at profit to each other. This is in keeping with the fact that they can be related, in an homogeneous fashion, with the absolute cross sections.

It is also of notice, that the spatial dimension, structure, non-locality and ω -dependence of the function (7.4.1) is expected to be rather different from that of the structure wavefunction of the Cooper pair, observed e.g. in high energy electron scattering or (p, pn) reactions, a question closely connected with linear response. While this concept has been, and continues to be quite useful in the study of many-body systems, it is a subtle one. In direct two-nucleon transfer reactions induced by both light and heavy (grazing) ion collisions, the contact between the two interacting nuclei is weak. Nonetheless, even a very low (normal) density overlap between target and projectile may induce important modifications in Cooper pairs. In particular, allow pairing correlated partner nucleons to profit from the enlarged volume compared to that available in the target or projectile nucleus. As a consequence the pair can expand, each of the partner nucleons be mainly in a different nucleus, and successive become the dominant transfer mechanism.

It can be stated that this picture is again an example of the fruitfulness of *linear response* to shed light on subtle questions regarding many-body systems. In the case under discussion, it allows the partners of the nuclear Cooper pair to correlate over dimensions larger than nuclear dimensions and in so doing make their intrinsic structure “observable”, almost free of the strong pressures of the mean field²⁰. The above discussion is schematically illustrated in Fig. 7.4.1. With it one comes back to the original question (Sect. 2.1). Which are the proper variables to be used in an attempt to describe the nuclear system? Elementary modes of excitation is a valid choice. But because these modes are in interaction, the above choice is not sufficient (unique). An operative definition requires that also the specific probe, reaction or decay process is specified. In fact, if one were to study Cooper pairs through electron scattering (two-nucleon correlations), one would likely obtain a picture of the system as that marked by the small ellipses in Fig. 7.4.1 (see also Fig. 3.6.3). Thus, rather different from the one which emerges from the specific two-nucleon transfer process (large ellipse, correlation length ξ), as testified by the observations made as a function of the width of the Josephson-like junction (distance of closest approach) transiently established in the reaction $^{116}\text{Sn}+^{60}\text{Ni}$ populating the **S-S** and **S-Q** channels.

²⁰Within this context, one can refer to the need of being able to distinguish between right and left weakly coupled (dioxide-layer separated) superconductors, to be able to measure gauge phase difference through the Josephson effect. See also Magierski et al. (2017).

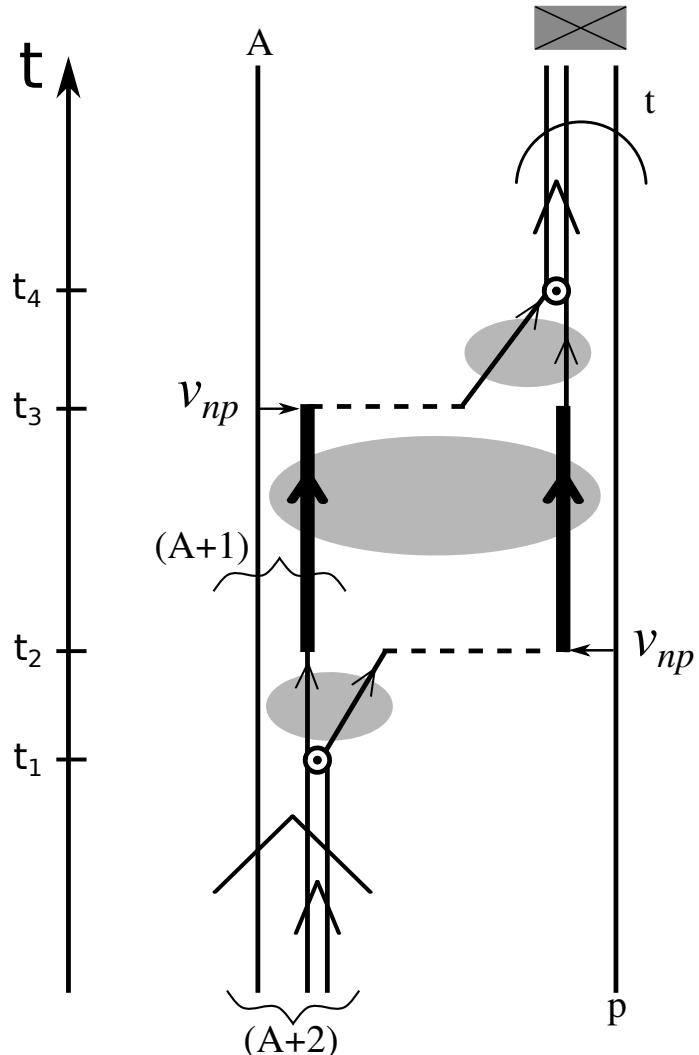


Figure 7.4.1: Diagram describing structure and reaction aspects of the main process through which a Cooper pair (di-neutron) tunnels from target to projectile in the reaction $(A + 2) + p \rightarrow A + t$. In order that the two-step process $(A + 2) + p \rightarrow (A + 1) + d \rightarrow A + t$ takes place, target and projectile have to be in contact at least in the time interval running between t_2 and t_3 . During this time, the two systems create, with local regions of ever so low nucleonic presence, a common density over which the non-local pairing field can be established, and the Cooper pair can be correlated. Even with regions in which the pairing interaction may be zero. Small ellipses (with linear dimensions of the order of the nuclear radius R_0) indicate situations in which the two neutron correlation is distorted by the external mean field of a single of the systems involved of the reaction, i.e. $A + 2$ in the entrance channel, t in the exit one (see e.g. Fig. 3.6.3). The large ellipse (with linear dimensions of the order of the correlation length ξ) indicate the region in which the two partners of the Cooper pair correlate over distances of the order of the correlation length ξ . It is this information that the outgoing particle of a Cooper pair transfer process brings to the detector. In other words, this is the observable Cooper pair in terms of its specific probe, and the reason why the neutrons are described, in the interval $\Delta t = t_3 - t_1$, in terms of bold face arrowed lines.

7.5 Closing the circle²¹

In one of the first references of this monograph²² entitled “Quantum mechanics of collision phenomena”, Born considers the elastic scattering of a beam consisting of N electrons which cross a unit area per unit time, scattered by a static potential. The stationary wavefunction describing the scattering process behaves asymptotically as,

$$e^{ikz} + f(\theta, \phi) \frac{e^{kr}}{r}, \quad \left(k = \frac{mv}{\hbar} \right). \quad (7.5.1)$$

The number of particles scattered into the solid angle $d\Omega = \sin \theta d\theta d\phi$ is given by $N|f(\theta, \phi)|^2 d\Omega$. To connect with Born notation one has to replace $f(\theta, \phi)$ by Φ_{mn} , where n denotes the initial state plane wave in the z -direction and m the asymptotic final state in which the waves move in the direction fixed by the angles (θ, ϕ) . Then Born writes that Φ_{mn} determines the probability for the scattering of the electron from the z - to the $(\theta\phi)$ -direction, adding a footnote in proof, as already mentioned, stating that a more precise consideration shows that the probability is proportional to the square of Φ_{mn} . In a second paper with the same title of the first²³ he states explicitly that the probability is to be connected with the modulus squared²⁴.

Within this context, the matrix element between the entrance and exit channel distorted waves of $F(\mathbf{r}_1, \mathbf{r}_2, \mathbf{r}_{Ap})$ is proportional to $f(\theta, \phi)$ and thus Φ_{mn} . The function F is not directly measurable, but the closer one can come of a theoretical construct connecting the Cooper pair ($s+r$) to experiment. For superfluid nuclei lying along the stability valley, this construct does not change much with the theory one uses to calculate the spectroscopic two-nucleon transfer amplitudes, provided they display off diagonal long range order (ODLRO)²⁵, a property closely related to phase coherence between Cooper pairs and deformation in gauge space. This is the reason why the associated absolute cross sections are rather stable, and α_0 provides a natural order parameter to measure deformation in gauge space.

On the other hand, in the case of halo, weakly bound exotic nuclei like ^{11}Li , the sensitivity of the absolute cross sections to the models used to calculate $F(\mathbf{r}_1, \mathbf{r}_2, \mathbf{r}_{Ap})$ is expected to be more relevant, and directly related to the conspicuous changes which renormalization produces in the radial shape of the single-particle wavefunctions of the halo neutrons forming the Cooper pair. This is in keeping with

²¹In this section we follow closely Pais (1986).

²²Born (1926a).

²³Born (1926b)

²⁴The motion of particles follows probability laws but the probability itself propagates according to the law of causality. And concerning the distinction between classical and quantal probabilities he states: “The classical theory introduces the microscopic coordinates which determine the individual processes only to eliminate them because of ignorance by averaging over their values; whereas the new theory get the same results without introducing them at all... We free the forces of their classical duty of determining directly the motion of particles and allow them instead to determine the probability of states”.

²⁵See footnote 42 Ch. 1. See also App. A, Potel et al. (2017) and refs. therein.

the fact that the nucleus is a finite many-body system, where surface effects play a central role, and light halo weakly bound exotic nuclei like ^{11}Be and ^{11}Li constitute paradigms of these leptodermic controlled entities. While the surface (S) to volume (V) ratio ($aS/V = 3a/R$, a being the diffusivity) is estimated to be 0.32 in the case of Sn, this value becomes 0.73 for Li.

Within this context, self-consistency in $(\text{NFT})_{\text{ren}}$ implies both the renormalization of single-particle energies and occupancies ($\tilde{\epsilon}_j, Z_j(\omega)$), as well as of radial wavefunctions ($\tilde{\phi}(r)^{(i)}$). For each $\tilde{\epsilon}_j$ there can be more than one radial function, depending on whether the nucleon is moving around the ground state ($i = \text{gs}$) or an excited state ($i = \text{coll}$) of the core. In keeping with the fact that $\tilde{\phi}(r)^{(i)}$ are the formfactors associated with transfer processes, self-consistency in $(\text{NFT})_{\text{ren}}$ is intimately related to unification of (s+r). In nuclei one has thus the unique possibility to study renormalization effects in terms of individual quantum states. Also pairing correlations resulting from bare plus induced 1S_0 interactions in systems with few (5-6) Cooper pairs (superfluid nuclei) down to the single Cooper pair (^{11}Li).

7.6 Summary

Nucleons moving in time reversal states lying close to the Fermi energy and interacting through the (1S_0 NN -potential) and the exchange of collective vibrations (induced pairing interaction), lead to $L = 0, S = 0$ nuclear Cooper pairs. Making use of the two-particle wavefunction describing them, it is found that the partner nucleons come closer to each other, as compared to the uncorrelated pair situation.

Cooper pairs were introduced in physics to describe low temperature superconductivity in metals, and constitute the building blocks of BCS. The BCS ground state wavefunction describes the condensation of these weakly bound (through the exchange of virtual photons, electron plasmons and lattice phonons), widely extended (dimensions much larger than typical distances between uncorrelated electrons), strongly overlapping (quasi) bosons (pair of electrons moving in time reversal states ($\nu, \tilde{\nu}$) with momentum and spin ($\mathbf{k} \uparrow, -\mathbf{k} \downarrow$)).

In presence of an external field, for example that associated with the tunneling interaction of a Josephson junction –involving energies smaller than the transition temperature and thus acting on the Cooper pairs center of mass– all pairs move in an identical (coherent) fashion, leading to a supercurrent across the junction of carriers of charge $q = 2e$. Its critical value being approximately equal to the (normal) single-electron current associated with a junction bias equal to the Cooper pair binding energy (2Δ) divided by the electron charge (e).

When the momentum imparted to the center of mass of the condensed Cooper pairs acquires a value of the order of the inverse of the mean square radius of the Cooper pairs (correlation length), the back-and-forth ($\ell = 0$) radial motion of electron partners (intrinsic ($\nu, \tilde{\nu}$) motion) cannot follow suit, getting out of step and resulting in the breaking (unbinding) of the pairs (critical current). As a consequence, supercurrent ceases and normal current ($q = e$) starts flowing.

The traditional view of nuclear Cooper pairs described above, seems to be partially at odds with that found at the basis of the BCS description of low temperature superconductivity. In particular concerning what can be considered the only (specific) coupling field able to detect Cooper pair (gauge) phase correlation, namely tunneling across a Josephson junction. In each of the two weakly coupled superconductors, pairs of electrons recede from each other as compared to the normal situation. In this way, they lower their relative kinetic energy of confinement, profiting best of the weak, attractive, (retarded) interaction. Being the electron pair phase correlated, with a correlation length much larger than the barrier thickness, although they tunnel one at a time, they do so as a single particle (of mass $2m_e$ and charge $2e$), and the probability of going through the junction is comparable to that of a single electron. It is like interference in optics, with phase coherent waves mixing.

In the transient, time dependent, Josephson-like junction established between two superfluid nuclei in a heavy ion collision about and below the Coulomb barrier, one finds that for bombarding energies leading to distances of closest approach smaller than, or of the order of the correlation length, quantity much larger than nuclear dimensions, the probability P_2 of pair transfer –process dominated by successive transfer where entangled nuclear Cooper pair partners are about a correlation length apart from each other– is about equal to that of one-nucleon transfer (P_1). For larger distances of closest approach, larger than the correlation length, the intrinsic back and forth relative motion characterizing the intrinsic structure of the transferred nuclear Cooper pair gets out of step, resulting in the transition from the superfluid-superfluid (S-S) transfer regime to the superfluid-quasiparticle (S-Q) one, as testified by the fact that P_1 becomes much larger than P_2 . Already for distances of closest approach about 0.6 fm larger than the nuclear correlation length, the absolute one-nucleon transfer cross section is about an order of magnitude larger than that associated with two-nucleon tunneling. The nuclear (S-S)→(S-Q) transition as a function of the distance of closest approach, where the Cooper pair is probed (stressed) beyond the correlation length, parallels that associated with the one undergoing by a low temperature superconductor where Cooper pairs are biased to move with a center of mass moment of the order of the inverse of the correlations length.

The gained homogeneity between the pictures of Cooper pairs in superconductors and in nuclei, is a consequence of fulfilling a basic requirement of quantum mechanics. Namely, that to specify the experimental setup in discussing a physical phenomena. In the present case one has tried to answer to the question what the structure of a nuclear Cooper pair is when probed in terms of its specific probe, namely that of the transfer (tunneling) of a pair of nucleons, comparing theory with experiment in terms of absolute cross sections.

Appendix 7.A One- and two-nucleon transfer computer codes

In this Appendix we provide a brief description of the numerical methods implemented in the computer codes **ONE** and **COOPER** written to evaluate the absolute differential cross sections associated with one- and two-nucleon transfer reactions.

COOPER²⁶: The two-nucleon transfer differential cross section is given by Eq. (6.1.3). The principal task consists in calculating the transfer amplitudes $T^{(1)}(\theta)$, $T_{succ}^{(2)}(\theta)$ and $T_{NO}^{(2)}(\theta)$ described in Eqs. 6.1.4a-6.1.4c, by numerically evaluating the corresponding integrals. The dimensionality of the integrals can be reduced by expanding in partial waves (eigenfunctions of the angular momentum operator) the distorted waves and wavefunctions present in the corresponding integrands. The resulting expressions are Eqs. (6.2.36) and (6.2.37) for $T^{(1)}(\theta)$, Eqs. (6.2.128), (6.2.129) and (6.2.130) for $T_{succ}^{(2)}(\theta)$, and Eqs. (6.2.154), (6.2.155) and (6.2.156) for $T_{NO}^{(2)}(\theta)$. The integrals are computed numerically with the method of Gaussian quadratures.

The one-dimensional (radial) functions (6.1.2) appearing in the integrands are defined in a spatial grid up to a given maximum radius r_{max} . The bound state wavefunctions are obtained by numerical integration of the radial Schrödinger equation for a Woods-Saxon potential with a spin-orbit term. Along with the optical potentials governing the relative nucleus-nucleus motion in the initial, intermediate, and final channels, they conform the basic structure input needed to define the calculation. They are generated by solving a one-body Schrödinger equation for a Woods-Saxon with standard geometry, where the radius and diffusivity parameters are given as an input. The depth of the potential is adjusted to reproduce the binding energy of the state under consideration, according to the prescription defined below. The resulting potential corresponding to the final (initial) nucleon bound state stands also for the interaction potential featured in the integrand in the prior (post) representation.

The binding energies of the single-particle wavefunctions determine the asymptotic behavior of the transfer formfactors, an essential ingredient for a quantitative account of the two-nucleon transfer cross section²⁷. According to the prescription followed in the calculations discussed in the present monograph, the orbital corresponding to the spin-parity of the $A + 1$ system will be bound by half the experimental two-neutron separation energy S_{2n} of the $A + 2$ nucleus. The energies of the remaining single-particle states relative to the $A + 1$ ground state are determined from a structure calculation. Let us illustrate this procedure with the $^{120}\text{Sn}(p, t)^{118}\text{Sn}$ reaction with a 21 MeV proton beam. In order to assess the impact of the binding energies of the single-particle orbitals in the absolute value of the transfer cross section, we also show a calculation in which all the single-particle wavefunctions are bound with half the two-neutron separation energy of ^{120}Sn . The

²⁶Potel (2012b).

²⁷For an example with a simple discussion of the problem see Fig. 6.4.4.

	^{120}Sn (I)	^{120}Sn (II)	B
$d_{5/2}$	-9.49	-7.79	0.41
$g_{7/2}$	-8.89	-7.79	0.57
$s_{1/2}$	-7.79	-7.79	0.41
$d_{3/2}$	-7.29	-7.79	0.66
$h_{11/2}$	-5.89	-7.79	1.03

Table 7.A.1: Energies (in MeV) of the single-particle orbitals involved in the description of the ground state of ^{120}Sn . The second column displays the energies of the single-particle orbitals according to the prescription used throughout this monograph (calculation (I)). The $s_{1/2}$ single-particle state, associated with the $1/2^+$ ground state of ^{119}Sn , has a binding energy of half the two-neutron separation energy of ^{120}Sn , $S_{2n}(^{120}\text{Sn})=15.59$ MeV. Calculation (II) (third column) corresponds to binding all the single-particle wavefunctions with half the two-neutron separation energy of ^{120}Sn . The last column displays the spectroscopic amplitudes B obtained from a BCS calculation of the ground state of ^{120}Sn (see Tab. 7.2.3).

binding energies used for the two calculations are collected in Table 7.A.1, and the corresponding results are shown in Fig. 7.A.1. When calculating the successive transfer amplitude, the energy of the Green's function of the intermediate channel $d + (A + 1)$ (Eq. 6.1.11) is taken at the experimental energy of each of the populated states of the $(A + 1)$ system. Off-shell effects due to the virtual population of the different $(A + 1)$ intermediate states are thus fully taken into account. In the study of ^{11}Li , a different approach have been followed concerning structure: the two-neutron wavefunction have been read from a two-dimensional numerical grid, generated from the NFT structure calculation.

The distorted waves are obtained by integrating the radial Schrödinger equation with positive energy from $r = 0$ to r_{max} , and matching the solution with the corresponding Coulomb wave function at a given $r = r_{match}$, big enough to lie outside of the range of the nuclear interaction. The Woods-Saxon optical potentials used to obtain the distorted waves consist on a real Coulomb term, a real and imaginary volume terms, an imaginary surface term, and a real and imaginary spin orbit terms. The parameters needed to specify all those terms are given as an input.

ONE²⁸: In order to obtain the one-particle transfer cross sections, one has to evaluate the amplitudes defined in Eq. (5.1.26). The corresponding numerical integrations are performed making use of Gaussian quadratures. The radial part of the single-particle wavefunctions (5.1.6) and of the partial waves (5.1.7) and (5.1.8), are generated in the way described above. The kinetic energies defining the distorting waves in the entrance and exit channels are defined by the bombarding energy in the center of mass frame, and the experimental Q -value of the reaction. The binding energies corresponding to the initial and final single-particle wavefunc-

²⁸Poté (2012a)

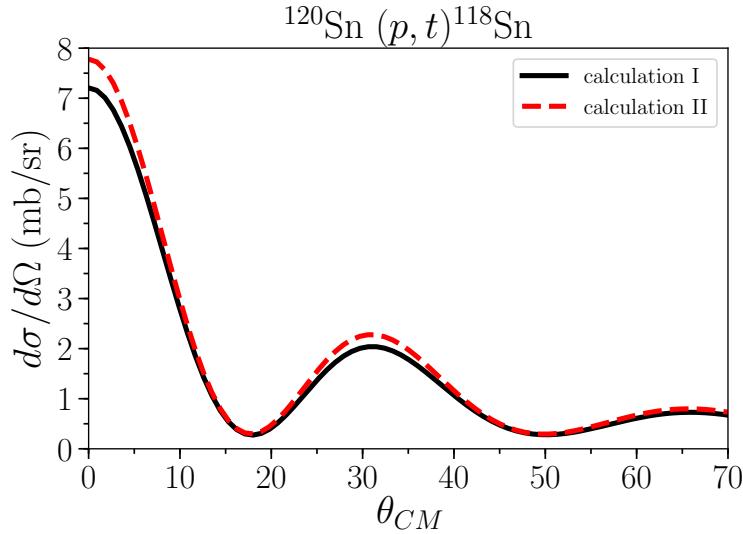


Figure 7.A.1: Numerical computation of the two-neutron transfer cross section corresponding to the reaction $^{120}\text{Sn}(p,t)^{118}\text{Sn}$. The two calculations shown in the figure correspond to the two different prescriptions for defining the binding energies of the single-particle orbitals described in the caption of Tab. 7.A.1.

tions are taken to be the experimental separation energy for each state. It is to note that the difference between the initial and final binding energy should be consistent with the reaction Q -value.

Appendix 7.B Statistics.

Consider two identical particles moving in a one-dimensional harmonic oscillator. It is assumed that one is in the ground state and the other is in the first excited state. According to the superposition principle

$$\Phi(x_1, x_2) = \lambda\phi_1(x_1)\phi_0(x_2) + \mu\phi_0(x_1)\phi_1(x_2). \quad (7.B.1)$$

We calculate the correlation of these particles, that is, the quantity

$$Corr = \frac{\langle x_1 x_2 \rangle - \langle x_1 \rangle \langle x_2 \rangle}{\sqrt{(\langle x_1^2 \rangle - \langle x_1 \rangle^2)(\langle x_2^2 \rangle - \langle x_2 \rangle^2)}} \quad (7.B.2)$$

Starting with

$$\begin{aligned}\langle x_1 x_2 \rangle &= \int dx_1 dx_2 (\lambda^* \phi_1^*(x_1) \phi_0^*(x_2) + \mu^* \phi_0^*(x_1) \phi_1^*(x_2)) \\ &\quad \times (x_1 x_2) (\lambda \phi_1(x_1) \phi_0(x_2) + \mu \phi_0(x_1) \phi_1(x_2)) \\ &= |\lambda|^2 \langle \phi_1 | x_1 | \phi_1 \rangle \langle \phi_0 | x_2 | \phi_0 \rangle + \lambda^* \mu \langle \phi_1 | x_1 | \phi_0 \rangle \langle \phi_0 | x_2 | \phi_1 \rangle \\ &\quad + \lambda \mu^* \langle \phi_0 | x_1 | \phi_1 \rangle \langle \phi_1 | x_2 | \phi_0 \rangle + |\mu|^2 \langle \phi_0 | x_1 | \phi_0 \rangle \langle \phi_1 | x_2 | \phi_1 \rangle.\end{aligned}\quad (7.B.3)$$

In keeping with the fact that

$$\langle \phi_1 | x | \phi_1 \rangle = \langle \phi_0 | x | \phi_0 \rangle = 0, \quad (7.B.4)$$

and

$$\langle \phi_0 | x | \phi_1 \rangle = \langle \phi_1 | x | \phi_0 \rangle = \sqrt{\frac{\hbar\omega}{2C}}, \quad (7.B.5)$$

one obtains

$$\langle x_1 x_2 \rangle = \left(\frac{\hbar\omega}{2C} \right)^2 \Re(\lambda^* \mu). \quad (7.B.6)$$

And

$$\sqrt{-} = \left(\frac{\hbar\omega}{2C} \right)^2, \quad (7.B.7)$$

for the denominator of Eq. (7.B.2). From the above results the correlation function between particle 1 and 2 is

$$Corr = 2\Re(\lambda^* \mu) = \begin{cases} 1 & (\lambda = +\mu = \frac{1}{\sqrt{2}}) \\ -1 & (\lambda = -\mu = \frac{1}{\sqrt{2}}) \end{cases} \quad (7.B.8)$$

It is of notice that, in quantum mechanics, average values imply the mean outcome of a large number of experiments. In this case, of the (simultaneous) measure of the position of the two particles²⁹.

Appendix 7.C Correlation length and quantality parameter.

The correlation length can be defined as³⁰

$$\xi = \frac{\hbar v_F}{\pi \Delta} \approx \frac{\hbar^2}{m} \frac{k_F}{\pi \Delta}, \quad (7.C.1)$$

²⁹Basdevant and Dalibard (2005).

³⁰See e.g. Annett (2013) p. 62.

where the Fermi momentum in the case of stable nuclei lying along the stability valley is

$$k_F \approx 1.36 \text{ fm}^{-1}. \quad (7.C.2)$$

Thus,

$$\xi = 40 \text{ MeV fm}^2 \times \frac{1.36}{\pi\Delta} \text{ fm}^{-1} \approx \frac{17}{\Delta} \text{ fm}, \quad (7.C.3)$$

and,

$$\xi \approx 14 \text{ fm}, \quad (\Delta \approx 1.2 \text{ MeV}). \quad (7.C.4)$$

Thus, the associated (generalized) quantality parameter is, in the present case,

$$q_\xi = \frac{\hbar^2}{2m\xi^2} \frac{1}{2\Delta} \approx 0.04. \quad (7.C.5)$$

That is, the two partner nucleons are, in the Cooper pair, rigidly correlated with each other.

We now consider $^{11}_3\text{Li}_8$, and calculate k_F (neutrons) with the help of the Thomas-Fermi model³¹

$$k_F = \left(3\pi^2 \frac{8}{\frac{4\pi}{3}(4.58)^3} \right)^{1/3} \text{ fm}^{-1} \approx \frac{(18\pi)^{1/3}}{4.58} \text{ fm}^{-1} \approx 0.8 \text{ fm}^{-1}. \quad (7.C.6)$$

The correlation length can, in the present case, be calculated in terms of the correlation energy ($E_{corr} \approx -0.5 \text{ MeV}$),

$$\xi \approx \frac{\hbar v_F}{\pi |E_{corr}|} \approx \frac{200 \text{ MeV} \times \text{fm} \times 0.16}{\pi \times 0.5 \text{ MeV}} \approx 20 \text{ fm}, \quad (7.C.7)$$

the resulting generalized quantality parameter being

$$q_\xi = \frac{\hbar^2}{2m\xi^2} \frac{1}{|E_{corr}|} \approx 0.1. \quad (7.C.8)$$

It is of notice that this result is but an alternative embodiment of the relation (4.4.8). One could argue that both (4.4.8) and (7.C.8) (as well as (7.C.5) for stable nuclei), are just a manifestation of (7.B.8). That there is more to it is forcefully expressed by the fact that, selecting the pure two-particle configuration $|s_{1/2}^2(0)\rangle$ ($|p_{1/2}^2(0)\rangle$) to describe the halo neutron Cooper pair of ^{11}Li leads to absolute two-particle transfer cross sections which are about one order of magnitude larger (smaller) than the observed cross section (see Fig. 7.1.2). The fact that the NFT result (7.1.1)–(7.1.3) reproduces observations within experimental errors, underscores the central role

³¹ Quantity which can be related to (v_F/c) according to $(v_F/c) = \hbar k_F/(mc) = (\hbar c/(mc^2))k_F \approx 0.2(k_F)\text{fm}^{-1}$. In the case in which $k_F \approx 0.8 \text{ fm}^{-1}$ (see Eq. (7.C.6)), $v_F/c \approx 0.16$.

played by structure on Cooper pair tunneling, through the emergent property of generalized pairing rigidity.

Summing up, both the bare NN - and the long range induced pairing interaction changes, under the circumstances that $E_{intrinsic} \ll 2\Delta$, $Q \ll \hbar/\xi$ and $D_0 \lesssim \xi$, the statistics of the elementary modes from fermionic to (quasi) bosonic ones³² and, at the same time, the value of the quantity parameter from $q \approx 1$ to $q_\xi \ll 1$. That is, from a regime of delocalized single nucleons to one of strongly overlapping, independent pair motion each being governed by the same phased wavefunction³³ $(U'_v + V'_v e^{-2i\phi} P_v^\dagger)|0\rangle$, and behaving as a single entity. The operator P_v^\dagger , being a product of two fermions, do not fulfill Bose statistics ($(P_v^\dagger)^2 = 0$). This property implies the presence of a pairing gap not only for breaking a pair, but also for making a pair move differently from the others³⁴. As a result one has long-range-order in the superfluid nuclear system, known as off-diagonal-long-range-order (ODLRO)³⁵. This effect leads to generalized gauge rigidity, the detailed renormalizing and dressing mechanisms ultimately deciding on the soundness and applicability of the description under discussion. The fact that in working out the reaction mechanism one uses, for practical reasons, a single-particle basis (second order DWBA corrected by non-orthogonality), reconstructing the pair correlations in term of sums over virtual states, is at the basis of the two-neutron transfer physical sum rule discussed in Sect. 7.2.2.

The extension of the validity of the picture of superconductivity in metals based on the motion of $\approx 10^6$ phase correlated Cooper pairs, down to few (5–8, ^{116}Sn) and eventually one Cooper pair (^{11}Li), seems short of a wonder. In particular in this last case, in which the binding of the Cooper pair is essentially due to a quantum-mechanical many-body retarded interaction resulting from the exchange of one of the most intriguing nuclear collective vibrations, at the edge between soft- $E1$ mode and Pygmy Dipole Resonance (see App. 7.D), the associated ZPF affecting, among other things, nuclear diffusivity and the neutron skin.

Within this context, the sense of wonder is common with the emergence of other unexpected quantum materials and macromolecules as a result of interactions resulting from ZPF, like the van der Waals interaction responsible, *inter alia*, for the van der Waals solid made out of C_{60} fullerene, as well as associated doped

³²Where $E_{intrinsic}(q)$ stands for the energy (momentum) given to the intrinsic (center of mass) motion of the transferred Cooper pair. The quantity D_0 is the distance of closest approach at which the two-nucleon transfer process takes place, ξ being the correlation length. Within this context see Sects. 3.4 and 4.7.1.

³³Since P_v^\dagger commutes for different v 's, $|BCS\rangle$ represents uncorrelated occupancy of the various pair states.

³⁴The BCS state can be written as $|BCS\rangle \sim \prod_{v>0} \tilde{\alpha}_v \alpha_v |0\rangle$, where $\alpha_v = U_v a_v - V_v a_v^\dagger$, is the annihilation quasiparticle operator (Bogoliubov (1958a); Valatin (1958); see App. G of Brink and Broglia (2005)). Thus, to make e.g. the pair associated with the state v' to behave differently from the rest of the pairs one has to avoid including in the product above the state v' , i.e. $\prod_{v \neq v'} \tilde{\alpha}_v \alpha_v |0\rangle$. The resulting state resembles a two-quasiparticle state.

³⁵See footnote 42 Ch. 1.

superconductors (Rb_3C_{60} , K_3C_{60} , $\text{Rb}_2\text{CsC}_{60}$, etc.)³⁶, let alone the Casimir interaction, and the related (strong)³⁷ hydrophobic force. Force which plays an essential role in the folding of the macromolecules responsible for metabolism³⁸, namely proteins (see App. 7.E).

Appendix 7.D Multipole pairing vibrations.

Although much work has been carried out concerning multipole pairing vibrations³⁹, i.e. modes with transfer quantum numbers $\beta = \pm 2$ and multipolarity and parity λ^π different from 0^+ , this remains a chapter essentially missing from the subject of pairing in nuclei. Arguably, with the partial exception of quadrupole pairing studied in the multiphonon pairing vibrational spectrum⁴⁰ around the closed shell ^{208}Pb , and in connection with strongly excited 0^+ pairing vibrational states in the actinide region⁴¹, and in the quadrupole and hexadecapole pairing vibrations in the multiplet spectrum⁴² of ^{209}Bi .

In what follows, we elaborate on the new insight on pairing vibrational modes, the studies of two-neutron pickup reactions on ^{11}Li have opened. As already explained, because of the small overlap existing between halo neutrons and core nucleons both the 1S_0 , NN -interaction and the symmetry-potentials become strongly screened, resulting in a subcritical value of the pairing strength and in a weak repulsion to separate protons from neutrons in the dipole channel. As a result, neither the $J^\pi = 0^+$ correlated neutron state (Cooper pair), nor the $J^\pi = 1^-$ one (vortex-like)⁴³ pair addition modes are bound (although both qualify to do so) to the $N = 6$ closed shell system (^9Li).

Being the bare NN -pairing interaction subcritical, the two neutrons can correlate their motion by exchanging vibrations of the medium in which they propagate, namely the halo. These modes can hardly be the $\lambda = 2^+, 3^-$ or 5^- surface vibrations found in nuclei lying along the stability valley. This is because the diffusivity of the halo is consistent, and it blurs the very definition of surface. It is of notice that

³⁶Hebard et al. (1991); Rosseinsky et al. (1991); Holczer et al. (1991); Fleming et al. (1991); Anderson (1991); Zhou et al. (1992); Hebard (1992); Gunnarsson (1995, 1997, 2004), see also Broglia et al. (2004) and refs. therein.

³⁷Chandler (2002).

³⁸Dyson (1999).

³⁹See Brink and Broglia (2005) Sect. 5.3 and refs. therein. See also Broglia et al. (1974), Ragnarsson and Broglia (1976), Broglia et al. (1971a), Broglia et al. (1971b), Bès and Broglia (1971b), Bès and Broglia (1971a), Flynn et al. (1971), Bès et al. (1972), Broglia (1981), Bohr and Mottelson (1980), Flynn et al. (1972a), Bortignon et al. (1976); see also Kubo et al. (1970) and references therein.

⁴⁰See Flynn et al. (1972a).

⁴¹Casten et al. (1972), Bès et al. (1972), Ragnarsson and Broglia (1976). It is of notice that β -vibrations and monopole pairing vibrations become mixed in quadrupole deformed nuclei.

⁴²Bortignon et al. (1976).

⁴³One of the effects of a vortex is that it allows rotation of a quantum system to take place about an axis of symmetry.

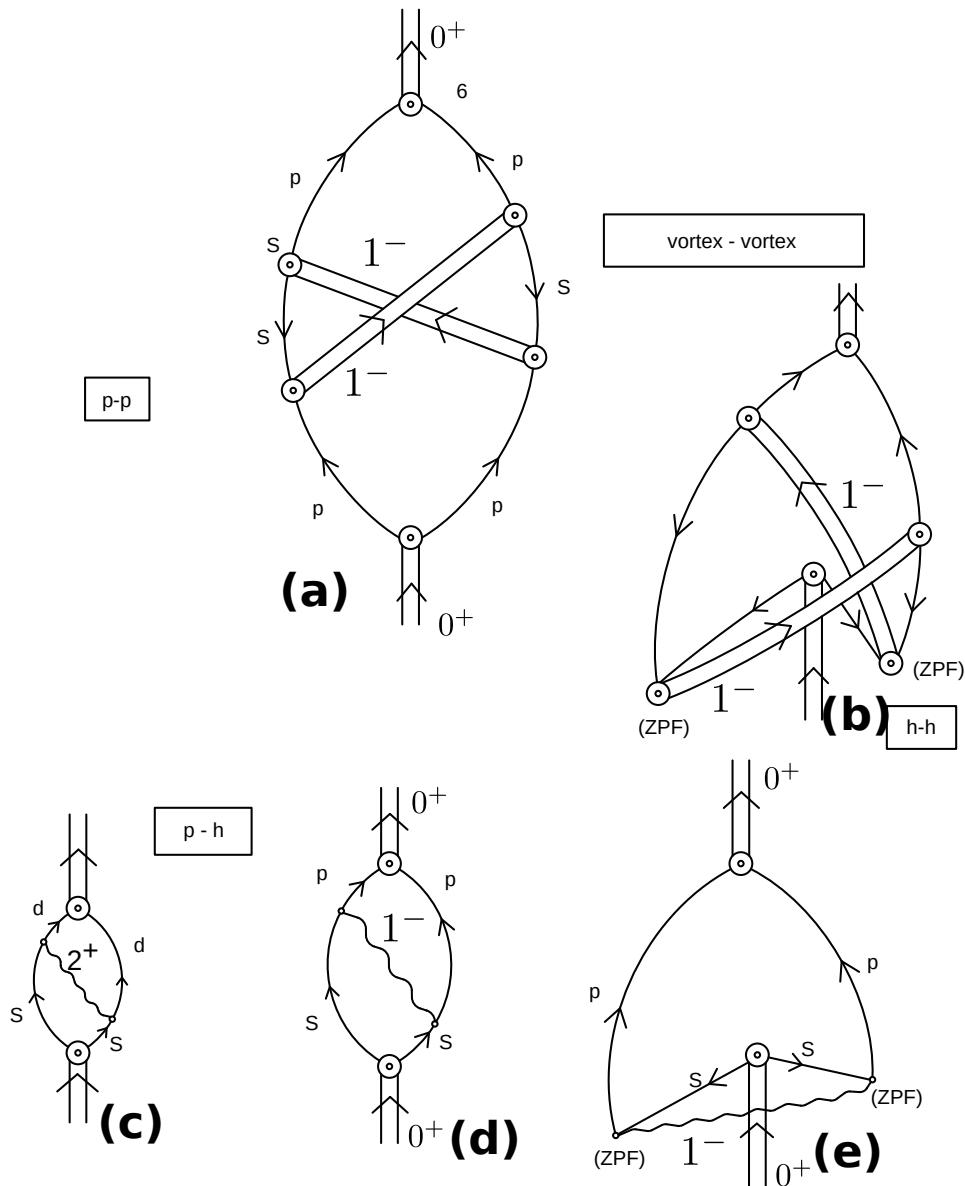


Figure 7.D.1: NFT-Feynman diagrams describing the interweaving between the neutron halo pair addition monopole and dipole modes (double arrowed lines labeled 0^+ and 1^- respectively). Above, the exchange of dipole modes binding the 0^+ pair addition mode through forwards going particle-particle p-p (h-h) components. Below, the assumption is made that the PDR of ^{11}Li can be viewed as a p-h (two quasiparticle), QRPA mode.

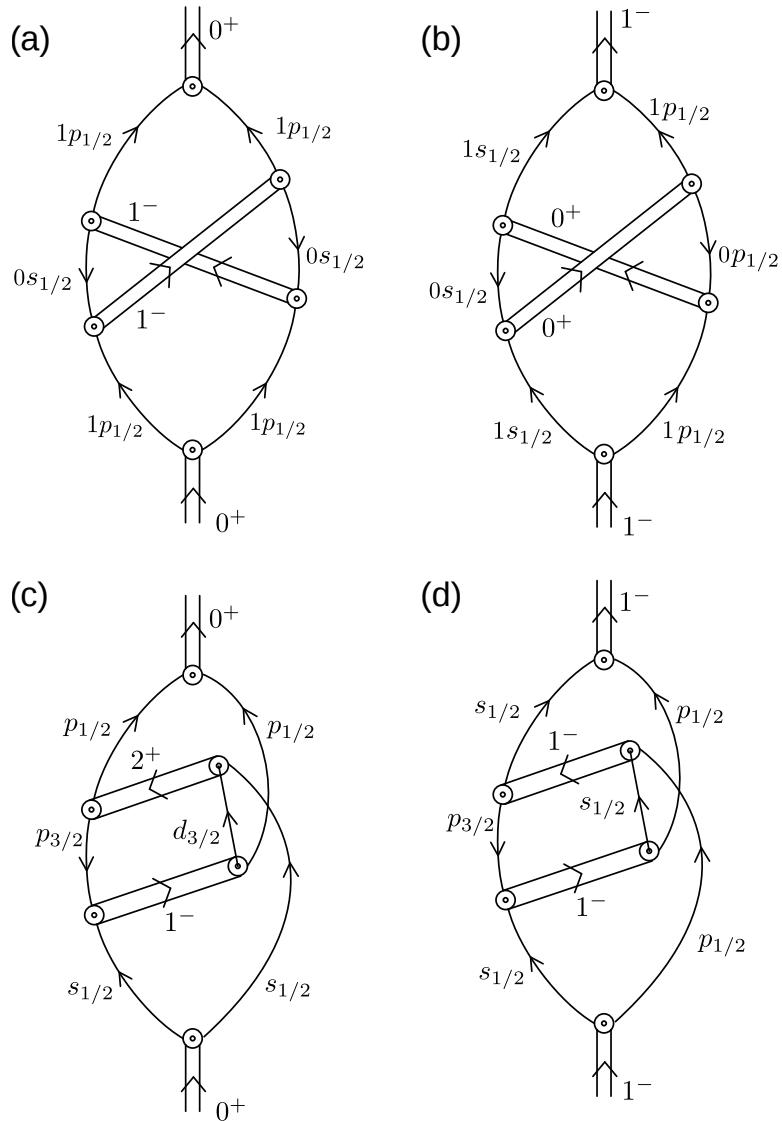


Figure 7.D.2: NFT-Feynman diagrams describing, (a,c) some of the particle-particle (pp), hh and ph processes binding the Cooper pair neutron halo and stabilizing ^{11}Li , as well as (b,d) giving rise to the PDR.

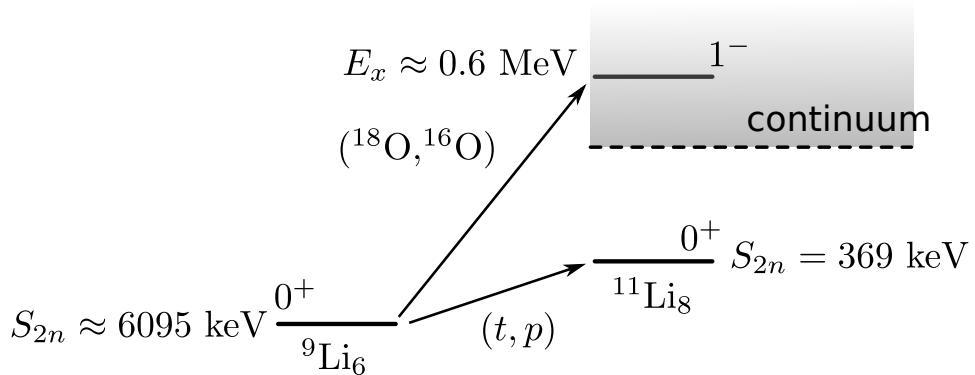


Figure 7.D.3: Schematic representation of levels of ^{11}Li which can be populated in two-nucleon transfer reactions, for example (t, p) , $(^{18}\text{O}^{16}\text{O})$ etc. Indicated in keV are the two-neutron separation energies S_{2n} . In labeling the different states, one has not considered the quantum numbers of the $p_{3/2}$ odd proton.

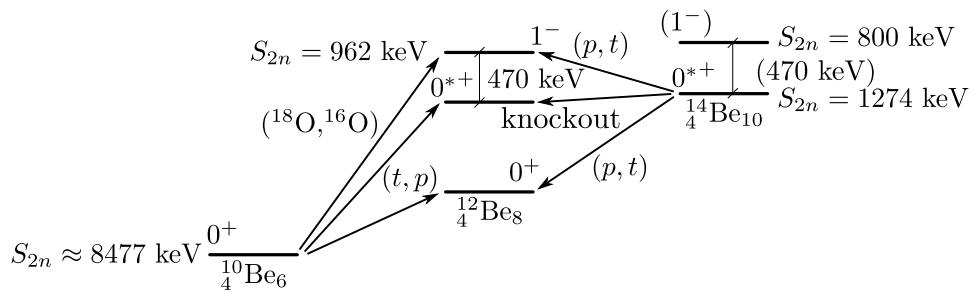


Figure 7.D.4: Levels of ^{12}Be expected to be populated in two-nucleon transfer and knockout processes. S_{2n} are the two-neutron separation energies.

those associated with the core (2^+ see Fig. 7.D.1 (c), and eventually also $3^-, 5^-$ etc.) provide some glue, but insufficient to bind the halo dineutron system.

As already mentioned, the next alternative is that of bootstrapping. Namely, that in which the two partners of the (monopole) Cooper pair exchange pairs of vortices (dipole Cooper pair), as well as one dipole Cooper pair and a quadrupole pair removal mode, while those of the vortex exchange pairs of Coopers pairs (monopole pairing vibrations), but also two dipole pairs, as shown in Figs. 7.D.1 and 7.D.2. In other words, by liaising with each other, the two dineutrons contenders at the role of ^{11}Li ground state settle the issue. As a result the Cooper pair becomes weakly bound ($S_{2n} = 380$ keV), the vortex state remaining barely unbound, $\approx 0.7\text{--}0.8$ MeV above the bound Cooper pair and thus few hundreds keV above threshold. There is, in principle, no physical reason why things could not have gone the other way. Within this context we refer to ^3He superfluidity, where condensation involve $S = 1$ pairs. It is of notice that we are not considering spin degrees of freedom in the present case, at least not dynamic ones.

For practical purposes, one can describe the 1^- as a two quasiparticle state and calculate it within the framework of QRPA⁴⁴, including the $p - h$, $p - p$ and $h - h$ channels and adjusting the strength of the dipole-dipole separable interaction to reproduce the experimental findings (Fig. 4.12.1). In this basis it is referred to as a Pygmy Dipole Resonance (PDR)⁴⁵. Exchanged between the two partners of the Cooper pair (Fig. 7.D.1(d)) leads to essentially the right value of dineutron binding to the ^9Li core. Within this context one can view the ^{11}Li neutron halo as a van der Waals Cooper pair (Fig. 7.D.1 (e); see also Sect. 3.7). The transformation between this picture and that discussed in connection with (a) and (b) as well as with Fig. 7.D.2 can be obtained expressing the PDR, QRPA wavefunction, in terms of particle creation and destruction operators (Bogoliubov-Valatin transformation) as seen from Fig. 7.D.1 (a) and (b). A vortex-vortex stabilised Cooper pair emerges.

Which of the two pictures is more adequate to describe the dipole mediated condensation is an open question, as each of them reflects important physical properties which characterise the PDR. In any case, both indicate the symbiotic character of the halo Cooper pair addition mode and of the pygmy dipole resonance built on top of, and almost degenerate with it. Insight into this question can be obtained by shedding light on the question of whether the velocity field of each of the symbiotic states is more similar to that associated with irrotational or vortex-like flow⁴⁶. Two-nucleon transfer reactions, specific probe of (multipole) pairing vibra-

⁴⁴Broglia et al. (2019a).

⁴⁵Barranco et al. (2001); Broglia et al. (2019c).

⁴⁶See Repko et al. (2013). Within this context, and making use of an analogy, one can mention that a consistent description of the GQR and of the GIQR is obtained assuming that the average eccentricity of neutron orbits is equal to the average eccentricity of the proton orbits (Bès et al. (1975b)), the scenario of neutron skin. The isoscalar quadrupole-quadrupole interaction is attractive. Furthermore, the valence orbitals of nuclei have, as a rule and aside from intruder states, homogeneous parity. These facts preclude the GQR to play the role of the PDR. In fact, there will always be a low-lying quadrupole vibration closely connected with the aligned coupling scheme and thus with nuclear plasticity. Within this context one can nonetheless posit that the GQR, related to neutron

tional modes, contain many of the answers to the above question (Fig. 7.D.3), to the extent that one is able, as one is, to calculate absolute differential cross sections within %10 accuracy. In fact, ground state correlations will play a very different role in the absolute value of the ${}^9\text{Li}(t,p){}^{11}\text{Li}$ (1^-) cross section, depending on which picture is correct. In the case in which it can be viewed as a vortex (pair addition dipole mode) it will lead to an increase of the two-particle transfer reaction (positive coherence). It will produce the opposite effect if the correct interpretation of the PDR is that of a more ($p - h$)-like excitation⁴⁷. Insight in the above question may also be obtained by studying the properties of a quantal vortex in a Wigner cell with parameters which approximately reproduce the halo of ${}^{11}\text{Li}$. Within this context, and for the solely purpose of providing an analogy, we refer to what have been done in the study of quantal vortices in the environment of neutron stars⁴⁸.

A possible test of the soundness of the physics discussed above, concerns the question of whether the first excited, 0^+ halo state ($E_x = 2.24$ MeV) of ${}^{12}\text{Be}$ can be viewed as the $|gs({}^{11}\text{Li})\rangle$ in a new environment. In other words, to consider the halo neutron pair addition mode a novel mode of elementary excitation: neutron halo pair addition mode of which the $|1^-({}^{12}\text{Be}) ; 2.71 \text{ MeV}\rangle$ is a fraction of its symbiotic PDR partner. One can gain insight concerning this question, by eventually measuring the E1-branching ratio $|1^-(2.71 \text{ MeV})\rangle \rightarrow |0^{+*}(2.24 \text{ MeV})\rangle$, and possibly finding other low-energy E1-transitions populating the 0^* state, as well as through two-nucleon stripping process, and two-nucleon pickup and knockout reactions (Fig. 7.D.4). A resumé of the picture discussed above is given in Fig. 4.12.1.

Appendix 7.E Vacuum fluctuations and interactions: the Casimir effect

In Fig. 7.E.1 (I) (a) an example is given of zero-point fluctuations (ZPF) of the nuclear vacuum (ground state) in which, a surface quantised vibration and an uncorrelated particle-hole mode get virtually excited for a short period of time. Adding a nucleon (odd system, Fig. 7.E.1 (I) (b)) leads, through the particle-vibration coupling mechanism to processes which contain the effect of the antisymmetry between the single-particle explicitly considered and the particles out of which the vibrations are built (Fig. 7.E.1 (I)(c)). Time ordering gives rise to the graph shown in Fig. 7.E.1 (I)(d). Processes I(c) and I(d), known as correlation (CO) and polarisation (PO) contributions to the mass operator, dress the particles and lead to physical nucleons whose properties can be compared with the experimental

skin, is closely associated with the aligned coupling scheme. Making a parallel, one can posit that the PDR is closely connected with vortical motion. Arguably, support for this picture is provided by the low-lying E1 strength of ${}^{11}\text{Li}$. It results from the presence of $s_{1/2}$ and $p_{1/2}$ orbitals almost degenerate and at threshold, leading to a low-lying Cooper pair coupled to angular momentum 1^- . (dipole pair addition mode). The scenario of vortical motion.

⁴⁷Broglia et al. (1971c).

⁴⁸Avogadro et al. (2007, 2008).

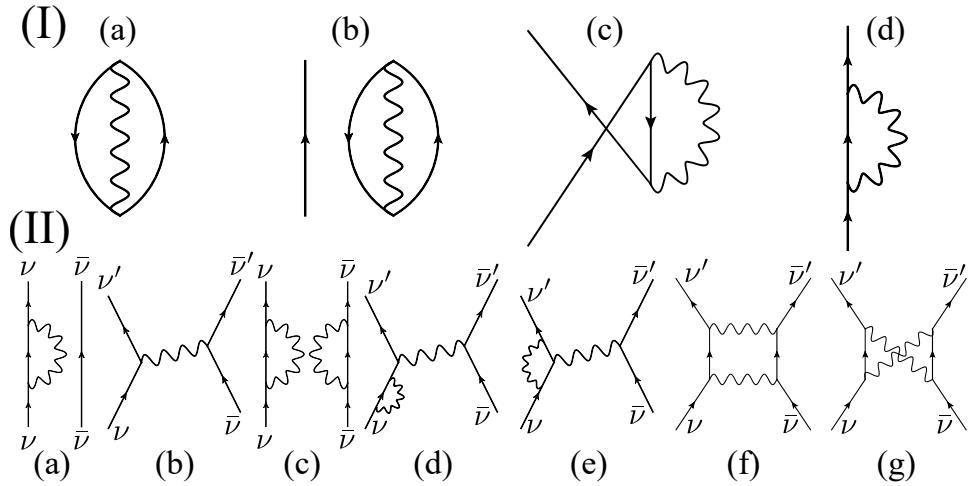


Figure 7.E.1: **(I)** **(a)** ZPF associated with (particle-hole) surface vibrations; **(b)** same process that in (a) but in the case of an odd system; **(c)** the antisymmetrization between the particles considered explicitly and those involved in the vibration; **(d)** time ordering of (c). Diagrams (c) and(d) lead to the clothing of single-particle motion in lowest order in the particle-vibration coupling vertex. **(II)** A dressed nucleon moving in a state ν in the presence of: **(a)** a bare nucleon moving in the time reversed state $\bar{\nu}$, **(b)** another dressed nucleon. Exchange of vibration in (a) leads to (b), NFT lowest-order contribution to the induced pairing interaction, in the particle-vibration coupling vertex. Exchange of vibrations in (c) leads to (d) self-energy, (e) vertex correction of the induced pairing interaction; (f) diagram contributing to the induced pairing interaction. The symmetrisation between the bosons displayed in (c) is shown in (g).

findings with the help of specific experimental probes. For example one-particle transfer reactions. Summing up, the processes shown in Fig. 7.E.1 (I) are examples of quantum field theory phenomena. They testify to the fact that the dressing of nucleons is at the basis of the quantal description of the atomic nucleus.

Nuclear superfluidity at large, and its incipience in the case of single Cooper pair like e.g. in ^{11}Li in particular, are among the most quantal of all the phenomena displayed by the nuclear many-body system. Even if the bare 1S_0 , NN -interaction was not operative, or was rendered subcritical by screening effects as in the case of ^{11}Li , Cooper binding will still be possible, as a result of the exchange of vibrations between pairs of physical (dressed) nucleons (Fig. 7.E.1 (II) (a), (c)) moving in time reversal states close to the Fermi energy (Fig. 7.E.1 (II)(b),(d)-(g)). This is a consequence of the ZPF associated with the nuclear vacuum (ground) state. *Arrived at this point, the reader may find it useful to read again Sect. 1.9, in particular the last paragraphs.*

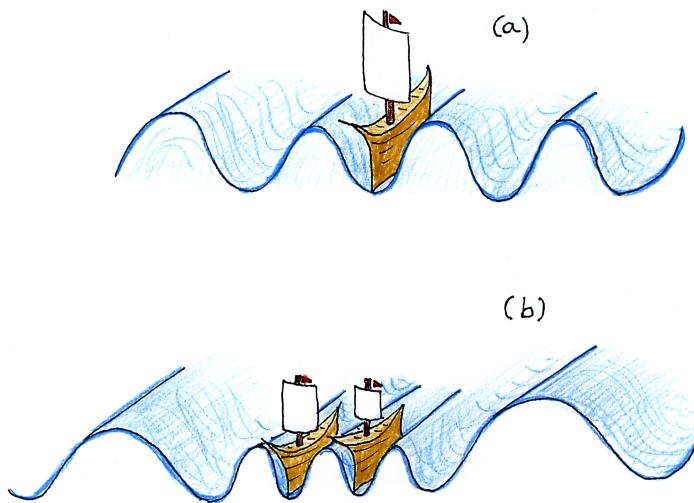


Figure 7.E.2: Schematic representation of the behaviour of an isolated ship at sea in a situation of no wind but of strong swell (a), and of two ships close by under similar conditions (b).

7.E.1 Measuring QED quantum fluctuations

The many-body character of the attractive interaction between nucleons arising from the nuclear (vacuum) ZPF discussed above, and important to describe phenomena occurring close to the surface of the Fermi sea, have a classical analogue. Fittingly, a classical phenomenon which emerges from maritime experience.

At sea, on a windless day in which the water surface can resemble a mirror, free floating ships singly or in groups do not do much, just stay put. The situation is quite different in a strong swell, still on a windless situation. In this case, single, isolated ships end up lying parallel to the wave crests (see Fig. 7.E.2(a)) and start rolling heavily. In the days of the clipper ships it was believed that under those circumstances, two vessels at close distance, attracted each other. This in keeping with the fact that the rigs of the rolling ships became often entangled leading to disaster. It was not until quite recently⁴⁹ that a quantitative understanding of the phenomenon (based on knowledge of similar quantal effects) was achieved, providing evidence that the old tale was true. Only waves with wavelength smaller or equal than the separation of the ships can exist between them. In the region of sea extending away from the ships to the horizon, waves of any wavelength can exist (see Fig. 7.E.2(b)). This fact results into an imbalance between the forces exerted by the internal (between ships) waves, in favour of that exerted by the external waves, leading to a net attraction. Quantum mechanical, such an effect is known as

⁴⁹Boersma (1996).

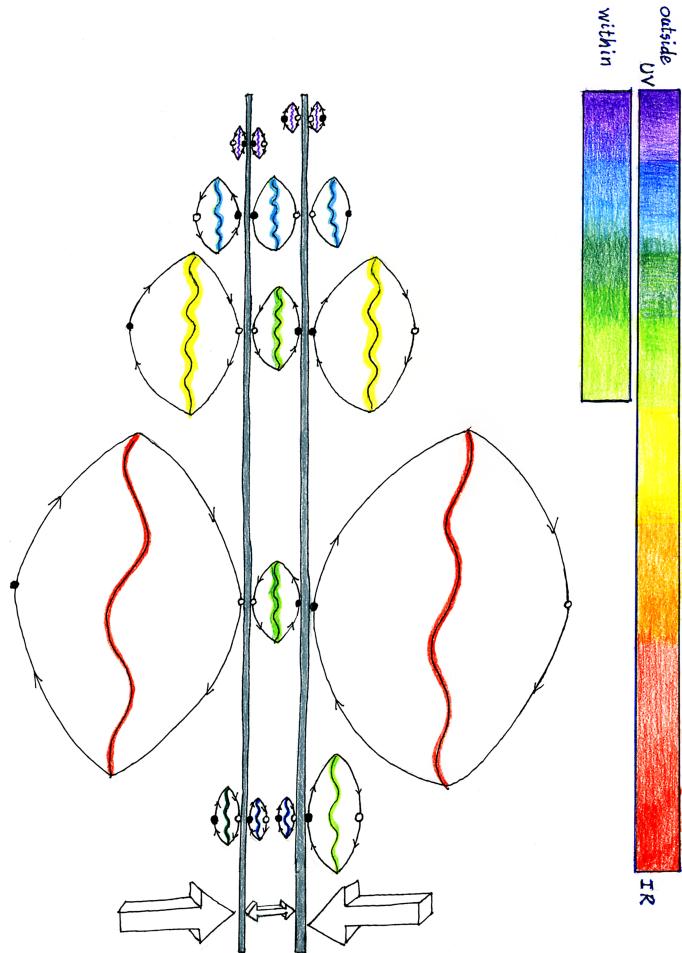


Figure 7.E.3: Casimir effect. Two metallic isolated, conducting plates (grey vertical sections) in vacuum attract each other when they are placed at very small distances (of the order of micron). This is known as the Casimir effect. The origin of such a force can be traced back to quantal zero point fluctuations (ZPF) of the electromagnetic vacuum. In the figure a cartoon of such processes is given. Virtual electrons (e^- , solid dots) and positrons (e^+ , open dots) pop up of the vacuum together with a photon, and travel for short distances and times. In their way, some of them hit the plates. Because the range of wavelengths allowed between the plates is smaller than the full spectrum allowed for the photons associated with the electromagnetic ZPF in the right and left unlimited halves, more fermions or bosons will be knocking the plates from outside than from the in between region, thus leading to an imbalance of the “quantal” pressure and consequently to an effective attractive force.

the Casimir effect⁵⁰.

Two conducting, neutral plates at very small distances, of the order of the micron ($1\mu = 10^{-6}$ cm), attract each other, due to the imbalance in electromagnetic field pressure exerted by the bombarding of the surface by photons, electrons and positrons, arising from the ZPF of the electromagnetic field⁵¹ (see Fig. 7.E.3). It is of notice that the Casimir effect, namely the attraction between two metallic, uncharged, plates which in Fig. 7.E.3 have been drawn as plane surfaces, can in principle have any shape.

Summing up, in a similar way in which one can state that there are neither bare fermions (e.g. electrons or nucleons) nor bare bosons (or quasi-bosons) (e.g. photons (surface vibrations)), one can posit that essentially all bare forces are, with varied degrees of complexity and strength, renormalized by many-body effects. This message comes also from the least suspected of all fundamental forces, the Coulomb interaction, responsible for all of chemistry and most of biophysics (see also Sect. 4.8.3).

7.E.2 The hydrophobic force

Water has three noteworthy properties: liquid water is heavier than ice, it has an exceptionally large specific heat⁵² –the amount of heat needed to raise the temperature of 1 kg of mass 1 degree Kelvin at room temperature (≈ 300 K)– aside from being an excellent solvent due to its dipole moment.

All these properties follow from the structure of the water molecule H₂O. This, in turn, is connected with the directional, anisotropic structure of the valence, electronic distribution of O. Oxygen is an open shell atom⁵³, having in its ground state only four electrons in the last occupied 2p-orbital which can host six ($|gs\rangle = |1s^2\rangle|2s^2\rangle|2p^4\rangle$). It can thus use at profit the electrons of the two H-atoms to dynamically become a closed shell system (²⁰Ne–noble gas–like electronic configuration, Fig. 7.E.4 (c)).

Hybridization between the four orbitals $|2s\rangle$, $|2p_x\rangle$, $|2p_y\rangle$ and $|2p_z\rangle$ (Fig. 7.E.4 (c)) leads to a tetrahedral correlation in which the four orbitals $|i\rangle$ point towards the corners of a tetrahedron with the oxygen atom at the center (Fig. 7.E.4 (a), 7.E.5). Because the electronic distribution has its charge center closer to the oxygen atom than to the two protons of the H-atoms, H₂O has a sizable dipole moment⁵⁴

⁵⁰Casimir (1948).

⁵¹As stated in the last sentences of the caption to Fig. 7.E.3, long wavelengths play the central role in the Casimir effect. A recurring property of the modes renormalizing the bare interaction or the associated collective variables (elementary modes of excitation) found in condensed matter (phonons of much lower frequency (ω_{ip}) than plasmons (ω_{ep}), nuclei (collective quadrupole and soft- E_1 (pygmy) modes ($\hbar\omega_2$, $\hbar\omega_{PDR}$) $\ll \hbar\omega_{GR}$), proteins (in this connection see Micheletti et al. (2004, 2001, 2002); Hamacher (2010); see also footnote 62 of Ch. 1).

⁵²Water absorbs at 25 °C, 4.18 Joules/(g K) of heat when the temperature of one gram is increased 1 degree celsius (°C). For the sake of comparison, it takes ≈ 2.0 Joules/(g K) of heat to raise 1 gram of acetic acid, acetone or ethanol 1 °C.

⁵³See e.g. Greiner (1998).

⁵⁴The dipole moment of a polar molecule is defined as $u = ql$, where l is the distance between the

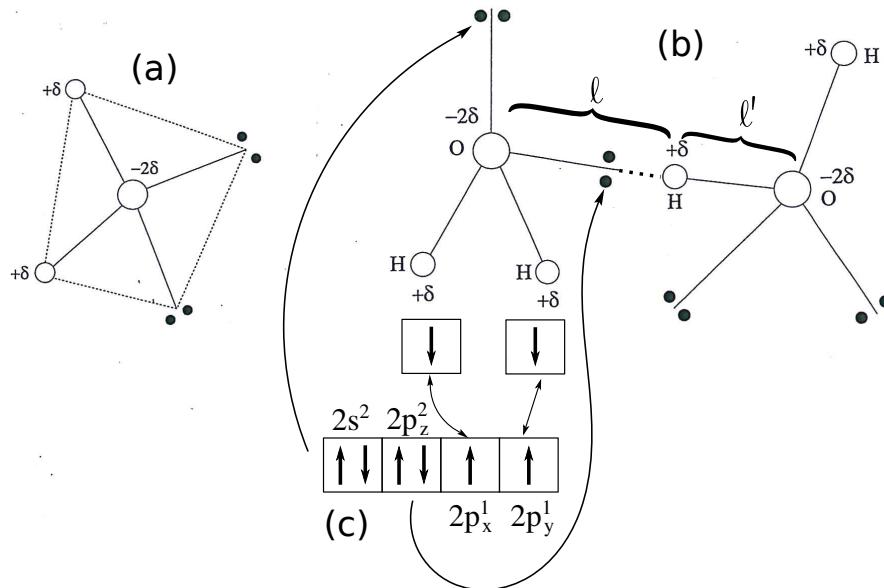


Figure 7.E.4: (a) To understand the behavior of water molecules one has to realize that in the covalent bonds, where the oxygen and the hydrogen atoms share a couple of electrons, the corresponding electronic cloud is somewhat more concentrated on the oxygen than on the hydrogen. The oxygen acquires a partial charge -2δ while each hydrogen gets a positive one $+2\delta$. (b) Such charges give rise to an attractive electric force between close lying molecules, in which the hydrogen atoms of a molecule point to the oxygen atom of the other molecule. One can view this attractive force as a type of chemical bond. It is known as hydrogen bond (dotted line). It is of notice that also in this type of bond there is a component which implies the sharing of electrons. Furthermore, the hydrogen atom in an hydrogen bond does not just stick indiscriminantly to the oxygen of another molecule. Being positively charged it goes where the electrons are. So the hydrogen bond is a bond between a hydrogen atom and a lone pair (two solid dots) its length ℓ being 1.83 \AA , while $\ell' = 1.04 \text{ \AA}$. This means that a water molecule can form four hydrogen bonds: the molecule's two hydrogens form two bonds with neighboring oxygens, while the molecule's two lone pairs interact with neighboring hydrogens. (c) Schematic electronic structure of H_2O is displayed below (b) to schematically indicate the origin of the lone pair of electrons, as well as the role the electrons of the two hydrogen atoms have in closing the $2p$ -shell of the oxygen atom.

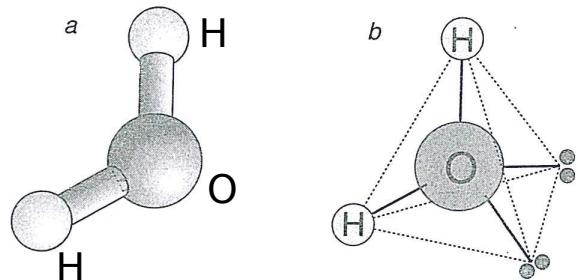


Figure 7.E.5: The water molecule is bent, with the two bonds between oxygen and hydrogen making an angle of 104.5° (a). To understand the structure of liquid water, we must also take into account the two “lone pairs” of electrons on the oxygen atom. The hydrogen atoms and the lone pairs sit more or less at the corners of a tetrahedron (b).

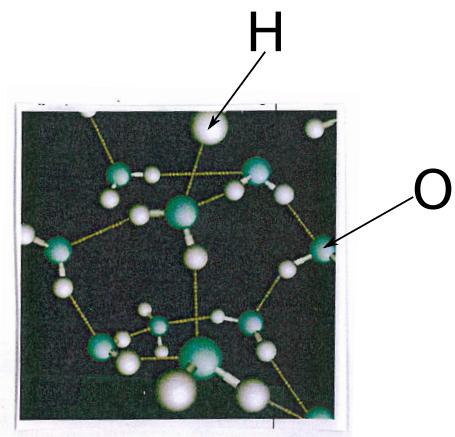


Figure 7.E.6: Stick and ball representation of a network of hydrogen bonds in bulk water.

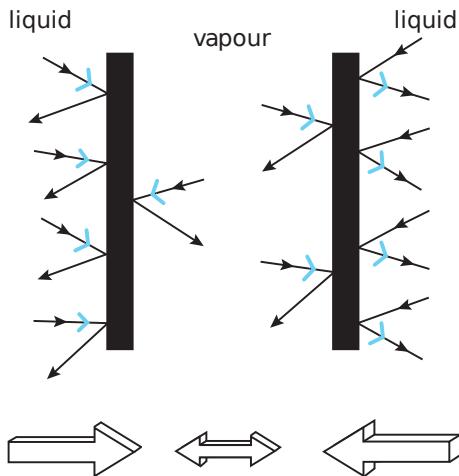


Figure 7.E.7: Schematic representation of water molecules (arrowed lines) impinging from the outside (liquid) and from the inter space (vapor) of two large parallel hydrophobic plates.

($\approx 0.68ea_0 \approx 0.6$ D, e being the electron charge, a_0 the Bohr radius and D the Debye unit). Water molecules can form four hydrogen bonds⁵⁵ (hb; see also Fig. 7.E.6) a special bond between molecules which is produced in situations when they share an hydrogen nucleus between them. The molecule's two hydrogen atoms form two bonds with neighboring oxygens, while the molecule's two lone pairs interact with neighboring hydrogens⁵⁶ (Figs. 7.E.4 (b) and 7.E.6).

Water and oil do not mix. The term hydrophobic (water-fearing) is commonly used to describe substances that, like oil, do not mix (dissolves) with (in) water. Although it may look as if water repels oil, these two types of molecules actu-

two charges $+q$ and $-q$. Thus, for two electronic charges $q = \pm e$ separated by $l = 0.1$ nm, the dipole moment is $u = 1.602 \times 10^{-19} \text{ C} \times 10^{-10} \text{ m} = 4.8 \text{ D}$, where the Debye=1D= $3.336 \times 10^{-30} \text{ Cm}$ is the unit of dipole moment. Small polar molecules have moments of the order of 1D (see e.g. Israelachvili (1985)).

⁵⁵Let us consider an H atom in a covalent bond with oxygen. When a second oxygen atom approaches the H atom, its nucleus, the proton, sees a potential with two minima, and tunnels through the corresponding barrier from one minimum to the other. In other words, the effective potential in which the proton moves becomes broader as compared to the single oxygen potential. Thus, the quantum mechanical confinement kinetic energy decreases by roughly a factor of two. This implies that the order of magnitude energy of an hydrogen bond between two oxygen atoms corresponds to the difference in the corresponding ZPF energies, i.e. ≈ 200 meV ($\approx 0.38 \text{ eV} - 0.19 \text{ eV} \approx 0.2 \text{ eV} \approx 4.6 \text{ kcal/mol}$ ($\approx 8 \text{ kT}$), where kT ($\approx 0.6 \text{ kcal/mol} \approx 25 \text{ meV}$) is the thermal energy at room temperature ($\approx 300 \text{ K}$)). For comparison a covalent bond corresponds to $\approx 96 \text{ kcal/mol} \approx 4 \text{ eV}$, while the van der Waals interaction energy between two H (Eqs. (3.B.18) and (3.B.19)) at a distance of 2.4 \AA , that is of the order of the hydrogen bond length of 1.83 \AA , is $\approx 20 \text{ meV}$ (see Povh and Rosina (2002))

⁵⁶A possible pedestrian explanation of this is to impersonate a water molecule. Quoting from Ball (2003): “Your hands are hydrogen atoms, your ankles are the lone pairs of electrons of oxygen. Stand legs apart... Twist 90° at the waist, stretch your arms and you’re H₂O. The way that water molecules join up has just one rule: hands can grab ankles, but nothing else. That grasp is an hydrogen bond”.

ally attract each other, e.g. through the van der Waals interaction, but not nearly as strongly as water attracts itself. Mixing enough oil (hydrophobic, non polar (NP) molecules) with water leads to a reduction in favorable bonding. Strong mutual attraction between water molecules induce segregation of NP molecules from water and results in an effective NP-NP (hydrophobic) attraction, as observed in surface force measurements⁵⁷. The loss of hydrogen bonds near the two extended hydrophobic surfaces depicted in Fig. 7.E.7 causes water to move away from those surfaces⁵⁸, producing thin vapour layers. Fluctuations in the interfaces formed in this way can destabilize and expel the remaining liquid contained between these surfaces. The resulting pressure imbalance will cause the surfaces to attract. If the liquid is close to coexistence with the vapor phase, as is the case for water at ambient (biological) conditions⁵⁹, this phenomenon occurs also for widely separated surfaces. The similarity with a (generalized) Casimir effect is apparent.

Within this context, and in keeping with the fact that the hydrophobic force plays a central role in the folding and, as a consequence, in the biological activity of proteins, one can point that water at physiological conditions ($T = 300\text{K}$, $\text{pH}=7$, etc.) can be viewed as the vacuum of the macromolecules responsible for metabolism, and thus one of the two origins of life on earth⁶⁰.

While the many-body basis of hydrophobicity cannot be denied (Fig. 7.E.6) the parallel with the zero-body Casimir effect (Fig. 7.E.3) can hardly be avoided. Neither with the nuclear phenomena discussed in Sect. 4.8 in connection with superconductivity in nuclei and in metals (see also Sect. 3.6 concerning pairing in ^{11}Li).

Within this scenario, one can also observe a (generalized) Nambu-tumbling effect. Namely the chain *hydrogen bond* → *weak and strong hydrophobicity* →

⁵⁷See e.g. Chandler (2002, 2005) and Lum et al. (1999) and references therein.

⁵⁸Hydrophobic molecules do not hydrogen bond to water, creating excluded regions where the density of water molecules vanishes. When these units are small enough, water can reorganize near them without loosing hydrogen bonds, building an ice-like cage around the NP molecule. The entropic cost of this structural change leads to low solubility for small apolar molecules in water. Said it differently, such a force, known as the weak hydrophobic force (WHF) leads only to metastable states of groups of nonpolar molecules (Chandler (2002)). On the other hand, close to a large hydrophobic object (Fig. 7.E.7), the persistence of an hydrogen bond network –four bond per water molecule–cannot persist, being geometrically impossible. In this case, water molecules near the hydrophobic cluster (surface) have typically three or fewer hydrogen bonds. This force, known as the strong hydrophobic force (SHF, Chandler (2002)) can be viewed as a *bona fide* hydrophobic interaction, stabilizing large aggregates of nonpolar molecules.

⁵⁹In particular at $T \approx 300\text{K} (\approx 27\text{C}^\circ)$. This is the reason why proteins unfold at low temperatures, in particular the monomer of the HIV-1-protease homodimer, which denaturizes at 8C° (Rösner et al. (2017)). In general this temperature lies below freezing point and one has to use particular techniques to be able to deal with supercool water (e.g. capilars, etc.).

⁶⁰To the question “what is life?” (Schrödinger (1944)) one is forced to answer that life is not one but two things (Dyson (1999)). Which ones? Replication and metabolism. The molecules of DNA and RNA are responsible for the first function (Watson and Crick (1953); Watson (1980)), proteins (i.e. polymers made out of the 20 commonly occurring amino acids in nature), for the second (Sanger (1952)). Because software (replication) is necessary a parasite of hardware (proteins), the becoming of a protein carries, to a large extent, the secret of life (Monod (1970)).

polypeptide chain evolution→*proteins (second order phase transition, spontaneous symmetry breaking in information space*⁶¹)→*protein folding (first order phase transition in conformation space)*→*metabolism.* And, at the basis of the first link of the chain, one finds the difference in the ZPF energy which an hydrogen nucleus feels when in presence of one (O) or two (O₂) oxygen nuclei, quantity closely connected with quantum fluctuation of confinement (Povh and Rosina (2002)).

⁶¹Broglia (2013).

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