

In App. A it was shown that $q \approx 0.5$ leads to particle delocalization and likely make the system amenable to a mean field description.

1.D. COOPER PAIRS (Fig. 1A.2; see however

1.D.1 independent particle motion

($q \approx 0.5$), mean field (see Fig. 1A.2)

$$\exp \sum_i \sigma(\sigma_i^\dagger(i))_{\text{exc}} = \sigma_{\text{exc}}$$

$$\frac{\sigma_{\text{exc}}}{\sigma_{\text{exc}} + \sigma_{\text{gs}}} \approx 0.03$$

(11)

(31)

provides end paragraph of App. A). In such a case Hartree-Fock approximation is tantamount to a self-consistent relation between density and potential, weighted by the nucleon-nucleon interaction v (eventual sum of the bare and the induced interaction), i.e.

(the Hartree-Fock ground state being

$$U(r) = \int dr' \rho(r') v(|r-r'|),$$

$$U_x(r, r') = - \sum_i \varphi_i^*(r') v(|r-r'|) \varphi_i(r'),$$

$$\rho(r) = \sum_i |\varphi_i(r)|^2; \int dr \rho(r) = N,$$

leading to

Hartree-Fock complete separation between occupied (i) and empty (k) states,

$$(U_v^2 + V_v^2) = 1 \varphi = \bar{a}_v^\dagger |0\rangle = (U_v + V_v \bar{a}_v^\dagger) |0\rangle; V_v^2 = \begin{cases} 1 & \epsilon_i \leq \epsilon_F \\ 0 & \epsilon_i > \epsilon_F \end{cases}$$

$$|HF\rangle = |\text{Nilsson}(52)\rangle_{\mathcal{K}} = \det(\varphi_v) = \Pi \bar{a}_v^\dagger |0\rangle = \Pi a_i^\dagger |0\rangle = \Pi a_i^\dagger a_i^\dagger |0\rangle.$$

$$|IKM\rangle \sim \int d\Omega D_{MK}^I(\Omega) |\text{Nilsson}(\Omega)\rangle; E_I = (\hbar^2/2I) I(I+1); I = I_{\text{rot}}$$

(A)-(A)

(B)-(B) from p. 11d

1.D.2 independent pair motion

Let us make use of the

constant matrix element approximation ($j_1 j_2 |0\rangle j_3 j_4\rangle = G$, The abnormal density mean field is related to pair operators (finite values), i.e.

$$\sum (a_{j_2}^\dagger a_{j_1}^\dagger) a_{j_3} a_{j_4} + \sum a_{j_2}^\dagger a_{j_1}^\dagger \langle a_{j_3} a_{j_4} \rangle; \varphi^{\text{COOPER}} = (U_j + V_j a_{jm}^\dagger a_{jm}^\dagger) |0\rangle,$$

(the ground state (mean field) wavefunction being the BCS wavefunction,

$$|BCS\rangle = \Pi_{jm>0} (U_j + V_j a_{jm}^\dagger a_{jm}^\dagger) |0\rangle; \alpha_0 = \langle BCS | \sum_{jm>0} a_{jm}^\dagger a_{jm}^\dagger |BCS\rangle.$$

pairing

Introducing the phony,

$$U_v = |U_v| = U'_v; V_v = e^{-2i\phi} V'_v (V'_v \equiv |V_v|) (v \equiv j, m); \bar{a}_v^\dagger = e^{-i\phi} a_v^\dagger;$$

here ϕ is the gauge angle, one can write for the intrinsic wavefunction,

$$|BCS(\phi)\rangle_{\mathcal{K}} = \Pi (U'_v + V'_v e^{-2i\phi} a_v^\dagger a_v^\dagger) |0\rangle = \Pi (U'_v + V'_v \bar{a}_v^\dagger a_v^\dagger) |0\rangle,$$

$$= |BCS(\phi=0)\rangle_{\mathcal{K}'} : \text{lab. system, } \mathcal{K}': \text{intr. system } \mathcal{K}'.$$

(two-nucleon systems with longitudinal are

The BCS number and gap equations, order parameter and

$$\alpha_0 = \alpha'_0 e^{-2i\phi}; \alpha'_0 = \sum_{v>0} U'_v V'_v; B_v = \langle BCS | [a_v^\dagger, \bar{a}_v^\dagger]_0 |BCS\rangle = (2v+1/2)^{1/2} U'_v V'_v.$$

Examples of Bv coefficients for the reaction $^{124}\text{Sn}(p, t)^{122}\text{Sn}(gs)$ are given in table 1.D.1 $\Delta = G\alpha_0; N_0 = 2 \sum_{v>0} V_v^2; \frac{1}{G} = \sum_{v>0} \frac{1}{2E_v} \left\| \frac{V'_v}{U'_v} \right\|^2 = \frac{1}{\sqrt{2}} \left(1 \mp \frac{1}{E_v} \right)^{1/2}$ // space

The wave function of the members of a pairing rotational band can be written as

$$|N_0\rangle \sim \int_0^{2\pi} d\phi |BCS(\phi)\rangle_{\mathcal{K}} \sim \left(\sum_{v>0} c_v a_v^\dagger a_v^\dagger \right)^{N_0/2} |0\rangle; E_N = (\hbar^2/2I) N^2 I \approx 2\hbar^2/G$$

In the case of a deformed nucleus, the system acquires not only a privileged orientation in gauge space, but also in 3D-space. Now, as summarized

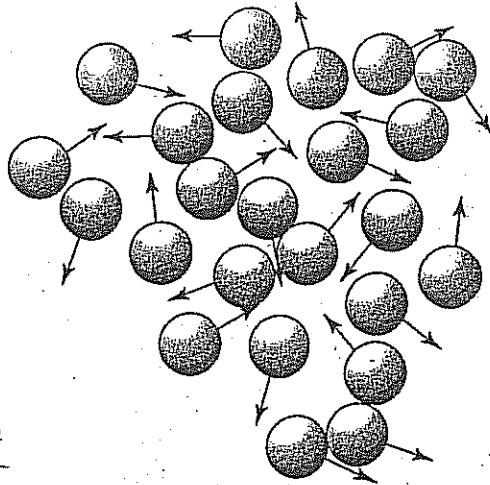
(a quadrupole

above, in a superfluid system, Cooper grains⁽¹⁹⁾
 and not single particles, are the building
 blocks of the system. But while the mean $\langle r^2 \rangle$
 square radius of a nucleon at the Fermi energy
 is $\langle r^2 \rangle^{1/2} \approx (3/5)^{1/2} R_0$ ($R_0 = 1.2 \text{ fm} \approx 6 \text{ fm}$ ($A \approx 120$)),
 that of a Cooper pair is determined by the
 correlation length $\xi \approx \hbar v_F / M_n \approx 36 \text{ fm}$ between
 the two nucleons forming the pair (within this
 context see Figs. 1.A.4 and 1.A.5). Consequently
 orienting the quadrupole deformed potential
 in different directions (angles Ω), will have a ~~much~~
 less pronounced effect on Cooper pairs
 than on independent particles. Within
 this context one can mention the fact that low-
 lying ^{nuclear} collective vibrations (and rotations)
 are not observed ^(essentially) at intrinsic excitation
 energies corresponding to temperatures of
 $\approx 1-2 \text{ MeV}$. In this case, this is because the
 surface is strongly fluctuating and thus
 not well defined, making it non-operative
 its anisotropic orientation in space.

In keeping with the fact that in
 FMB ^{and due to the important role played by quantal fluctuations} deformations are exhibited through
 rotational bands, superfluid nuclei display
 well defined pairing rotational bands
 (e.g. the ground state of the superfluid
 Sn-isotopes). The moment of inertia is directly
 See Fig. 1.3.

related to the effective pairing interaction, $(11)_6$ eventually, sum of the bare and of the induced interactions. These rotational bands are specifically excited in two-nucleon transfer reactions [cf. Figs. 1.3, 1.4]

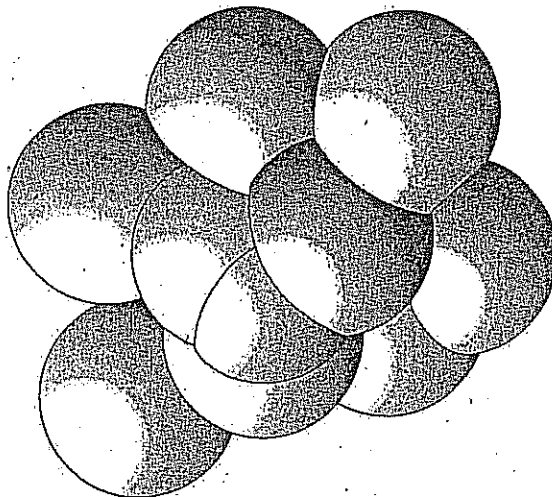
(33)



1.A.4

Figure 1.4. A system of independent Cooper pairs (Schafer pairs). This situation corresponds to the incoherent solution of the many Cooper pair problem, the so called Fock state.

In cold gases the system after the Feshbach resonance leading to BEC



1.A.5

Figure 1.5. There are about 10^{18} Cooper pairs per cm^3 in a superconducting metal. A Cooper pair has a spatial extension of about 10^{-4} cm. Thus a given Cooper pair will overlap with 10^6 other Cooper pairs, leading to strong pair-pair correlation, as schematically shown. This solution corresponds to the coherent solution of the many Cooper pair problem (coherent state), also valid in atomic nuclei.

flucts (paring vibration, the right physics)

(11) C

or at least not only

(E) From this vantage point one can posit that it is not so much the nuclear superfluid state which is abnormal, but the normal (closed shell) system (see App. D).

p. 4 version June 6/14

(E)

(34)

(A)-(A) p. 11

Fig. 5 WSPC Spont. breaking II

in the case in which $\langle Q_{2M} \rangle = 0$ the system can display quadrupole as a spectrum of quadrupole vibration the associated ZPF leads to dynamic violations

~~The above self consistent~~

(where $\langle Q_{2M} \rangle \neq 0$) of quadrupole rotational invariance

carrying & paring (vibrations) or entire nuclei $\beta=0$ in keeping with the quadrupole - Goldstone character

To be solved, the above self consistent equation have to be given boundary conditions. Namely in particular whether the system has a spherical ^{for example} or a quadrupole deformed shape. That is, whether $\langle HF | Q_2 | HF \rangle$ is zero or different or has a finite value ~~at the origin~~ (cf. App. H)

$$Q_{2M} = \sum_{\beta=0, \pm 1, \pm 2} \langle r_2 | r^2 Y_{2M} | r_1 \rangle [a_{\beta}^\dagger a_{\beta}]_{2M}$$

being the quadrupole operator. in the case in which $\langle Q_{2M} \rangle \neq 0$, the $|HF\rangle$ state is

known as the Nilsson state, $|Nilsson(52)\rangle_K$ defining a privileged orientation in 3D space and thus an intrinsic, body-fixed system of reference K' which makes an angle Ω (Euler angle) with the laboratory frame K . The fluctuation associated

with the fluctuations associated of the angle Ω become then is no restoring force associated with the different orientations, (different fluctuation in Ω diverge leading to a restoration of symmetry ($C \rightarrow 0$, D finite in the case of low-lying collective, large amplitude vibrations mentioned above) leading to a rotational band displaying a rigid moment of inertia and whose members are the states

(relation between collectivity and correl. of r.o.s & vibs)

type by E. V. S.

p. 11

(A) p. 11 (version 6/14)

(B) Similar dynamic and static spontaneous symmetry breaking phenomena takes place in connection with particle-particle ^($B=+2$) and hole-hole ^($B=-2$) correlations, that is ~~with~~ gauge space, (see ~~Table~~ Fig. 1.D.1), ~~These~~ subjects discussed in App. E (~~dynamic~~ : parring vibrations) and below (static : parring rotations), ~~to~~

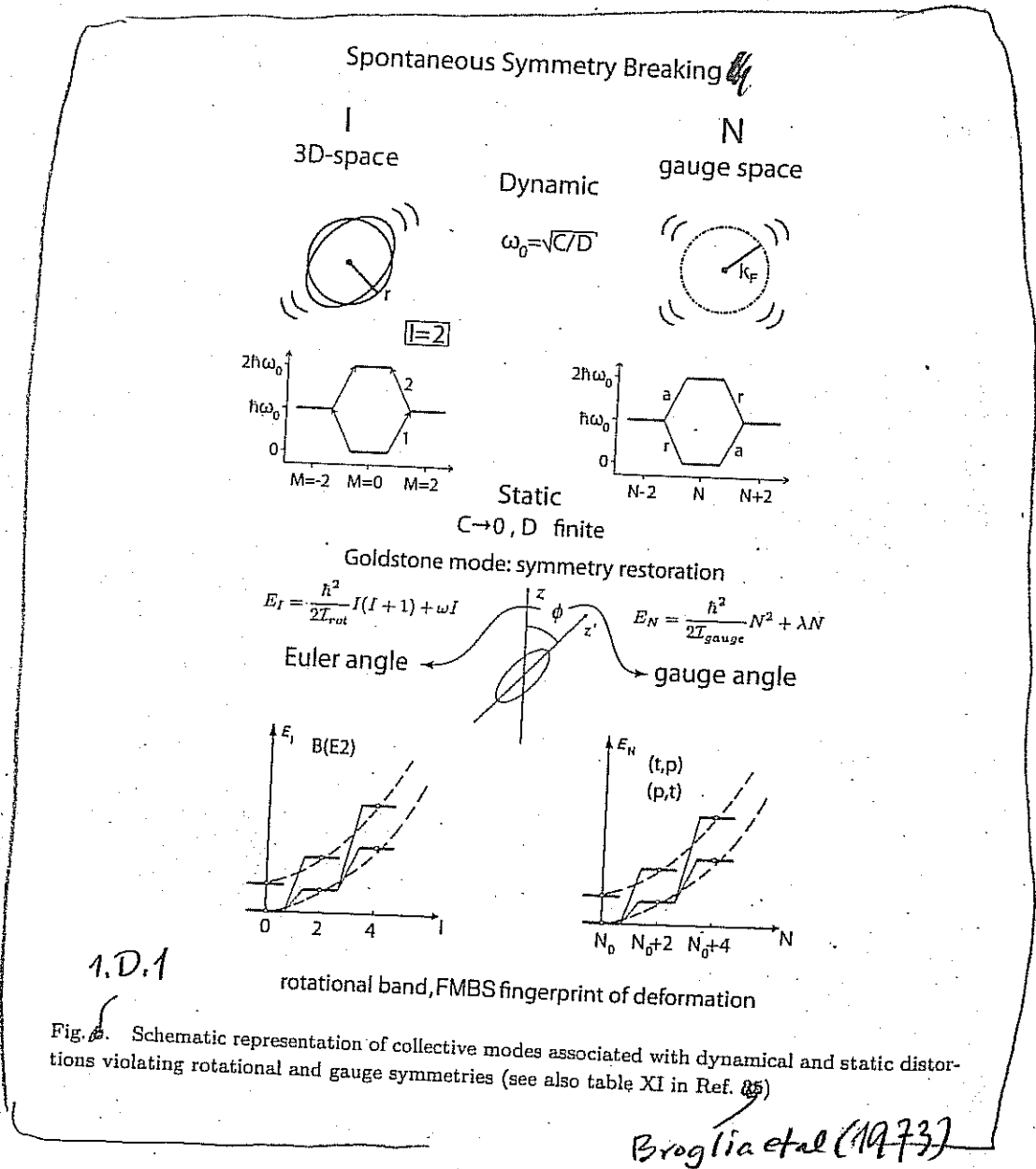
(see also Figs. 1.1 and 1.3)

(B)

p. 11 version
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(35)

the successive contribution to the two-particle transfer cross section is the dominant one, non-orthogonality canceling much of the already weak, simultaneous contribution. Of notice that similar issues were debated in connection with the proposal of Josephson²⁹ concerning the possibility of observing a supercurrent across a dioxide layer separating two superconductors, and Bardeen's objection that the pairing gap is zero inside the layer.³⁰ The answer to such an objection is to be found in the fact that it is $\alpha_0 (= \langle P^\dagger \rangle)$ which controls tunneling and not Δ , a fact that emerges naturally from Gorkov's formulation of superconductivity (see contribution of Potel and Broglia to the present volume).



1.D.3 Two-nucleon spectroscopic amplitudes associated with transitions between members of a pairing rotational band

(19)C

Parity change part. vib.

Explain in this appendix pairing rots + p. vibs.
(refer Advance, two-entry

Table 1, A.1

Charge Transfer
multishell
wave function
Covello

use all three w.f.s
p. 3, 1

Pair transfer in
a multishell Charge 1



Two-entry
Table WSPC Table 5
p. 652

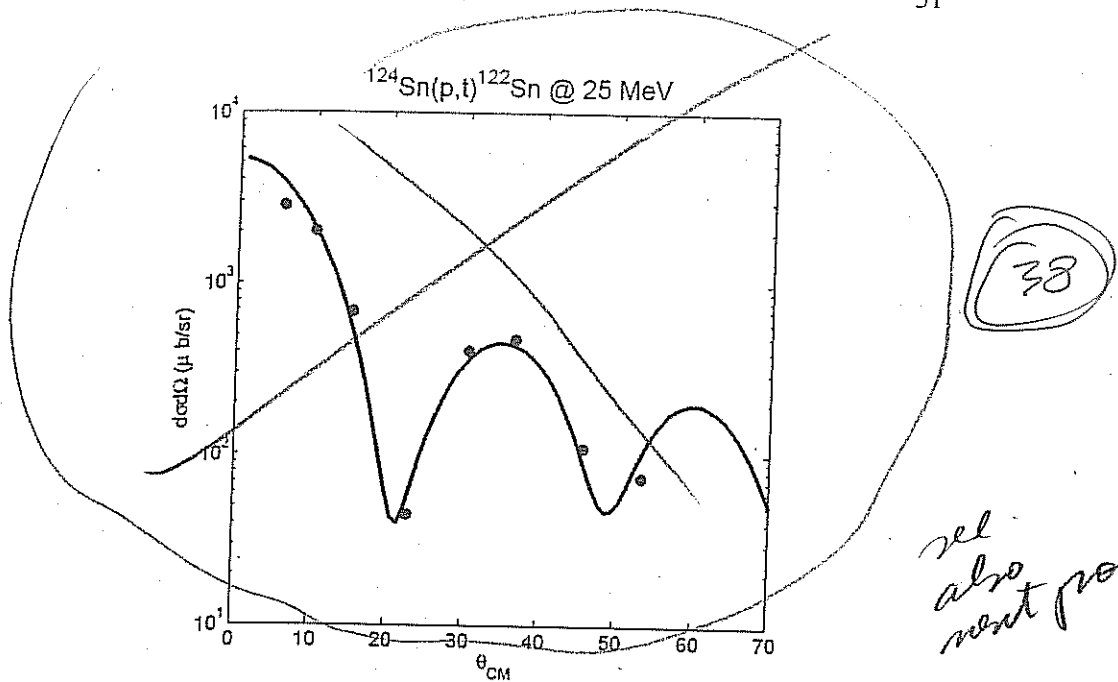
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talk
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(37)



rel.
also
next page

Table 1.D.1

Figure 1.A.3:

nlj^a	2n spectr. ampls. $^{124}\text{Sn}(p,t)^{122}\text{Sn}(\text{gs})$		
	BCS ^b	NuShell ^c	V_{low-k}^d
$1g_{7/2}$	0.44	0.63	-1.1073
$2d_{5/2}$	0.35	0.60	-0.7556
$2d_{3/2}$	0.58	0.72	-0.4825
$3s_{1/2}$	0.36	0.52	-0.3663
$1h_{11/2}$	1.22	-1.24	-0.6647

Table 1.A.1: a) quantum numbers of the two-particle configurations $(nlj)_{J=0}^2$ coupled to angular momentum $J = 0$. b) $\langle \text{BCS} | P_v | \text{BCS} \rangle = \sqrt{2j_v + 1} U_v(A) V_v(A+2)$ ($A+2 = 124$) where $P_v = a_v^\dagger a_v$ ($v \equiv nlj$) (cf. Potel et al. (2011, 2013a,b)). c) Two-neutron overlap functions obtained making use of the shell-model wavefunctions for the ground state of ^{122}Sn and ^{124}Sn and the code NuShell (Brown and Rae, 2007) (cf. also ?). 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).

Objection

What about $v_{\text{pairing}} (= G)$ becoming zero, e.g. between the two nuclei?

(Covello)

spectroscopic
amplitude

$$\sqrt{2j_v + 1} U_v(A) V_v(A+2)$$

↑ check

is it $\sqrt{2} S_v$
or $\sqrt{S_v}$?

$\sqrt{2} S_v$

Sney Table IV
Covello

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Table 1. Two-neutron overlap function for $(^{122}\text{Sn}|^{124}\text{Sn})$.

$1g_{7/2}^2$	0.62944
$2d_{5/2}^2$	0.59927
$2d_{3/2}^2$	0.71913
$3s_{1/2}^2$	0.51892
$1h_{11/2}^2$	-1.24399

part of
Table 1. P. 1

(29)

of $(0g_{7/2}, 1d_{5/2}, 1d_{3/2}, 2s_{1/2}, 0h_{11/2})$ for neutrons with the code NuShell.²⁴ The model-space two-body matrix elements are those used in Refs. 29 and 30. They were obtained starting with a G matrix derived from the CD-Bonn³¹ nucleon-nucleon interaction. The harmonic oscillator basis was employed for the radial wave functions with an oscillator energy $\hbar\omega = 7.87$ MeV. The effective interaction for the above shell-model space is obtained from the Q-box method and includes all non-folded diagrams through third-order in the interaction G to sum up the folded diagrams to infinite order.^{32,33} The single-particle energies were adjusted to reproduce the observed states in ^{131}Sn .

The inputs to the reaction code are the two-nucleon spectroscopic amplitudes (TNA) of Table 1. A center of mass correction³⁴ equal to $[A/(A-2)]^{2n+l}$ for the TNA has been applied, where $A = 124$. Our sign convention is that the radial wave functions are positive at the origin. The sequential process was calculated by a single intermediate state for each of these orbits connected by a product of one-nucleon spectroscopic amplitudes that are equal to the center-of-mass corrected TNA multiplied by $\sqrt{2}$ that takes into account the normalization of the two-particle amplitude. Future calculations should also take into account the TNA obtained from the mixing of neutron pairs for orbitals outside of the model space.

We use the triton potential of Li,³⁵ the deuteron potential of Daehnick,³⁶ and the proton potential of Chapel Hill 89.³⁷ All the two-neutron wave functions are constructed within the half-separation-energy prescription. For a triton wave function we use the pure s^2 configuration found by the product of eigenstates at the half-separation energy (4.24 MeV) in a Woods-Saxon potential with $V = 77.83$ MeV, $R = 0.95$ fm, and $a = 0.65$ fm (the results are not sensitive to these values). The Sn wave functions shown in Table 1 are found at the half-separation energy (7.219 MeV) in a WS potential with $r = 1.17$ fm, and $a = 0.75$ fm that has the fixed spin-orbit component $V_{so} = 6.2$ MeV, $r = 1.01$ fm, and $a = 0.75$ fm.

The complete cross section prediction is shown in Fig. 1, compared with the experimental data of Guazzoni *et al.*²⁸ Now we see that, with the shell-model overlaps and proper finite-range and sequential contributions, the unhappiness factors are much closer to unity. A better agreement between theory and experiment has already been published,¹⁹ but in the present calculations there are still questions about the angular oscillations which are in not so good agreement with experiment. Note that Guazzoni *et al.*²⁸ took the better agreement of the *simultaneous* transfer

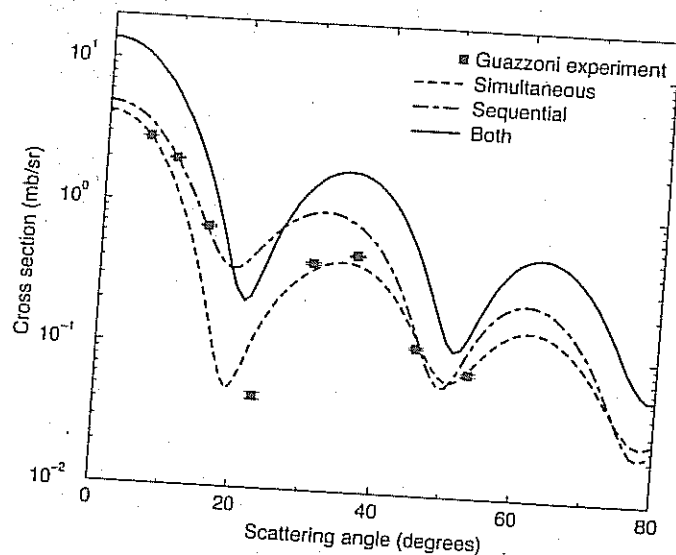


Fig. 1. Simultaneous (short dash), sequential (dot-dash) and simultaneous + sequential (solid line) cross sections for the reaction $^{124}\text{Sn}(p,t)^{122}\text{Sn}$ at 25 MeV, in comparison with the experimental data of Guazzoni *et al.*²⁸

curve (dashed line) to indicate small effects for sequential transfers, but this is not correct since we do know that sequential transfers occur, and can calculate them with good accuracy in this model (dot-dashed line).

To see the importance of the nonorthogonality terms, and hence of choosing 'prior-post' couplings if nonorthogonality terms are to be avoided, Fig. 2 plots the different sequential cross sections for all possible combinations of post and prior for the two steps. The prior-post solid curve is the dot-dashed curve in Fig. 1. The other curves are all different from this one, and cannot be simply added to the simultaneous amplitude to get the correct result. This also implies that no complete calculation with only zero-range couplings is possible.

Finally, it is instructive to look at the interference effects between the various simultaneous and sequential contributions. To display these coherence effects, I choose to plot the scattering amplitude at zero degrees for the non-spin-flip amplitude $m_p = m_t = 1/2$ (the only nonzero amplitude at this angle). Figure 3 plots all the simultaneous and sequential contributions from the different components listed in Table 1, along with their coherent sums. We see that all the contributions to the simultaneous transfer are constructively coherent, as are all the contributions to the total sequential amplitude. This constructive coherence follows from the signs of the amplitudes in Table 1, and reflects the significant pairing enhancement in ^{124}Sn . The total sequential and simultaneous amplitudes are not uniformly coherent with each other, however. A uniform 90° angle between the simultaneous and sequential amplitudes in Fig. 3 would indicate an incoherent summation of the two

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~~Guazzoni~~
~~Edwards~~
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✓ Appendix 1.E Two-nucleon spectroscopic amplitudes associated with pairing vibrational modes in closed shell systems: the ^{208}Pb case.

The solution of the pairing Hamiltonian

$$H = H_{sp} + H_p,$$

where

$$H_{sp} = \sum_v \epsilon_v a_v^\dagger a_v$$

and

$$H_p = -GP^\dagger P,$$

with

$$P^\dagger = \sum_{v>0} a_v^\dagger a_v^\dagger,$$

for the case of closed shell systems and in the harmonic approximation (RPA) leads to pair addition (a) pair removal (r) two-particle, two-hole correlated modes, the associated creation and annihilation operator being

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

and

$$\Gamma_r^\dagger(n) = \sum_i X_n^r(i) \Gamma_i^\dagger + \sum_k Y_n^r(k) \Gamma_k,$$

with

$$\sum X^2 - Y^2 = 1$$

and

$$\Gamma_k^\dagger = a_k^\dagger a_k^\dagger, \quad (\epsilon_k > \epsilon_F),$$

and

$$\Gamma_i^\dagger = a_i a_i, \quad (\epsilon_i \leq \epsilon_F).$$

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, while n labels the roots of the corresponding dispersion relations in increasing order of energy, ~~(1, 2, 3, ...)~~

$$\frac{1}{G(\pm 2)} = \sum_k \frac{(\Omega_k/2)}{2\epsilon_k \mp W_n(\pm 2)} + \sum_i \frac{(\Omega_i/2)}{2\epsilon_i \pm W_n(\pm 2)},$$

Pair vibration

1.E. (TWO-NUCLEON SPECTROSCOPIC AMPLITUDES ASSOCIATED WITH PAIRING VIBRATION) 1

orbit	ϵ_j	$\epsilon_{p1/2} - \epsilon_j \equiv \epsilon_j - \epsilon_{p1/2} $
$0h_{9/2}$	-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
$\epsilon_F = -5.825 \text{ keV}$		$\epsilon_k - \epsilon_{g9/2} \equiv \epsilon_{g9/2} - \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 1.E.1: Valence single particle levels of ^{208}Pb . In the upper part the occupied levels ($\epsilon_i \leq \epsilon_F$) are shown while in the lower part the empty levels ($\epsilon_k > \epsilon_F$). Of notice that $\epsilon_{p1/2} - \epsilon_{g9/2} = 3.41 \text{ MeV}$, is the single-particle gap associated associated with $N = 126$ shell closure.

where $\Omega_j = j + 1/2$ is the pair degeneracy.

For the case of the (neutron) pair addition and pair subtraction modes of ^{208}Pb the above equation can be solved graphically (see Fig 1.E.1). The minimum of the dispersion relation defining the Fermi energy. This is in keeping with the fact that in the case in which $W_1(\beta = +2) = W_1(\beta = -2) = 0$, corresponding to the phase transition between normal and superfluid phases, in which case the Fermi energy value is well defined, the BCS λ variational parameter coincides with ϵ_F . It is of notice that, as a rule the Fermi energy of closed shell nuclei is defined as half the energy difference between the last occupied and the first empty single particle state (cf. e.g. ?). One then obtains

$$\begin{cases} E_{\text{corr}}(+2) = B(208) + B(210) - 2B(209) = 1.248 \text{ MeV} , \\ E_{\text{corr}}(-2) = B(208) + B(206) - 2B(207) = 0.630 \text{ MeV} , \end{cases}$$

$W_1(+2) + W_1(-2) = (B(208) - B(206)) - (B(210) - B(208)) = 14.11 - 9.115 = 4.995 \text{ MeV}$. It is of notice that in the above derivations all energies are > 0 . In particular, (see Table 1.E.1)

$$\epsilon_i < \epsilon_F \Rightarrow \epsilon_F - \epsilon_i = -|\epsilon_F| + |\epsilon_i| = |\epsilon_i| - |\epsilon_F| > 0 ,$$

and

$$\epsilon_k > \epsilon_F \Rightarrow \epsilon_k - \epsilon_F = -|\epsilon_k| + |\epsilon_F| = |\epsilon_F| - |\epsilon_k| > 0 ,$$

of the system under study (^{208}Pb in the present case).

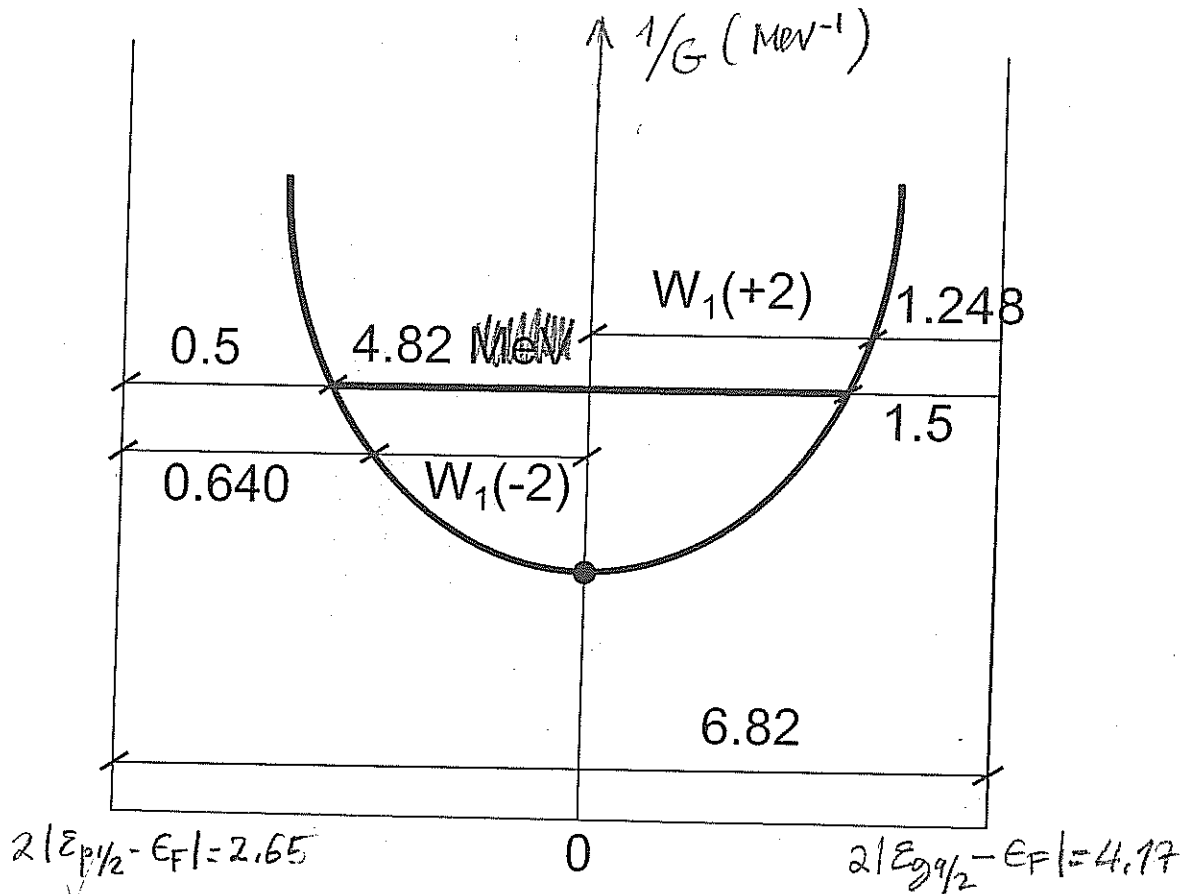
It is of notice that

The energy value at which the dispersion relation touches for the first time the corresponding axis, coincides with the

One obtains

difference

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✓ Figure 1.E.1: The right hand side of the RPA pairing vibrational dispersion relation for neutrons in the case of the closed shell system ^{208}Pb (cf. 1.7) in the region between the two neighboring shells ($p_{1/2}$ and $g_{9/2}$). All quantities are in MeV. For each 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 pairing vibrational state $(2p - 2h)$ of ^{208}Pb .

(Bes and
Broglia,
1966)

(44)

1.E. TWO-NUCLEON SPECTROSCOPIC AMPLITUDES ASSOCIATED WITH PAIRING VIBRATIONAL

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. 1.E.1 one can write,

$$2(\epsilon_{g_{9/2}} - \epsilon_F) = W_1(+2) + E_{corr}(+2),$$

and thus

$$4.17 \text{ MeV} = 2(-3.74 \text{ MeV} - (-5.825) \text{ MeV}) = W_1(+2) + 1.248 \text{ MeV}.$$

Consequently,

$$W_1(-2) = 2.01 \text{ MeV}, \quad W_1(+2) + W_1(-2) = 4.93 \text{ MeV}.$$

Let us make a rigid shift of energies setting $\epsilon_F = 0$. Thus,

$$W_1(+2) = 2\epsilon_{g_{9/2}} - E_{corr}(+2) \quad W_1(-2) = -2\epsilon_{p_{1/2}} - E_{corr}(-2).$$

1.E.1 Pair removal mode

In Fig. 1.E.2 the graphical representation of the forwards going RPA amplitude of the pair removal mode is shown. Its expression is

$$X_1^r(i) = \frac{\frac{1}{2}\Omega_i^{1/2}\Lambda(-2)}{2(\epsilon_F - \epsilon_i) - W_1(-2)}$$

where

$$\begin{aligned} 2(\epsilon_F - \epsilon_i) - W_1(-2) &= 2(\epsilon_F - \epsilon_i) - 2(\epsilon_F - \epsilon_{p_{1/2}}) - E_{corr}(-2) \\ &= 2(\epsilon_{p_{1/2}} - \epsilon_i) + E_{corr}(-2) = 2(|\epsilon_i| - |\epsilon_{p_{1/2}}|) + E_{corr}(-2). \end{aligned}$$

Thus,

$$X_1^r(i) = \frac{\frac{1}{2}\Omega_i^{1/2}\Lambda(-2)}{2(|\epsilon_i| - |\epsilon_{p_{1/2}}|) + E_{corr}(-2)}.$$

Making use of the empirical value of $E_{corr}(-2)$ worked out above one obtains,

$$X_1^r(i) = \frac{\frac{1}{2}\Omega_i^{1/2}\Lambda(-2)}{2(|\epsilon_i| - |\epsilon_{p_{1/2}}|) + 0.64 \text{ MeV}} \quad 0.640$$

In Fig. 1.E.3 we display the graphical process associated with the backwards going RPA amplitude,

$$Y_1^r(k) = \frac{\frac{1}{2}\Omega_k^{1/2}\Lambda(-2)}{2(\epsilon_k - \epsilon_F) + W_1(-2)}.$$

Making use of

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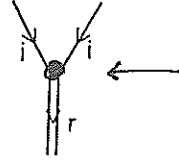


Figure 1.E.2: NFT representation of the forwards going RPA amplitude of the pair removal mode (double downward going arrowed line) describing a two correlated hole state (single downward going arrowed line for each hole with quantum numbers collectively labeled i).

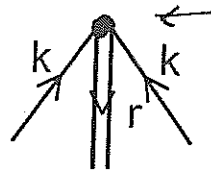


Figure 1.E.3: Same as Fig. 1.E.2 but for the backwards going amplitudes.

$$\begin{aligned} 2(\epsilon_k - \epsilon_F) + 2(\epsilon_F - \epsilon_{p_{1/2}}) - E_{corr}(-2) &= 2(\epsilon_k - \epsilon_{p_{1/2}}) - E_{corr}(-2) \\ &= 2(|\epsilon_{p_{1/2}}| - |\epsilon_k|) - E_{corr}(-2) = 2(|\epsilon_{p_{1/2}}| - |\epsilon_{g_{9/2}}|) + 2(|\epsilon_{g_{9/2}}| - |\epsilon_k|) - E_{corr}(-2), \end{aligned}$$

One can write,

$$Y_1^r(k) = \frac{\frac{1}{2}\Omega_k^{1/2}\Lambda(-2)}{2(|\epsilon_{g_{9/2}}| - |\epsilon_k|) + 2(|\epsilon_{p_{1/2}}| - |\epsilon_{g_{9/2}}|) - E_{corr}(-2)}$$

to finally obtain, making use of $2(|\epsilon_{p_{1/2}}| - |\epsilon_{g_{9/2}}|) - E_{corr}(-2) = 6.82\text{MeV} - 0.640\text{MeV} = 6.18\text{MeV}$, the expression

$$Y_1^r(k) = \frac{\frac{1}{2}\Omega_k^{1/2}\Lambda(-2)}{2(|\epsilon_{g_{9/2}}| - |\epsilon_k|) + 6.18\text{MeV}}.$$

Because (see Fig. 1.A.1)

The above calculated expressions of $X_1^r(i)$ and $Y_1^r(k)$ contain the experimental values of the 2-hole correlation energies (0.640 MeV). In keeping with the fact that 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 intermediate value and use the resulting $E_{corr}(-2)$ (≈ 0.5 MeV) and $E_{corr}(+2)$ (≈ 1.5 MeV), and calculate with these values the corresponding X, Y amplitudes for both the lowest removal and lowest addition pairing modes. Thus, making use of from

$$X_1^r(i) = \frac{\frac{1}{2}\Omega_i^{1/2}\Lambda(-2)}{2(|\epsilon_i| - |\epsilon_{p_{1/2}}|) + E_{corr}(-2)}; \quad Y_1^r(k) = \frac{\frac{1}{2}\Omega_k^{1/2}\Lambda(-2)}{2(|\epsilon_{g_{9/2}}| - |\epsilon_k|) + 2(|\epsilon_{p_{1/2}}| - |\epsilon_{g_{9/2}}|) - E_{corr}(-2)}$$

and of and

$$2(|\epsilon_{p_{1/2}}| - |\epsilon_{g_{9/2}}|) = 6.82\text{MeV}; \quad 2(|\epsilon_{p_{1/2}}| - |\epsilon_{g_{9/2}}|) - E_{corr} = (6.82 - 0.5)\text{MeV} = 6.32\text{MeV},$$

(Tamm-Dancoff; TD)

It is of notice that the coupling strength with which the pair removal mode couples to the single-particle (hole) states is calculated by normalizing the amplitudes: $1) \sum_i A_i^2 = 1.5549 \text{ MeV}^{-2}$ and thus $\Lambda(-2) = 0.802 \text{ MeV}$, $\sum_i (X_i^r(i))^2_{TD} = 1$; 2) RPA, $\Lambda_1^2(-2) (\sum_i A_i^2(i) - \sum_k B^2(k)) = \Lambda_1^2(-2) \times 1.45073$.

1.E. TWO-NUCLEON SPECTROSCOPIC AMPLITUDES ASSOCIATED WITH PAIRING VIBRATIONAL

units		MeV	MeV ⁻¹	RPA	TD
nlj	Ω_i	$ \epsilon_i - \epsilon_{p1/2} $	$A(i) = \frac{\frac{1}{2} \Omega_i^{1/2} \Lambda(-2)}{2(\epsilon_i - \epsilon_{p1/2}) + 0.5 \text{ MeV}}$	$X_1^r(i)$	$X_1^r(i)$
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 1.E.2: Forwards going RPA amplitudes of the pair removal mode of ²⁰⁸Pb (i.e., ²⁰⁶Pb) state), cf. Table XVI Broglia et al. (1973).

units		MeV	MeV ⁻¹	
nlj	Ω_k	$ \epsilon_{g9/2} - \epsilon_k $	$B(k) = \frac{\frac{1}{2} \Omega_k^{1/2} \Lambda(-2)}{2(\epsilon_{g9/2} - \epsilon_k) + 6.23 \text{ MeV}}$	$Y_1^r(k)$
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 1.E.3: Same as Table 1.E.2 but for the backwards amplitude

Thus $\Lambda(-2) = 0.83$. The above results show that there is a few percent difference between the two values of Λ (TD and RPA) as well as for the corresponding X-amplitudes. Nonetheless, ground state corre-

$$\frac{0.830}{0.802} = \frac{0.028}{0.028} \approx 4\%$$

Integrated relations as expressed by the Y-amplitudes, give rise to a 50% increase in the ²⁰⁶Pb(t,p) ²⁰⁸Pb(gs) section, from 0.34 mb to 0.52 mb, to be compared with the experimental data ($\sigma = 0.68 \pm 24 \text{ mb}$), see Fig. 1.E.4

Fig. sem. Gregory

one can write,

²⁰⁶Pb(gs)

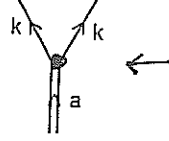
$$\begin{cases} X_1^r(i) = \frac{\frac{1}{2} \Omega_i^{1/2} \Lambda(-2)}{2(|\epsilon_i| - |\epsilon_{p1/2}|) + 0.5 \text{ MeV}} \\ Y_1^r(k) = \frac{\frac{1}{2} \Omega_k^{1/2} \Lambda(-2)}{2(|\epsilon_{g9/2}| - |\epsilon_k|) + 6.23 \text{ MeV}} \end{cases}$$

Tables 1.E.2 and 1.E.3 contain the amplitudes of the pair removal mode of ²⁰⁸Pb ($\Gamma_1^+(1) = \sum_i X_1^r(i) \Gamma_i^+ + \sum_k Y_1^r(k) \Gamma_k^+$) that is the two neutron (hole) correlated state describing ²⁰⁶Pb = $\Gamma_1^+(1)|0\rangle$.

It is of notice that while the harmonic (RPA) description of the pair vibrational mode of ²⁰⁸Pb provides a fair description 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 calculated (cf. for example Flynn, Bortignon, Clark).

Let us conclude this Appendix by noting that

to p. 19



(47)

Figure 1.E.4. Same as Fig. 1.E.2 but for the pair addition mode

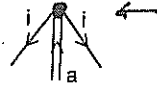


Figure 1.E.5. Same as Fig. 1.E.3 but for the pair addition mode

1.E.2 Pair addition modes

In Figs. 1.E.4 and 1.E.5 the X and Y amplitudes of the pair addition mode are given. The associated expressions are given below:

$$X_1^a(k) = \frac{\frac{1}{2}\Omega_k^{1/2}\Lambda_1(+2)}{2(\epsilon_k - \epsilon_F) - W_1(+2)},$$

which, making use of

$$\begin{aligned} 2(\epsilon_k - \epsilon_F) - W_1(+2) &= 2(\epsilon_k - \epsilon_F) - 2(\epsilon_{g_{9/2}} - \epsilon_F) - E_{corr}(+2) \\ &= 2(\epsilon_k - \epsilon_{g_{9/2}}) + E_{corr}(+2) = 2(|\epsilon_{g_{9/2}}| - |\epsilon_k|) + E_{corr}(+2) \end{aligned}$$

can be written as

$$X_1^a(k) = \frac{\frac{1}{2}\Omega_k^{1/2}\Lambda_1(+2)}{2(|\epsilon_{g_{9/2}}| - |\epsilon_k|) + W_1(+2)}.$$

Similarly,

$$Y_1^a(i) = \frac{\frac{1}{2}\Omega_i^{1/2}\Lambda_1(+2)}{2(\epsilon_F - \epsilon_i) + W_1(+2)}.$$

which with the help of the relation,
Making use of the relation

$$\begin{aligned} 2(\epsilon_F - \epsilon_i) + W_1(+2) &= 2(\epsilon_F - \epsilon_i) - 2(\epsilon_{g_{9/2}} - \epsilon_F) - E_{corr}(+2) \\ &= 2(\epsilon_{p_{1/2}} - \epsilon_i) + 2(\epsilon_{g_{9/2}} - \epsilon_{p_{1/2}}) - E_{corr}(+2) \\ &= 2(|\epsilon_i| - |\epsilon_{p_{1/2}}|) + 2(|\epsilon_{p_{1/2}}| - |\epsilon_{g_{9/2}}|) + E_{corr}(+2), \end{aligned}$$

can be written 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 $E_{corr}(+2) = 1.5 \text{ MeV}$ (cf. Fig. 1.E.1) and

$$\Delta\epsilon_{sp} = 2(|\epsilon_{p_{1/2}}| - |\epsilon_{g_{9/2}}|) = 6.28 \text{ MeV},$$

(48)

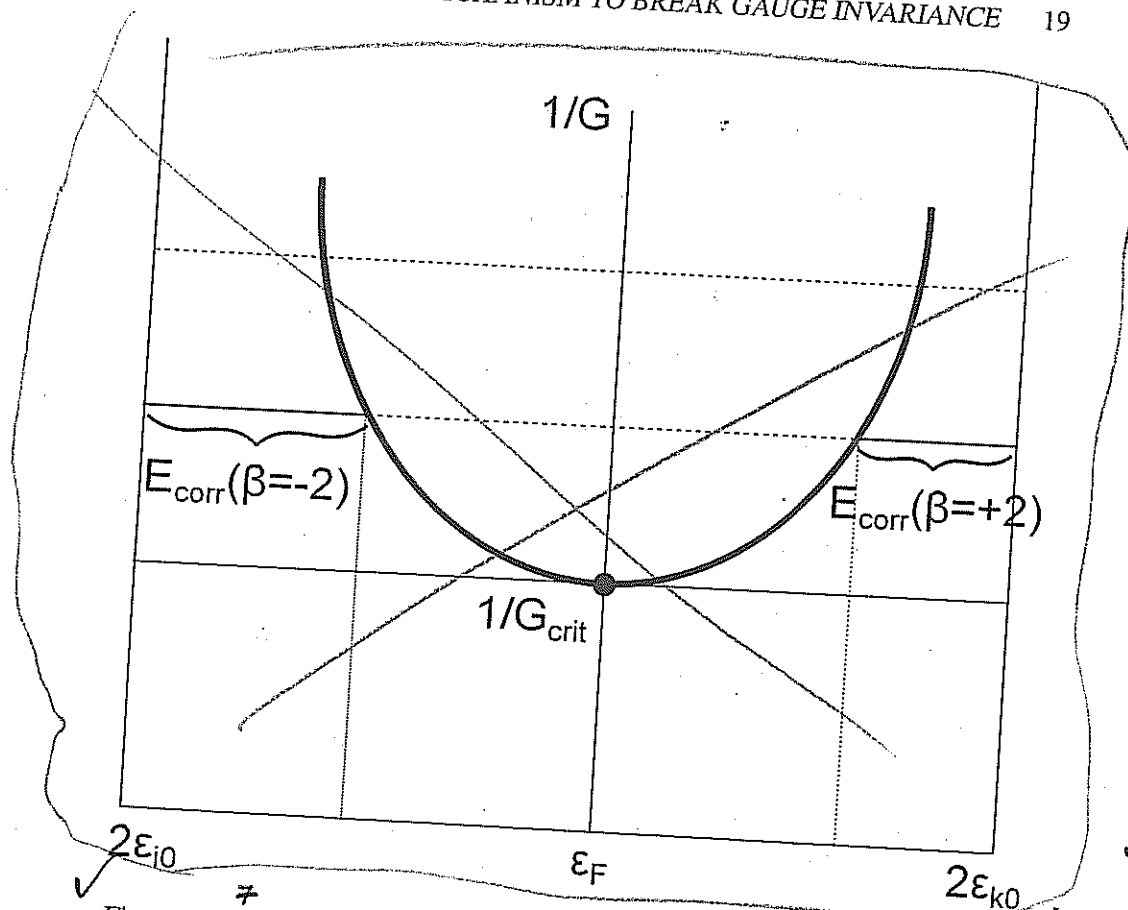


Figure 1.E.6: Same as Fig. 1.E.1 but schematically emphasizing the different elements characterizing the pairing vibration dispersion relation.

one thus obtains $2\Delta\epsilon_{sp} - E_{corr} = (6.82 - 1.5) \text{ MeV} = 5.32 \text{ MeV}$, leading to

$$\begin{cases} X_1^a(k) = \frac{\frac{1}{2}\Omega_k^{1/2}\Lambda(-2)}{2(|\epsilon_{q9/2}| - |\epsilon_k|) + 1.5 \text{ MeV}} \\ Y_1^a(i) = -\frac{\frac{1}{2}\Omega_i^{1/2}\Lambda(+2)}{2(|\epsilon_i| - |\epsilon_{p1/2}|) + 5.32 \text{ MeV}} \end{cases}$$

The corresponding numerical values are displayed in Tables 1.E.4 and 1.E.5 (see also Fig. 1.E.7).

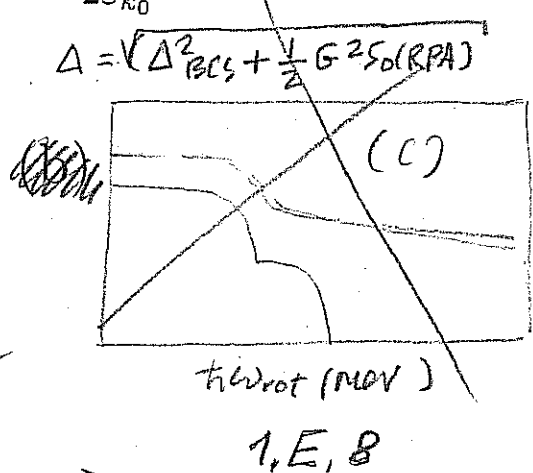
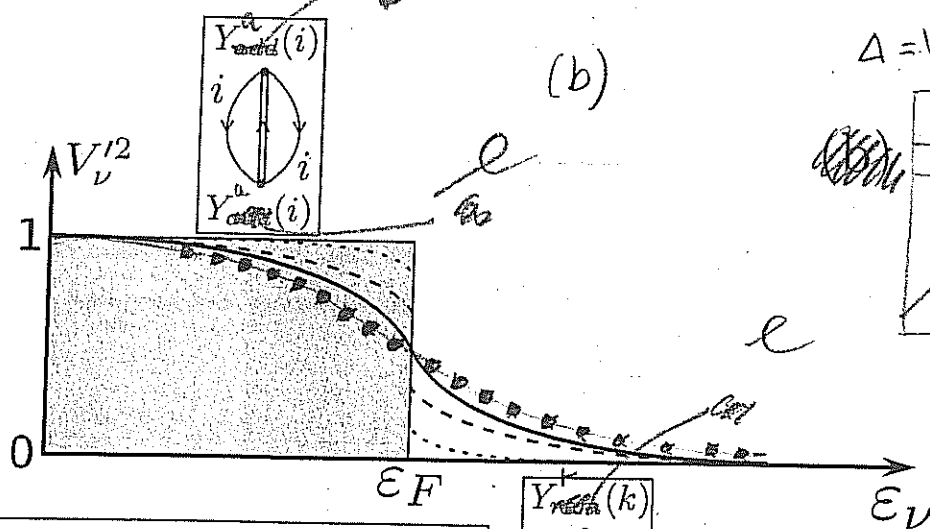
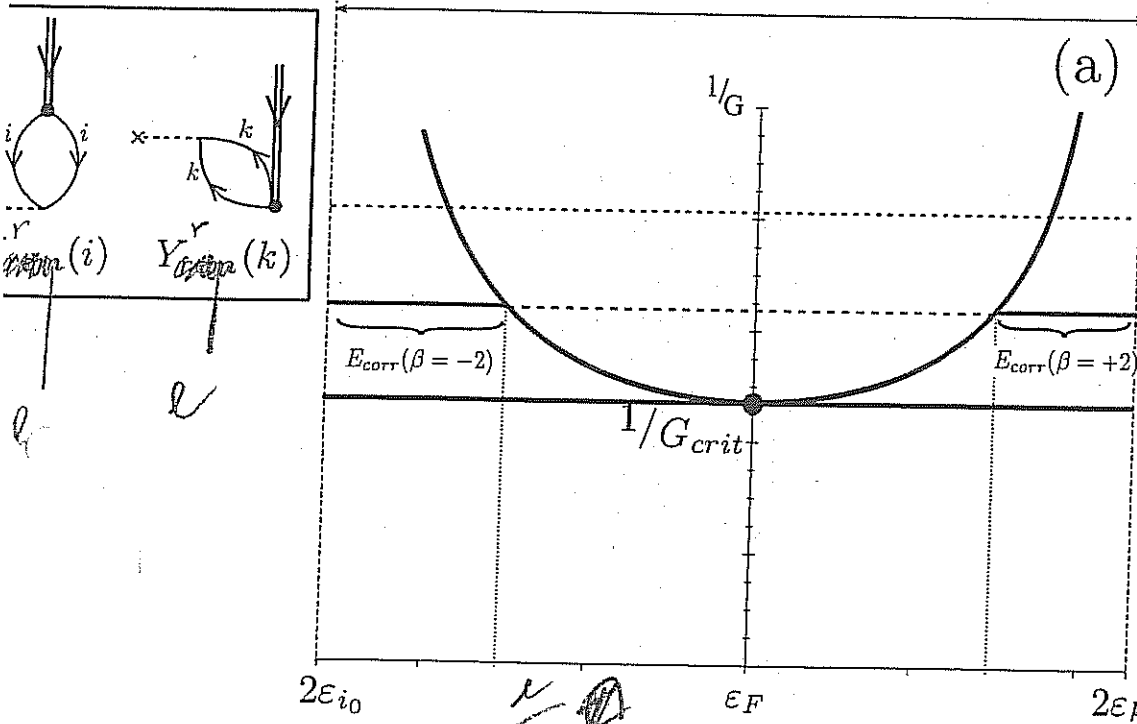
Appendix 1.F Microscopic mechanism to break gauge invariance

Pairing is intimately connected with particle number violation and thus spontaneous breaking of gauge invariance, as testified by the order parameter $\langle BCS | P^\dagger | BCS \rangle = v_0$. Now, in the nuclear case and at variance with condensed matter, dynamical

Within the framework of Fig. 1.D.1 (parallel between 3D and gauge phenomena of spontaneous symmetry breaking), we display in Fig. 1.E.8, the relative importance of dynamic and static pairing distortions, in comparison with the corresponding quantities in the case of quenching of nuclear distortions in 3-D space.

These results underscore the major role pairing vibrations play in nuclei around closed shells.

(49)



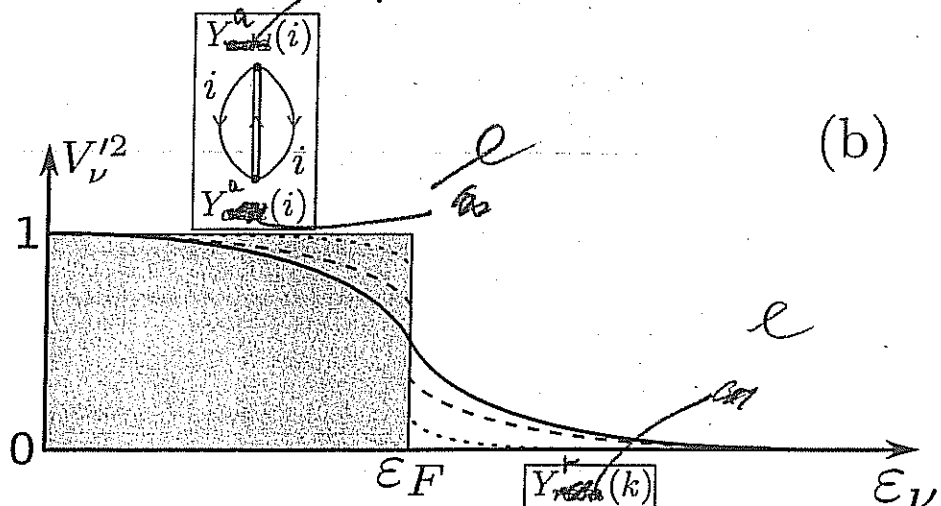
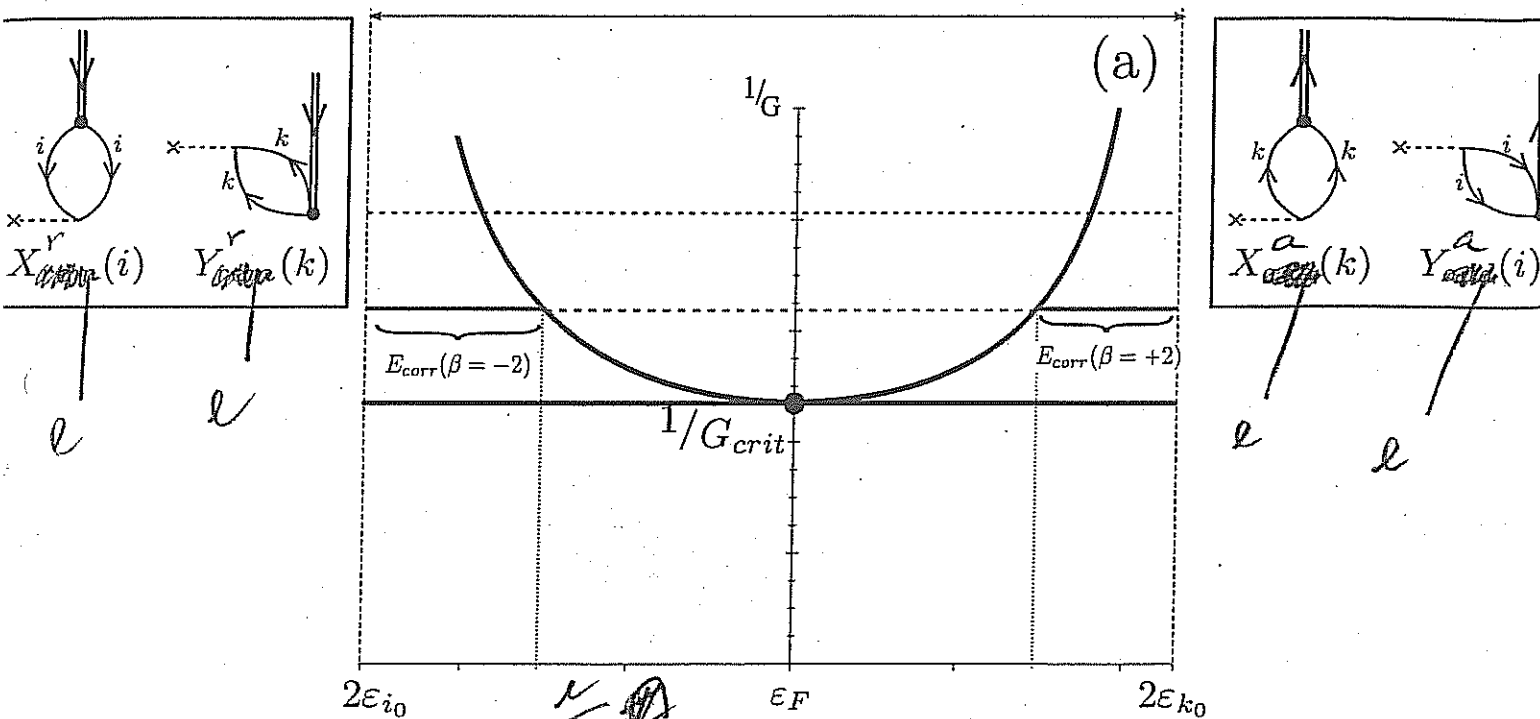
$(\langle 0 | P^\dagger P | 0 \rangle)^{1/2} = \begin{cases} \text{dynamic} & \cdots \\ \text{static} & - \end{cases}$

1.E.7
Fig. 1.1

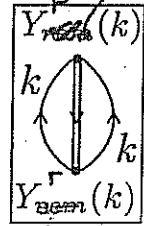
1.E.8
Fig. 1.2

Fig. 1.3
 $\chi = \frac{2G\Omega_c}{D} = 6N(0)$
 $\frac{d\chi}{d\mu}$
 Fig. 2

(50)



$$(\langle 0 | P^\dagger P | 0 \rangle)^{1/2} = \begin{cases} \text{dynamic} & \cdots \cdots \\ \text{static} & \text{---} \end{cases}$$



~~Fig. 2.1.2~~

Fig. 2

$$P^\dagger = \sum_{\nu > 0} a_\nu^\dagger a_\nu^\dagger$$

$$x = \frac{2G\Omega}{D} = GN(0)$$

$$x > 1$$

$$\alpha_0 = \langle P^\dagger \rangle = \frac{\Delta}{G} \approx 7$$

$$x < 1$$

$$\alpha_{dyn} = \frac{1}{G} \frac{\langle PP^\dagger \rangle^{1/2} + \langle P^\dagger P \rangle^{1/2}}{2}$$

$$\approx \frac{1}{2} \left(\frac{E_{corr}(A+2)}{G} + \frac{E_{corr}(A-2)}{G} \right) \approx 10$$

$$\frac{\alpha_0}{\alpha_{dyn}} \approx 0.7$$

$$\frac{\beta_2}{(\beta_2)_{dyn}} \approx 3 - 6$$

1.E.8

Fig. 1.E.8

Relative importance of dynamic pairing distortion around closed shell and for pairing deformed nuclei, as compared with similar quantities for the case of quadrupole surface degree of freedom.

(51)

(52)

units		MeV	MeV ⁻¹	
nlj	Ω_k	$ \epsilon_{g9/2} - \epsilon_k $	$C(k) = \frac{\frac{1}{2}\Omega_k^{1/2}}{2(\epsilon_{g9/2} - \epsilon_k) + 1.5 \text{ MeV}}$	$X_1^a(k)$
$1g_{9/2}$	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 1.E.4: Forwards going RPA amplitudes associated with the pair addition mode of ^{208}Pb (cf Table XVI Broglia et al (1973))

a) $\sum_k C^2(k) = 0.903$

units		MeV	MeV ⁻¹	
nlj	Ω_i	$ \epsilon_i - \epsilon_{p1/2} $	$D(i) = \frac{\frac{1}{2}\Omega_i^{1/2}}{2(\epsilon_i - \epsilon_{p1/2}) + 5.32 \text{ MeV}}$	$Y_1^a(i)$
$2p_{1/2}$	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 1.E.5: Same as Table 1.E.4 but for the backwards going amplitude.

a) $\sum_i D^2(i) = 0.079$; Thus $\Lambda^2(+2)(\sum_k C^2(k) - D^2(i)) = \Lambda^2(+2)(0.903 - 0.079) \text{ MeV}^{-2}$
 $= 0.824 \text{ MeV}^{-2}$; $\Lambda(+2) = \text{MeV}(0.824)^{-1/2} = 1.102 \text{ MeV}$;
 $\Lambda(+2) = 1.102 \text{ MeV}$.

53

breaking of gauge symmetry is equally important (pairing vibrations around closed shell nuclei, cf. Fig. ??). The fact that the average single-particle field acts as an external potential (like e.g. magnetic field in metallic superconductors) is at the basis of the existence of a critical value of the pairing strength G to bind Cooper pairs in nuclei. In fact, spatial quantization in finite systems at large and in nuclei in particular, intimately connect with the tantamount role the surface has in these systems, is at the basis of the existence of a critical G value. Also of the fact that in nuclei an important fraction (30-50%) of Cooper pair binding is due to the exchange of collective vibrations between the partners of the pair, the rest being associated with the bare NN interaction in the 1S_0 channel (cf. Fig. ??).

Now, there are situations in which spatial quantization seems, essentially completely, the NN -interaction. This happens in the case in which the nuclear valence orbitals s , p -states at threshold (pairing anti halo effect). Examples of situations of this type are provided by $N = 6$ (parity inversion) isotones. In particular, by ^{11}Li , in which case the strongly renormalized $s_{1/2}$ and $p_{1/2}$ valence orbitals are a virtual and a resonant state lying at ≈ 0.1 and 0.6 MeV in the continuum, respectively. In keeping with the fact that the binding provided to a pair of fermions moving in time reversal states by a contact pairing interaction (δ -force) is (cf. e.g. Eq. (2.12) Brink and Broglia (2005)) $E_0 = (2j+1)/2V_0J(j) \approx \frac{(2j+1)}{2}V_0\frac{3}{R^3}$, the ratio

$$r = \frac{2}{(2j+1)} \left(\frac{R_0}{R} \right)^3,$$

where $R_0 = 1.2A^{1/3}\text{fm} = 2.7\text{fm}$ ($A = 11$), and $R = \sqrt{\frac{5}{3}}\langle r^2 \rangle_{^{11}\text{Li}}^{1/2} = \sqrt{\frac{5}{3}}3.74\text{ fm} = 4.6\text{ fm}$ are the radius of a stable nucleus of mass $A = 11$ (systematics), while R is the measured one, while j is the angular momentum representative for a nucleus of mass $A = 11$ ($j \sim k_F R_0 \approx 3-4$), one obtains $r = 0.06$. Making use of the multipole expansion of a general interaction

$$v(|\mathbf{r}_1 - \mathbf{r}_2|) = \sum_{\lambda} V_{\lambda}(r_1, r_2) P_{\lambda}(\cos \theta_{12}).$$

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 $\lambda = 0$ term is 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 $^{11}\text{Li}(\text{gs})$ we are thus forced to accept the need for 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.

1.F.1 Bootstrap Cooper pair binding

Within the s, p subspace, the most natural low wavelength vibration is the dipole mode. From systematics, the centroid of these vibrations is $\hbar\omega_{GDR} \approx 100 \text{ MeV}/R$, R being the nuclear radius. Thus, in the case of ^{11}Li , one expects the centroid of the Giant Dipole Resonance carrying $\approx 100\%$ of the energy weighted sum rule (EWSR) at $\hbar\omega_{GDR} \approx 100 \text{ MeV}/2.7 \approx 37 \text{ MeV}$. Now, 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 of a rather sharp dipole state with centroid at $\approx 1 \text{ MeV}$ and carrying $\approx 10\%$ of the EWSR. The existence of this "pigmy resonance" which can be viewed as a simple consequence of the existence of a low-lying particle hole state associated with the transition $s_{1/2} \rightarrow p_{1/2}$, arguably, testifies to the coexistence of two states with rather different radii in the ground state¹. One, closely connected with the ^9Li core, ($\approx 2.7 \text{ fm}$), the second with the diffuse halo ($\approx 4.6 \text{ fm}$). Because the overlap between them is small ($\approx (2.7/4.6)^3 \approx 0.2$), one can posit that a bona fide dipole pigmy resonance is a GDR based on an exotic, unusually extended state as compared to systematics ($A \approx (4.6/1.2)^3 \approx 60$), i.e. to system with an effective A mass number about 5 times than predicted by systematics.

Let us try to shed some light 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 radius, one may write

$$\langle r^2 \rangle_{^{11}\text{Li}} \approx \frac{3}{5} R_{eff}^2(^{11}\text{Li}).$$

with

$$R_{eff}^2(^{11}\text{Li}) = \left(\frac{9}{11} R_0^2(^9\text{Li}) + \frac{2}{11} \left(\frac{\xi}{2} \right)^2 \right),$$

where

$$R_0(^9\text{Li}) = 2.5 \text{ fm}$$

is the ^9Li radius ($R_0 = r_0 A^{1/3}$, $r_0 = 1.2 \text{ fm}$), where ξ is the correlation length of the Cooper pair neutron halo. An estimate of this quantity is provided by the relation

$$\xi = \frac{\hbar v_F}{2E_{corr}} \approx 20 \text{ fm},$$

in keeping with the fact that in ^{11}Li , $(v_F/c) \approx 0.1$ and $E_{corr} \approx 0.5 \text{ MeV}$. Consequently, $\langle r^2 \rangle_{^{11}\text{Li}}^{1/2} \approx 3.74 \text{ fm}$ ($R_{eff}^{11}\text{Li} \approx 4.83 \text{ fm}$), in overall agreement with the experimental value $\langle r^2 \rangle_{^{11}\text{Li}}^{1/2} = 3.55 \pm 0.1 \text{ fm}$ (Kobayashi et al., 1989).

We now proceed to the calculation of the centroid of the dipole pigmy resonance of ^{11}Li . Making use of the dispersion relation given in Eq. (3.30) p.55 of ?, and of the fact that $\epsilon_{v_k} - \epsilon_{v_i} = \epsilon_{p_{1/2}} - \epsilon_{s_{1/2}} \approx 0.5 \text{ MeV}$ (see Fig. 11.1 p.264 Brink

¹Within this context the dipole strength found at $\approx 10 \text{ MeV}$ in neutron skin rich nuclei can hardly be considered pigmy resonance, but the long tail of the GDR.

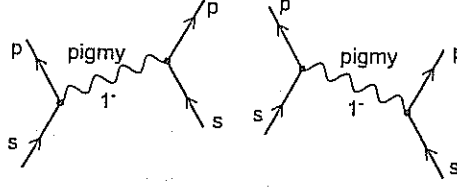


Figure 1.F.1:

and Broglia (2005)), and that the EWSR associated with the ^{11}Li pigmy resonance is $\approx 10\%$ of the total Thomas–Reiche–Kuhn sum rule one can write,

$$0.1 \frac{\hbar^2 A}{2M} = \frac{1}{\kappa_1} [(0.5\text{MeV})^2 - (\hbar\omega_{\text{pigmy}})^2],$$

and thus

$$(\hbar\omega_{\text{pigmy}})^2 = (0.5\text{MeV})^2 + 0.1 \frac{\hbar^2 A}{2M},$$

where (see ?)

$$\kappa_1 = \frac{5V_1}{A(\xi/2)^2} \left(\frac{2}{11} \right) = -\frac{125\text{MeV}}{A100\text{fm}^2} \left(\frac{2}{11} \right) \approx -\frac{2.5}{A^2} \text{fm}^{-2} \text{MeV},$$

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. One then obtains,

$$-0.1 \frac{\hbar^2 A}{2M} \kappa_1 = 0.1 \times 20\text{MeV fm}^2 A \times \frac{2.5}{A^2} \text{fm}^{-2} \text{MeV} \approx 0.45\text{MeV} \approx (0.7\text{MeV})^2$$

consequently

$$\hbar\omega_{\text{pigmy}} = \sqrt{(0.5)^2 + (0.7)^2} \text{MeV} \approx 1\text{MeV},$$

in overall agreement with the experimental findings (?). It is of notice that the centroid of the pigmy resonance calculated in the RPA with the help of a separable interaction is $\approx (0.8\text{MeV} + 2.0\text{MeV})/2 \approx 1.4\text{MeV}$ (see Fig. 11.3(a) p.269, Brink and Broglia (2005). Let us now estimate the binding which the exchange of the pigmy resonance between two neutron of the Cooper pair halo of ^{11}Li can provide. The associated particle vibration coupling $\Lambda \left(\frac{\partial W(E)}{\partial E} \right)_{\hbar\omega_{\text{pigmy}}}^{-1/2}$, where $W(E)$ is the dispersion relation used to determine $\hbar\omega_{\text{pigmy}}$ (cf. e.g. Brink and Broglia (2005) Eq. (8.42) p.189; note the use of a dimensionless dipole single particle field $F' = F/\langle r^2 \rangle_{^{11}\text{Li}}$)

$$W(E) = \frac{2(\epsilon_k - \epsilon_i) |\langle \tilde{i} | F / \langle r^2 \rangle_{^{11}\text{Li}} | k \rangle|^2}{(\epsilon_k - \epsilon_i)^2 - E^2}$$

(56)

24 CHAPTER 1. STRUCTURE AND PAIR TRANSFER IN A NUTSHELL

One then obtains

$$\begin{aligned}
 \Lambda^2 &= \left\{ 2\hbar\omega_{pigm} \frac{0.1(TRK)/\langle r^2 \rangle_{^{11}\text{Li}}}{[(\epsilon_{p_{1/2}} - \epsilon_{s_{1/2}})^2 - (\hbar\omega_{pigm})^2]^2} \right\}^{-1} \\
 &= \left\{ 2\text{Mev} \frac{0.1(\hbar^2 A/2M)(1/\langle r^2 \rangle_{^{11}\text{Li}})}{[(0.5)^2 - (1\text{MeV})^2]^2 \text{MeV}^4} \right\}^{-1} \\
 &= \left(\frac{0.75}{1.57} \right)^2 = 0.48 \text{MeV}^2
 \end{aligned}$$

leading to $\Lambda = 0.7\text{MeV}$. The value of induced interaction matrix elements is then given by

$$M_{ind} = -\frac{\Lambda^2}{\hbar\omega_{pigm}} = -0.5\text{MeV},$$

and the same contribution for the other time ordering. Assuming the halo neutrons to spend the same amount of time in the $|s_{1/2}^2(0)\rangle(\epsilon_{s_{1/2}} = 0.1 \text{ MeV})$ than in the $|p_{1/2}^2(0)\rangle(\epsilon_{p_{1/2}} = 0.6 \text{ MeV})$ configuration, the correlation energy is $E_{corr} = |2(\epsilon_{s_{1/2}} + \epsilon_{p_{1/2}})/2 + 2M_{ind}| = 0.3 \text{ MeV}$, in overall agreement with the findings (0.380 MeV).

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