

Let us examine the term

$$\sum_M (-1)^M \langle K \Lambda - M | P | \mu - M \rangle \left[Y^{l_1}(\hat{\mathbf{r}}_{A1}) Y^{l_1}(\hat{\mathbf{r}}_{B1}) \right]_{\Lambda}^K \left[Y^{l_2}(\hat{\mathbf{r}}'_{A2}) Y^{l_2}(\hat{\mathbf{r}}'_{B2}) \right]_{\mu-M}^P. \quad (7.2.167)$$

Making use of the relation

$$\langle l_1 l_2 m_1 m_2 | L M_L \rangle = (-1)^{l_2 - m_2} \sqrt{\frac{2L+1}{2l_1+1}} \langle L l_2 - M_L m_2 | l_1 - m_1 \rangle, \quad (7.2.168)$$

the expression (7.2.168) is equivalent to

$$(-1)^K \sqrt{\frac{2P+1}{2\Lambda+1}} \left\{ \left[Y^{l_1}(\hat{\mathbf{r}}'_{A2}) Y^{l_2}(\hat{\mathbf{r}}'_{B2}) \right]^P \left[Y^{l_1}(\hat{\mathbf{r}}_{A1}) Y^{l_1}(\hat{\mathbf{r}}_{B1}) \right]^K \right\}_{\mu}^{\Lambda}. \quad (7.2.169)$$

We now recouple *the term* ←

$$\left[Y^{l_a}(\hat{\mathbf{r}}'_{aA}) Y^{l_b}(\hat{\mathbf{r}}_{aA}) \right]_0^0 \left[Y^{l_b}(\hat{\mathbf{r}}_{bB}) Y^{l_b}(\hat{\mathbf{r}}_{bB}) \right]_0^0, \quad (7.2.170)$$

arising from the partial wave expansion of the incoming and outgoing distorted waves *to have,*

$$((l_a l_b)_0 (l_b l_b)_0 | (l_a l_b)_{\Lambda} (l_a l_b)_{\Lambda})_0 \left\{ \left[Y^{l_a}(\hat{\mathbf{r}}'_{aA}) Y^{l_b}(\hat{\mathbf{r}}_{bB}) \right]^{\Lambda} \left[Y^{l_a}(\hat{\mathbf{r}}_{aA}) Y^{l_b}(\hat{\mathbf{r}}_{bB}) \right]^{\Lambda} \right\}_0^0. \quad (7.2.171)$$

The only term which does not vanish upon integration is

$$\frac{(-1)^{\Lambda+\mu}}{\sqrt{(2l_a+1)(2l_b+1)}} \left[Y^{l_a}(\hat{\mathbf{r}}'_{aA}) Y^{l_b}(\hat{\mathbf{r}}_{bB}) \right]_{-\mu}^{\Lambda} \left[Y^{l_a}(\hat{\mathbf{r}}_{aA}) Y^{l_b}(\hat{\mathbf{r}}_{bB}) \right]_{\mu}^{\Lambda}. \quad (7.2.172)$$

Again, the only term surviving

$$\left\{ \left[Y^{l_1}(\hat{\mathbf{r}}'_{A2}) Y^{l_2}(\hat{\mathbf{r}}'_{B2}) \right]^P \left[Y^{l_1}(\hat{\mathbf{r}}_{A1}) Y^{l_1}(\hat{\mathbf{r}}_{B1}) \right]^K \right\}_{\mu}^{\Lambda} \left[Y^{l_a}(\hat{\mathbf{r}}'_{aA}) Y^{l_b}(\hat{\mathbf{r}}_{bB}) \right]_{-\mu}^{\Lambda} \quad (7.2.173)$$

is

$$\frac{(-1)^{\Lambda+\mu}}{\sqrt{2\Lambda+1}} \left[\left[Y^{l_1}(\hat{\mathbf{r}}'_{A2}) Y^{l_2}(\hat{\mathbf{r}}'_{B2}) \right]^P \left[Y^{l_1}(\hat{\mathbf{r}}_{A1}) Y^{l_1}(\hat{\mathbf{r}}_{B1}) \right]^K \right]_{\mu}^{\Lambda} \left[Y^{l_a}(\hat{\mathbf{r}}'_{aA}) Y^{l_b}(\hat{\mathbf{r}}_{bB}) \right]_{-\mu}^{\Lambda}. \quad (7.2.174)$$

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We now couple this last term with the term $[Y^{l_c}(\hat{\mathbf{r}}_{cC})Y^{l_c}(\hat{\mathbf{r}}_{cC})]_0^0$, arising from the partial wave expansion of the Green function. That is,

$$\begin{aligned}
 & \left[\left\{ [Y^{l_f}(\hat{\mathbf{r}}'_{A2})Y^{l_a}(\hat{\mathbf{r}}'_{b2})]^P [Y^{l_f}(\hat{\mathbf{r}}_{A1})Y^{l_b}(\hat{\mathbf{r}}_{b1})]^K \right\}^\Lambda [Y^{l_a}(\hat{\mathbf{r}}'_{aA})Y^{l_b}(\hat{\mathbf{r}}_{bB})]^\Lambda \right]_0^0 [Y^{l_c}(\hat{\mathbf{r}}_{cC})Y^{l_c}(\hat{\mathbf{r}}_{cC})]_0^0 \\
 &= ((l_a l_b)_\Lambda (l_c l_c)_0 (l_b l_c)_P (l_b l_c)_K)_\Lambda \left\{ \left[[Y^{l_f}(\hat{\mathbf{r}}'_{A2})Y^{l_a}(\hat{\mathbf{r}}'_{b2})]^P [Y^{l_f}(\hat{\mathbf{r}}_{A1})Y^{l_b}(\hat{\mathbf{r}}_{b1})]^K \right]^\Lambda \right\}^\Lambda \\
 & \left\{ [Y^{l_a}(\hat{\mathbf{r}}_{aA})Y^{l_c}(\hat{\mathbf{r}}_{cC})]^P [Y^{l_b}(\hat{\mathbf{r}}_{bB})Y^{l_c}(\hat{\mathbf{r}}_{cC})]^K \right\}_0^0 = ((l_a l_b)_\Lambda (l_c l_c)_0 (l_b l_c)_P (l_b l_c)_K)_\Lambda \\
 & \times ((PK)_\Lambda (PK)_\Lambda (PP)_0 (KK)_0)_0 \left\{ \left[[Y^{l_f}(\hat{\mathbf{r}}'_{A2})Y^{l_a}(\hat{\mathbf{r}}'_{b2})]^P [Y^{l_a}(\hat{\mathbf{r}}'_{aA})Y^{l_c}(\hat{\mathbf{r}}_{cC})]^P \right]_0^0 \right\}^0 \\
 & \times \left\{ \left[[Y^{l_f}(\hat{\mathbf{r}}_{A1})Y^{l_b}(\hat{\mathbf{r}}_{b1})]^K [Y^{l_b}(\hat{\mathbf{r}}_{bB})Y^{l_c}(\hat{\mathbf{r}}_{cC})]^K \right]_0^0 \right\}^\Lambda = ((l_a l_b)_\Lambda (l_c l_c)_0 (l_b l_c)_P (l_b l_c)_K)_\Lambda \\
 & \times \sqrt{\frac{2\Lambda + 1}{(2K + 1)(2P + 1)}} \left\{ \left[[Y^{l_f}(\hat{\mathbf{r}}'_{A2})Y^{l_a}(\hat{\mathbf{r}}'_{b2})]^P [Y^{l_a}(\hat{\mathbf{r}}'_{aA})Y^{l_c}(\hat{\mathbf{r}}_{cC})]^P \right]_0^0 \right\}^0 \\
 & \times \left\{ \left[[Y^{l_f}(\hat{\mathbf{r}}_{A1})Y^{l_b}(\hat{\mathbf{r}}_{b1})]^K [Y^{l_b}(\hat{\mathbf{r}}_{bB})Y^{l_c}(\hat{\mathbf{r}}_{cC})]^K \right]_0^0 \right\}^\Lambda.
 \end{aligned}
 \tag{7.2.175}$$

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

$$\begin{aligned}
 T_\mu^{succ} &= (-1)^{j_f + j_b} \frac{2048\pi^5 \mu_{Cc}}{\hbar^2 k_{Aa} k_{Bb} k_{Cc}} \sqrt{\frac{(2j_{i1} + 1)}{(2\Lambda + 1)(2j_f + 1)}} \sum_{K,P} ((l_f \frac{1}{2})_{j_f} (l_{i2} \frac{1}{2})_{j_a} (l_f l_{i2})_P (\frac{1}{2} \frac{1}{2})_0)_P \\
 & \times ((l_f \frac{1}{2})_{j_f} (l_{i1} \frac{1}{2})_{j_b} (l_f l_{i1})_K (\frac{1}{2} \frac{1}{2})_0)_K ((j_{i1} j_f)_K (j_{i1} j_{i2})_\Lambda (j_{i1} j_{i1})_0 (j_f j_{i2})_P)_P \\
 & \times \frac{(-1)^K}{(2K + 1) \sqrt{2P + 1}} \sum_{l_c, l_a, l_b} ((l_a l_b)_\Lambda (l_c l_c)_0 (l_a l_c)_P (l_b l_c)_K)_\Lambda e^{i(\sigma_a^{l_a} + \sigma_f^{l_b})} i^{l_a - l_b} \\
 & \times (2l_c + 1)^{3/2} [Y^{l_a}(\hat{\mathbf{k}}_{aA})Y^{l_b}(\hat{\mathbf{k}}_{bB})]_\mu^\Lambda S_{K,P,l_a,l_b,l_c},
 \end{aligned}
 \tag{7.2.176}$$

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 (7.2.132))

$$\begin{aligned}
 S_{K,P,l_a,l_b,l_c} &= \int r_{Cc}^2 dr_{Cc} r_{b1}^2 dr_{b1} \sin \theta d\theta v(r_{b1}) u_{l_f}(r_{C1}) u_{l_i}(r_{b1}) \\
 & \times \frac{S_{P,l_a,l_c}(r_{Cc})}{r_{Cc}} \frac{F_{l_b}(r_{Bb})}{r_{Bb}} \\
 & \times \sum_M \langle l_c 0 l_b M | K M \rangle [Y^{l_f}(\hat{\mathbf{r}}_{C1})Y^{l_b}(\theta + \pi, 0)]_M^K Y_{-M}^{l_b}(\hat{\mathbf{r}}_{Bb}),
 \end{aligned}
 \tag{7.2.177}$$

and

$$\begin{aligned}
 s_{P, l_a, l_c}(r_{Cc}) &= \int r_{Cc}'^2 dr_{Cc}'^2 r_{A2}'^2 dr_{A2}'^2 \sin \theta' d\theta' v(r_{c2}') u_{l_f}(r_{A2}') u_{l_i}(r_{c2}') \\
 &\times \frac{F_{l_a}(r_{Aa}')}{r_{Aa}'} \frac{f_{l_c}(k_{Cc}, r_c)}{r_{Cc}'} P_{l_c}(k_{Cc}, r_c) \\
 &\times \sum_M \langle l_c 0 l_a M | P M \rangle \left[Y^{l_f}(r_{A2}') Y^{l_a}(r_{c2}') \right]_M^P Y_{-M}^{l_a}(r_{Aa}').
 \end{aligned} \quad (7.2.178)$$

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

$$\frac{d\sigma}{d\Omega}(\hat{k}_{bB}) = \frac{k_{bB}}{k_{aA}} \frac{\mu_{aA} \mu_{bB}}{(2\pi\hbar^2)^2} \left| \sum_{\mu} \langle \Lambda \mu J_f M_i - \mu | J_i M_i \rangle T_{\mu} \right|^2. \quad (7.2.179)$$

For a non polarized incident beam,

$$\frac{d\sigma}{d\Omega}(\hat{k}_{bB}) = \frac{k_{bB}}{k_{aA}} \frac{\mu_{aA} \mu_{bB}}{(2\pi\hbar^2)^2} \frac{1}{2J_i + 1} \sum_{M_i} \left| \sum_{\mu} \langle \Lambda \mu J_f M_i - \mu | J_i M_i \rangle T_{\mu} \right|^2. \quad (7.2.180)$$

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 .

$$\frac{d\sigma}{d\Omega}(\hat{k}_{bB}) = \frac{k_{bB}}{k_{aA}} \frac{\mu_{aA} \mu_{bB}}{(2\pi\hbar^2)^2} \frac{1}{2J_i + 1} \sum_{\mu} |T_{\mu}|^2 \sum_{M_i, M_f} |\langle \Lambda \mu J_f M_f | J_i M_i \rangle|^2. \quad (7.2.181)$$

The sum over M_i, M_f of the Clebsh-Gordan coefficients gives $(2J_i + 1)(2\Lambda + 1)$ (see ??) One then gets,

$$\frac{d\sigma}{d\Omega}(\hat{k}_{bB}) = \frac{k_{bB}}{k_{aA}} \frac{\mu_{aA} \mu_{bB}}{(2\pi\hbar^2)^2} \frac{1}{(2\Lambda + 1)} \sum_{\mu} |T_{\mu}|^2, \quad (7.2.182)$$

where one can write

$$\begin{aligned}
 T_{\mu} &= \sum_{l_a, l_b} C_{l_a, l_b} \left[Y^{l_a}(\hat{k}_{aA}) Y^{l_b}(\hat{k}_{bB}) \right]_{\mu}^{\Lambda} \\
 &= \sum_{l_a, l_b} C_{l_a, l_b} i^{l_a} \sqrt{\frac{2l_a + 1}{4\pi}} \langle l_a l_b 0 \mu | \Lambda \mu \rangle Y_{\mu}^{l_b}(\hat{k}_{bB}).
 \end{aligned} \quad (7.2.183)$$

Note that (7.2.182) 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 (7.2.40)

$$\frac{d\sigma}{d\Omega}(\hat{k}_{bB}) = \frac{k_{bB}}{k_{aA}} \frac{\mu_{aA} \mu_{bB}}{(2\pi\hbar^2)^2} \frac{1}{2(2\Lambda + 1)} \sum_{\mu} |A_{\mu}|^2 + |B_{\mu}|^2. \quad (7.2.184)$$

note different factors in Eq. (7.2.40)

(245) App A

gregory

??

7.2

~~parte de las correcciones~~

parte de las correcciones están en la versión

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Appendix

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7.A. ZPF AND PAULI PRINCIPLE AT THE BASIS OF MEDIUM POLARIZATION EFFECTS: SELF-ENERGY, VER

Appendix 7.A ZPF and Pauli principle at the basis of medium polarization effects: self-energy, vertex corrections and induced interaction

In keeping with a central objective of the formulation of quantum mechanics, namely that the basic concepts on which it is based relate directly to experiment (Heisenberg), elementary modes of nuclear excitation (single-particle, collective vibrations and rotations), are solidly anchored on observation (inelastic and Coulomb excitation, one- and two-particle transfer reactions). Of all quantal phenomena, zero point fluctuations (ZPF), closely connected with virtual states, are likely to be most representative of the essential difference existing between quantum and classical mechanics. In fact, ZPF are intimately connected with the complementary principle (Bohr), and thus with the indeterminacy (Heisenberg) and non-commutative (Born, Jordan) relations, and with the probabilistic interpretation (Born) of the (modulus squared) of the wavefunctions, solution of Schrödinger's or Dirac's equations.

Pauli principle 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 excitation (see Figs. 7.J.2 and 7.J.3); within the present context, see also Schrieffer (1964).

parcialmente corregido

Appendix 7.B Coherence and effective formfactors

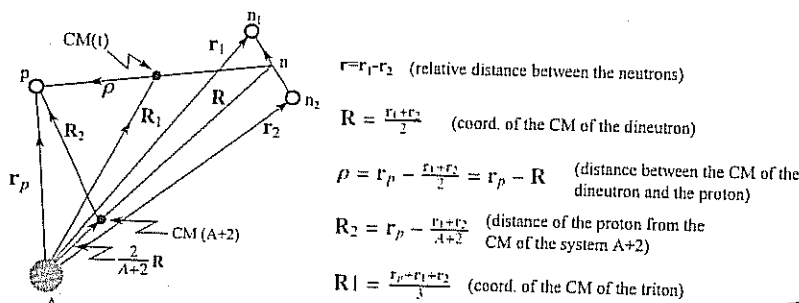


Figure 7.B.1

coordinate system used in the calculation of the transfer amplitude

In what follows we shall work on a simplified derivation of the simultaneous two-nucleon transfer amplitude, within the framework of first order DWBA specially suited to discuss correlation aspects of pair transfer in general, and of the associated effective formfactors in particular.

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

$$\begin{aligned}\psi_\alpha &= \psi_{M_i}^{J_i}(\xi_A) \sum_{s s_f} [\chi^s(\sigma_\alpha) \chi^{s_f}(\sigma_\beta)]_{M_i}^{s_f} \phi_i^{L=0}(\sum_{i < j} |\vec{r}_i - \vec{r}_j|) \\ &= \psi_{M_i}^{J_i}(\xi_A) \sum_{M_i M_i'} (s M_i' s_f' M_i' | s_i M_i) \chi_{M_i}^s(\sigma_\alpha) \chi_{M_i'}^{s_f'}(\sigma_\beta) \\ &\quad \times \phi_i^{L=0}(\sum_{i < j} |\vec{r}_i - \vec{r}_j|)\end{aligned}\quad (7.B.1)$$

while

$$\begin{aligned}\psi_\beta &= \psi_{M_f}^{J_f}(\xi_{A+2}) \chi_{M_f}^{s_f'}(\sigma_\beta) \\ &= \sum_{\substack{n_1 l_1 j_1 \\ n_2 l_2 j_2}} B(n_1 l_1 j_1, n_2 l_2 j_2; J J_i' J_f) [\phi^J(j_1 j_2) \phi^{J_i'}(\xi_A)]_{M_f}^{J_f} \\ &\quad \times \chi_{M_f}^{s_f'}(\sigma_\beta)\end{aligned}\quad (7.B.2)$$

Making use of the above equation one can define the spectroscopic amplitude B as

$$\begin{aligned}B(n_1 l_1 j_1, n_2 l_2 j_2; J J_i' J_f) \\ = \langle \psi^{J_i'}(\xi_{A+2}) | [\phi^J(j_1 j_2) \phi^{J_i'}(\xi_A)]^{J_i'} \rangle,\end{aligned}\quad (7.B.3)$$

where

$$\phi^J(j_1 j_2) = \frac{[\phi_{j_1}(\vec{r}_1) \phi_{j_2}(\vec{r}_2)]^J - [\phi_{j_1}(\vec{r}_2) \phi_{j_2}(\vec{r}_1)]^J}{\sqrt{1 + \delta(j_1, j_2)}},\quad (7.B.4)$$

is an antisymetrized, normalized wave function of the two transferred particles. The function $\chi_{M_f}^{s_f'}(\sigma_\beta)$ appearing both in eq. (7.B.1) and (7.B.2) is the spin wave function of the proton while

$$\chi^s(\sigma_\alpha) = [\chi^{s_1}(\sigma_{n_1}) \chi^{s_2}(\sigma_{n_2})]^s, \quad (7.B.5)$$

is the spin function of the two-neutron system.

A convenient 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.

$$\begin{aligned}\phi_i^{L=0}(\sum_{i < j} |\vec{r}_i - \vec{r}_j|) &= N_i e^{l(r_1 - r_2)^2 + (r_1 - r_p)^2 + (r_2 - r_p)^2} \\ &= \phi_{000}(\vec{r}) \phi_{000}(\vec{\rho}),\end{aligned}\quad (7.B.6)$$

$$\phi_{000}(\vec{r}) = R_{nl}(v^{1/2} r) Y_{lm}(\hat{r})$$

The coordinate $\vec{\rho}$ is the radius vector which measures the distance between the center of mass of the dineutron and the proton, while the vector \vec{r} is the dineutron relative coordinate (cf. Fig. 7.B.1.)

$$\vec{r} = \vec{r}_1 - \vec{r}_2 \quad (\text{relative distance between the neutrons}) \quad (7.B.7)$$

$$\begin{aligned}\vec{R} &= \frac{\vec{r}_1 + \vec{r}_2}{2} \quad (\text{coord. of the CM of the dineutron}) & (7.B.7b) \\ \vec{\rho} &= \vec{r}_p - \frac{\vec{r}_1 + \vec{r}_2}{2} \quad (\text{distance between the CM of the dineutron and the proton}) & (7.B.7c) \\ \vec{R}_2 &= \vec{r}_p - \frac{\vec{r}_1 + \vec{r}_2}{A+2} \quad (\text{distance of the proton from the CM of the system } A+2) & (7.B.7d) \\ \vec{R}_1 &= \frac{\vec{r}_p + \vec{r}_1 + \vec{r}_2}{3} \quad (\text{coord. of the CM of the triton}) & (7.B.7e)\end{aligned}$$

To obtain the DWBA cross section we have to calculate the integral

$$T = \int d\xi_A d\vec{r}_1 d\vec{r}_2 d\vec{r}_p \chi_p^{(-)}(\vec{R}_2) \psi_p^*(\xi_{A+2}, \sigma_p) V_p \psi_a(\xi_A, \sigma_a, \sigma_p) \psi_i^{(+)}(\vec{R}_1) \quad (7.B.8)$$

where the final state effective interaction $V_p(\rho)$ is assumed to depend only on the distance ρ between the center of mass of the di-neutron and of the proton. Instead of integrating over $\xi_A, \vec{r}_1, \vec{r}_2$ and \vec{r}_p we would integrate over ξ_A, \vec{r}, \vec{r}' and \vec{r}_p . The Jacobian of the transformation is equal to 1, i.e. $\partial(\vec{r}_1, \vec{r}_2)/\partial(\vec{r}, \vec{r}') = 1$. *such*

To carry out the integral (7.B.8) we transform the wave function (7.B.4) into center of mass and relative coordinates. If we assume that both $\phi_{j_1}(\vec{r}_1)$ and $\phi_{j_2}(\vec{r}_2)$ are harmonic oscillator wave functions, this transformation can easily 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 \vec{r}_1 and \vec{r}_2 and $|nl, NL; \lambda \mu\rangle$ is the corresponding one depending on \vec{r} and \vec{R} , we can write

$$\begin{aligned}|n_1 l_1, n_2 l_2; \lambda \mu\rangle &= \left(\sum_{nl, NL} |nl, NL; \lambda \mu\rangle \langle nl, NL; \lambda \mu| \right) |n_1 l_1, n_2 l_2; \lambda \mu\rangle \\ &= \sum_{nl, NL} |nl, NL; \lambda \mu\rangle \langle nl, NL; \lambda \mu | n_1 l_1, n_2 l_2; \lambda \rangle \quad (7.B.9)\end{aligned}$$

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

$$\begin{aligned}2n_1 + l_1 + 2n_2 + l_2 &= 2n + l + 2N + L, \\ (-1)^{l_1+l_2} &= (-1)^{l+L}.\end{aligned} \quad (7.B.10)$$

The coefficients $\langle nl, NL, L | n_1 l_1, n_2 l_2, L \rangle$ are tabulated and were first discussed by ~~Moshinsky~~ Moshinsky in Nucl. Physics, 13 (1959) 104.

With the help of eq.(7.B.9) we can write the wave function $\psi_{M_i}'(\xi_{A+2})$ as

Moshinsky, 1959

(used as a basis to expand the Saxon-Woods single-particle wave functions)

$$\begin{aligned}
\psi_{M_f}^{J_f}(\xi_{A+2}) &= \sum_{\substack{n_1 l_1 j_1 \\ n_2 l_2 j_2 \\ J J_i}} B(n_1 l_1 j_1, n_2 l_2 j_2; J J_i' J_f) [\phi^J(j_1 j_2) \phi^{J_i'}(\xi_A)]_{M_f}^{J_f} \\
&= \sum_{\substack{n_1 l_1 j_1 \\ n_2 l_2 j_2}} \sum_{J J_i} B(n_1 l_1 j_1, n_2 l_2 j_2; J J_i' J_f) \\
&\quad \times \sum_{M_f M_{J_i}} \langle J M_f J_i' M_{J_i} | J_f M_{J_f} \rangle \psi_{M_{J_i}}^{J_i'}(\xi_A) \\
&\quad \times \sum_{LS'} \langle S' L J | j_1 j_2 J \rangle \sum_{M_L M_S'} \langle L M_L S' M_S' | J M_J \rangle \chi_{M_S'}^{S'}(\sigma_\alpha) \\
&\quad \times \sum_{n l N \Lambda} \langle n l, N \Lambda, L | n_1 l_1, n_2 l_2, L \rangle \\
&\quad \times \sum_{m_l M_\Lambda} \langle l m_l M_\Lambda | L M_L \rangle \phi_{n l m_l}(\vec{r}) \phi_{N \Lambda M_\Lambda}(\vec{R}) .
\end{aligned} \tag{7.B.11}$$

Integration over \vec{r} gives

$$\langle \phi_{n l m_l}(\vec{r}) | \phi_{000}(\vec{r}) \rangle = \delta(l, 0) \delta(m_l, 0) \Omega_n , \tag{7.B.12}$$

where

$$\Omega_n = \int R_{n0}(v_1^{1/2} r) R_{00}(v_2^{1/2} r) r^2 dr . \tag{7.B.13}$$

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 v_2 which is different from the one that the two neutrons are subjected to, when moving in the system A .

Integration over ξ_A and multiplication of the spin functions gives

$$\begin{aligned}
V(\rho) & \quad \langle \psi_{M_{J_i}}^{J_i}, V_{\beta}'(\rho) \psi_{M_{J_i}'}^{J_i'} \rangle = \delta(J_i, J_i') \delta(M_{J_i}, M_{J_i}') V_{\beta}(\rho) , \\
& \quad \langle \chi_{M_{S_i}}^S(\sigma_\alpha), \chi_{M_{S_i}'}^{S'}(\sigma_\alpha) \rangle = \delta(S, S') \delta(M_{S_i}, M_{S_i}') , \\
& \quad \langle \chi_{M_{S_f}}^{S_f}(\sigma_\beta), \chi_{M_{S_f}'}^{S_f'}(\sigma_\beta) \rangle = \delta(S_f, S_f') \delta(M_{S_f}, M_{S_f}') .
\end{aligned} \tag{7.B.14}$$

The integral (7.B.8) can now be written as *then*

$$\begin{aligned}
T &= \sum_{\substack{n_1 l_1 j_1 \\ n_2 l_2 j_2}} \sum_{J M_J} \sum_{n N} \sum_S B(n_1 l_1 j_1, n_2 l_2 j_2; J J_i' J_f) \\
&\quad \times \langle J M_J J_i M_{J_i} | J_f M_{J_f} \rangle \langle S L J | j_1 j_2 J \rangle \\
&\quad \times \langle L M_L S M_S | J M_J \rangle \langle n 0, N L, L | n_1 l_1, n_2 l_2, L \rangle \\
&\quad \times \langle S M_S S_f M_{S_f} | S_i M_{S_i} \rangle \Omega_n \\
&\quad \times \int d\vec{R} d\vec{r}_p \chi_i^{(+)*}(\vec{R}_1) \phi_{N L M_L}^*(\vec{R}) V_{\beta}'(\rho) \phi_{000}(\vec{R}) \chi_i^{(+)}(\vec{R}_1) ,
\end{aligned} \tag{7.B.15}$$

We now define the effective two-nucleon transfer form factor as

where we have approximated V_{β}' by an effective interaction depending only on $\rho = |\vec{p}|$.

(non-orthogonality effect)

$V(\rho)$

$$u_{LSJ}^{JJ_f}(R) = \sum_{n_1 l_1 j_1, n_2 l_2 j_2; J J_i J_f} B(n_1 l_1 j_1, n_2 l_2 j_2; J J_i J_f) \langle S L J | j_1 j_2 J \rangle \langle n_0, NL, L | n_1 l_1, n_2 l_2; L \rangle \Omega_n R_{nL}(R) \quad (7.B.16)$$

We can now rewrite eq. (7.B.15) as

$$T = \sum_J \sum_L \sum_S (J M_J J_i M_{J_i} | J_f M_{J_f}) (S M_S S_f M_{S_f} | S_i M_{S_i}) (L M_L S M_S | J M_J) \times \int d\vec{R} d\vec{r}_p \chi_p^{*(-)}(\vec{R}_2) u_{LSJ}^{JJ_f}(R) Y_{LM_L}^*(\vec{r}) V(\rho) \phi_{000}(\vec{r}) \chi_i^{(+)}(\vec{R}_1) \quad (7.B.17)$$

Because the di-neutron has $S = 0$, we have that

$$(L M_L 00 | J M_J) = \delta(J, L) \delta(M_L, M_J) \quad (7.B.18)$$

and the summations over S and L disappear from eq. (7.B.17). Let us now make also here, as done in App. for one particle transfer reactions, the zero range approximation, that is,

$$V(\rho) \phi_{000}(\vec{r}) = D_0 \delta(\vec{r}) \quad (7.B.19)$$

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.F.1)*

From eqs. (??) we obtain

$$\vec{R} = \vec{R}_1 = \vec{r} \quad \leftarrow \quad (7.B.20)$$

$$\vec{R}_2 = -\frac{A}{A+2} \vec{R} \quad \leftarrow$$

Then eq. (7.B.15) can be written as

$$T = D_0 \sum_L (L M_L J_i M_{J_i} | J_f M_{J_f}) \times \int d\vec{R} \chi_p^{*(-)}\left(-\frac{A}{A+2} \vec{R}\right) u_L^{JJ_f}(R) Y_{LM_L}^*(\vec{R}) \chi_i^{(+)}(\vec{R}) \quad (7.B.21)$$

From eq. (7.B.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 (i, p) and (p, i) reactions are in zero-range approximation are,

$$\Delta S = 0$$

$$\Delta J = \Delta L = L$$

$$\Delta\pi = (-1)^L \quad \text{for} \quad (7.B.22)$$

i.e. only normal parity states are excited.

The integral appearing in eq. (7.B.21) has the same structure as the DWBA integral appearing in *Eq. (6.F.16)* derived in the case of one-nucleon transfer reactions.

The difference between the two processes manifest itself through the different structure of the two form factors. While $u_i(r)$ appearing in Equation (6.F.1) is a single-particle bound state wave function, $u_L^{JJ_f}$ is a coherent summation over the center of mass states of motion of the two transferred neutrons. In other words, an effective quantity (function), It is of notice that this difference essentially vanishes, when one considers dressed particles resulting from the coupling to collective motion, and leading to ω -dependent effective masses.

Examples of two-nucleon transfer form factors are given in Figs. 7.B.2, 7.B.3 and 7.B.4

Brogia and Riedel (1967)

lower case

$\Delta\pi$

E

Then eq. (7.B.15) can be written as

$$T = D_0 \sum_L (L M_L J_i M_{J_i} | J_f M_{J_f}) \times \int d\vec{R} \chi_p^{*(-)}\left(-\frac{A}{A+2} \vec{R}\right) u_L^{JJ_f}(R) Y_{LM_L}^*(\vec{R}) \chi_i^{(+)}(\vec{R}) \quad (7.B.21)$$

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among other things,

lower case

(i) a pure shell-model configuration $(p_{1/2})^{-2}$ for the ground state of ^{206}Pb , (ii) a configuration mixing caused by pairing a residual interaction and (iii) the same model as in (ii) including ground state correlations.

The results show clearly that different assumptions about the structure of one definite nuclear state lead to almost the same shape of the f_L functions inside the nucleus. The

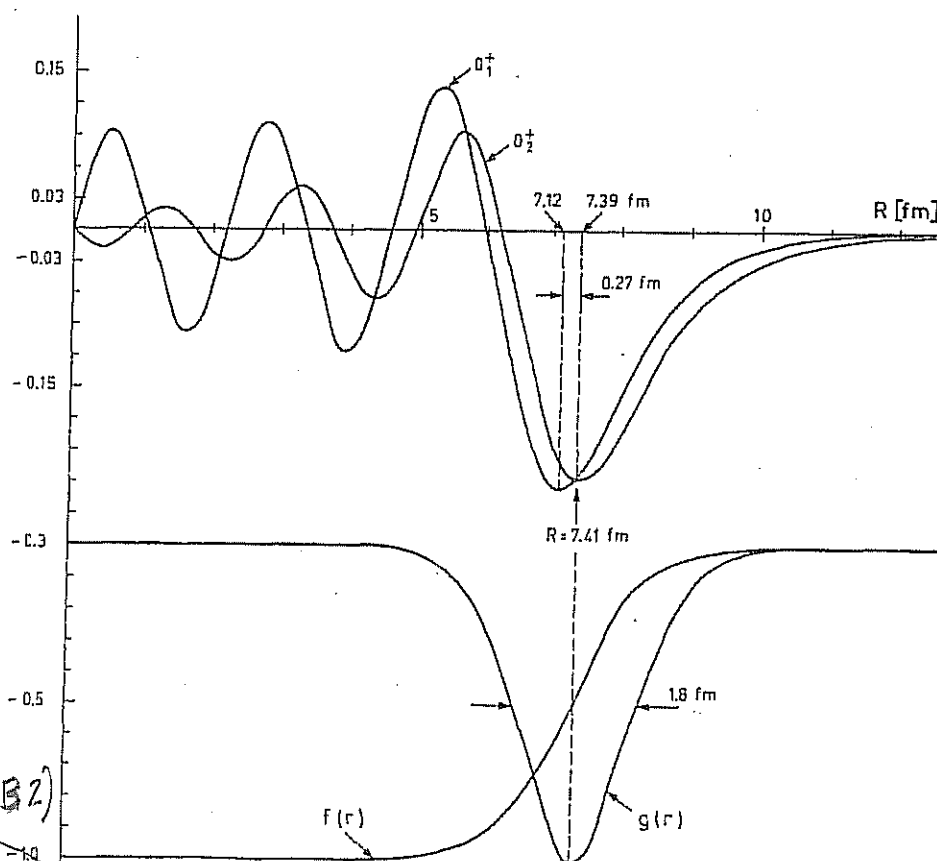


Fig. 7.13.2. The upper part of the figure shows the modified form factor for the transition to the ground state ($0^+_{1/2}$) and the pairing vibrational state ($0^+_{3/2}$) at 4.87 MeV. Both curves are matched with appropriate Hankel functions. In the lower part the form factors of the real ($f(r)$) and the imaginary ($g(r)$) part of the optical potential are given in the same scale for the radius.

form factor calculated by means of the shell-model wave functions of True and Ford¹³⁾ is very similar to that of model (ii). The main difference of the three bound states is the magnitude of the maximum around the nuclear surface. Then, coherence effects only affect the degree of collectiveness of the form factor but not its shape.

Let us now discuss some details of the asymptotic behaviour of the form factors.

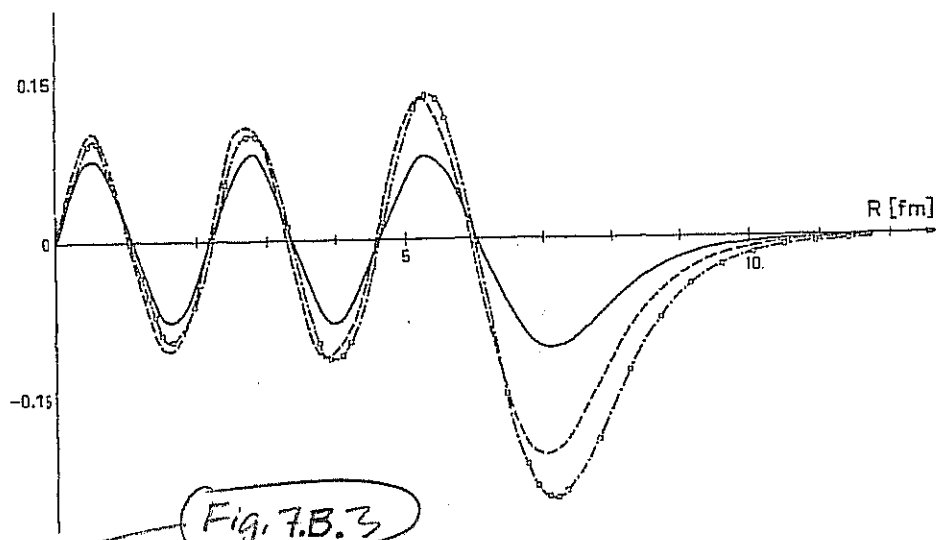


Fig. 5. Modified form factor for the transition to the ground state calculated in different spectroscopic models (pure shell-model configuration —, shell model plus pairing residual interaction — — —, including ground state correlations — o — o —).

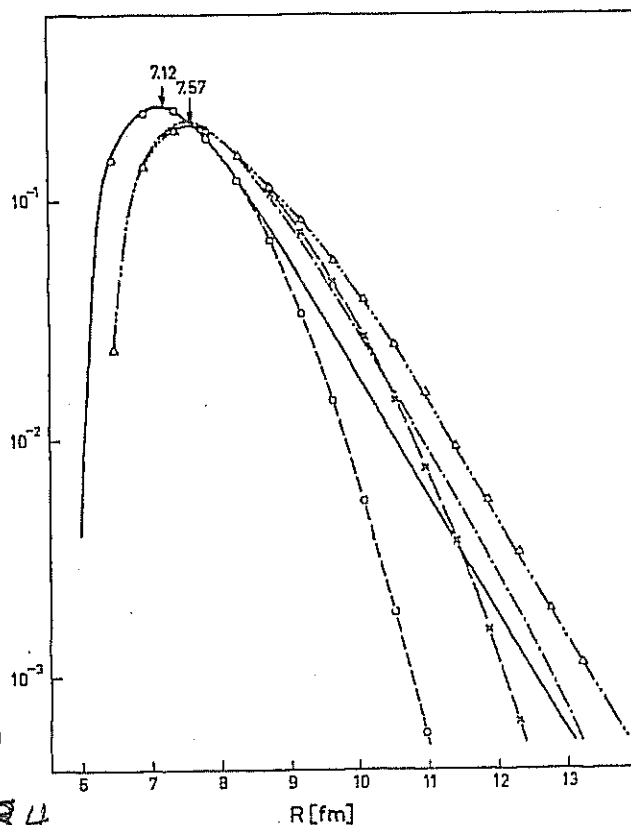


Fig. 6. Asymptotic behaviour of the modified form factor for the ground state transition for oscillator plus Hankel wave functions (—), oscillator wave functions alone (— o — o), using Woods-Saxon wave functions according to method I (— x — x —); (the same curve matched with a Hankel function — — Δ — — Δ — —) and method II (— · — · —).