Structure and reactions of N=7 isotones: parity inversion and transfer.

F. Barranco
Sevilla University
R.A. Broglia
The Niels Bohr Institute
G. Potel
MSU,USA
E. Vigezzi
INFN Milano
NSD2019, Venice (Italy)



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Spirit of the talk: Mean Field at the limits

Mean Field is probably the most genuine Nuclear Structure concept. Intimately linked to the collective behavior of nuclei, it is expected to emerge as a useful tool when the number of nucleons increases.

On the other hand, Nuclear Field Theory (NFT) is the systematic perturbative procedure for improving the zero order Mean Field description towards the exact description of nuclei, by incorporating higher and higher order many body effects.

In this talk we explore how things are going in the region of very light, one valence neutron N=7 isotones (from Z=3 to Z=6), and find that very good description of low-energy spectra and associated reaction processes is attainable as far as the NFT procedure is pursued up to third order terms. It will be shown that a consistent description of

O spectra,

O one neutron transfer absolute $d\sigma/dEd\Omega$

O dipole transitions and

O isotopic rms charge radius change

It is obtained by including

+ up to "3p-2h" configurations, and

+ coupling to the **single-particle continuum**.

It will be stressed that antisymmetrization (**Pauli principle**) between the valence neutron and core, plays a crucial role.

By incorporating coupling to the **single-particle continuum**, a **common framework for the study of structure and reactions**, based on the Nuclear Field Theory emerges. A formulation in terms of **Generalized Coupled Channels allows for the proper inclusion of Pauli principle**.



N=7 isotnes, from stability valley to drip line





¹³C (1st exc.: 1/2+)

• 1d5/2 2s1/2 1p1/2

Ij continuum



¹³C (2nd exc.: 5/2+)

Ij continuum

1d5/2 2s1/2 1p1/2 1p3/2 1s1/2



¹³C (gs: 1/2-)



¹¹Be (gs: 1/2-)



Z=4

¹¹Be (1st exc.: 1/2+)



Z=4

¹¹Be (2nd exc.: 5/2+)



Z=4





HOW TO IMPROVE SIMPLE (SPHERICAL STATIC) ZERO ORDER MEAN FIELD DESCRIPTION?

Include quadrupole deformed mean field potential

+Static/rigid and strong coupling limt (A.Macchiavelli)

+Static/rigid and weak coupling (F.Nunes; A. Moro)

+Dynamic and weak/medium coupling (This work, Esbensen&Sagawa)

HOW TO IMPROVE SIMPLE ZERO ORDER MEAN FIELD DESCRIPTION?

Nuclear Fierld Theory (NFT)

shows how to construct an *EXACT* solution of the many body problem, for a given two body force, in terms of a **rapidly convergent series expansion** for each observable, similar to the Feynman's QED diagrams expansion, and **based on the use of realistic (coherent-collective) core excitations** rather than independent 1p-1h, 2p-2h,... configurations, that is in the **dynamical mean field response rather than in the static (zero order) one.**

- 1. Fermions: Neutron moving in a HF mean field
- 2. Bosons: Core realistic excited states; RPA Vibrational like
- 3. Fermion-Boson vertex: Microscopically evaluated from RPA and HF
- 4. Expansion parameter: inverse of effective degeneracy, Ω .

Empirical NFT: fix Fermions, Vibrations and Vertex from experiment in order to give a unified description of "ALL(II)" data of "ALL" nuclei

PVC: Vibrational Core (even-even) + One particle (neutron)

$$H = H_{sp} + H_{coll} + H_{PVC}$$

$$H_{sp} = -d/d \vec{r} \hbar^{2}/2 \mu(r) d/d \vec{r} + V(r) + V_{ls}(r)$$

$$H_{coll} = \sum_{\lambda \mu \nu} \hbar \omega_{\lambda,\nu} [\Gamma^{+}_{\lambda \mu,\nu} \Gamma_{\lambda \mu,\nu} + 1/2]$$
Dynamic deformed nucleus
$$H_{PVC} = \sum_{\lambda \mu \nu} \delta V_{\lambda\nu}(r) Y_{\lambda\mu}(\hat{r}) [\Gamma^{+}_{\lambda \mu,\nu} + (-1)^{\mu} \Gamma_{\lambda \mu,\nu}]$$
where $\delta V_{\lambda\nu}(r) \approx -rdV/dr \beta_{\lambda,\nu}$

$$\mu(r) = \mu_{0} - \frac{0.2M}{1 + \exp(r - R)/a}; \mu_{0} = \frac{A}{(A+1)}M$$

Where $\lambda\mu$ is the excited core angular momentum and z-projection, v labels the different states with the same $\lambda\mu$. For the collective states the form factor in PVC is the semiclassic one.







V0=-72MeV ; R=2.14fm ; a= 0.72











Charge Radii



 $B_{bare}(E1)=0.29 e^2 fm^2 --->0.11 e^2 fm^2 (exp.0.10 e^2 fm^2)$



Data: Schmitt, PRC88,064612(2013); Winfield,NPA683,48(2001) DWBA; Optical potential: Koening-Delaroche/Han et al.



N=7 isotones results



Bare mean field potential for N=7 isotones

		$\hbar\omega_{2^+}$ (MeV)	β_2^n	V_{WS} (MeV)	V_{ls}	a_{WS} (fm)	R_{WS} (fm)		
	¹⁰ Li	3.37	0.68	64.0	0.50	0.75	2.10		
	$^{11}\mathrm{Be}$	3.37	0.71	72.0 (73.8)	0.58(0.54)	0.72(0.78)	2.14		
	$^{12}\mathrm{B}$	3.80	0.57	76.8	0.65	0.78(0.76)	2.18		
	¹³ C	4.4	0.46(0.43)	82(83.0)	0.76(0.7)	0.73(0.72)	2.23 (2.25)		
New simple parametrization: V_{ws} =-82+-54(N-Z)/A <i>MeV</i>									
0					a = 0.75	5 <i>fm;</i> R _{ws} =	=0.99A ^{1/3} /	^r m	
-10	¹⁰ Li ¹¹ Be	$-\frac{10}{10}$ $V_{LS} = 0.00$)82V _{ws}		
-20	— ¹² B — ¹³ C		4.8				•		
\$ ⁻³⁰			4.6		•	0.8-	0		
≥40- >			A.2			500	•		
-50-	///	/	C) 4 - 9 4 - 9 4 -		0	- 0.0-	0	•	
-00-			3.6- 3.4-	•		0.4-		0	
-70			3.2			0.2		<u> </u>	
0.0	1	2 3 4 R (fn	n]	2 3 4 Z	5 6	1 2	3 4 5 Z	0	

Odd Z nuclei

Effect of coupling to 1p3/2 proton hole in 12B







FIG. 2. Spectral functions of the low-lying negative (a) and positive (b) parity states in ¹⁰Li.

d(⁹Li,p)¹⁰Li @ E_{lab} 100MeV ; θ_{CM} [5.5°,16°]



Theor.: *ArXiV1812.01761;* see also A.Moro PLB793,13(2019) Data: Cavallaro et al. PRL118:012701, 2017

d(⁹Li,p)¹⁰Li @ E_{lab} 21MeV ; θ_{CM} [98^o,134^o]



Theor.: ArXiV1812.01761; see also A.Moro PLB793,13(2019),S.Orrigo PLB633,469(2006)

Data: Jeppesen et al. PLB 642:449, 2006

We plan to produce an universal bare mean field to be used together with quadrupole (eventually also octupole) coupling (as generalisation of the classical B&M static one) that can be applied to the whole mass table from A about 10 and from proton to neutron dripline in the unified study of low lying spectroscopic observables.

Results in the Sn isotopes region already exist.

A QRPA based protocol for calculating realistic quadrupole properties exists.

We plan to produce an universal bare mean field to be used together with quadrupole (eventually also octupole) coupling (as generalisation of the classical B&M static one) that can be applied to the whole mass table from A about 10 and from proton to neutron dripline in the unified study of low lying spectroscopic observables.

Results in the Sn isotopes region already exist.

A QRPA based protocol for calculating realistic quadrupole properties exists.

Thank you!



FIG. 5: Main processes contributing to the dipole transition between the first excited state and the ground state of ${}^{11}Be$. The vertex correction associated with the low-lying 1⁻ strength (incipient GDPR) is estimated to be small.

One has to consider, however, that one expects other important contributions from manybody processes. The two time orderings associated with the most important ones are shown in the right part of Figs. 5. They interfere in a destructive way with the leading contribution, leading to

$$B(E1; I_i \to I_f) = \frac{e_{eff}^2}{2I_i + 1} a_{1/2^+}^2 a_{1/2^-}^2 |M^0 + M^{1(a)} + M^{1(b)}|^2 = \frac{16}{121} \frac{1}{2} 0.66 |\sqrt{\frac{1}{2\pi}} \times 5 - 0.25 - 0.19|^2 \quad e^2 \text{fm}^2$$

$$B_{\text{exp}}(\text{E1}) = 0.10 \ e^2 \ \text{fm}^2 = 0.11 \quad e^2 \text{fm}^2. \tag{16}$$

$$B_{\text{bare}}(\text{E1}) = 0.29 \ e^2 \ \text{fm}^2$$

Charge Radii

$$\langle r^{2} \rangle_{11\text{Be}} = \left(\langle r^{2} \rangle_{10\text{Be}} + \left(\frac{\langle r^{2} \rangle_{1s1/2}^{1/2}}{11} \right)^{2} \right) \times S^{2} + (1 - S^{2}) \times \left(\langle r^{2} \rangle_{10\text{Be}} \left(1 + \frac{2}{4\pi} \beta_{\pi}^{2} \right) + \left(\frac{\langle r^{2} \rangle_{d5/2\,coll}^{1/2}}{11} \right)^{2} \right) = .$$

$$\langle r^{2} \rangle_{10\text{Be}} + \left(\frac{\langle r^{2} \rangle_{1s1/2}^{1/2}}{11} \right)^{2} \times S^{2} + (1 - S^{2}) \times \left(\left(\frac{\langle r^{2} \rangle_{d5/2\,coll}^{1/2}}{11} \right)^{2} + \langle r^{2} \rangle_{10\text{Be}} \frac{2}{4\pi} \beta_{\pi}^{2} \right)$$

$$\Delta < r^2 > \frac{1}{2} = 0.11 \text{ fm} / 0.27 \text{ fm}$$
Another standard strategy is mixing more complex configurations, including 1p-1h, 2p-2h,... core excitations. For example for the 1/2+



The amplitudes are obtained diagonalizing the nuclear many body hamiltonian , in a more or less extended basis. But this method is rather slowly convergent.







B&M I

surface thickness parameter a is taken to be A independent.

$$R = r_0 A^{1/3} \qquad r_0 = 1.27 \text{ fm}$$

$$a = 0.67 \text{ fm}$$
(2-181)

The potential strengths include a term depending on the neutron excess, in order to describe approximately the potential acting on a single neutron,

$$V = \left(-51 + 33 \frac{N-Z}{A} \quad \text{MeV} \right)$$

$$V_{1s} = -0.44V = \left(22 - 14 \frac{N-Z}{A}\right) \quad \text{MeV}$$
(2-182)

and the neutron excess for each A has been chosen to correspond to the





Core GS is not simply the HF (zero order) lower levels filled state

$$\sum_{\lambda} 1/2 \, \hbar \, \omega_{\lambda}$$

This is the zero point motion associated to the normal modes in the harmonic approximation. Its theoretical value depends on the used model for calculating the frequencies w_L. In fact the GSC energy can be calculated as (c.f. *Ring&Schuck*)

$$E_{corr} = \sum_{\lambda,\nu} 1/2 \hbar \omega_{\lambda,\nu}^{RPA} - 1/2 \hbar \omega_{\lambda,\nu}^{HF}$$

This is a negative energy indicating that the RPA Ground State is more bound than the HF one,

What in turn means that in the core GS time by time fluctuations happen around the HF (zero order core GS) (c.f. Bohr&Mottelson Vol.II)



Heyde, Shell Model book The neutron proton interaction $V_{np}(r_n, r_p) = F_0 \,\delta(r_n - r_p) \,(1 - a + a \,\sigma_n, \sigma_p)$ Proton frozen on the 1p_{3/2} orbital No neutron mixing with other lj_'s An additional $J_{nn}(Ij)_n$ dependent neutron potential $\delta V_{J np(lj) n}(r_{n}) = F_{0} |\Phi_{1p3/2}(r_{p})|^{2} (1-a+a \sigma_{n} \sigma_{p})_{J np(lj) n}$ to be added to the bare potential; (finally for simplicity the form factor is taken the same that for the bare potential) $\delta V_{J np(lj) n}(\mathbf{r}_{n}) = F'_{WS0} WSff(\mathbf{r}_{n})(1-a+a \sigma_{n}, \sigma_{p})_{J np(lj) n}$ = -2 MeVa=0.5 and F'_{WS0} (1-a+a σ_n . σ_p)_{J np=1(p1/2) n}

Single-particle energy correction due to GSC Pauli rearrangement

The contribution of a given p-h configuration to the GS Correlation Energy is (*B&MII*)

nк, lb, jb h, la, ja
$$\delta E = \frac{(-h_{ai,bk,\lambda}\sqrt{(2j_a+1)})^2}{0 - (E_{ai} + E_{bk} + \hbar \omega_{\lambda})} < 0$$

The presence of a new neutron (scattering- or bound-like) inhibits some of these correlations, producing an energy modification of the core state...

$$-\frac{\delta E}{2j_a+1} > 0$$

This is the meaning/value of the NFT self energy diagram (B&MII, eq.6.225)

$$(-1)\frac{\left(-h_{ai,bk,\lambda}\sqrt{2j_{a}+1}\right)^{2}\langle\left(\left(j_{a1},j_{a2}\right)J=0,j_{a3};j_{a}|\left(j_{a1},j_{a3}\right)J=0,j_{a2};j_{a}\right)\rangle}{E_{ai}-\left(2E_{ai}+E_{bk}+\hbar\omega_{\lambda}\right)}$$
nk,lb,jb ni,la,ja,ma
$$(-1)\frac{\left(-h_{ai,bk,\lambda}\sqrt{2j_{a}+1}\right)^{2}\langle\left(\left(j_{a1},j_{a2}\right)J=0,j_{a3};j_{a}|\left(j_{a1},j_{a3}\right)J=0,j_{a2};j_{a}\right)\rangle}{E_{ai}-\left(2E_{ai}+E_{bk}+\hbar\omega_{\lambda}\right)}$$





Including higher orders

This is the non-crossing phonon aproximation to the many phonons full solution. It excludes the vertex correction diagrams, that should be incorporated independently



These equations must be solved iteratively





Transfer form factors









$$\begin{split} \widetilde{|1/2^+\rangle} &= \sqrt{0.83} |s_{1/2}\rangle + \sqrt{0.17} |(d_{5/2} \otimes 2^+)_{1/2^+}\rangle + \sqrt{0.01} |(s_{1/2}^2(0) \otimes 0^{(-)})_{0^+} s_{1/2}\rangle \tag{1} \\ \widetilde{|1/2^-\rangle} &= \sqrt{0.81} |p_{1/2}\rangle + \sqrt{0.15} |(p_{1/2}, 1p_{3/2})_{2^+}^{-1} \otimes 2^+)_{0^+} |p_{1/2}\rangle + \sqrt{0.02} |(d_{5/2} \otimes 3^-)_{1/2^-}\rangle \\ &+ \sqrt{0.04} |(p_{1/2}^2(0) \otimes 0^{(-)})_{0^+} p_{1/2}\rangle \tag{2} \\ |0^{(-)}\rangle &= |^8 \text{Be}(\text{gs}, 0^+)\rangle \quad \text{monopole pair removal mode of the closed shell system } {}^{10}_4 \text{Be}_6 \tag{3} \\ |\widetilde{5/2^+\rangle} &= \sqrt{0.34} |d_{5/2}\rangle + \sqrt{0.32} |(s_{1/2} \otimes 2^+)_{5/2^+}\rangle + \sqrt{0.34} |(d_{5/2} \otimes 2^+)_{5/2^+}\rangle \tag{4} \\ &+ \sqrt{0.003} |(d_{5/2}(0)^2 \otimes 0^{(-)})_{0^+} d_{5/2}\rangle \end{split}$$



FIG. 2: (a) $5/2^+$ (NFT)_{ren} phase shifts as a function of energy. (b) $d_{5/2}$ phase shifts calculated with the bare potential.

Taking care of the Pauli principle.1st part



Project out the occupied waves in HF GS

$$R_{a}^{x}(r) = \int dr' [\delta(r-r') - \sum_{k \in occ.} R_{ak}(r) R_{ak}(r')] R_{a}^{x}(r') = \int dr' [\delta(r-r') - P(r, r')] R_{a}^{x}(r')$$

Similarly for R_{b}^{C}

The CC equation becomes non-local (integro-differential): must be solved in R^2

$$\begin{bmatrix} \left[-\hbar^{2}/2\mu \ d^{2}/d \ r^{2}+V_{a}(r)\right] & \Xi_{a,b\lambda}(-\beta_{\lambda}rdV/dr) \\ \Xi_{a,b\lambda}(-\beta_{\lambda}rdV/dr) & \left[-\hbar^{2}/2\mu \ d^{2}/d \ r^{2}+V_{b}(r)+\hbar\omega\right] \end{bmatrix} \begin{bmatrix} R_{a}^{x}(r) \\ R_{b}^{C}(r) \end{bmatrix} = \tilde{E} \begin{bmatrix} R_{a}^{x}(r) \\ R_{b}^{C}(r) \end{bmatrix}$$



Alternative method: Expand using only *non-occupied* waves

$$\begin{split} R_{a}^{x}(r) &= \sum_{i=1}^{N} x_{ai} R_{ai}^{WS}(r) \quad i \in non - occ. \qquad R_{b}^{C}(r) = \sum_{i=1}^{N} C_{bi} R_{bi}^{WS}(r) \quad i \in non - occ. \\ & \left[\begin{bmatrix} -\hbar^{2}/2\mu \ d^{2}/d \ r^{2} + V_{a}(r) \end{bmatrix} \quad \Xi_{a,b\lambda}(-\beta_{\lambda} r dV/dr) \\ \Xi_{a,b\lambda}(-\beta_{\lambda} r dV/dr) \quad \begin{bmatrix} -\hbar^{2}/2\mu \ d^{2}/d \ r^{2} + V_{b}(r) + \hbar \omega \end{bmatrix} \right] \begin{bmatrix} R_{a}^{x}(r) \\ R_{b}^{C}(r) \end{bmatrix} = \tilde{E} \begin{bmatrix} R_{a}^{x}(r) \\ R_{b}^{C}(r) \end{bmatrix} \\ & \left(\frac{e_{al}}{2} - \frac{h_{al,bl\lambda}}{h_{a2,bl\lambda}} + \frac{h_{al,bl\lambda}}{h_{a2,bl\lambda}} + \frac{h_{al,bl\lambda}}{h_{a2,bl\lambda}} + \frac{h_{al,bl\lambda}}{2} + V_{b}(r) + \hbar \omega \end{bmatrix} \left| \begin{pmatrix} x_{al} \\ x_{a2} \dots \\ C_{bl} \\ C_{bl} \\ C_{bl} \end{pmatrix} \right| = \tilde{E} \begin{pmatrix} x_{al} \\ x_{a2} \dots \\ C_{bl} \\ C_{bl} \\ C_{b2} \dots \end{pmatrix} \end{split}$$

FULL HAMILTONIAN MATRIX (Block-Arrowhead, energy independent, matrix), where

 $h_{ai,bj\lambda} = \Xi_{a,b\lambda} \beta_{\lambda} \int R_{ai}^{WS}(r) (-r dV / dr) R_{bj}^{WS}(r) dr$

Energy dependent Self-energy

$$\begin{array}{l} R_{a}^{x}(r) = \sum_{i} x_{ai} R_{ai}^{WS}(r) \quad i \in non - occ. \\ R_{b}^{C}(r) = \sum_{i} C_{bi} R_{bi}^{WS}(r) \quad i \in non - occ. \\ \begin{pmatrix} e_{al} & 0 \dots & h_{al,bl\lambda} & h_{al,b2\lambda} \dots \\ 0 & e_{a2} \dots & h_{a2,bl\lambda} & h_{a2,b2\lambda} \dots \\ h_{al,bl\lambda} & h_{a2,b2\lambda} \dots & 0 & e_{b2} + \hbar \omega \dots \\ h_{al,b2\lambda} & h_{a2,b2\lambda} \dots & 0 & e_{b2} + \hbar \omega \dots \\ \end{pmatrix} \begin{pmatrix} x_{al} \\ x_{a2} \dots \\ C_{bl} \\ C_{b2} \dots \end{pmatrix} = \tilde{E} \begin{pmatrix} x_{al} \\ x_{a2} \dots \\ C_{bl} \\ C_{b2} \dots \end{pmatrix} \\ \begin{array}{c} e_{al} + \sum_{11} (\tilde{E}) & \sum_{12} (\tilde{E}) \dots \\ \sum_{21} (\tilde{E}) & e_{a2} + \sum_{22} (\tilde{E}) \dots \end{pmatrix} \begin{pmatrix} x_{al} \\ x_{a2} \dots \end{pmatrix} = \tilde{E} \begin{pmatrix} x_{al} \\ x_{a2} \dots \\ C_{bl} \\ C_{b2} \dots \end{pmatrix} \\ \end{array}$$

 $\Sigma_{ij}(\tilde{E}) = \sum_{bk} \frac{h_{ai,bk,\lambda}h_{aj,bk,\lambda}}{\tilde{E} - (E_{i,i} + \hbar \omega_{\lambda})}$ aj Where the so called basic vertex is $h_{ai,bk\lambda} = \Xi_{a,b\lambda} \beta_{\lambda} \int R_{ai}^{WS}(r) (-rdV/dr) R_{bk}^{WS}(r) dr$ bk To be noted that this Feynman Diagram contains all the Often, for bound states, a 1 s-p state richness of the CC equation, approximation is used and in fact it is completely ai $e_{al} + \Sigma_{11}(\tilde{E}) = \tilde{E}$ equivalent to it, eventual open inelastic channels included. In this approximation the radial dependence is blocked. Only the amplitudes are determined

Optical Potential as the r,r'-E-dependent Self-energy

.. obtained by means of the standard algebraic manipulations:

$$C_{bk} = \frac{\sum_{ai} x_{ai} h(a_i, b_k \lambda)}{\tilde{E} - (E_{bk} + \hbar \omega_\lambda)} = \frac{\Xi_{a,b\lambda} \int dr' R_{bk}^{WS}(r') \delta \rho(r') R_{ai}^{x}(r')}{\tilde{E} - (E_{bk} + \hbar \omega_\lambda)}$$
$$h_{ai,bk\lambda} = \Xi_{a,b\lambda} \beta_\lambda \int R_{ai}^{WS}(r) (-rdV/dr) R_{bk}^{WS}(r) dr$$
$$R_a^{x}(r) = \sum_i x_{ai} R_{ai}^{WS}(r); E_{ai} > 0$$

$$R_{b}^{C}(r) = \sum_{k} C_{bk} R_{bk}^{WS}(r) = \sum_{k} R_{bk}^{WS}(r) \frac{\Xi_{a,b\lambda} \int dr' R_{bk}^{WS}(r') \delta\rho(r') R_{ai}^{x}(r')}{\tilde{E} - (E_{bk} + \hbar \omega_{\lambda})}; E_{bk} > 0$$

$$\left(\frac{H_{p} - e_{F}}{\Xi_{a,b\lambda} f(r)} H_{p} - e_{F} + \hbar \omega \right) \left(\frac{R_{a}^{x}}{R_{b}^{C}} \right) = \tilde{E} \left(\frac{R_{a}^{x}}{R_{b}^{C}} \right)$$

$$(H_{p}(r)-e_{F})R_{ai}^{x}(r)+\int dr'\sum_{bk>0}\frac{\Xi_{a,b\lambda}^{2}R_{bk}^{WS}(r')R_{bk}^{WS}(r)\delta\rho(r')\delta\rho(r')}{\tilde{E}-(E_{bk}+\hbar\omega_{\lambda}-i\eta)}R_{ai}^{x}(r')=\tilde{E}R_{ai}^{x}(r')$$

Microscopic optical potentials for high enough E's

 $1/2\hbar\omega_{\lambda}$

This is the zero point motion associated to the normal modes in the harmonic approximation. Its theoretical value depends on the used model for calculating the frequencies w_L. In fact the GSC energy can be calculated as (c.f. *Ring&Schuck*)

$$E_{corr} = \sum_{\lambda,\nu} 1/2 \hbar \omega_{\lambda,\nu}^{RPA} - 1/2 \hbar \omega_{\lambda,\nu}^{HF}$$

This is a negative energy indicating that the RPA Ground State is more bound than the HF one, what in turn is due to the Correlations Included in the RPA (c.f. Bohr&Mottelson Vol.II)

Matrix elements due to GSC Pauli rearrangement

The contribution of a given p-h configuration to the GS Correlation Energy is (*B&MII*)

The presence of a new neutron (scattering- or bound-like) inhibits some of these correlations, producing an energy modification of the core state...



This is the meaning/value of the NFT self energy diagram (B&MII, eq.6.225)

Meaning of the auxiliary GSCPR channel.

$$u_{a}^{x}(r) = \sum_{i} x_{ai} R_{ai}^{WS}(r); e_{ai} > e_{F}$$

$$u_{b}^{C}(r) = \sum_{i} C_{bi} R_{bi}^{WS}(r); e_{bi} > e_{F}$$

$$v_{c}^{D}(r) = \sum_{i} D_{ci} R_{ci}^{WS}(r); e_{ci} < e_{F}$$

GSC Pauli rearr.: An auxiliary Coupled Channel

$$\Psi_{jama} = \left[\psi_{jama}^{x} + \left[\psi_{jb}^{C} \cdot \Gamma_{\lambda}^{+}\right]_{jama}\right] \Phi_{GS}$$

but if

$$\Psi_{GS} = \left[1 + \varepsilon \left[\psi_{ja}^{x} \cdot \left[\psi_{jc}^{-1} \cdot \Gamma_{\lambda}^{+}\right]_{ja}\right]\right]_{J=0} \Phi_{GS}^{HF}$$

a new term can 1 must be added

$$\Psi_a = [\psi_a^x + [\psi_b^C \cdot \Gamma_\lambda^+]_{j_a} - [\psi_c^D \cdot \Gamma_\lambda]_{j_a} + \dots] \Phi_{GS}$$

with the hole – phonon anihilator

 $[\psi_c^D \cdot \Gamma_{\lambda}]_{j_a} = (v_c^D(r)/r) [\Theta_{j_c} \cdot \Phi_{\lambda}^{-1}]_{j_a},$

$$v_{c}^{D}(r) = \sum_{i} D_{ci} R_{ci}^{WS}(r); e_{ci} < e_{F}$$

p-h + phonon "virtual"excitations with respect to the HF GS.

Creation of a neutron in an "occupied" level jc levels + phonon anihilation may/will give a non nule contribution

$$\begin{pmatrix} H_p - e_F & \Xi_{a,b\lambda} f(r) & \Xi_{a,c\lambda} f(r) \\ \Xi_{a,b\lambda} f(r) & H_p - e_F + \hbar \omega & 0 \\ \Xi_{a,c\lambda} f(r) & 0 & (H_p - e_F) - \hbar \omega \end{pmatrix} \begin{pmatrix} u_a^x \\ u_b^C \\ -v_c^D \end{pmatrix} = \tilde{E} \begin{pmatrix} u_a^x \\ u_b^C \\ -v_c^D \end{pmatrix}$$

PVC: Vibrational Core (even-even) + One particle (neutron)

$$H = H_{coll} + H_{sp} + H_{PVC}$$

$$H_{coll} = \sum_{\lambda,\mu\nu} \hbar \omega_{\lambda,\nu} [\Gamma^{+,}_{\lambda\mu,\nu} - \Gamma_{\lambda\mu,\nu} + 1/2]$$

$$H_{sp} = -\hbar^{2}/2 \mu d^{2}/d \vec{r}^{2} + V(r) + V_{ls}(r)$$

$$H_{PVC} = \sum_{\lambda,\mu\nu} \delta V_{\lambda\nu}(r) Y_{\lambda\mu}(\hat{r}) [\Gamma^{+,}_{\lambda\mu,\nu} + (-1)^{\mu} - \Gamma_{\lambda\mu,\nu}]$$

$$\delta V_{\lambda\nu}(r) \approx -rdV/dr \beta_{\lambda,\nu}$$

$$Extended \begin{bmatrix} [-\hbar^{2}/2 \mu d^{2}/d r^{2} + V_{a}(r)] \\ \equiv_{a,b\lambda}(-\beta_{\lambda}rdV/dr) \\ \equiv_{a,b\lambda}(-\beta_{\lambda}rdV/dr) \end{bmatrix} \begin{bmatrix} R_{a}^{*}(r) \\ R_{b}^{*}(r) \end{bmatrix} = \tilde{E} \begin{bmatrix} R_{a}^{*}(r) \\ R_{b}^{*}(r) \end{bmatrix}$$

$$Extended Coupled \\ Channels \\ Equations with = \sum_{a,b\lambda} = \langle \Theta_{j,m_{e}} \sum_{\mu} Y_{\lambda\mu} [\Gamma^{+,}_{\lambda\mu} + (-1)^{\mu} - \Gamma_{\lambda\mu}] [\Theta_{j,b} - \Phi_{\lambda}]_{j,m_{e}} \rangle \sim \langle j_{b}, 1/2; \lambda 0| j_{a}, 1/2 \rangle$$

$$1. Here the zero point energy \sum_{\lambda} 1/2 \hbar \omega_{\lambda}$$
 is assigned to the GS. Later we will discuss a more dynamical role of this term $\sum_{\lambda} 1/2 \hbar \omega_{\lambda}$ is assigned to the GS. Later we will discuss a more dynamical role of this term $\sum_{\lambda} 1/2 \hbar \omega_{\lambda}$ is assigned to the GS. Later we will discuss a more dynamical role of this term $\sum_{\lambda} 1/2 \hbar \omega_{\lambda}$ is assigned to the GS. Later we will discuss a more dynamical role of this term $\sum_{\lambda} 1/2 \hbar \omega_{\lambda}$ is assigned to the GS. Later we will discussed later.
3. The core is supposed to be closed shell. Open shell (superfluid) cores will be discussed later.
4. (n,d) and (n,t) channels will be discussed later

RESULTADOS poner lo del PRL

Treatment of the neutron-proton interaction.

simple but quantitative treatment of the effects related to the presence of the odd proton in 12 B, that can be directly extrapolated to case of 10 Li.

The lowest states observed in the energy spectrum of the odd-odd nucleus 12B are two doublets with angular momentum and parity (1+; 2+) and (1-; 2-) which can be naturally interpreted as arising from the coupling of the odd proton lying on the 1p3=2 orbital with the odd neutron lying on the 1p1=2 $(1p3=2 \ 1p1=2)$ or on the 2s1=2 $(1p3=2 \ 2s1=2)$ weakly bound orbitals. The doublets display a similar energy splitting equal to approximately 0.95 MeV. The 1+ and the 2- states lie lowest in the two multiplets. The relative position of the centroids of the (1+; 2+) (Ecent = 0:59 MeV) and (1-; 2-)doublets (Ecent = 2:02) in 12B indicates a dierence in the position of the 1p1=2or on the 2s1=2 neutron states equal to ps = -1.43 MeV, to be compared with with the value ps = -3.09 MeV associated with the relative position of the States 1/2- (g.s.) and 1/2+ (3.09 MeV) in 13C.

PHYSICAL REVIEW C 69, 041302(R) (2004)

Parity inversion and breakdown of shell closure in Be isotopes

G. Gori,^{1,2} F. Barranco,³ E. Vigezzi,² and R. A. Broglia^{1,2,4}

¹Dipartimento di Fisica, Università degli Studi di Milano, via Celoria 16, 20133 Milano, Italy

²INFN, Sezione di Milano, via Celoria 16, 20133 Milano, Italy

³Departamento de Fisica Aplicada III, Escuela Superior de Ingenieros, Camino de los Descubrimientos s/n, 41092 Sevilla, Spain

⁴The Niels Bohr Institute, University of Copenhagen, Blegdamsvej 17, 2100 Copenhagen Ø, Denmark

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Why? What for?

NFT has been mostly used in the study of the structure of well bound nuclei:

Single-particle and multiplet states: Hamamoto Phys.Rep.(73), Mahaux,Bortignon, Broglia and Dasso, Phys.Rep.(85) Density distribution: F.B. and Broglia, PLB(85) Width of the Giant Resonances: Colo', Bortignon and Broglia, NPA(87) Superfluidity properties: Idini et al.,PRC(2012,2015) Studies of the **structure** of nuclei close to the **neutron drip line** like **11-Li, 11-Be and 12-Be** led us naturally to the need of **including the continuum** in the model basis for describing **ground state** and bound excited states properties (*Barranco et a.,EPJ(2001*) (nicely confirmed in 11-Li by the Tanihata et al. (p,t) data) (*Potel et al. PRL(2010*))

The study of **resonant states** is now under development, the formal **connection with scattering and reaction** processes has appeared naturally.

These formal advances find simple expression in the **Coupled Channels** formalism. (*Parallel developments have been pursuid by Mizuyama et. al. and by Blanchon et al. in terms of microspical optical potentials; connections will be established*)

Using NFT as the (exact!) starting point allows in principle the exact treatment of some interesting aspects of the reaction side, as the **antisymmetrization in n-N reactions.**

Summary/Message

Small nuclei mean field/PVC based description must consider several levels of correlations, but **still a useful tool,** in order to give a **global** description of low energy spectroscopy data.

Nice laboratories for many body physics.





Nuclear Physics A 683 (2001) 48-78



www.elsevier.nl/locate/npe

Single-neutron transfer from ¹¹Be_{gs} via the (p, d) reaction with a radioactive beam

J.S. Winfield^{a,b,*}, S. Fortier^a, W.N. Catford^b, S. Pita^a, N.A. Orr^c, J. Van de Wiele^a, Y. Blumenfeld^a, R. Chapman^d, S.P.G. Chappell^e, N.M. Clarke^f, N. Curtis^{b,1}, M. Freer^f, S. Galès^a, H. Langevin-Joliot^a, H. Laurent^a, I. Lhenry^a, J.M. Maison^a, P. Roussel-Chomaz^g, M. Shawcross^b, K. Spohr^d, T. Suomijärvi^a, A. de Vismes^g

Abstract

The ¹¹Be(p, d)¹⁰Be reaction has been performed in inverse kinematics with a radioactive ¹¹Be beam of E/A = 35.3 MeV. Angular distributions for the 0⁺ ground state, the 2⁺, 3.37 MeV state and the multiplet of states around 6 MeV in ¹⁰Be were measured at angles up to 16^o_{Cm} by detecting the ¹⁰Be in a dispersion-matched spectrometer and the coincident deuterons in a silicon array. Distorted wave and coupled-channels calculations have been performed to investigate the amount of 2⁺ core excitation in ¹¹Be_{gs}. The use of "realistic" ¹¹Be wave functions is emphasised and bound-state form factors have been obtained by solving the particle-vibration coupling equations. This calculation admixture. Cross sections calculated with these form factors are in good agreement with the present data. The Separation Energy prescription for the bound-state wave function also gives satisfactory fits to the data, but leads to a significantly larger [2⁺ \otimes 1*d*] component in ¹¹Be_{gs}. \otimes 2001 Elsevier Science B.V. All rights reserved.

Appendix A. Transfer form factors in a vibrational coupling approach

It has long been established that in the presence of core-polarisation admixtures created by an interaction Hamiltonian H_{int} , radial form factors of one-nucleon transfer reactions should no more be approximated by the product of a spectroscopic amplitude and a singleparticle wave function U_{lj}^{SE} deduced from the standard separation-energy procedure, but have to be determined by solving the Pinkston–Satchler coupled equations [71]. In the present analysis of the ¹¹Be(p, d)¹⁰Be reaction, these coupled equations were solved in the framework of the particle-vibration coupling model [72,73], with the program CCVIB [55]. The interaction Hamiltonian was chosen as:

$$H_{\rm int} = -R \frac{\mathrm{d}V}{\mathrm{d}r} \sum_{LM} \frac{\mathrm{i}^{-L} \beta_L}{\sqrt{2L+1}} \left(c_{LM}^{\dagger} + (-1)^{L+M} c_{L-M} \right) Y_{LM}^*(\theta, \phi), \tag{A.1}$$

where c_{LM}^{\dagger} and c_{L-M} are operators for phonon creation and annihilation and β_L is the deformation parameter. Here the term dV/dr corresponds to the deformation of the central part V(r) of the potential, with no action on the spin-orbit part. Coupled-channel calculations were performed for both $1/2^+$ ground state and $1/2^-$ first excited state at 0.32 MeV, with the additional constraint that the same well depth V_0 reproduced the experimental separation energies relative to the ground state of ${}^{10}\text{Be}$, thus ensuring the right asymptotic behaviour of the wave functions. This reproduction of the separation energies could be obtained by adjusting either the value of the β_2 parameter or the strength of the spin-orbit part of the potential. The configuration space was truncated to the 0_{gs}^+ and 2_1^+ core states in ${}^{10}\text{Be}$ (first-order vibrational coupling), coupled with one neutron in the $1p_{3/2}$ or $1p_{1/2}$, and $2s_{1/2}$ or $1d_{5/2}$ orbitals, for the $1/2^-$ and $1/2^+$ states, respectively. Renormalization processes acting on s-, p- and d-waves Coupling to pair vibrations included

Parameters are varied around a standard parameterization of the neutron-core potential (Bohr-Mottelson)

$$V_0 = -51 + \frac{33(N-Z)}{A} \text{ MeV}$$
; $a = 0.67 \text{ fm};$; $R = 1.27A^{1/3},$

 $(\boldsymbol{\beta}_2)_0^{\text{nuc}} = 0.67$ $(\boldsymbol{\beta}_2)_0^{\text{Coul}} = 1.15$

$E(p_{1/2})$	$\tilde{E}(1/2^{-})$	$A(1/2^{-})$	$E(s_{1/2})$	$\tilde{E}(1/2^{+})$	$A(1/2^{+})$	$E_{res}(d_{5/2})$	$\tilde{E}(5/2^{+})$
-2.22	-0.19	0.15	0.18	-0.18	0.07	3.5	1.21

$\beta_2/(\beta_2)_0$	$eta_3/(eta_3)_0$	$V_{WS}/(V_{WS})_0$	$V_{ls}/(V_{ls})_0$	a	$m_k(0)/m$
1.3	0.9	1.0	1.0	0.74	0.9



Initial single-particle wavefunctions and Final many-body wavefunctions R_{box} = 30 fm



Determination of phase shifts, $R_{box} = 40$ fm




TRANSFER FORM FACTORS





Admixture of $5/2^+$ state with 2^+ phonon





(n,nn) channel: Coupling to pair modes

There are Ground State Correlations apart from a particle-hole plus a surface vibration (a correlated p-h)



In fact: 2 holes plus a (correlated) 2 particles-state (pair addition) and 2 particles plus a (correlated) 2 holes-state (pair removal)





...(n,np) channel: Coupling to T=0 pairs



 Φ_{npL} is the deuteron-mode wave with angular momentum L (up to 8 @ 1MeV), scattering ones included!!...(n,d) can be calculated

...(n,np) channel: Coupling to T=0 pairs



Note: The T=0 pair correlations continue to be a Nuclear Structure challenge.

The direct connection to reaction processes opens new windows for experimentally studying it.

One notable difference with T=1 pairing is the low energy GSC viability, which for T=1 is always perfect, while for T=0, due to the asymmetry between proton and neutron mean field, is much more dificult.

At higher energies instead the strongest n-p interaction favours the T=0 channel

(n,t) channel



The interaction between the T=0 pair and the incoming neutron will lead to bound structures (tritium) whose scattering waves can be analyzed to obtain the transfer amplitudes

(p,t) channel



The interaction between the T=1 pair and the incoming proton will lead to bound structures (tritium) whose scattering waves can be analyzed to obtain the transfer amplitudes

(p,t) using 2nd order DWBA

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Unified description of structure and reactions: implementing the nuclear field theory program

R A Broglia^{1,2}, P F Bortignon^{1,3}, F Barranco⁴, E Vigezzi^{3,8}, A Idini⁵ and G Potel⁶⁷ ¹Dipartimento di Fisica, Università degli Studi Milano, Via Celoria 16, I-20133 Milano, Italy ²The Niels Bohr Institute, University of Copenhagen, DK-2100 Copenhagen, Denmark ³INFN Sezione di Milano, Via Celoria 16, I-20133 Milano, Italy ⁴Departamento de Fisica Aplicada III, Escuela Superior de Ingenieros, Universidad de Sevilla, Camino de los Descubrimientos, Sevilla, Spain ⁵Department of Physics, University of Jyvaskyla, FI-40014 Jyvask



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THANK YOU

In the previous calculation, many-phonon contributions have been neglected

Configuration of the dressed 5/2⁺ state E ~1.4 MeV :

 $0.9 \, 1d_{5/2} + 0.4 \, 2s_{1/2} \times 2+$, where the $2s_{1/2}$ state is not bound

However, our dressed $1/2^+$ configuration is bound, with $1/2^+ = 0.95 \text{ s}_{1/2} + 0.3 \text{ d}_{5/2} \times 2^+$

By iterating, in the calculation of the dressed $5/2^+$ state we take into account the 2-phonon contribution and the fact that the $1/2^+$ state is localized.

Empirical renormalization : instead of iterating, take empirical states as intermediate states, and check the consistency of the final solution

The energy of the 5/2+ state can be raised reducing the depth of the potential for I=2 waves

Add the admixture

$V(d_{5/2})/(V_{WS})_0$	$E(p_{1/2})$	$\tilde{E}(p_{1/2})$	$A(p_{1/2})$	$E(s_{1/2})$	$\tilde{E}(s_{1/2})$	$A(s_{1/2})$	$E_{res}(d_{5/2})$	$\tilde{E}(d_{5/2})$
0.5	-1.97	-0.21	0.15	0.18	-0.51	0.16	-	1.65
0.6	-1.97	-0.21	0.15	0.18	-0.51	0.16	-	1.42
0.7	-1.97	-0.21	0.15	0.18	-0.51	0.16	-	1.13
0.8	-2.22	-0.35	0.12	0.18	-0.31	0.13	-	1.15
0.9	-1.98	-0.15	0.12	0.22	-0.44	0.12	6.5	0.66
1.0	-1.98	-0.02	0.12	0.22	-0.18	0.13	4.2	0.09
1.1	-1.98	-0.02	0.13	0.22	-0.17	0.12	2.6	-0.71

Admixture of 5/2⁺ state with 2⁺ phonon doubles with empirical renormalization

 $V(d5/2) = (V(d5/2))_{0}$

No confined intermediate s1/2 x2⁺ state

 $V(d5/2) = 0.7 (V(d5/2))_{0}$

Confined intermediate s1/2 x 2+ state at 3.3 - 0.5 = 2.8 MeV

Single particle resonance at 3.5 MeV

No single particle resonance





Lay's Waves...



Optical Potential as the r,r'-E-dependentSelf-energy (c.f. Idini and Mizuyama)

$$\begin{split} u_a^x(r) &= \sum_i x_{ai} R_{ai}^{WS}(r); E_{ai} > 0 \\ u_b^C(r) &= \sum_i C_{bi} R_{bi}^{WS}(r); E_{bi} > 0 \\ v_c^D(r) &= \sum_i D_{ci} R_{ci}^{WS}(r); E_{ci} < 0 \\ H_p - e_F &\equiv_{a,b\lambda} f(r) &\equiv_{a,c\lambda} f(r) \\ \Xi_{a,b\lambda} f(r) & H_p - e_F + \hbar \omega & 0 \\ \Xi_{a,c\lambda} f(r) & 0 & (H_p - e_F) - \hbar \omega \\ \int \left(\frac{u_a^C}{-v_c^D} \right) &= \tilde{E} \begin{pmatrix} u_a^x \\ u_b^C \\ -v_c^D \end{pmatrix} \\ C_{bk} &= \frac{\Xi_{a,b\lambda} \int dr' R_{bk}^{WS}(r') f(r') u_{ai}^x(r')}{\tilde{E} - (E_{bk} + \hbar \omega_{\lambda})} \\ D_{ck} &= \frac{-\Xi_{a,b\lambda} \int dr' R_{ck}^{WS}(r') f(r') u_{ai}^x(r')}{\tilde{E} + (E_{ck} + \hbar \omega_{\lambda})} \end{split}$$

$$(H_{p}(r)-e_{F})u_{ai}^{x}(r)+\int dr' \left[\sum_{bk>0} \frac{\Xi_{a,b\lambda}^{2}R_{bk}^{WS}(r')R_{bk}^{WS}(r)f(r')f(r)}{\tilde{E}-(E_{bk}+\hbar\omega_{\lambda})}+\sum_{ck<0} \frac{\Xi_{a,b\lambda}^{2}R_{ck}^{WS}(r')R_{ck}^{WS}(r)f(r')f(r)}{\tilde{E}+(E_{ck}+\hbar\omega_{\lambda})}\right]u_{ai}^{x}(r')$$

$$=\tilde{E}u_{ai}^{x}(r)$$

Microscopic optical potentials

In HPVC only linear terms are/must be included. Higher orders will be discussed here.

$$C_{bk} = \frac{\Xi_{a,b\lambda} \int dr' R_{bk}^{WS}(r') f(r') u_{ai}^{x}(r')}{\tilde{E}_{ai} - (E_{bk} + \hbar \omega_{\lambda})} \longrightarrow \frac{\Xi_{a,b\lambda} \int dr' u_{bk}^{x}(r') f(r') u_{ai}^{x}(r')}{\tilde{E}_{ai} - (\tilde{E}_{bk} + \hbar \omega_{\lambda})} \left(\tilde{E}_{bk}(part.) > 0\right)$$

$$D_{ck} = \frac{-\Xi_{a,b\lambda} \int dr' R_{ck}^{WS}(r') f(r') u_{ai}^{x}(r')}{\tilde{E}_{ai} + (E_{ck} + \hbar \omega_{\lambda})} \longrightarrow -\frac{\Xi_{a,c\lambda} \int dr' v_{ck}^{x}(r') f(r') u_{ai}^{x}(r')}{\tilde{E}_{ai} + (\tilde{E}_{ck} + \hbar \omega_{\lambda})} \left(\tilde{E}_{ck}(hole) > 0\right)$$

This is the non-crossing phonon aproximation to the many phonons full solution. It excludes the vertex correction diagrams, that should be incorporated independently

(a)

$$\Sigma(\mathbf{k}, i\omega_n) = \underbrace{\sum_{k=1}^{(a)} (\mathbf{k}, i\omega_n)}_{\mathbf{k}, (\mathbf{k}, i\omega_n)} = \underbrace{\sum_{k=1}^{(a)} (\mathbf{k}, i\omega_n)}_{\mathbf{k}, (\mathbf{k}, i\omega_n)}_{\mathbf{k}, (\mathbf{k}, i\omega_n)} = \underbrace{\sum_{k=1}^{(a)} (\mathbf{k}, i\omega_n)}_{\mathbf{k}, (\mathbf{k$$

These equations must be solved iteratively or $= \tilde{E}_{ai} u_{ai}^{x}(r)$ Making a good ansatz for the final solution->experimental info: Empirical Renormalization

No free parameter !







Structure and scattering from (Gorkov-type) self-energies Andrea Idini

Towards consistent approaches for nuclear structure and reactions 09 Jun 2016

Partial Waves Contribution



C. MAHAUX

Institute of Physics B5, University of Liège, Sart Tilman, B-4000 Liège 1, Belgium

P.F. BORTIGNON

Istituto di Fisica Galileo Galilei, I-35100 Padova, Italy

R.A. BROGLIA

Niels Bohr Institutet, University of Copenhagen, DK-2100 Copenhagen Ø, Denmark

and



NORDITA, Blegdamsvej 17, DK-2100 Copenhagen Ø, Denmark





Fig. 4.45a. Energy dependence of the quantity $\Delta V_{2g9/2}(E)$ defined by eq. (3.18a), for the energy-averaging interval $\Delta = 1$ MeV. The straight dashed line represents $E - \varepsilon_{2g9/2}^{(0)}$, where $\varepsilon_{2g9/2}^{(0)} = -3.3$ MeV is derived from the Skyrme III-Hartree-Fock approximation. The approximation scheme is that of ref. [283].

Fig. 4.45b. Energy dependence of the quantity $W_{2g9/2}(E)$ defined by eq. (3.18b), for $\Delta = 1$ MeV. The location of the Skyrme III-Hartree-Fock approximation to the single-particle energy is indicated by an arrow. The large peak of $W_{2g9/2}(E)$ at $E \approx 0.5$ MeV is due to the coupling to the $\{1j_{15/2}\}_*3^-\}$ configuration.

(n,nn) channel: Coupling to pair modes

There are Ground State Correlations apart from a particle-hole plus a surface vibration (a correlated p-h)



In fact: 2 holes plus a (correlated) 2 particles-state (pair addition) and 2 particles plus a (correlated) 2 holes-state (pair removal)





No open channels since dineutron is not stable, but...

...(n,np) channel: Coupling to T=0 pairs



 Φ_{npL} is the deuteron-mode wave with angular momentum L (up to 8 @ 1MeV), scattering ones included!!

...(n,np) channel: Coupling to T=0 pairs



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The interaction between the T=1 pair and the incoming proton will lead to bound structures (tritium) whose scattering waves can be analyzed to obtain the transfer amplitudes

(p,t) using 2nd order DWBA

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...and beyond

Totally especulative:

Heavy ions collisions.... TDHF... TDNFT

Open shell (superfluid) nucleii.1 Spinor CC equations

$$\begin{pmatrix} H_{p}-e_{F} & \Xi_{a,b\lambda}f(r) & \Xi_{a,c\lambda}f(r) \\ \Xi_{a,b\lambda}f(r) & H_{p}-e_{F}+\hbar\omega & 0 \\ \Xi_{a,c\lambda}f(r) & 0 & (H_{p}-e_{F})-\hbar\omega \end{pmatrix} \begin{pmatrix} u_{a}^{x} \\ u_{b}^{C} \\ -v_{c}^{D} \end{pmatrix} = \tilde{E} \begin{pmatrix} u_{a}^{x} \\ u_{b}^{C} \\ -v_{c}^{D} \end{pmatrix}$$

$$= \tilde{E} \begin{pmatrix} v_{a}^{x} \\ u_{b}^{C} \\ -v_{c}^{D} \end{pmatrix}$$

$$= \tilde{E} \begin{pmatrix} v_{a}^{x} \\ v_{b}^{C} \\ v_{b}^{C} \\ -u_{c}^{D} \end{pmatrix} = \tilde{E} \begin{pmatrix} v_{a}^{x} \\ v_{b}^{C} \\ -u_{c}^{D} \end{pmatrix}$$
Holes

Using the Pair-Spin notation, these six equation can be written as

$$\begin{pmatrix} H_{p} - e_{F} & 0\\ 0 & -(H_{p} - e_{F}) \end{pmatrix} \begin{pmatrix} u_{a}^{x}\\ v_{a}^{x} \end{pmatrix} + \Xi_{a,b\lambda} f(r) \sum_{k} \begin{pmatrix} C_{bk} & 0\\ 0 & -C_{bk} \end{pmatrix} \begin{pmatrix} u_{bk}\\ v_{bk} \end{pmatrix} + \Xi_{a,c\lambda} f(r) \sum_{k} \begin{pmatrix} 0 & -D_{ck}\\ -D_{ck} & 0 \end{pmatrix} \begin{pmatrix} u_{ck}\\ v_{ck} \end{pmatrix} = \tilde{E} \begin{pmatrix} u_{a}^{x}\\ v_{a}^{x} \end{pmatrix}$$

$$\Xi_{a,b\lambda} f(r) \begin{pmatrix} u_{a}^{x}\\ v_{a}^{x} \end{pmatrix} + \sum_{k} (E_{ck} + \hbar \omega) \begin{pmatrix} 0 & -D_{ck}\\ 0 & -C_{bk} \end{pmatrix} \begin{pmatrix} u_{ck}\\ v_{ck} \end{pmatrix} = \tilde{E} \sum_{k} \begin{pmatrix} 0 & -D_{ck}\\ 0 & -C_{bk} \end{pmatrix} \begin{pmatrix} u_{ck}\\ v_{ck} \end{pmatrix}$$

$$\Xi_{a,b\lambda} f(r) \begin{pmatrix} u_{a}^{x}\\ v_{a}^{x} \end{pmatrix} + \sum_{k} (E_{ck} + \hbar \omega) \begin{pmatrix} 0 & -D_{ck}\\ -D_{ck} & 0 \end{pmatrix} \begin{pmatrix} u_{ck}\\ v_{ck} \end{pmatrix} = \tilde{E} \sum_{k} \begin{pmatrix} 0 & -D_{ck}\\ -D_{ck} & 0 \end{pmatrix} \begin{pmatrix} u_{ck}\\ v_{ck} \end{pmatrix}$$

Open shell (superfluid) cores 2. Adding pairing field

$$\begin{pmatrix} H_{p} - e_{F} & \Delta(r) \\ \Delta(r) & -(H_{p} - e_{F}) \end{pmatrix} \begin{pmatrix} u_{a}^{x}(r) \\ v_{a}^{x}(r) \end{pmatrix} + \sum_{b} \Xi_{a,b\lambda} f(r) \sum_{k} \begin{pmatrix} C_{bk} & -D_{bk} \\ -D_{bk} & -C_{bk} \end{pmatrix} \begin{pmatrix} u_{bk}(r) \\ v_{bk}(r) \end{pmatrix} = \tilde{E} \begin{pmatrix} u_{a}^{x}(r) \\ v_{a}^{x}(r) \end{pmatrix}$$

$$\Xi_{a,b\lambda} f(r) \begin{pmatrix} u_{a}^{x}(r) \\ v_{a}^{x}(r) \end{pmatrix} + \sum_{k} (E_{bk} + \hbar \omega) \begin{pmatrix} C_{bk} & -D_{bk} \\ -D_{bk} & -C_{bk} \end{pmatrix} \begin{pmatrix} u_{bk}(r) \\ v_{bk}(r) \end{pmatrix} = \tilde{E} \sum_{k} \begin{pmatrix} C_{bk} & -D_{bk} \\ -D_{bk} & -C_{bk} \end{pmatrix} \begin{pmatrix} u_{bk}(r) \\ v_{bk}(r) \end{pmatrix}$$

$$C_{bk} = \frac{\Xi_{a,b\lambda}}{\tilde{E}_{ai} - (E_{bk} + \hbar \omega_{\lambda})} \langle (u_{bk}(r), -v_{bk}(r)) f(r) \begin{pmatrix} u_{ai}^{x}(r) \\ v_{ai}^{x}(r) \end{pmatrix} \rangle$$
$$D_{bk} = \frac{-\Xi_{a,b\lambda}}{\tilde{E}_{ai} + (E_{bk} + \hbar \omega_{\lambda})} \langle (v_{bk}(r), u_{bk}(r)) f(r) \begin{pmatrix} u_{ai}^{x}(r) \\ v_{ai}^{x}(r) \end{pmatrix} \rangle$$

Open shell (superfluid) cores 3. Self-energy

$$\int dr' \left(\begin{pmatrix} H_p(r,r') - e_F & \Delta(r,r') \\ \Delta(r,r') & -(H_p(r,r') - e_F) \end{pmatrix} + \begin{pmatrix} \Sigma_a^{11}(r,r';\tilde{E}) & \Sigma_a^{12}(r,r';\tilde{E}) \\ \Sigma_a^{21}(r,r';\tilde{E}) & \Sigma_a^{22}(r,r';\tilde{E}) \end{pmatrix} \right) \begin{pmatrix} u_a^x(r') \\ v_a^x(r') \end{pmatrix} = \tilde{E} \begin{pmatrix} u_a^x(r) \\ v_a^x(r) \end{pmatrix}$$

$$\begin{split} \Sigma_{a}^{11}(r,r';\tilde{E}) &= \sum_{bk} \Xi_{a,b\lambda}^{2} f(r) f(r') \left[\frac{u_{bk}(r)u_{bk}(r')}{\tilde{E} - (E_{bk} + \hbar \omega_{\lambda})} + \frac{v_{bk}(r)v_{bk}(r')}{\tilde{E} + (E_{bk} + \hbar \omega_{\lambda})} \right] \\ \Sigma_{a}^{12}(r,r';\tilde{E}) &= \sum_{bk} \Xi_{a,b\lambda}^{2} f(r) f(r') \left[\frac{-u_{bk}(r)v_{bk}(r')}{\tilde{E} - (E_{bk} + \hbar \omega_{\lambda})} + \frac{v_{bk}(r)u_{bk}(r')}{\tilde{E} + (E_{bk} + \hbar \omega_{\lambda})} \right] \\ \Sigma_{a}^{21}(r,r';\tilde{E}) &= \sum_{bk} \Xi_{a,b\lambda}^{2} f(r) f(r') \left[\frac{-v_{bk}(r)u_{bk}(r')}{\tilde{E} - (E_{bk} + \hbar \omega_{\lambda})} + \frac{u_{bk}(r)v_{bk}(r')}{\tilde{E} + (E_{bk} + \hbar \omega_{\lambda})} \right] \\ \Sigma_{a}^{22}(r,r';\tilde{E}) &= \sum_{bk} \Xi_{a,b\lambda}^{2} f(r) f(r') \left[\frac{v_{bk}(r)v_{bk}(r')}{\tilde{E} - (E_{bk} + \hbar \omega_{\lambda})} + \frac{u_{bk}(r)u_{bk}(r')}{\tilde{E} + (E_{bk} + \hbar \omega_{\lambda})} \right] \end{split}$$

c.f. Van der Sluys et al. NPA551.

Open shell (superfluid) cores 4. Self-energy

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Fragmentation of single-particle strength in spherical open-shell nuclei: Application to the spectral functions in ¹⁴²Nd

V. Van der Sluys, D. Van Neck¹, M. Waroquier² and J. Ryckebusch¹ Institute for Nuclear Physics and Laboratory for Theoretical Physics, Proefluinstraat 86, B-9000 Ghent, Belgium

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$$\left[\begin{pmatrix} E_a & 0\\ 0 & -E_a \end{pmatrix} + \begin{pmatrix} \Sigma_{11}(E) & \Sigma_{12}(E)\\ \Sigma_{12}(E) & \Sigma_{22}(E) \end{pmatrix} \right] \begin{pmatrix} x_0\\ y_0 \end{pmatrix} = E \begin{pmatrix} x_0\\ y_0 \end{pmatrix}.$$
 (24)

The energy-dependent self-energy matrix elements $\Sigma_{ii}(E)$ stand for:

$$\Sigma_{11}(E) = \sum_{bJ\nu} \left(\frac{|V(abJ\nu)|^2}{E - (E_b + E_{J\nu})} + \frac{|W(abJ\nu)|^2}{E + (E_b + E_{J\nu})} \right),$$

$$\Sigma_{22}(E) = \sum_{bJ\nu} \left(\frac{|W(abJ\nu)|^2}{E - (E_b + E_{J\nu})} + \frac{|V(abJ\nu)|^2}{E + (E_b + E_{J\nu})} \right),$$

$$\Sigma_{12}(E) = -\sum_{bJ\nu} V(abJ\nu) W(abJ\nu) \left(\frac{1}{E - (E_b + E_{J\nu})} - \frac{1}{E + (E_b + E_{J\nu})} \right).$$
 (25)




Tadpole, Triangle, Multiplet-C and Butterfly Diagrams



Part.-Phon.-Int.: Triangle, Multiplet-C and Butterfly Diagrams



Tadpole, Triangle, Multiplet-C and Butterfly Diagrams



Part.-Phon.-Vertex and Two-Phonon Energy Butterfly Diagrams QRPA estimate (using up1/2=vp1/2=1/sqrt(2); vp3/2=sqrt(3)up3/2=sqrt(3/4)









MAIN PVC EFFECTS

PVC/NFT: Vibrational Core (even-even) + One particle (neutron)

$$H = H_{c} + H_{p} + H_{PVC}$$

$$H_{c} = \sum_{\lambda\mu\nu} \hbar \omega_{\lambda,\nu} [\Gamma_{\lambda\mu,\nu}^{+} \Gamma_{\lambda\mu,\nu} + 1/2]$$

$$H_{p} = -\hbar^{2}/2 \mu d^{2}/d \vec{r}^{2} + V(r) + V_{ls}(r)$$

$$H_{PVC} = -rdV/dr \sum_{\lambda\mu\nu} Y_{\lambda\mu}(\hat{r})\beta_{\lambda,\nu} [\Gamma_{\lambda\mu,\nu}^{+} + (-1)^{\mu}\Gamma_{\lambda\mu,\nu}]$$

$$\begin{bmatrix} -\hbar^{2}/2 \mu d^{2}/d r^{2} + V_{a}(r)] \qquad \Xi_{a,b\lambda}(-\beta_{\lambda}rdV/dr)$$

$$\Xi_{a,b\lambda}(-\beta_{\lambda}rdV/dr) \qquad [-\hbar^{2}/2 \mu d^{2}/d r^{2} + V_{b}(r) + \hbar\omega] \begin{bmatrix} R_{a}^{x}(r) \\ R_{b}^{C}(r) \end{bmatrix} = \tilde{E} \begin{bmatrix} R_{a}^{x}(r) \\ R_{b}^{C}(r) \end{bmatrix}$$
Channels
Equations with $\Xi_{a,b\lambda} = \langle \Theta_{j_{a}m_{a}} \sum_{\lambda\mu} Y_{\lambda\mu} [\Gamma_{\lambda\mu}^{+} + (-1)^{\mu}\Gamma_{\lambda\mu}] [\Theta_{j_{b}} \cdot \Phi_{\lambda}]_{j_{a}m_{a}} \rangle \sim \langle j_{b}, 1/2; \lambda 0 | j_{a}, 1/2 \rangle$

Other comments:

- 1- For r>10fm we have Bessel functions for positive E's → Phase shifts (S-matrix) may be extracted and thus cross sections (differential or integrated)
- 2. The theory input are as usual: V(r), VIs(r), the beta's and the hw's.
- 3. We aim to use a unique V(r) and VIs(r) functions: independent of Ij, of E, local (non locality will be discussed)

Matrix elements due to GSC Pauli rearrangement. 2



In HPVC only linear terms are/must be included. Higher orders will be discussed here.

$$C_{bk} = \frac{\Xi_{a,b\lambda} \int dr' R_{bk}^{WS}(r') f(r') u_{ai}^{x}(r')}{\tilde{E}_{ai} - (E_{bk} + \hbar \omega_{\lambda})} \longrightarrow \frac{\Xi_{a,b\lambda} \int dr' u_{bk}^{x}(r') f(r') u_{ai}^{x}(r')}{\tilde{E}_{ai} - (\tilde{E}_{bk} + \hbar \omega_{\lambda})} \left(\tilde{E}_{bk}(part.) > 0\right)$$

$$D_{ck} = \frac{-\Xi_{a,b\lambda} \int dr' R_{ck}^{WS}(r') f(r') u_{ai}^{x}(r')}{\tilde{E}_{ai} + (E_{ck} + \hbar \omega_{\lambda})} \longrightarrow -\frac{\Xi_{a,c\lambda} \int dr' v_{ck}^{x}(r') f(r') u_{ai}^{x}(r')}{\tilde{E}_{ai} + (\tilde{E}_{ck} + \hbar \omega_{\lambda})} \left(\tilde{E}_{ck}(hole) > 0\right)$$

This is the non-crossing phonon aproximation to the many phonons full solution. It excludes the vertex correction diagrams, that should be incorporated independently

(a)

$$\Sigma(\mathbf{k}, i\omega_n) = \underbrace{\sum_{k=1}^{(a)} (\mathbf{k}, i\omega_n)}_{\mathbf{k}, (\mathbf{k}, i\omega_n)} = \underbrace{\sum_{k=1}^{(a)} (\mathbf{k}, i\omega_n)}_{\mathbf{k}, (\mathbf{k}, i\omega_n)}_{\mathbf{k}, (\mathbf{k}, i\omega_n)} = \underbrace{\sum_{k=1}^{(a)} (\mathbf{k}, i\omega_n)}_{\mathbf{k}, (\mathbf{k$$

These equations must be solved iteratively or $= \tilde{E}_{ai} u_{ai}^{x}(r)$ Making a good ansatz for the final solution->experimental info: Empirical Renormalization

The phonon's calculation will be discused here

The butterfly diagram in the TLM (two level model), after NPA260,1(76)

$$\begin{array}{c} & \underset{m=-j,\ldots,j;}{\overset{\alpha}{=}} 2\Omega = 2j+1 \\ \text{m=-j},\ldots,j; 2\Omega = 2j+1 \\ \text{TDA:} \begin{cases} \omega = \varepsilon - 2V\Omega \\ X = \frac{\Lambda}{\varepsilon - \hbar\omega} = \frac{1}{\sqrt{2\Omega}} \\ \Lambda = V\sqrt{2\Omega} \\ \Lambda = V\sqrt{2\Omega} \end{cases}$$

Part.-Phon.-Vertex and Two-Phonon Energy Butterfly Diagrams QRPA estimate (using up1/2=vp1/2=1/sqrt(2); vp3/2=sqrt(3)up3/2=sqrt(3/4)





EBUTT= 2 x (1.50MeV) = 3.00MeV (without square)

Part.-Phon.-Int.: Butterfly Diagrams; sx2+



Part.-Phon.-Int.: Butterfly Diagrams: dx2+



The factor of 2 is for the two internal affected verteces



 $\langle (2\ 5/2)\ 1/2\ 2\ 5/2|2(5/2\ 2)\ 1/2\ 5/2\rangle = \sqrt{(2\ x\ 1/2+1)(2\ x\ 1/2+1)}(-1)^{2+5/2+2+5/2} \begin{cases} 2 & 5/2 & 1/2 \\ 2 & 5/2 & 1/2 \\ \end{cases} = 2/30$

 $(e_s + \hbar \omega - e_d - 2\hbar \omega) \approx -5.2 \text{MeV}; \quad since \quad \frac{V^2(s; 2d)}{5.2^2} \approx 0.17 \rightarrow \frac{V^2(s; 2d)}{-5.2 \text{MeV}} \approx -1.0 \text{MeV}$

thus the diagram is -2/30 MeV = -0.066 MeV

Tadpole

This is an energy independent diagram. Its effect can be incorpored in static mean fields. The monopole part of V' is totally democratic and thus leads to a democratic (state independent correction), which is included in our common empirical mean filed.

nlj Being the phonon of Quadrupole type (and thus of p1/2,p3/2^-1 nature), the quadrupole component of V' acts only in the nlj= 1p1/2, and thus with large Pauli inhibition.

In any case the tadpole is a small part of....

Charge Radii

$$\langle r^{2} \rangle_{11\text{Be}} = \left(\langle r^{2} \rangle_{10\text{Be}}^{1/2} + \left(\frac{\langle r^{2} \rangle_{151/2}^{1/2}}{11} \right)^{2} \right) \times S^{2} + (1 - S^{2}) \times \left(\langle r^{2} \rangle_{10\text{Be}}^{10\text{Be}} \left(1 + \frac{2}{4\pi} \beta_{\pi}^{2} \right) + \left(\frac{\langle r^{2} \rangle_{d5/2\,coll}^{1/2}}{11} \right)^{2} \right)^{2} \right) = .$$

$$\langle r^{2} \rangle_{10\text{Be}} + \left(\frac{\langle r^{2} \rangle_{151/2}^{1/2}}{11} \right)^{2} \times S^{2} + (1 - S^{2}) \times \left(\left(\frac{\langle r^{2} \rangle_{d5/2\,coll}^{1/2}}{11} \right)^{2} + \langle r^{2} \rangle_{10\text{Be}} \frac{2}{4\pi} \beta_{\pi}^{2} \right)$$

$$(2.47)^{2} = (2.36)^{2} + (7.10/11)^{2} \times S^{2} + (1 - S^{2}) \times (2.36^{2} 2/4\pi\beta^{2} + (3/11)^{2})$$

$$(2.47)^{2} = (2.36)^{2} + (7.10/11)^{2} \times S^{2} + (1 - S^{2}) \times (2.36^{2} 2/4\pi\beta^{2} + (3/11)^{2})$$

$$(2.47)^{2} = (2.36)^{2} + (7.10/11)^{2} \times S^{2} + (1 - S^{2}) \times (2.36^{2} 2/4\pi\beta^{2} + (3/11)^{2})$$

$$(3.10 = 5.57 + 0.42 S^{2} + (1 - S^{2})(0.89\beta^{2} + 0.07)$$

$$\text{ with } \beta = 1.2 \quad 6.10 = 5.57 + 0.42 S^{2} + (1 - S^{2})(1.35 \to S^{2}) = 0.88$$

$$\text{ with } S^{2} = 0.83 \quad 6.10 = 5.57 + 0.35 + 0.17 \times (0.89\beta^{2} + 0.07) \to \beta = 1.06$$



These are diagrams belonging to the Rainbow series, effectively included in the diagonalization.

