Nuclear Reactions II: Applications to Nuclear Structure

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Oak Ridge, June 28 2019



Outline

Why do reactions? Probing nuclear structure

- Elastic scattering
- Inelastic scattering
- One-particle transfer

2 An advanced example: 2-neutron transfer and pairing

- Pairing correlations and successive transfer
- Reaction and structure models
- A quantitative measure of pairing correlations

3 Moving forward: integrating structure and reactions

- (d, p) reactions: a unified approach
- Choosing the potential: examples of calculations

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Elastic scattering



FIG. 10. Elastic scattering for ${}^{6}\text{He} + {}^{12}\text{C}$ at 38.3 MeV/nucleon in comparison with the OM results given by the real folded potential (obtained with the CDMSY6 interaction and the Gaussian ga density for ${}^{6}\text{He}$). The dashed curve is obtained with the unrenormalized folded potential only. The solid curve is obtained by adding a complex surface polarization potential to the real folded potential. Its parameters, and those of the imaginary part, are explained in the text. The dotted line is obtained by folding the CDM3Y6 interaction with the compact Gaussian density r_{O} .

[Lapoux et al, PRC 66 (02) 034608]

traditionally used to extract optical potentials, rms radii, density distributions.

Inelastic scattering



One-particle transfer

populates states with strong single-particle content shape: angular momentum. magnitude: spectroscopic factor (single-particle strength).



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2-neutron transfer and reactions



- Reaction $A + a \longrightarrow (a 2) + (A + 2)$.
- Measure of the pairing correlations between the transferred nucleons.
- Need to correctly account for the correlated wavefunction.

Delocalization of the pair transfer process



Let's remember the Born series

$|\phi\rangle = |\phi_0\rangle + G_0 V |\phi\rangle$

$|\phi\rangle = |\phi_0\rangle + G_0 V (|\phi_0\rangle + G_0 V |\phi\rangle)$

$|\phi\rangle = |\phi_0\rangle + G_0 V (|\phi_0\rangle + G_0 V [|\phi_0\rangle + G_0 V |\phi\rangle])$

$$|\phi\rangle = |\phi_0\rangle + G_0 V \phi_0 + G_0 V G_0 V \phi_0 + G_0 V G_0 V G_0 V \phi_0 + \dots$$



Two-nucleon transfer: stick to second order

In order to account for the successive transfer of two nucleons, we stick to 2^{nd} order, $A + a \rightarrow (A + 1) + (a - 1) \rightarrow (A + 2) + (a - 2)$



- $V \rightarrow$ one-nucleon transfer
- $G_0 \rightarrow$ Green's function of each one of the many intermediate states (A+1)+(a-1)

Reaction and structure models

Structure:

$$\Phi_{i}(\mathbf{r}_{1},\sigma_{1},\mathbf{r}_{2},\sigma_{2}) = \sum_{j_{i}} B_{j_{i}} \left[\psi^{j_{i}}(\mathbf{r}_{1},\sigma_{1})\psi^{j_{i}}(\mathbf{r}_{2},\sigma_{2}) \right]_{0}^{0}$$

$$\Phi_{f}(\mathbf{r}_{1},\sigma_{1},\mathbf{r}_{2},\sigma_{2}) = \sum_{j_{f}} B_{j_{f}} \left[\psi^{j_{f}}(\mathbf{r}_{1},\sigma_{1})\psi^{j_{f}}(\mathbf{r}_{2},\sigma_{2}) \right]_{0}^{0}$$



mean field potentials

Reaction:



radial wave functions $u^{j_i}(r)$



radial wave functions $u^{j_f}(r)$

$$T_{2NT} = \sum_{j_f j_i} B_{j_f} B_{j_i} \left(T^{(1)}(j_i, j_f) + T^{(2)}_{succ}(j_i, j_f) - T^{(2)}_{NO}(j_i, j_f) \right)$$

Introducing $T^{(1)}(j_i, j_f)$, $T^{(2)}_{succ}(j_i, j_f)$ and $T^{(2)}_{NO}(j_i, j_f)$

very schematically, the first order (simultaneous) contribution is

 $T^{(1)} = \langle \beta | V | \alpha \rangle,$

while the second order contribution can be separated in a *successive* and a *non-orthogonality* term

$$T^{(2)} = T^{(2)}_{succ} + T^{(2)}_{NO}$$

= $\sum_{\gamma} \langle \beta | \mathbf{V} | \gamma \rangle \mathbf{G} \langle \gamma | \mathbf{V} | \alpha \rangle - \sum_{\gamma} \langle \beta | \gamma \rangle \langle \gamma | \mathbf{V} | \alpha \rangle.$

If we sum over a complete basis of intermediate states γ , we can apply the closure condition and $T_{NO}^{(2)}$ cancels $T^{(1)}$

the transition potential being single particle, two-nucleon transfer is a second order process.

Two particle transfer in 2-step DWBA

$$\overset{\mathbf{a}}{\underset{|\alpha\rangle}{}} \overset{\mathbf{A}}{\underset{|\gamma\rangle}{}} \overset{\mathbf{f}}{\underset{|\gamma\rangle}{}} \overset{\mathbf{F}}{\underset{|\gamma\rangle}{}} \overset{\mathbf{b}}{\underset{|\beta\rangle}{}} \overset{\mathbf{b}}{\underset{|\beta\rangle}{}} \overset{\mathbf{B}}{\underset{|\beta\rangle}{}} \overset{\mathbf{Potel et al., PRL 107 092501 (2011)}{Potel et al., PRL 105 172502 (2010)}$$

$$T_{2NT} = \sum_{j_f j_i} B_{j_f} B_{j_i} \left(T^{(1)}(j_i, j_f) + T^{(2)}_{succ}(j_i, j_f) - T^{(2)}_{NO}(j_i, j_f) \right)$$

Simultaneous transfer

$$T^{(1)}(j_i, j_f) = 2 \sum_{\sigma_1 \sigma_2} \int d\mathbf{r}_{fF} d\mathbf{r}_{b1} d\mathbf{r}_{A2} [\Psi^{j_f}(\mathbf{r}_{A1}, \sigma_1) \Psi^{j_f}(\mathbf{r}_{A2}, \sigma_2)]_0^{0*} \chi_{bB}^{(-)*}(\mathbf{r}_{bB}) \\ \times v(\mathbf{r}_{b1}) [\Psi^{j_i}(\mathbf{r}_{b1}, \sigma_1) \Psi^{j_i}(\mathbf{r}_{b2}, \sigma_2)]_0^0 \chi_{aA}^{(+)}(\mathbf{r}_{aA})$$

Two particle transfer in 2-step DWBA

$$\begin{aligned} & \overset{a}{\underset{|\alpha\rangle}{}} \overset{A}{\underset{|\gamma\rangle}{}} \overset{f}{\underset{|\gamma\rangle}{}} \overset{F}{\underset{|\gamma\rangle}{}} \overset{b}{\underset{|\beta\rangle}{}} \overset{b}{\underset{|\beta\rangle}{}} \overset{B}{\underset{|\beta\rangle}{}} \overset{O}{\underset{|\beta\rangle}{}} \overset{Potel et al., PRL 107 092501 (2011)}{}_{Potel et al., PRL 105 172502 (2010)} \\ & T_{2NT} = \sum_{j_{f}j_{i}} B_{j_{i}} \left(T^{(1)}(j_{i}, j_{f}) + T^{(2)}_{succ}(j_{i}, j_{f}) - T^{(2)}_{NO}(j_{i}, j_{f}) \right) \\ & \text{Successive transfer} \\ & T^{(2)}_{succ}(j_{i}, j_{f}) = 2 \sum_{K,M} \sum_{\substack{\sigma_{1}\sigma_{2} \\ \sigma_{1}'\sigma_{2}'}} \int d\mathbf{r}_{fF} d\mathbf{r}_{b1} d\mathbf{r}_{A2} [\Psi^{j_{f}}(\mathbf{r}_{A1}, \sigma_{1})\Psi^{j_{f}}(\mathbf{r}_{A2}, \sigma_{2})]_{0}^{0*} \\ & \times \chi^{(-)*}_{bB}(\mathbf{r}_{bB}) v(\mathbf{r}_{b1}) [\Psi^{j_{f}}(\mathbf{r}_{A2}, \sigma_{2})\Psi^{j_{i}}(\mathbf{r}_{b1}, \sigma_{1})]_{M}^{K} \\ & \times \int d\mathbf{r}'_{fF} d\mathbf{r}'_{b1} d\mathbf{r}'_{A2} G(\mathbf{r}_{fF}, \mathbf{r}'_{fF}) [\Psi^{j_{f}}(\mathbf{r}'_{A2}, \sigma_{2}')\Psi^{j_{i}}(\mathbf{r}'_{b1}, \sigma_{1}')]_{M}^{K} \\ & \times \frac{2\mu_{fF}}{\hbar^{2}} v(\mathbf{r}'_{f2}) [\Psi^{j_{i}}(\mathbf{r}'_{b2}, \sigma_{2}')\Psi^{j_{i}}(\mathbf{r}'_{b1}, \sigma_{1}')]_{0}^{0} \chi^{(+)}_{aA}(\mathbf{r}'_{aA}) \end{aligned}$$

Two particle transfer in 2-step DWBA

Non-orthogonality term

$$\begin{split} \mathcal{T}_{NO}^{(2)}(j_{i},j_{f}) &= 2\sum_{K,M}\sum_{\substack{\sigma_{1}\sigma_{2} \\ \sigma_{1}'\sigma_{2}'}} \int d\mathbf{r}_{fF} d\mathbf{r}_{b1} d\mathbf{r}_{A2} [\Psi^{j_{f}}(\mathbf{r}_{A1},\sigma_{1})\Psi^{j_{f}}(\mathbf{r}_{A2},\sigma_{2})]_{0}^{0*} \\ &\times \chi_{bB}^{(-)*}(\mathbf{r}_{bB}) v(\mathbf{r}_{b1}) [\Psi^{j_{f}}(\mathbf{r}_{A2},\sigma_{2})\Psi^{j_{i}}(\mathbf{r}_{b1},\sigma_{1})]_{M}^{K} \\ &\times \int d\mathbf{r}_{b1}' d\mathbf{r}_{A2}' [\Psi^{j_{f}}(\mathbf{r}_{A2}',\sigma_{2}')\Psi^{j_{i}}(\mathbf{r}_{b1}',\sigma_{1}')]_{M}^{K} \\ &\times [\Psi^{j_{i}}(\mathbf{r}_{b2}',\sigma_{2}')\Psi^{j_{i}}(\mathbf{r}_{b1}',\sigma_{1}')]_{0}^{0}\chi_{aA}^{(+)}(\mathbf{r}_{aA}') \end{split}$$











Transfer driven by single-particle potential (mean field) \rightarrow essentially a successive process!

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Probing pairing with 2-transfer: ¹¹²Sn(p,t)¹¹⁰Sn @ 26 MeV



Experimental data and shell model wavefunction from Guazzoni *et al.* PRC **74** 054605 (2006)

experiment very well reproduced with mean field (BCS) wavefunctions

Examples of calculations





good results obtained for halo nuclei, population of excited states, superfluid nuclei, normal nuclei (pairing vibrations), heavy ion reactions...

Potel *et al.*, Rep. Prog. Phys. **76** (2013) 106301

Absolute cross sections reproduced

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Extracting the structure information: a standard approach



- Factorization of structure and reactions.
- Can suffer from inconsistency between the two schemes.
- Extracted spectroscopic factor $S_i^2 = \sigma / \tilde{\sigma}$ problematic.

(d, p) reactions: a unified approach

- Obscribe the structure of the 2-body subsystems in some given framework of choice.
- Employ the same quantum many-body methods to work out the interactions U_{An}, U_{Ap}, U_{pn} (in general, non-local and energy-dependent).
- Solution Write down the resulting 3-body Hamiltonian H.
- Obtain cross sections from *H* using controlled approximations.

3-body Hamiltonian

$$H = T + U_{An}(r_n, r'_n, E_n, J_n, \pi_n) + U_{Ap}(r_p, r'_p, E_p, J_p, \pi_p) + U_{pn}(r_{pn}, r'_{pn}, E_{pn}, J_{pn}, \pi_{pn})$$

Disclaimer

Still not the end of the story! 3-body forces U_{Anp} not taken into account at this stage.

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(d, p) reaction as a 2-step process



U_{An} potential and neutron states



Dispersive Optical Model (DOM): Calcium isotopes



GP et al., Eur. Phys. J. A 53 (2017) 178.

- DOM used to compute (d, p) cross sections on Ca isotopes.
- Both bound and continuum neutron states described.
- DOM can be extrapolated to unknown territory (⁶⁰Ca).

Dispersive Optical Model (DOM): Calcium isotopes



Absolute transfer cross sections without spectroscopic factors.









Coupled-Cluster (CC): Ca isotopes



Coupled-Cluster (CC): Ca isotopes



structure calculated within ab initio coupled cluster framework by J. Rotureau (MSU)

