Nuclear ab intio methods

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Introduction



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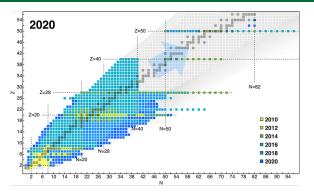


Figure: H. Hergert, Front. Phys. (2020)

Ab initio methods: Methods for solving the quantum mechanical many-body problem for all constituent nucleons, which

- describe the atomic nucleus from the bottom up with the bare nuclear forces compatible with the symmetries of QCD.
- This is done either exactly for very light nuclei or by employing certain well-controlled approximations for heavier nuclei.

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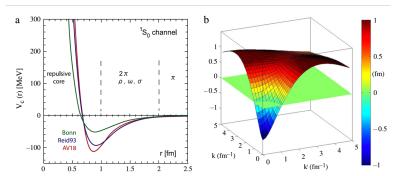


Figure: Several phenomenological NN potentials in the 1S_0 channel and momentumspace matrix elements of the Argonne v18 (AV18) 1S_0 potential after Fourier (Bessel) transformation. S.K. Bogner et al. Prog. Part. Nucl. Phys. 65 (2010) 94.

Nuclear structure calculations are complicated due to the coupling of low to high momenta by these potentials.

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The Brueckner-Hartree-Fock (BHF) theory

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The Brueckner-Hartree-Fock (BHF) theory



The G matrix is given by:

$$\langle ab|G(W)|cd\rangle = \langle ab|V|cd\rangle + \sum_{mn}\langle ab|V|mn\rangle \frac{Q(m,n)}{W-\varepsilon_m-\varepsilon_n}\langle mn|G(W)|cd\rangle$$

$$\stackrel{a}{\downarrow} G \stackrel{b}{\downarrow} = \stackrel{a}{\downarrow} V \stackrel{b}{\downarrow} + \stackrel{a}{\downarrow} - \stackrel{b}{\downarrow} - \stackrel{b}{\downarrow} - \stackrel{b}{\downarrow} - \stackrel{b}{\downarrow} - \stackrel{b}{\downarrow} - \stackrel{b}{\downarrow} - \stackrel{b}{\downarrow}$$

- W is the starting energy, its value depends on the position of G-matrix in the diagram.
- \bullet ϵ_m , ϵ_n are the HF single-particle energies.
- Q is the Pauli operator which forbids the states being scattered below Fermi surface.

K. Brueckner, C. Levinson, H. Mahmoud, Phys. Rev. 95, 217 (1954)

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The Brueckner-Hartree-Fock (BHF) theory



The HF equation is given by:

$$\sum_{j} (T + U)_{ij} D_{ja} = \varepsilon_{a} D_{ia},$$

where the one-body mean-field potential is

$$U_{ij} = \sum_{c=1}^{A} \langle ic | \bar{G}(W) | jc \rangle.$$

Rel. BHF: S.H. Shen et al., Prog. Part. Nucl. Phys. 109, 103713 (2019).

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Procedure of the (R)BHF calculation

- 1. Initial single-particle basis $\{|i\rangle\}$ trial for RBHF final solution $\{|a\rangle\}$
- $\bar{G}_{aba'b'}(W) = \bar{V}_{aba'b'} + \frac{1}{2} \sum_{i} \frac{\bar{V}_{abcd}Q(c,d)}{W \varepsilon_c \varepsilon_d} \bar{G}_{cda'b'}(W),$ 2. Bethe-Goldstone equation Solving with matrix inversion method M. Haftel and F. Tabakin, NPA 158, 1 (1970)
- $U_{ab} = \sum_{i=1}^{A} \langle ac|\bar{G}(W)|bc\rangle.$ Single-particle potential
- $\sum_{i} (T_{ij} + U_{ij}) D_{ji'} = \varepsilon_{i'} D_{ii'},$ 4. RHF iteration If converged $\{|i'\rangle\} = \{|i\rangle\} = \{|a\rangle\}$, RBHF iteration finishes.
- $\mbox{Basis transformation} \quad \bar{V}_{k'l'm'n'} = \sum D^*_{kk'} D^*_{ll'} D_{mm'} D_{nn'} \bar{V}_{klmn}.$ Go back to step 2.

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The no-core shell model



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The Monte-Carlo methods

The Monte-Carlo method



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The many-body perturbation theory

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Spontaneous symmetry breaking



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The coupled-cluster theory

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The in-medium similarity renormalization group

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Spontaneous symmetry breaking



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