

# Short range unitary two-body correlations

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Due to the short-ranged repulsive core in the nucleon-nucleon potential  $V$  the many-body state is depleted as a function of the relative distance  $x_{ij} = |\vec{x}_i - \vec{x}_j|$  for each pair  $(ij)$  when they are close to each other. These short range correlations cannot be described by shell model states. The most common procedure to remedy this problem is Brueckner's G-matrix method which replaces the bare  $V$  by an effective interaction  $G$ . The second method is the Jastrow approach where the correlated ground state of the nucleus is assumed to be of the form  $\prod_{i<j} f(x_{ij})|\Phi\rangle$  where  $f(x_{ij})$  is a correlation factor which diminishes the probability to find two nucleons at small distances  $x_{ij}$ .

We propose a third new method where the correlated state  $|\Psi\rangle$  is obtained by a unitary transformation  $e^{-i\mathbf{S}}$  (not to be confused with the expS method [1])

$$|\Psi\rangle = e^{-i\mathbf{S}}|\Phi\rangle,$$

where  $\mathbf{S}$  is a hermitean two-body operator which depends in general on the relative distance  $\vec{x}_{12}$ , the relative momentum  $\vec{p}_{12}$ , the spins and isospins of the two nucleons. The aim of  $e^{-i\mathbf{S}}$  is to push two nucleons away from each other whenever they get too close. The most simple ansatz which does that is

$$\mathbf{S} = \frac{1}{2} \sum_{i<j} \left( \vec{p}_{ij} \vec{V}_s(\mathbf{x}_{ij}) + \vec{V}_s(\mathbf{x}_{ij}) \vec{p}_{ij} \right).$$

$|\vec{V}_s|(x)$  is roughly speaking the distance which the particles are moved away from each other by  $e^{-i\mathbf{S}}$  if they are found at a distance  $x$  in  $|\Phi\rangle$ .  $|\vec{V}_s|(x)$  is largest if  $x$  lies inside the hard core and  $|\vec{V}_s|(x) \rightarrow 0$  if  $x$  is outside the repulsive interaction.

Fig. 1 displays the radial dependence of the correlated and uncorrelated deuteron wave function together with the Afnan Tang S3 potential [2]. The functional form of  $s(x)$  is adjusted such that the correlated state  $|\Psi_d\rangle$  for short distances equals the exact solution.

Once the correlator  $e^{-i\mathbf{S}}$  is adjusted to reproduce the two-body system at low energies (long wave length) we calculate

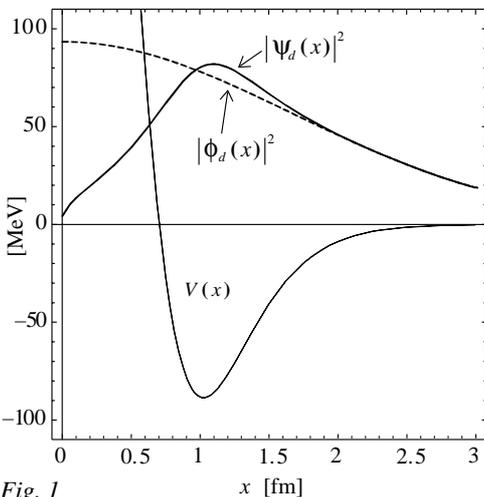


Fig. 1

the ground state energies of nuclei with many particles by minimizing

$$E = \langle \Psi | \mathbf{H} | \Psi \rangle = \langle \Phi | e^{i\mathbf{S}} \mathbf{H} e^{-i\mathbf{S}} | \Phi \rangle$$

with respect to  $|\Phi\rangle$  which is a single Slater determinant composed of localized Gaussians (FMD [3]). Unlike  $\mathbf{H} = \mathbf{T} + \mathbf{V}$

$$\begin{aligned} \mathbf{H}_{eff} &\equiv e^{i\mathbf{S}} \mathbf{H} e^{-i\mathbf{S}} = e^{i\mathbf{S}} \mathbf{T} e^{-i\mathbf{S}} + e^{i\mathbf{S}} \mathbf{V} e^{-i\mathbf{S}} \\ &= \mathbf{T} + \mathbf{T}_{eff}^{[2]} + \mathbf{V}_{eff}^{[2]} + \mathbf{T}_{eff}^{[3]} + \mathbf{V}_{eff}^{[3]} + \dots \end{aligned}$$

is not a one- plus two-body operator but contains three-body and higher interactions.

We calculate  $\mathbf{T}_{eff}^{[2]}$  and  $\mathbf{V}_{eff}^{[2]}$ , which are then functionals of  $s(x)$ , analytically and approximate the energy by neglecting three-body and higher terms in the cluster expansion. This turns out to be a very good approximation at typical nuclear densities. Estimations of the three-body terms give corrections less than 5% of binding energy for the  $\alpha$ -particle. Fig. 2 compares uncorrelated (left column) an correlated (right column) energies. The correlated potential energy  $\langle \Phi | \mathbf{V}_{eff}^{[2]} | \Phi \rangle$  (grey bars at negative values) is about twice the uncorrelated  $\langle \Phi | \mathbf{V} | \Phi \rangle$  in all nuclei. This gain in binding is counteracted by an increase in the kinetic energies (grey bars at positive values). Both together yield binding energies (black bars) which are within 8% deviation from results of Yakubovskii calculations [4] for  ${}^4\text{He}$  and FHNC calculations [5] for  ${}^{16}\text{O}$  and  ${}^{40}\text{Ca}$ . This supprises since  $\mathbf{H}_{eff}$  is the same for all nuclei and not density dependent. In addition one may easily conceive improved shell model states  $|\Phi\rangle$ .

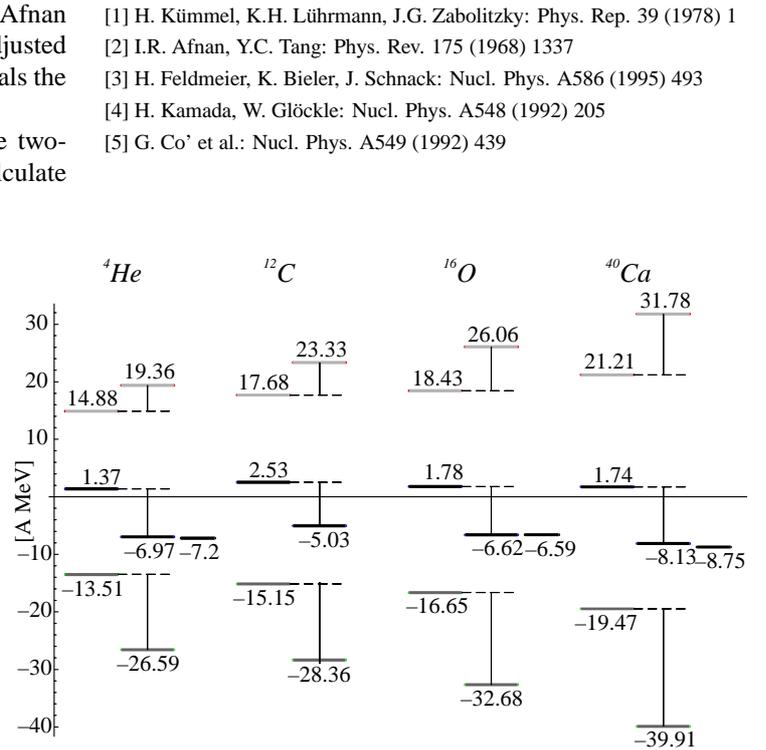


Fig. 2

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