One- and two-nucleon knock-out in neutrino-nucleus scattering: Nuclear mean-field approaches

Kajetan Niewczas







Kaietan Niewczas

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Nuclear response in the quasielastic and Δ regions



Outline

Lecture 1. the general framework of the nuclear mean-field model

- (1) Independent-particle model
- (2) Nucleon in a central potential
- (3) Mean-field nuclear potential

Lecture 2. *one- and two-nucleon knock-out in lepton-nucleus scattering*

Let's model a nucleus

Basic property of the nucleus-binding

$$M(Z,N) = ZM_p + NM_n - B$$

Nuclear packing fraction:

- $\rightarrow~$ for nuclei: 0.07 $< \rm NFP < 0.42$
- $\rightarrow~$ for hard spheres: ≈ 0.74
- $\rightarrow~$ for liquid argon: ≈ 0.032

Nucleus is like a dense quantum liquid





Fig. 3.1 Average binding energy B/A in Mev per nucleon for the naturally occurring nuclides (and Be^{*}), as a function of mass number A. Note the change of magnification in the A scale at A = 30. The Pauli four-shells in the lightest nuclei are evident. For $A \ge 16$, B/A is roughly constant; hence, to a first approximation, B is proportional to A.

R. Evans, The Atomic Nucleus (1955)

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Liquid-drop model

Bethe-Weiszäcker mass formula





Fig. 3.5 Summary of the semiempirical liquid-drop-model treatment of the average binding-energy curve from Fig. 3.1 of Chap. 9. Note how the decrease in surface energy and the increase in coulomb energy conspire to produce the maximum observed in B/A at $A \sim 60$. For these curves, the constants used in the semiempirical mass formula are given in the last line of Table 3.3.



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Independent-particle model

Basic assumptions

Elementary model of nuclear physics:

- nonrelativistic,
- nucleons are explicit degrees of freedom,
- described by the following Hamiltonian

$$\hat{H} = \sum_i^A \hat{T}_i + \sum_{i < j}^A \hat{V}_{ij} + \sum_{i < j < k}^A \hat{V}_{ijk} + ...$$

- two-body potential obtained from
 - \rightarrow phenomenology,
 - $\rightarrow~$ one-boson exchange models,
 - $\rightarrow~$ using χEFT ;

- three-body potential obtained from
 - \rightarrow phenomenology,
 - \rightarrow using χEFT ;

Nucleon-nucleon interaction

K. A. Olive et al. (Particle Data Group), Chin.Phys. C 38 (2014), 090001



Two-body potentials (e.g. Argonne v_{18}) use **angular momentum** and **isospin** operators of the form {1, L · S, $\sigma_1 \cdot \sigma_2$, S₁₂, L², (L · S)², L² $\sigma_1 \cdot \sigma_2$ }, {1, $\tau_1 \cdot \tau_2$ }

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Nucleon-nucleon interaction

Nuclear force:

- Short range
- $\circ \ \ {\rm Repulsive \ core}$
- Charge symmetry and independence

• Spin dependence

Isospin symmetry:

- $\circ~isospin~T=1/2$
 - \rightarrow neutron $T_z = -1/2$
 - \rightarrow proton $T_z = +1/2$
 - \rightarrow nucleus $T_z = 1/2(Z N)$

Approximately conserved in nuclei



Nucleon-nucleon interaction





FIG. 11. The deuteron S- and D-wave function components divided by r.

R. B. Wiringa, V. G. J. Stoks, and R. Schiavilla, Phys.Rev. C 51 (1995), 38–51

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Independent-particle model

What do we know so far:

- nuclei are made of nucleons,
- **binding per nucleon** is relatively small ($\simeq 7.5$ MeV for ¹²C),
- **distance between particles** larger than the nucleon radius ($\simeq 1 2$ fm),

Probability for a particle to propagate over a distance x with **no interactions** is

$$P(x) = \frac{1}{\lambda} \exp(-x/\lambda)$$

where $\lambda = (\rho \sigma)^{-1}$ is the mean free path, while ρ is target density and σ is interaction cross section

For nucleons inside nuclei:

$$\tilde{\lambda} \ll d < \lambda < R$$

where $\tilde{\lambda}$ is the de Broglie wavelength, d is the distance between targets, and R is the nuclear radius

 \rightarrow nucleus can be modeled as a system of independent, quasifree nucleons

Independent-particle model

General characteristics:

• **discrete energy levels** of a particle in a potential well

$$\mathsf{E}_{\mathfrak{i}}=\mathsf{T}_{\mathfrak{i}}-\mathsf{U}(\mathfrak{r}_{\mathfrak{i}})<\mathfrak{0},$$

 \rightarrow nuclear binding

$$B = \sum_{i}^{A} (T_i - U(r_i)),$$

 \rightarrow separation energy

$$E_s = T_{max} - U(r),$$

• Coulomb barrier for protons



Fermi gas model

Let's assume a gas of nucleons:

- nucleons are **fermions**,
 - \rightarrow wave functions are **antisymmetric**

 $\psi(...,x_a,...,x_b,...)=-\psi(...,x_b,...,x_a,...)$

- degeneracy pressure from Pauli principle,
- no interactions between nucleons,
- everything immersed in an infinite potential well

$$\frac{-\hbar^2}{2m}\nabla^2\psi(x,y,z)=E\psi(x,y,z)$$

 \rightarrow stationary Schrödinger equation



Infinitely deep potential well

The wave functions:

$$\Psi_{n_x,n_y,n_z}(x,y,z) = \psi_{n_x}(x)\psi_{n_y}(y)\psi_{n_z}(z) = \sin(k_x x)\sin(k_y y)\sin(k_z z)$$

The **energies**:

$$\mathsf{E}_{\mathsf{n}_x,\mathsf{n}_y,\mathsf{n}_z} = \frac{\hbar^2 p_x^2}{2\mathfrak{m}} + \frac{\hbar^2 p_y^2}{2\mathfrak{m}} + \frac{\hbar^2 p_z^2}{2\mathfrak{m}} = \frac{\hbar^2 \pi^2}{2\mathfrak{m}} \left(\frac{\mathfrak{n}_x^2}{\mathsf{L}_x^2} + \frac{\mathfrak{n}_y^2}{\mathsf{L}_y^2} + \frac{\mathfrak{n}_z^2}{\mathsf{L}_z^2} \right)$$

The number of states up to the Fermi momentum:

$$p_x^2 + p_y^2 + p_z^2 < p_F^2 \implies n_x^2 + n_y^2 + n_z^2 < \frac{p_F^2 L^2}{\pi^2 \hbar^2}$$

We calculate the number of occupied of states:

$$n = 2\frac{1}{8}\frac{4}{3}\pi \left(\frac{p_F L}{\pi \hbar}\right)^3 = \frac{1}{3}\pi \left(\frac{p_F}{\pi \hbar}\right)^3 V = \frac{1}{3}\pi \left(\frac{p_F}{\pi \hbar}\right)^3 \frac{4}{3}\pi r_A^3$$

(we took only 1/8 of the total sphere ($n_x > 0$, $n_y > 0$, $n_z > 0$), but with 2 spin states)

Finally, using $r_A = r_0 A^{1/3}$, we obtain the Fermi momenta for protons and neutrons:

$$p_{F} = \frac{\hbar}{r_{0}} \sqrt[3]{\frac{9\pi}{4}} \sqrt[3]{\frac{Z}{A}} \text{ and } p_{F} = \frac{\hbar}{r_{0}} \sqrt[3]{\frac{9\pi}{4}} \sqrt[3]{\frac{A-Z}{A}}$$

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(1)

(2)

(3)

(4)

(5)

- Fermi gas model... why not?
 - analytical model for efficient computations,
 - nucleon transition from below the Fermi sea to above

$$\begin{split} \theta(k_F - |\vec{k}|) &\rightarrow \theta(|\vec{k} + \vec{q}| - k_F) \\ |\vec{k}|^2 / 2M - V &\rightarrow (|\vec{k} + \vec{q}|^2 + M^2)^{1/2} \end{split}$$

- \rightarrow final nucleon is a plane wave;
- captures general features of quasielastic peak
 - \rightarrow **Fermi momentum** controls the spread,
 - $\label{eq:controls} \begin{array}{l} \rightarrow \mbox{ average interaction energy controls the shift;} \\ (\bar{\varepsilon} = \langle E \rangle \mbox{ ? } \bar{\varepsilon} \neq V) \end{array}$





...maybe better not



Artur Ankowski

R. Whitney et al., Phys.Rev. C 9 (1974), 2230

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Fermi gas model

- general assumptions are unclear
 - \rightarrow taken limits are inconsistent;
- fails to predict proper **energy levels**
 - $\rightarrow~$ unreliable for exclusive processes;
- lack of nucleon-nucleon interactions
 - \rightarrow overestimates the inclusive data;
- local Fermi gas is more robust
 - \rightarrow but makes even less sense;



J. Mougey, Nucl. Phys. A 335 (1980) 35

Problem 1.

Let's take the ^{12}C nucleus with $k_F=221$ MeV/c and $\bar{\varepsilon}=25$ MeV.

- (1) What are the general properties of this Fermi gas (E_F, V, E_s) ?
 - \rightarrow what is the average nucleon energy?
- (2) How does the spectral function of the Fermi gas model look like?
 - $\rightarrow~$ what is the energy-momentum relation?
- (3) How does the spectral function of a local Fermi gas looks like?
 - $\rightarrow~$ how can we parametrize k_F as a function of density $\rho(r)?$

Nuclear density and nucleon distribution for Carbon



http://discovery.phys.virginia.edu/research/groups/ncd/index.html

Let's consider a nucleon in a **central nuclear potential**

- $\rightarrow \ V = V(r) \ only$
- $\rightarrow~{\rm angular}~{\rm momentum}$ is conserved
 - harmonic oscillator

$$V_{\rm HO}(\mathbf{r}) = \frac{1}{2}\mathfrak{m}\omega^2\mathbf{r}^2 - V_1$$

• Woods-Saxon potential

$$V_{WS}(r) = -V_0 \frac{1}{1 + \exp((r - R)/a)}$$

with $V_0\simeq 50$ MeV and $\alpha\simeq 0.60$



The Schrödinger equation:

$$-\frac{\hbar^2}{2m}\nabla^2\psi - V(r)\psi = E\psi$$
(1)

 \rightarrow in carthesian coordinates:

$$\nabla^2 = \frac{\partial^2}{\partial x^2} + \frac{\partial^2}{\partial y^2} + \frac{\partial^2}{\partial z^2}$$

 \rightarrow in spherical coordinates:

$$\nabla^{2} = \frac{1}{r^{2}} \frac{\partial}{\partial r} \left(r^{2} \frac{\partial}{\partial r} \right) + \frac{1}{r^{2} \sin \theta} \frac{\partial}{\partial \theta} \left(\sin \theta \frac{\partial}{\partial \theta} \right) + \frac{1}{r^{2} \sin^{2} \theta} \left(\frac{\partial^{2}}{\partial^{2} \varphi} \right)$$

In spherical coordinates we separate variables:

$$\psi(\mathbf{r}, \theta, \phi) = \mathsf{R}(\mathbf{r})\mathsf{Y}(\theta, \phi)$$

We obtain two equations:

$$\frac{1}{R}\frac{d}{dr}\left(r^{2}\frac{dR}{dr}\right) - \frac{2mr^{2}}{\hbar^{2}}\left[V(r) - E\right] = l(l+1)$$

$$1\left(1 - \frac{\partial}{\partial r}\left(r + \frac{\partial}{\partial Y}\right) + \frac{1}{2} - \frac{\partial^{2}Y}{\partial r^{2}}\right) = l(l+1)$$
(5)

$$\frac{1}{Y}\left\{\frac{1}{\sin\theta}\frac{\partial}{\partial\theta}\left(\sin\theta\frac{\partial 1}{\partial\theta}\right) + \frac{1}{\sin^2\theta}\frac{\partial}{\partial\phi^2}\right\} = -l(l+1)$$
(6)

(2)

(3)

(4)

Let's consider the **angular part** and use $Y(\theta, \phi) = \Theta(\theta)\Phi(\phi)$:

$$\frac{1}{\Theta} \left[\sin \theta \frac{d}{d\theta} \left(\sin \theta \frac{d\Theta}{d\theta} \right) \right] + l(l+1) \sin^2 \theta = m^2$$

$$\frac{1}{\Phi} \frac{d^2 \Theta}{d\phi^2} = -m^2$$
(8)

First we solve the latter:

$$(\phi) = e^{im\phi}$$

 \rightarrow applying the condition $\Phi(\varphi) = \Phi(2\pi + \varphi)$ we must have $m = 0, \pm 1, \pm 2, ...$

Then, we solve the remaining:

$$\Theta(\theta) = AP_{l}^{m}(\cos\theta) \tag{10}$$

 \rightarrow where $P_l^m(\cos \theta)$ are the associate Legendre polynomials, and l = 0, 1, 2, ... for m = -l, -l + 1, ..., l - 1, 2

Φ

 $Y(\theta, \varphi) = \Theta(\theta) \Phi(\varphi)$ are the **spherical harmonics**:

$$Y(\theta, \phi) = (-1)^{m} \sqrt{\frac{(2l+1)}{4\pi} \frac{(l-m)!}{(l+m)!}} e^{im\phi} P_{l}^{m}(\cos\theta)$$
(11)

(9)

The spherical harmonics are the angular solution to **any central potential problem**

The shape of the potential V(r) only affects the radial **part** of the wave function

We have:

$$L^{2}Y_{lm}(\theta,\varphi) = l(l+1)\hbar^{2}Y_{lm}(\theta,\varphi)$$
(12)

$$L_z Y_{lm}(\theta, \phi) = m\hbar Y_{lm}(\theta, \phi)$$
(13)

Angular momentum is quantized:

- \rightarrow the allowed values of l are 0, 1, 2, ...
- \rightarrow sometimes we use the letters s, p, d, f, ...
- \rightarrow the allowed values of m are 0, ±1, ..., ±1
- $\rightarrow \$ the eigenvalues of L^2, L_z are $l(l+1)\hbar^2$ and $m\hbar$



Now, let us come back to the radial part:

$$\frac{1}{R}\frac{d}{dr}\left(r^{2}\frac{dR}{dr}\right) - \frac{2mr^{2}}{\hbar^{2}}\left[V(r) - E\right] = l(l+1)$$
(14)

We introduce R(r) = u(r)/r:

$$-\frac{\hbar}{2m}\frac{d^{2}u(r)}{dr^{2}} + \left(\frac{l(l+1)\hbar^{2}}{2mr^{2}} + V(r)\right)u(r) = Eu(r),$$
(15)

where

$$u(\infty) = 0, \quad u(0) = 0, \quad \int_0^\infty u^2(r) dr = 1$$
 (16)

E.g., for the harmonic oscillator of $U(r) = \frac{1}{2}m\omega^2 r^2$:

$$u_{k,l}(r) = \left(\frac{m\omega}{\hbar}\right)^{1/2+1/2} e^{-\frac{m\omega}{2\hbar}} r^{l+1} L_k^{l+1/2} (\frac{m\omega}{\hbar} r^2)$$
(17)

with energy levels

$$E_{k,l} = \hbar \omega (2k + l + 3/2) = \hbar \omega (N + 3/2)$$
(18)

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Harmonic oscillator energy spectrum is degenerate

Energy levels are quantized:

- \rightarrow major oscillator quantum number: N = 0, 1, 2, ...
- \rightarrow orbital quantum number: l = N, N 2, ..., 1, 0
- \rightarrow radial quantum number: k = (N l)/2

Magic numbers appear in the spectrum



Spin-orbit coupling

So far we worked with the following Hamiltonian:

$$H_0 = \sum_{i}^{A} (T_i + U(r_i)) = \sum_{i}^{A} h_0(i), \quad E_0 = \sum_{i}^{A} \varepsilon_{a_i}$$
(19)

Let's introduce a **spin-orbit term**:

$$= \mathbf{h}_0 + \zeta(\mathbf{r})\mathbf{l} \cdot \mathbf{s}$$

🕗 Nobel Prize in Physics 1963, E.P. Wigner, M. Goeppert Meyer, J.H.D. Jensen

So far, both parallel and antiparallel orientations have the same energies:

$$\langle \mathbf{nlj}, \mathbf{m} | \mathbf{h}_0 | \mathbf{nlj}, \mathbf{m} \rangle = \boldsymbol{\varepsilon}_{\mathbf{nlj}}^{(0)}, \quad \langle \mathbf{r}, \sigma | \mathbf{nlj}, \mathbf{m} \rangle = \frac{\mathbf{u}_{\mathbf{nl}}(\mathbf{r})}{\mathbf{r}} \left[\mathbf{Y}_1(\boldsymbol{\theta}, \boldsymbol{\varphi}) \otimes \boldsymbol{\chi}^{1/2}(\sigma) \right]_{\mathbf{m}}^{(j)}$$
(21)

We can express the spin-orbit term as $\zeta(r)\frac{1}{2}(j^2-l^2-s^2)$ and obtain $\varepsilon_{nlj} = \varepsilon_{nlj}^{(0)} + \Delta \varepsilon_{nlj}$ with

h

$$\Delta \epsilon_{\mathfrak{nlj}} = \langle \mathfrak{nlj}, \mathfrak{m} | \zeta(\mathfrak{r}) \mathbf{l} \cdot \mathbf{s} | \mathfrak{nlj}, \mathfrak{m} \rangle, \quad \Delta \epsilon_{\mathfrak{nlj}} = \frac{D}{2} \left[\mathfrak{j}(\mathfrak{j}+1) - \mathfrak{l}(\mathfrak{l}+1) - \frac{3}{4} \right]$$
(22)

Finally, defining $D = \int u_{nl}^2(r)\zeta(r)dr$ and $\zeta(r) = V_{ls}r_0^2 \frac{1}{r} \frac{\partial U(r)}{\partial r}$, we get:

$$\Delta \epsilon_{\mathfrak{nl} j=\mathfrak{l}+1/2} = (D/2) \cdot \mathfrak{l}, \quad \Delta \epsilon_{\mathfrak{nl} j=\mathfrak{l}-1/2} = -(D/2) \cdot (\mathfrak{l}+1)$$
(23)

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(20)

Spin-orbit coupling



K. Heyde, The Nuclear Shell Model (1990)

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Woods-Saxon potential





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Problem 2.

Let's take the nucleon in a spherical potential well:

$$\frac{\mathrm{d}R(r)}{\mathrm{d}r} + \frac{2}{r}\frac{\mathrm{d}R(r)}{\mathrm{d}r} + \left[\frac{2m}{\hbar}(E - V(r)) - \frac{l(l+1)}{r^2}\right]R(r) = 0$$

- (1) What are the radial solutions to this problem?
 - \rightarrow Handbook of Mathemathical Functions..., M. Abramowitz, I. A. Stegun, Eq. 10.1.1
- (2) What are the energy levels for nucleons?
 - \rightarrow what are the roots of the solution and their relation to $k = \hbar p$?
- (3) What is the average nucleon energy as confronted with a Fermi gas?
 - \rightarrow what is the depth of the potential using the same separation energy as before?

Mean-field nuclear potential

Mean-field nuclear picture



\rightarrow let's try to use a realistic nucleon-nucleon potential to derive the central nuclear potential

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Mean-field potential

Single-particle radial Schrödinger equation:

$$(T + U(r)) \varphi_{\alpha}(r) = \varepsilon_{\alpha} \varphi_{\alpha}(r)$$

 $[\mathbf{1}, (\mathbf{n})] = [\mathbf{1}, (\mathbf{n})]$

Nuclear Hamiltonian:

$$H_0 = \sum_{i=1}^A \left(T_i + U(r_i) \right) = \sum_{i=1}^A h_0(i), \quad E_0 = \sum_{i=1}^A \varepsilon_{\alpha_i}(r_i)$$

Nuclear wave function is a Slater determinant:

$$\Phi_{\alpha_1,...,\alpha_A}(\mathbf{r}_1,...,\mathbf{r}_A) = \frac{1}{\sqrt{A!}} \begin{vmatrix} \phi_{\alpha_1}(\mathbf{r}_1) & \dots & \phi_{\alpha_1}(\mathbf{r}_A) \\ \vdots & \ddots & \vdots \\ \phi_{\alpha_A}(\mathbf{r}_1) & \dots & \phi_{\alpha_A}(\mathbf{r}_A) \end{vmatrix}$$

Let's restrict ourselves to two-body interactions only and evaluate the mean-field potential:

$$H = \sum_{i=1}^{A} T_i + \frac{1}{2} \sum_{i,j=1}^{A} V_{ij}$$
(4)

$$H = \sum_{i=1}^{A} (T_i + U(r_i)) + \left(\frac{1}{2} \sum_{i,j=1}^{A} V_{ij} - \sum_{i=1}^{A} U(r_i)\right) = H_0 + H_{res} = \sum_{i=1}^{A} h_0(i) + H_{res},$$
(5)

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(1)

(2)

(3)

Hartree-Fock methods

Let's consider a density in terms of the **occupied single-particle states**:

$$\rho(\textbf{r}) = \sum_b \varphi_b^*(\textbf{r}) \varphi_b(\textbf{r})$$

The Hartree potential at a given point generated by the two-body interaction:

$$U_{H}(\textbf{r}) = \sum_{b} \int \varphi_{b}^{*}(\textbf{r}') V(\textbf{r},\textbf{r}') \varphi_{b}(\textbf{r}') d\textbf{r}'$$

The Schrödinger equation becomes:

$$-\frac{\hbar^{2}}{2m}\nabla^{2}\phi_{i}(\mathbf{r}) + \sum_{b}\int\phi_{b}^{*}(\mathbf{r}')V(\mathbf{r},\mathbf{r}')\phi_{b}(\mathbf{r}')d\mathbf{r}'\cdot\phi_{i}(\mathbf{r}) -\sum_{b}\int\phi_{b}^{*}(\mathbf{r}')V(\mathbf{r},\mathbf{r}')\phi_{b}(\mathbf{r})\phi_{i}(\mathbf{r}')d\mathbf{r}' = \varepsilon_{i}\phi_{i}(\mathbf{r}) -\frac{\hbar^{2}}{2m}\nabla^{2}\phi_{i}(\mathbf{r}) + U_{H}(\mathbf{r})\phi_{i}(\mathbf{r}) - \int U_{F}(\mathbf{r},\mathbf{r}')\phi_{i}(\mathbf{r}')d\mathbf{r}' = \varepsilon_{i}\phi_{i}(\mathbf{r})$$
(9)

where the exchange term is driven by the Fock potential:

$$U_{\mathsf{F}}(\mathbf{r}) = \sum_{\mathsf{b}} \phi_{\mathsf{b}}^*(\mathbf{r}') V(\mathbf{r}, \mathbf{r}') \phi_{\mathsf{b}}(\mathbf{r})$$
(10)

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(6)

(7)

The iterative Hartree-Fock method

- $\rightarrow\,$ start with an initial guess for the average field or the wave functions
- $\rightarrow \mbox{ using the nucleon-nucleon potential } V(r,r')$ solve the equation

$$-\frac{\hbar^2}{2m}\nabla^2\varphi_{\mathfrak{i}}(\textbf{r})+U_{H}(\textbf{r})\varphi_{\mathfrak{i}}(\textbf{r})-\int U_{F}(\textbf{r},\textbf{r}')\varphi_{\mathfrak{i}}(\textbf{r}')d\textbf{r}'=\varepsilon_{\mathfrak{i}}\varphi_{\mathfrak{i}}(\textbf{r})$$

 $\rightarrow \,$ determine new values of $U_{H}(\textbf{r}),\,U_{F}(\textbf{r},\textbf{r}'),\,\varphi_{i}(\textbf{r}),\,\varepsilon_{i}$

 $\rightarrow\,$ at convergence: the final field $U_H(\textbf{r}),$ wave function $\varphi_i(\textbf{r}),$ and single-particle energy ε_i

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Nucleons in the mean-field potential



W. H. Dickhoff, D. Van Neck, Many-body Theory Exposed! (2005)

 \rightarrow nucleon lines are dressed according to the Hartree-Fock procedure

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Charge densities from the mean-field framework



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Charge densities from the mean-field framework



Fig. 3.18. The nuclear density distribution for the least bound proton in ²⁰⁶Pb. The shell-model predicts the last $(3s_{1/2})$ proton in ²⁰⁶Pb to have a sharp maximum at the centre, as shown at the left-hand side. On the right-hand side the nuclear charge density difference $\rho_c (^{206}\text{Pb}) - \rho_c (^{205}\text{Tl}) = \varphi_{3s_{1/2}}^2(r)$ is given [taken from (Frois 1983) and Doe 1983)]

K. Heyde, The Nuclear Shell Model (1990)

Relativistic mean-field

All of this can be also done in a **relativistic** framework:

- $\circ~$ Schrödinger equation \rightarrow Dirac equation,
- Wave functions \rightarrow **Dirac spinors**,
- Spin-orbit term comes for free!

$$\begin{split} \tilde{(E}\gamma_0 - \vec{p}\cdot\vec{\gamma} - \tilde{M}) \psi &= 0 \\ \tilde{E} &= E - V(r) \\ \tilde{M} &= M - S(r) \end{split}$$

