

The Nuclear Shell Model

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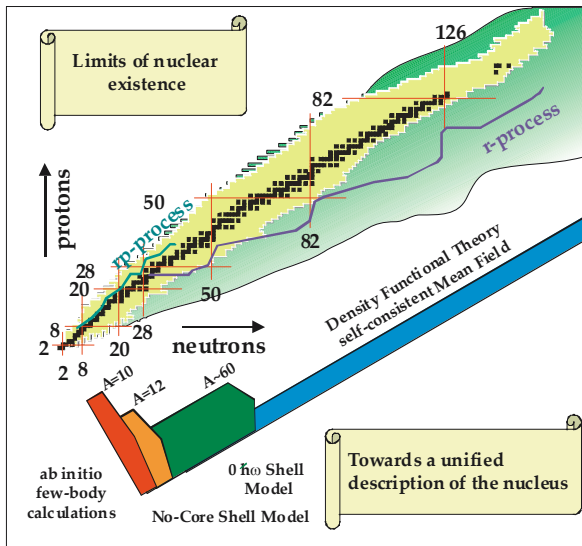
Lecture of A. P. Zuker "Beyond the Shell Model" on Friday afternoon

Structure of complex nuclei

- Nuclear charge (matter) density distribution $\rho_{ch}(\vec{r})$ ($\rho_m(\vec{r})$) with sharp radius $R \simeq r_0 A^{1/3}$.
- **Empirical evidence on the existence of an average potential and the corresponding shell structure**
(From masses, nucleon separation energies, low-energy spectra, etc.)
- **Independent-particle motion near Fermi-level.**
(From nucleon transfer reactions: nuclear mean-field with strong spin-orbit splitting and large shell gaps)
- **Pairing (superfluid behavior) at low excitation energy.**
(S_n versus S_{2n} ; two-nucleon spectra, comparison of spectra of even-even and even-odd, odd-even nuclei, etc.)
- **Low-lying multipole (quadrupole) modes, vibrational or rotational energy structures.**
(Coulomb excitation; scattering of charge particles; heavy-ion fusion-evaporation reactions, etc)

The aim of the microscopic theory is to describe these motions starting from a NN force.

Nuclear many-body problem



Nuclear many-body problem

Non-relativistic Hamiltonian for A nucleons

$$\hat{H} = \sum_{i=1}^A \frac{\vec{p}_i^2}{2m} + \sum_{i<j=1}^A W(\vec{r}_i - \vec{r}_j)$$
$$\hat{H} = \underbrace{\sum_{i=1}^A \left[\frac{\vec{p}_i^2}{2m} + U(\vec{r}_i) \right]}_{\hat{H}^{(0)}} + \underbrace{\sum_{i<j=1}^A W(\vec{r}_i - \vec{r}_j) - \sum_{i=1}^A U(\vec{r}_i)}_{\hat{V}},$$

Mean-field theories

Search for the most optimum mean-field potential starting from a given two-body interaction + correlations

Shell model

Schematic average potential + residual interaction

Variational approach to the nuclear many-body problem

Starting point: antisymmetric product wave function

$$\Psi(1, 2, \dots, A) = \frac{1}{\sqrt{A!}} \begin{vmatrix} \phi_{\alpha_1}(1) & \phi_{\alpha_1}(2) & \dots & \phi_{\alpha_1}(A) \\ \phi_{\alpha_2}(1) & \phi_{\alpha_2}(2) & \dots & \phi_{\alpha_2}(A) \\ \vdots & \vdots & \ddots & \vdots \\ \phi_{\alpha_A}(1) & \phi_{\alpha_A}(2) & \dots & \phi_{\alpha_A}(A) \end{vmatrix}$$

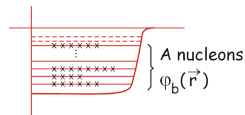
The best wave function is determined via a variational principle:

$$\delta \langle \Psi | H | \Psi \rangle = \langle \delta \Psi | H | \Psi \rangle = 0 \quad \text{with} \quad \int |\phi_{\alpha_i}(\vec{r})|^2 d\vec{r} = 1.$$

Self-consistent mean-field potential

Hartree-Fock equations

$$\left(-\frac{\hbar^2}{2m} \Delta + U_H(\vec{r}) \right) \phi_i(\vec{r}) + \int U_F(\vec{r}, \vec{r}') \phi_i(\vec{r}') d\vec{r}' = \varepsilon_i \phi_i(\vec{r})$$



Direct (Hartree) term

$$U_H(\vec{r}) = \sum_{b \in F} \int \phi_b^*(\vec{r}') W(\vec{r}, \vec{r}') \phi_b(\vec{r}') d\vec{r}'$$

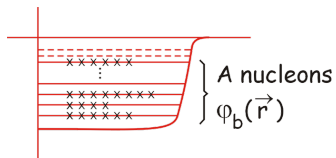
Exchange (Fock) term

$$U_F(\vec{r}, \vec{r}') = \sum_{b \in F} \phi_b^*(\vec{r}) W(\vec{r}, \vec{r}') \phi_b(\vec{r}')$$

Iterative solution of Hartree-Fock equations:

$$\begin{array}{ccccccc}
 \phi_i^{(0)}(\vec{r}) & & \phi_i^{(1)}(\vec{r}), \varepsilon_i^{(1)} & & \dots & & \phi_i^{(HF)}(\vec{r}), \varepsilon_i^{(HF)} \\
 \downarrow & & \downarrow & & \downarrow & & \\
 U_H^{(0)}(\vec{r}), U_F^{(0)}(\vec{r}, \vec{r}') & \nearrow & U_H^{(1)}(\vec{r}), U_F^{(1)}(\vec{r}, \vec{r}') & \nearrow & \dots & \nearrow & U^{(HF)}(\vec{r})
 \end{array}$$

Self-consistent mean-field potential



Direct (Hartree) term

$$U_H(\vec{r}) = \sum_{b \in F} \int \phi_b^*(\vec{r}') W(\vec{r}, \vec{r}') \phi_b(\vec{r}') d\vec{r}'$$

$$U_H(\vec{r}) = \int \rho(\vec{r}') W(\vec{r}, \vec{r}') d\vec{r}'$$

$$\rho(\vec{r}) = \sum_{b \in F} |\phi_b(\vec{r})|^2$$

Example: a delta-force

$$W(\vec{r}, \vec{r}') \propto \delta(\vec{r} - \vec{r}')$$

$$U_H(\vec{r}) \propto \rho(\vec{r})$$

Hartree-Fock ground state properties

From the product HF wave function, a number of ground-state properties can be calculated:

$$\Psi_{HF}(1, 2, \dots, A) = \frac{1}{\sqrt{A!}} \mathcal{A} \prod_{i=1}^A \phi_{\alpha_i}^{HF}(i)$$

$$\langle \Psi_{HF} | \hat{H} | \Psi_{HF} \rangle = E_0$$

$$\langle \Psi_{HF} | \sum_{i=1}^A \hat{r}_i^2 | \Psi_{HF} \rangle = \langle r^2 \rangle$$

$$\langle \Psi_{HF} | \sum_{i=1}^A \hat{\rho}(\vec{r}_i) | \Psi_{HF} \rangle = \rho(\vec{r})$$

Application to open-shell nuclei and calculation excitation spectra require inclusion of correlations (pairing, etc) \rightarrow beyond mean-field techniques (lecture of Ph. Quentin).

Shell model: energy matrix diagonalization

Non-relativistic Hamiltonian for A nucleons

$$\hat{H} = \sum_{i=1}^A \underbrace{\left[\frac{\vec{p}_i^2}{2m} + U(\vec{r}_i) \right]}_{\hat{h}_i} + \underbrace{\sum_{i<j=1}^A W(\vec{r}_i - \vec{r}_j) - \sum_{i=1}^A U(\vec{r}_i)}_{\hat{V}}$$

Construction of a basis from single-particle states

$$\hat{h}\phi_\alpha(\vec{r}) = \varepsilon_\alpha\phi_\alpha(\vec{r}) \quad \rightarrow \quad \{\varepsilon_\alpha, \phi_\alpha(\vec{r})\}$$

Spherical potential $U(\vec{r}) = U(r)$: $\alpha = \{n_\alpha, l_\alpha, j_\alpha, m_\alpha\}$

$$\phi_{nljm}(\vec{r}) = \frac{R_{nlj}(r)}{r} \underbrace{\left[Y_l(\theta, \varphi) \times \chi_{\frac{1}{2}} \right]}_{\sum_{m_l m_s} (l m_l \frac{1}{2} m_s | jm) Y_{l m_l}(\theta, \varphi) \chi_{\frac{1}{2} m_s}^{(j)}}$$

Single-particle wave functions

Radial differential equation

$$-\frac{\hbar^2}{2m}R''(r) + \frac{\hbar^2}{2m}\frac{l(l+1)}{r^2}R(r) + [U(r)R(r) + a_{ls}f_{ls}(r)]R(r) = \varepsilon R(r)$$

Normalization condition

$$\int |\phi_{nlsjm}(\vec{r})|^2 d\vec{r} = \sum_{m_l m_s m'_l m'_s} (l m_l \frac{1}{2} m_s | j m) (l m'_l \frac{1}{2} m'_s | j m) \int Y_{l m_l}(\theta, \phi) Y_{l m'_l}(\theta, \phi) d\Omega \\ \times \langle \chi_{\frac{1}{2} m_s} | \chi_{\frac{1}{2} m'_s} \rangle \int_0^\infty |R_{nl}(r)|^2 dr = \int_0^\infty |R_{nl}(r)|^2 dr = 1$$

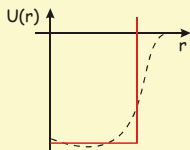
Parity

$$\hat{P}\phi_{nlsjm}(\vec{r}) = \hat{P}\phi_{nlsjm}(-\vec{r}) = (-1)^l \phi_{nlsjm}(\vec{r})$$

Examples of spherically-symmetric potentials

- Square-well potential + strong spin-orbit term

$$\rightarrow J_{l+1/2}(kr)$$



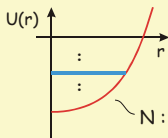
$$\underline{s} \quad l=0 \quad (100)$$

$$\underline{d} \quad l=2 \quad (93)$$

ex. $N = 4$

$$\underline{g} \quad l=4 \quad (75.5)$$

- Harmonic oscillator potential + orbital + spin-orbit term



$$\rightarrow (vr)^l e^{-\frac{v^2 r^2}{2}} L_{n-1}^{l+1/2}(v^2 r^2)$$

$$(v = \sqrt{\frac{m\omega}{\hbar}})$$

N : degenerate in n, l

Harmonic oscillator potential

$$U(r) = \frac{m\omega^2 r^2}{2} + \alpha \vec{l} \cdot \vec{l} + \beta \vec{l} \cdot \vec{s}$$

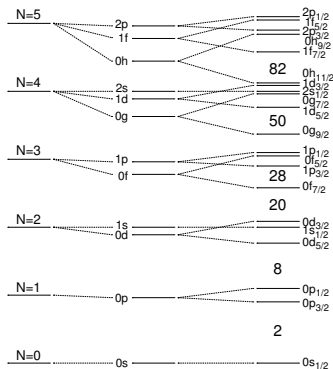
$$\varepsilon_N = \hbar\omega \left(2n + l + \frac{3}{2} \right) = \hbar\omega \left(N + \frac{3}{2} \right),$$

$$N = 0, 1, 2, \dots,$$

$$l = N, N - 2, \dots, 1 \text{ or } 0$$

$$n = (N - l)/2.$$

Harmonic oscillator potential possesses many **symmetry properties** which make it a preferable choice as a basis.



M. Mayer (1949)

O. Haxel, H. Jensen, H.E. Suess (1949)

Isospin

The idea of W.Heisenberg: proton and neutron are considered as two states of a nucleon.

$$V_{\pi\pi} \approx V_{\nu\nu} \approx V_{\pi\nu} \qquad \pi = \begin{pmatrix} 0 \\ 1 \end{pmatrix}, \nu = \begin{pmatrix} 1 \\ 0 \end{pmatrix}$$

Isospin operators (in analogy with the Pauli matrices):

$$\vec{t} = \frac{1}{2}\vec{\tau}, \quad \tau_x = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}, \quad \tau_y = \begin{pmatrix} 0 & -i \\ i & 0 \end{pmatrix}, \quad \tau_z = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}$$

Single-particle wave functions

$$\begin{aligned} \phi_\nu(r) &= \phi(\vec{r}) \begin{pmatrix} 1 \\ 0 \end{pmatrix} = \phi(\vec{r}) \theta_{t=1/2, m_t=1/2}, \\ \phi_\pi(r) &= \phi(\vec{r}) \begin{pmatrix} 0 \\ 1 \end{pmatrix} = \phi(\vec{r}) \theta_{t=1/2, m_t=-1/2}. \end{aligned}$$

Isospin and classification of nuclear states

$$\hat{T} = \sum_{i=1}^A \hat{t}_i, \quad \hat{T}_z = \sum_{i=1}^A \hat{t}_{zi}.$$

Charge independent Hamiltonian: $[\hat{H}, \hat{T}] = 0$.

$$M_T = \frac{1}{2}(N - Z), \quad \frac{1}{2}(N - Z) \leq T \leq \frac{A}{2}.$$

Realistic situation

$$m_p \approx m_n; \quad \hat{V}_{Coulomb} = \sum_{i < j = 1}^Z \frac{e^2}{|\vec{r}_i - \vec{r}_j|}$$

$$\hat{V}_{Coulomb} = \hat{V}^{(T=0)} + \hat{V}^{(T=1)} + \hat{V}^{(T=2)}.$$

$$E(T, M_T) = a(T) + b(T)M_T + c(T)M_T^2,$$

Two-particle wave function for identical fermions

Angular-momentum coupled state ($j_a \neq j_b$)

$$\Phi_{j_a(1)j_b(2);JM}(\vec{r}_1, \vec{r}_2) = \sum_{m_a m_b} (j_a m_a j_b m_b | J M) \phi_{j_a m_a}(\vec{r}_1) \phi_{j_b m_b}(\vec{r}_2) = [\phi_{j_a}(\vec{r}_1) \times \phi_{j_b}(\vec{r}_2)]_M^{(J)}$$

$j_a \equiv (n_a l_a j_a)$, $J = |j_a - j_b|, |j_a - j_b| + 1, \dots, j_a + j_b$, while $M = -J, -J + 1, \dots, J - 1, J$.

Not antisymmetric with respect to permutation of two identical fermions !

Normalized and antisymmetric state

$$\Phi_{j_a j_b; JM}(\vec{r}_1, \vec{r}_2) = \frac{1}{\mathcal{N}} \sum_{m_a m_b} (j_a m_a j_b m_b | J M) [\phi_{j_a m_a}(\vec{r}_1) \phi_{j_b m_b}(\vec{r}_2) - \phi_{j_b m_b}(\vec{r}_1) \phi_{j_a m_a}(\vec{r}_2)]$$

Since the Clebsch-Gordan coefficients have the following property:

$$(j_a m_a j_b m_b | J M) = (-1)^{j_a + j_b - J} (j_b m_b j_a m_a | J M)$$

$$\Phi_{j_a j_b; JM}(\vec{r}_1, \vec{r}_2) = \frac{1}{\mathcal{N}} \sum_{m_a m_b} [(j_a m_a j_b m_b | J M) \phi_{j_a m_a}(\vec{r}_1) \phi_{j_b m_b}(\vec{r}_2) - (-1)^{j_a + j_b - J} (j_b m_b j_a m_a | J M) \phi_{j_b m_b}(\vec{r}_1) \phi_{j_a m_a}(\vec{r}_2)]$$

Two-particle wave function for identical fermions

Normalized and antisymmetric state

$$j_a \neq j_b$$

$$\Phi_{j_a j_b; JM}(\vec{r}_1, \vec{r}_2) = \frac{1}{\sqrt{2}} \left\{ [\phi_{j_a}(\vec{r}_1) \times \phi_{j_b}(\vec{r}_2)]_M^{(J)} - (-1)^{j_a + j_b + J} [\phi_{j_b}(\vec{r}_1) \times \phi_{j_a}(\vec{r}_2)]_M^{(J)} \right\}$$

$$j_a \equiv (n_a l_a j_a), J = |j_a - j_b|, |j_a - j_b| + 1, \dots, j_a + j_b, \text{ while } M = -J, -J + 1, \dots, J - 1, J.$$

$$j_a = j_b = j$$

$$\Phi_{j^2; JM}(\vec{r}_1, \vec{r}_2) = \frac{1 + (-1)^J}{2} [\phi_j(\vec{r}_1) \times \phi_j(\vec{r}_2)]_M^{(J)}$$

Important consequence: if $j_a = j_b = j$, then $J = 0, 2, 4, \dots, 2j - 1$

$$(\nu 0 d_{5/2})^2 : J = 0, 2, 4$$

Two-particle wave function for protons and neutrons

$$j_a \neq j_b$$

$$\Phi_{j_a j_b; JM T M_T}(\vec{r}_1, \vec{r}_2) = \left\{ [\phi_{j_a}(\vec{r}_1) \times \phi_{j_b}(\vec{r}_2)]_M^{(J)} + (-1)^{j_a + j_b + J + T} [\phi_{j_b}(\vec{r}_1) \times \phi_{j_a}(\vec{r}_2)]_M^{(J)} \right\} \frac{\Theta_{TM T}}{\sqrt{2}}$$

$$\Theta_{1,1} = \theta_{1/2,1/2}(1)\theta_{1/2,1/2}(2),$$

$$\Theta_{1,-1} = \theta_{1/2,-1/2}(1)\theta_{1/2,-1/2}(2),$$

$$\Theta_{1,0} = [\theta_{1/2,1/2}(1)\theta_{1/2,-1/2}(2) + \theta_{1/2,-1/2}(1)\theta_{1/2,1/2}(2)] / \sqrt{2},$$

$$\Theta_{0,0} = [\theta_{1/2,1/2}(1)\theta_{1/2,-1/2}(2) - \theta_{1/2,-1/2}(1)\theta_{1/2,1/2}(2)] / \sqrt{2},$$

$$j_a = j_b = j$$

$$\Phi_{j^2; JM T M_T}(\vec{r}_1, \vec{r}_2) = \frac{1 - (-1)^{J+T}}{2} [\phi_j(\vec{r}_1) \times \phi_j(\vec{r}_2)]_M^{(J)} \Theta_{TM T}$$

Remark: if $j_a = j_b = j$, then $(J + T)$ is odd!

$$(\nu 0 d_{5/2})^2 : J = 0, 2, 4$$

$$(\nu 0 d_{5/2} \pi 0 d_{5/2}) : J = 0, 2, 4 (T = 1); J = 1, 3, 5 (T = 0)$$

Many-particle wave function: $J(T)$ -coupled states.

J -coupled state

- Consider N identical fermions in a single- j shell. We construct a totally antisymmetric and coupled to good J N -nucleon wave function from a set of totally antisymmetric $(N - 1)$ -nucleon wave functions coupled to all possible J' :

$$\Phi_{\chi JM}^{j(N)}(\vec{r}_1, \dots, \vec{r}_N) = \sum_{\chi' J'} [j^{N-1}(\chi' J') j | \} j^N \chi J] \Phi_{\chi' J' M'}^{j(N-1)}(\vec{r}_1, \dots, \vec{r}_{N-1}) \phi_{jm}(\vec{r}_N),$$

where

$$[j^{N-1}(\chi' J') j | \} j^N \chi J]$$

are one-particle *coefficients of fractional parentage (cfp's)*

- Repeat this procedure for N' particles in j' orbital and so on. Construct thus basis states by consecutive coupling of angular momenta and antisymmetrization.

Many-particle wave function: m -scheme basis

Slater determinants

$$\Phi_{\alpha}(1, 2, \dots, A) = \frac{1}{\sqrt{A!}} \begin{vmatrix} \phi_{\alpha_1}(\vec{r}_1) & \phi_{\alpha_1}(\vec{r}_2) & \dots & \phi_{\alpha_1}(\vec{r}_A) \\ \phi_{\alpha_2}(\vec{r}_1) & \phi_{\alpha_2}(\vec{r}_2) & \dots & \phi_{\alpha_2}(\vec{r}_A) \\ \vdots & \vdots & \ddots & \vdots \\ \phi_{\alpha_A}(\vec{r}_1) & \phi_{\alpha_A}(\vec{r}_2) & \dots & \phi_{\alpha_A}(\vec{r}_A) \end{vmatrix}$$

where $\alpha_i = (n_i, l_i, j_i, m_i)$ and α stores a set of single-particle configurations $\{\alpha_1, \alpha_2, \dots, \alpha_A\}$

$$M = \sum_{i=1}^A m_i$$

Projection on J can be performed.

Exercise 1

Basis construction

Consider 2 neutrons in $0f_{7/2}$ orbital.

- Write down the basis in J -coupled form.
- Write down the basis in m -scheme.
- How many different J -states exist in this model space?

Solution to exercise 1

Basis construction

- Basis in J -coupled scheme:

$$|(0f_{7/2})^2; J, T = 1\rangle \quad J = 0, 2, 4, 6$$

Solution to exercise 1 (continued)

Basis construction

- Basis in m -scheme: $|m_1 m_2; M\rangle$:

$$\begin{array}{l} | \frac{7}{2}, \frac{5}{2}; 6 \rangle \\ | \frac{7}{2}, \frac{3}{2}; 5 \rangle \\ | \frac{7}{2}, \frac{1}{2}; 4 \rangle \quad | \frac{5}{2}, \frac{3}{2}; 4 \rangle \\ | \frac{7}{2}, -\frac{1}{2}; 3 \rangle \quad | \frac{5}{2}, \frac{1}{2}; 3 \rangle \\ | \frac{7}{2}, -\frac{3}{2}; 2 \rangle \quad | \frac{5}{2}, -\frac{1}{2}; 2 \rangle \quad | \frac{3}{2}, \frac{1}{2}; 2 \rangle \\ | \frac{7}{2}, -\frac{5}{2}; 1 \rangle \quad | \frac{5}{2}, -\frac{3}{2}; 1 \rangle \quad | \frac{3}{2}, -\frac{1}{2}; 1 \rangle \\ | \frac{7}{2}, -\frac{7}{2}; 0 \rangle \quad | \frac{5}{2}, -\frac{5}{2}; 0 \rangle \quad | \frac{3}{2}, -\frac{3}{2}; 0 \rangle \quad | \frac{1}{2}, -\frac{1}{2}; 0 \rangle \\ | -\frac{7}{2}, \frac{5}{2}; -1 \rangle \quad | -\frac{5}{2}, \frac{3}{2}; -1 \rangle \quad | -\frac{3}{2}, \frac{1}{2}; -1 \rangle \\ | -\frac{7}{2}, \frac{3}{2}; -2 \rangle \quad | -\frac{5}{2}, \frac{1}{2}; -2 \rangle \quad | -\frac{3}{2}, -\frac{1}{2}; -2 \rangle \\ | -\frac{7}{2}, \frac{1}{2}; -3 \rangle \quad | -\frac{5}{2}, \frac{1}{2}; -3 \rangle \\ | -\frac{7}{2}, \frac{1}{2}; -4 \rangle \quad | -\frac{5}{2}, \frac{3}{2}; -4 \rangle \\ | -\frac{7}{2}, \frac{3}{2}; -5 \rangle \\ | -\frac{7}{2}, \frac{5}{2}; -6 \rangle \end{array}$$

- There are 4 different J -states.

Solution of a many-body Schrödinger equation

Construct a basis in the valence space for each J

$$\Phi_k^{(J)} = \left\{ (j_a)_{J_a}^{n_a} (j_b)_{J_b}^{n_b} \dots \right\}_k^{(J)}, \quad \hat{H}^{(0)} \Phi_k = E_k^{(0)} \Phi_k$$

Expand unknown wave function in terms of basis functions

$$\Psi_p = \sum_{k=1}^n a_{kp} \Phi_k \quad \Rightarrow \quad \hat{H} \Psi_p = E_p \Psi_p \quad (\hat{H} = \hat{H}^{(0)} + \hat{V})$$

Multiplying by $\langle \Phi_l |$, we get a system of equations

$$\sum_{k=1}^n H_{lk} a_{kp} = E_p a_{lp}$$

\Rightarrow diagonalization of the matrix

$$H_{lk} = \langle \Phi_l | \hat{H} | \Phi_k \rangle = E_k^{(0)} \delta_{lk} + V_{lk}$$

Calculate Hamiltonian matrix $H_{ij} = \langle \Phi_j | \hat{H} | \Phi_i \rangle$

— Diagonalize to obtain eigenvalues

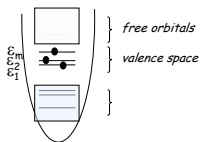
$$\begin{pmatrix} H_{11} & H_{12} & \dots & H_{1N} \\ H_{21} & H_{22} & & \\ \vdots & & \ddots & \\ H_{N1} & & \dots & H_{NN} \end{pmatrix} \rightarrow \begin{array}{c} \text{=====} \\ \text{=====} \\ \text{=====} \\ \text{=====} \end{array}$$

Basis dimension and choice of the model space

Basis dimension grows quickly

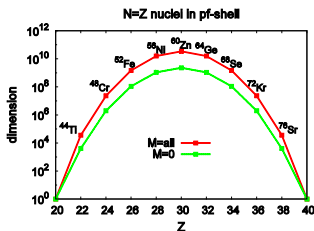
$$\dim \approx \begin{pmatrix} \Omega_{\pi} \\ N_{\pi} \end{pmatrix} \begin{pmatrix} \Omega_{\nu} \\ N_{\nu} \end{pmatrix}$$

Model space: a few valence orbitals beyond the closed-shell core.



Example: $^{60}\text{Zn}_{30}$ in *pf*-shell

$$\begin{aligned} \dim(^{60}\text{Zn}) &= \begin{pmatrix} 20 \\ 10 \end{pmatrix} \begin{pmatrix} 20 \\ 10 \end{pmatrix} \\ &= \frac{20!}{10!10!} \frac{20!}{10!10!} \approx 3.4 \times 10^{10}. \end{aligned}$$



Practical shell-model for ^{18}O in sd -shell

$$\hat{H} = \hat{H}^{(0)} + \hat{V} = \hat{h}(1) + \hat{h}(2) + \hat{V}$$

Single-particle energies:

$$\varepsilon(0d_{5/2}) = E_B(^{17}\text{O}_9) - E_B(^{16}\text{O}_8) = -4.143 \text{ MeV}$$

$$\varepsilon(1s_{1/2}) = \varepsilon(0d_{5/2}) + E_{\text{ex}}(^{17}\text{O}; 1/2_1^+) = -3.273 \text{ MeV}$$

$$\varepsilon(0d_{3/2}) = \varepsilon(0d_{5/2}) + E_{\text{ex}}(^{17}\text{O}; 3/2_1^+) = 0.942 \text{ MeV}$$

Basis of states for each (JT) denoted as $|j_a j_b\rangle_{JT}$:

$$0^+, T = 1 : \begin{aligned} |\Phi_1(0^+)\rangle &\equiv |d_{5/2}^2\rangle_{01} \\ |\Phi_2(0^+)\rangle &\equiv |s_{1/2}^2\rangle_{01} \\ |\Phi_3(0^+)\rangle &\equiv |d_{3/2}^2\rangle_{01} \end{aligned} \Rightarrow \begin{pmatrix} H_{11} & H_{12} & H_{13} \\ H_{21} & H_{22} & H_{23} \\ H_{31} & H_{32} & H_{33} \end{pmatrix}$$

$$1^+, T = 1 : \begin{aligned} |\Phi_1(1^+)\rangle &\equiv |d_{5/2} d_{3/2}\rangle_{11} \\ |\Phi_2(1^+)\rangle &\equiv |s_{1/2} d_{3/2}\rangle_{11} \end{aligned} \Rightarrow \begin{pmatrix} H_{11} & H_{12} \\ H_{21} & H_{22} \end{pmatrix}$$

and so on for $J = 2, 3, 4$.

Exercise 2: Energies of 0^+ states in ^{18}O

Diagonal Two-body matrix elements (TBMEs)

$$H_{11} = 2\varepsilon(d_{5/2}) + \underbrace{\langle d_{5/2}^2 | V | d_{5/2}^2 \rangle_{01}}_{-2.82 \text{ MeV}}$$

$$H_{22} = 2\varepsilon(s_{1/2}) + \underbrace{\langle s_{1/2}^2 | V | s_{1/2}^2 \rangle_{01}}_{-2.12 \text{ MeV}}$$

$$H_{33} = 2\varepsilon(d_{3/2}) + \underbrace{\langle d_{3/2}^2 | V | d_{3/2}^2 \rangle_{01}}_{-2.18 \text{ MeV}}$$

Non-diagonal TBMEs

$$H_{12} = H_{21} = \underbrace{\langle d_{5/2}^2 | V | s_{1/2}^2 \rangle_{01}}_{-1.32 \text{ MeV}}$$

$$H_{23} = H_{32} = \underbrace{\langle s_{1/2}^2 | V | d_{3/2}^2 \rangle_{01}}_{-1.08 \text{ MeV}}$$

$$H_{13} = H_{31} = \underbrace{\langle d_{5/2}^2 | V | d_{3/2}^2 \rangle_{01}}_{-3.19 \text{ MeV}}$$

Practical shell-model for ^{18}O in sd -shell

Eigenvalues (g.s. and two excited 0^+ states):

$$\begin{array}{ll} E(0_1^+) = -12.602 \text{ MeV} & \Rightarrow E_{gs}(0_1^+) = 0 \text{ MeV} \\ E(0_2^+) = -8.097 \text{ MeV} & E_{ex}(0_2^+) = 4.505 \text{ MeV} \\ E(0_3^+) = 0.622 \text{ MeV} & E_{ex}(0_3^+) = 13.224 \text{ MeV} \end{array}$$

Eigenstates:

$$|\Psi(0_1^+)\rangle = a_{11}|d_{5/2}^2\rangle_{01} + a_{21}|s_{1/2}^2\rangle_{01} + a_{31}|d_{3/2}^2\rangle_{01}$$

$$|\Psi(0_2^+)\rangle = a_{12}|d_{5/2}^2\rangle_{01} + a_{22}|s_{1/2}^2\rangle_{01} + a_{32}|d_{3/2}^2\rangle_{01}$$

$$|\Psi(0_3^+)\rangle = a_{13}|d_{5/2}^2\rangle_{01} + a_{23}|s_{1/2}^2\rangle_{01} + a_{33}|d_{3/2}^2\rangle_{01}$$

$$\sum_k a_{kp}^2 = 1$$

The full spectrum:

Repeat the same procedure for $J^\pi = 1^+, 2^+, 3^+, 4^+$.

Shell-model codes

m-scheme codes

- ANTOINE (E. Caurier)
http://www.iphc.cnrs.fr/nutheo/code_antoine/menu.html
- NuShellX@MSU (W. Rae, B. A. Brown)
- MSHELL (T. Mizusaki)
- REDSTICK (W. E. Ormand, C. Johnson)
- ...

J-coupled codes

- NATHAN (E.Caurier, F.Nowacki)
- DUPSM (Novoselsky, Vallières)
- Ritsschil (Zwarts)
- ...

Features

- Matrix dimension: $\sim 10^{10}$ and beyond
- Lanczos diagonalization algorithm
- Calculation of the matrix elements on-the-fly

Lanczos algorithm

Creation of a tri-diagonal matrix:

$$\begin{aligned}\hat{H}|1\rangle &= E_{11}|1\rangle + E_{12}|2\rangle \\ \hat{H}|2\rangle &= E_{21}|1\rangle + E_{22}|2\rangle + E_{23}|3\rangle \\ &\dots\end{aligned}$$

Matrix elements:

$$\begin{aligned}E_{11} &= \langle 1|\hat{H}|1\rangle \\ E_{12}|2\rangle &= (\hat{H} - E_{11})|1\rangle \\ E_{21} &= E_{12}, E_{22} = \langle 2|\hat{H}|2\rangle \\ E_{23}|3\rangle &= (\hat{H} - E_{22})|2\rangle - E_{21}|1\rangle \\ &\dots\end{aligned}$$

How to get the lowest states converged:

$$\begin{pmatrix} E_{11} & E_{12} \\ E_{21} & E_{22} \end{pmatrix} \Rightarrow \begin{pmatrix} E_{11} & E_{12} & 0 \\ E_{21} & E_{22} & E_{23} \\ 0 & E_{32} & E_{33} \end{pmatrix} \Rightarrow \begin{pmatrix} E_{11} & E_{12} & 0 & 0 \\ E_{21} & E_{22} & E_{23} & 0 \\ 0 & E_{32} & E_{33} & E_{34} \\ 0 & 0 & E_{43} & E_{44} \end{pmatrix} \Rightarrow \dots$$

You may need only a few iterations to get the lowest state of a $(10^3 \times 10^3)$ matrix converged!

Calculation of observables

General scheme:

- Construct the basis: $|\Phi_k\rangle$
- Expand the wave function: $|\Psi_p\rangle = \sum_k a_{kp} |\Phi_k\rangle$ and compute the Hamiltonian matrix $\{H_{lk}\}$.
- Solution of the Schrödinger equation by Hamiltonian matrix diagonalization: $\{H_{lk}\} \Rightarrow E_p, |\Psi_p\rangle$ (coefficients a_{kp})
- Calculation of matrix elements of the operators

$$T_{fi} \propto |\langle \Psi_f | \hat{O} | \Psi_i \rangle|^2$$

Electroweak operators:

$$\hat{O}(E, LM) = \sum_{k=1}^A e(k) r^L(k) Y_{LM}(\hat{r}(k))$$

$$\hat{O}(M, 1M) = \sum_{k=1}^A \mu_n(g_s(k) \vec{s}(k) + g_l(k) \vec{l}(k))$$

$$\hat{O}(F) = \sum_{k=1}^A \tau_{\pm}(k), \quad \hat{O}(GT) = \sum_{k=1}^A \vec{\sigma}(k) \tau_{\pm}(k) \quad \dots$$

Exercise 3

Electromagnetic transitions in ^{17}O

Calculate the $B(E2; 1/2_1^+ \rightarrow 5/2_{g.s.}^+)$ in ^{17}O modeled as a valence neutron in a $1s0d$ shell beyond the ^{16}O closed-shell core.

We take:

$$\int_0^{\infty} R_{1s_{1/2}}(r)r^2R_{0d_{5/2}}(r)dr \approx 12 \text{ fm}^2.$$

Compare your result to experimental value $B_{exp}(E2) = 6.3 \text{ e}^2 \cdot \text{fm}^4$. What can you conclude?

Solution to exercise 3

Electromagnetic transitions in ^{17}O

The single-particle electric multipole operator reads

$$\hat{O}(E, 2M) = er^2 Y_{2M}(\theta, \phi),$$

The reduced probability of the $\mathcal{E}L$ -transition from the initial to the final state is

$$B(EL; J_i \rightarrow J_f) = \frac{1}{2J_i + 1} |\langle J_f || O(EL) || J_i \rangle|^2.$$

In our case, there is one valence particle: $J_i = (1s_{1/2})$, $J_f = (0d_{5/2})$.

$$\begin{aligned} B(E2; 1/2^+ \rightarrow 5/2^+) &= \frac{1}{2} |\langle 0d_{5/2} || \hat{O}(E2) || 1s_{1/2} \rangle|^2 = \\ &= \frac{1}{2} |\langle n_f=0, l_f=2, s_f=\frac{1}{2}, j_f=\frac{5}{2} || er^2 Y_2(\theta, \phi) || n_i=1, l_i=0, s_i=\frac{1}{2}, j_i=\frac{1}{2} \rangle|^2 = \\ &= \frac{1}{2} e^2 \underbrace{\left(\int R_{0d_{5/2}}^*(r) r^2 R_{1s_{1/2}}(r) dr \right)^2}_{\langle r^2 \rangle^2} \left| \langle 2\frac{1}{2}; \frac{5}{2} || Y_2(\theta, \phi) || 0\frac{1}{2}; \frac{1}{2} \rangle \right|^2. \end{aligned}$$

$$B(EL; j_i \rightarrow j_f) = e^2 \frac{1}{4\pi} \langle r^2 \rangle^2 (2j_f + 1)(2l_i + 1)(2L + 1) (l_i 0 l 0 | l_f 0)^2 \left\{ \begin{matrix} \frac{1}{2} & l_f & j_f \\ L & j_i & l_i \end{matrix} \right\}^2$$

Solution to exercise 3

Electromagnetic transitions in ^{17}O (continued)

$$B(E2; 1/2^+ \rightarrow 5/2^+) = e^2 \frac{1}{4\pi} \langle r^2 \rangle^2 6 \times 5(0020|20)^2 \left\{ \begin{matrix} 1/2 & 2 & 5/2 \\ 2 & 1/2 & 0 \end{matrix} \right\}^2 = 34.4 e^2 \cdot \text{fm}^4.$$

Experimental value $B_{\text{exp}}(E2; 1/2^+ \rightarrow 5/2^+) = 6.3 e^2 \cdot \text{fm}^4$.

This means that the neutron should have an effective charge: $\tilde{e}_n \approx 0.43 e$. because we work in a severely restricted model space (one valence nucleon !).

Standard effective $E2$ and $M1$ operators

$$\tilde{e}_\pi \approx 1.5 e, \quad \tilde{e}_\nu \approx 0.5 e$$

$$\tilde{g}_s(\pi) \approx 0.7 \underbrace{g_s(\pi)}_{5.586}, \quad \tilde{g}_l(\pi) = g_l(\pi) = 1$$

$$\tilde{g}_s(\nu) \approx 0.7 \underbrace{g_s(\nu)}_{-3.826}, \quad \tilde{g}_l(\nu) = g_l(\nu) = 0$$

Nuclear many-body problem

Non-relativistic Hamiltonian for A nucleons

$$\hat{H} = \sum_{i=1}^A \frac{\hat{p}_i^2}{2m} + \sum_{i<j=1}^A \hat{W}(\vec{r}_i - \vec{r}_j)$$
$$\hat{H} = \underbrace{\sum_{i=1}^A \left[\frac{\vec{p}_i^2}{2m} + U(\vec{r}_i) \right]}_{\hat{H}^{(0)}} + \underbrace{\sum_{i<j=1}^A W(\vec{r}_i - \vec{r}_j) - \sum_{i=1}^A U(\vec{r}_i)}_{\hat{V}}$$

The residual interaction is assumed to have a two-body form

$$\hat{V} = \sum_{i<j=1}^A \hat{V}(\vec{r}_i - \vec{r}_j)$$

One and two-body Hamiltonian

$$\hat{H} = \sum_{i=1}^A \hat{h}(\vec{r}_i) + \sum_{i<j=1}^A \hat{V}(\vec{r}_i - \vec{r}_j)$$

Occupation-number representation (second quantization)

Creation and annihilation operators

$$|\alpha\rangle = a_{\alpha}^{\dagger}|0\rangle \quad \langle\alpha| = \langle 0|a_{\alpha}$$

Wave function of a fermion in a quantum state α in coordinate space:

$$\langle\vec{r}|\alpha\rangle = \phi_{\alpha}(\vec{r})$$

(Anti-)commutation relations:

$$\begin{aligned}\{a_{\alpha}^{\dagger}, a_{\beta}\} &= a_{\alpha}^{\dagger}a_{\beta} + a_{\beta}a_{\alpha}^{\dagger} = \delta_{\alpha\beta} \\ \{a_{\alpha}^{\dagger}, a_{\beta}^{\dagger}\} &= \{a_{\alpha}, a_{\beta}\} = 0\end{aligned}$$

Normalized and antisymmetric A -fermion state:

$$|\alpha_1\alpha_2\dots\alpha_A\rangle = a_{\alpha_A}^{\dagger}a_{\alpha_{A-1}}^{\dagger}\dots a_{\alpha_2}^{\dagger}a_{\alpha_1}^{\dagger}|0\rangle$$

One-body operators

$$\hat{O} = \sum_{k=1}^A \hat{O}(\vec{r}_k)$$

$$\langle \alpha | \hat{O} | \beta \rangle = \int \phi_{\alpha}^*(\vec{r}) \hat{O}(\vec{r}) \phi_{\beta}(\vec{r}) d\vec{r}$$

Second-quantized form of the one-body operator \hat{O} :

$$\hat{O} = \sum_{\alpha\beta} \langle \alpha | \hat{O} | \beta \rangle a_{\alpha}^{\dagger} a_{\beta}$$

For example, the number operator reads

$$\hat{N} = \sum_{\alpha\beta} \langle \alpha | \hat{1} | \beta \rangle a_{\alpha}^{\dagger} a_{\beta} = \sum_{\alpha} a_{\alpha}^{\dagger} a_{\alpha}$$

Operators in the occupation-number formalism

Symmetric two-body operator acting on an A -fermion system

$$\hat{T} = \sum_{j < k=1}^A \hat{T}(\vec{r}_k, \vec{r}_j)$$

$$\langle \alpha\beta | \hat{T} | \gamma\delta \rangle = \int \phi_\alpha^*(\vec{r}_1) \phi_\beta^*(\vec{r}_2) \hat{T}(\vec{r}_1, \vec{r}_2) (1 - \hat{P}_{12}) \phi_\gamma(\vec{r}_1) \phi_\delta(\vec{r}_2) d\vec{r}_1 d\vec{r}_2,$$

Second-quantized form of the two-body operator \hat{T} :

$$\hat{T} = \frac{1}{4} \sum_{\alpha\beta\gamma\delta} \langle \alpha\beta | \hat{T} | \gamma\delta \rangle a_\alpha^\dagger a_\beta^\dagger a_\delta a_\gamma$$

Nuclear many-body Hamiltonian in the occupation-number formalism

Non-relativistic Hamiltonian for A nucleons

$$\hat{H} = \sum_{i=1}^A \hat{h}(\vec{r}_i) + \sum_{i < j=1}^A \hat{V}(\vec{r}_i - \vec{r}_j) = \hat{H}^{(0)} + \hat{V}$$

↓

$$\hat{H} = \underbrace{\sum_{\alpha} \varepsilon_{\alpha} a_{\alpha}^{\dagger} a_{\alpha}}_{\text{one-body term}} + \underbrace{\frac{1}{4} \sum_{\alpha\beta\gamma\delta} \langle \alpha\beta | V | \gamma\delta \rangle a_{\alpha}^{\dagger} a_{\beta}^{\dagger} a_{\delta} a_{\gamma}}_{\text{two-body term}},$$

Two-body term in a JT -coupled form

$$\hat{V} = -\frac{1}{4} \sum_{j_{\alpha} j_{\beta} j_{\gamma} j_{\delta}} \langle j_{\alpha} j_{\beta} | V | j_{\gamma} j_{\delta} \rangle_{JT} \sqrt{(1 + \delta_{\alpha\beta})(1 + \delta_{\gamma\delta})} \left[a_{j_{\alpha}}^{\dagger} a_{j_{\beta}}^{\dagger} \right]^{(JT)} \left[\tilde{a}_{j_{\gamma}} \tilde{a}_{j_{\delta}} \right]^{(JT)} \quad (00)$$

Necessary ingredients

Single-particle energies

$$\varepsilon_{\alpha}$$

from experimental spectra of A_{core} plus a neutron or a proton

Two-body matrix elements (TBMEs)

$$\langle j_{\alpha} j_{\beta} | V | j_{\gamma} j_{\delta} \rangle_{JT}$$

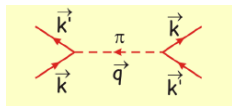
from theory ?

Bare nucleon-nucleon (NN) interaction

The NN interaction between two nucleons in the vacuum:
NN scattering data, deuteron bound states properties.

Elastic scattering in momentum space (Yukawa)

$$V_{\pi NN}(1, 2) = \frac{g_{\pi NN}^2}{4M^2} \frac{(\vec{\sigma}_1 \cdot \vec{q})(\vec{\sigma}_2 \cdot \vec{q})}{\vec{q}^2 + m_\pi^2}$$



Potential (Fourier transform) in coordinate space

$$V_{\pi NN}^{OPEP}(1, 2) = \frac{g_{\pi NN}^2}{4M^2} \frac{m_\pi^3}{12} \left\{ \vec{\sigma}_1 \cdot \vec{\sigma}_2 + \left(1 + \frac{3}{m_\pi r} + \frac{3}{(m_\pi r)^2} \right) (3\vec{\sigma}_1 \cdot \vec{r} \vec{\sigma}_2 \cdot \vec{r} - \sigma_1 \cdot \sigma_2) \right\} \frac{e^{-m_\pi r}}{m_\pi r}$$

Meson-exchange theories of NN potential: high-precision potentials (CD-Bonn, AV18, etc)

Concept of effective interaction (operators)

Effective in-medium nucleon-nucleon interaction

- In-medium effects (renormalization of the hard core)
- Truncated model space

$$\hat{H}\Psi = (\hat{H}^{(0)} + \hat{V})\Psi = E\Psi ,$$

True wave function:

$$\Psi = \sum_{k=1}^{\infty} a_k \Phi_k .$$

Wave function in a model space:

$$\Psi' = \sum_{k=1}^M a_k \Phi_k .$$

$$\langle \Psi' | \hat{H}_{eff} | \Psi' \rangle = \langle \Psi | \hat{H} | \Psi \rangle = E$$

$$\langle \Psi' | \hat{O}_{eff} | \Psi' \rangle = \langle \Psi | \hat{O} | \Psi \rangle$$

Approaches to the problem: **phenomenological** or **microscopic**.

Practical approaches to effective interaction

- **Schematic** interaction (parametrized interaction between two nucleons in a nuclear medium)
- **Phenomenological** interaction (Fit of TBME's to energy levels of nuclei to be described within the chosen model space)
- **Microscopic** interaction (derived from a bare NN-force)

Schematic interaction

Some examples:

$$V(1,2) = -V_0 \exp(\mu r)/(\mu r)$$

$$V(1,2) = -V_0 \delta(\vec{r}_1 - \vec{r}_2)$$

$$V(1,2) = -V_0 \delta(\vec{r}_1 - \vec{r}_2)(1 + \alpha \vec{\sigma}_1 \cdot \vec{\sigma}_2)$$

$$V(1,2) = -V_0 \delta(\vec{r}_1 - \vec{r}_2) \delta(r_1 - R)$$

...

$$V(1,2) = \chi Q \cdot Q \quad (Q = r^2 Y_{2\mu}(\Omega_r))$$

...

A few parameters (interaction strengths) are fitted to reproduce energy levels in a certain region of (a few) neighboring nuclei \Rightarrow local description only!

Exercise 4: TBMEs of the δ -force

Multipole expansion of the delta-function

$$V(1,2) = -V_0\delta(\vec{r}_1 - \vec{r}_2),$$

$$\delta(\vec{r}_1 - \vec{r}_2) = \sum_k \frac{\delta(r_1 - r_2)}{r_1 r_2} \frac{2k+1}{4\pi} P_k(\cos\theta_{12})$$

Diagonal TBMEs between normalized and antisymmetric states

$$\langle j_1 j_2 | V | j_1 j_2 \rangle_{JT} = I(2j_1 + 1)(2j_2 + 1) \begin{pmatrix} j_1 & j_2 & J \\ \frac{1}{2} & -\frac{1}{2} & 0 \end{pmatrix}^2 \frac{1 + (-1)^{l_1 + l_2 + J}}{2}$$

$$\langle j^2 | V | j^2 \rangle_{JT} = I(2j + 1)^2 \begin{pmatrix} j & j & J \\ \frac{1}{2} & -\frac{1}{2} & 0 \end{pmatrix}^2, \quad \text{if } j_1 = j_2 = j$$

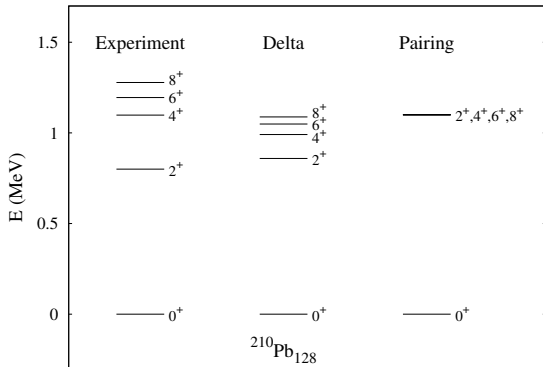
$$I = \frac{1}{4\pi} \int_0^\infty \frac{1}{r^2} [R_{n_1 l_1}(r) R_{n_2 l_2}(r)]^2 dr$$

Example 1: ^{210}Pb in $(\nu 0h_{9/2})^2$

$$V_{\text{delta}}(1,2) = -V_0 \delta(\vec{r}_1 - \vec{r}_2)$$

$$V_{\text{pairing}}(1,2) = -G \hat{S}_+ \cdot \hat{S}_-$$

$$\langle j_a^2 | V_{\text{pairing}}(1,2) | j_b^2 \rangle_{01} = -(-1)^{l_a+l_b} \frac{1}{2} G \sqrt{(2j_a+1)(2j_b+1)}$$

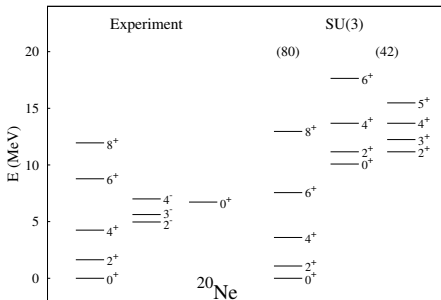


Example 2: ^{20}Ne and SU(3) model of Elliott

$$\hat{H} = \sum_{i=1}^A \left[-\frac{p_i^2}{2m} + \frac{1}{2}m\omega^2 r_i^2 \right] - \chi Q \cdot Q$$

Q is an algebraic quadrupole operator (Q_μ, L_μ are SU(3) generators) *J.P.Elliott (1958)*

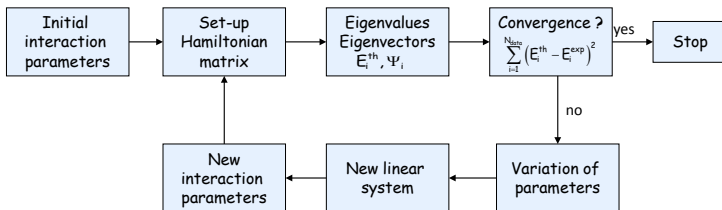
Group-theoretical classification of nuclear states (analytical solution) — see *lecture of H. Molique*.



Rotational classification of nuclear states as mixing of many spherical configurations

Empirical interaction (least-squares-fit method)

All TBME's $\langle j_a j_b | V | j_c j_d \rangle_{JT}$ are considered as free parameters !



Examples:

- $0p$ -shell: ${}^4\text{He}$ - ${}^{16}\text{O}$ (15 TBME's) *Cohen, Kurath (1965)*
- $1s0d$ -shell: ${}^{16}\text{O}$ - ${}^{40}\text{Ca}$ (63 TBME's) *Brown, Wildenthal (1988)*
- $1p0f$ -shell: ${}^{40}\text{Ca}$ - ${}^{80}\text{Zr}$ (195 TBME's) *Honma et al (2002, 2004)*

Microscopic effective interaction

A bare NN potential (CD-Bonn, AV18, chiral N3LO, etc) requires regularization and modification to be applied for many-body calculations in a restricted model space.

Renormalization schemes (*see lecture of L. Bonneau*)

- G matrix followed by the many-body perturbation theory
- V_{low-k}
- SRG (IM-SRG)
- Okubo-Lee-Suzuki transformation

Successful, but still lack precision of the empirical interactions, mainly due to behavior of centroids. One of possible reasons: absence of 3N forces (A. Poves, A.P. Zuker, 1981; A.P. Zuker, 2003)

Interacting shell model

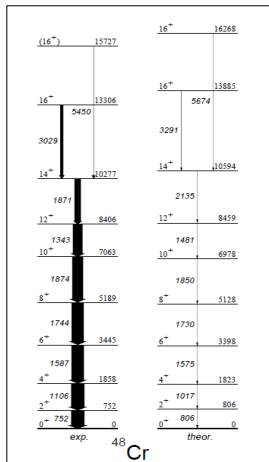
Oscillator-based shell model with accurate realistic interactions formulated in one or two harmonic-oscillator shells model spaces (large-scale diagonalization).

- Detailed information on individual states and transitions at low energies
- Conservation of principal symmetries

Numerous applications to nuclear structure, weak interaction and astrophysics

E. Caurier et al, Rev. Mod. Phys. 77, 427 (2005)

State-of-the-art calculations: backbending in ^{48}Cr



$J < 10$: collective rotation
 $J = 10-12$: backbending phenomenon
 (competition between rotation and
 alignment of $Of_{7/2}$ particles)
 $J > 12$: spherical states

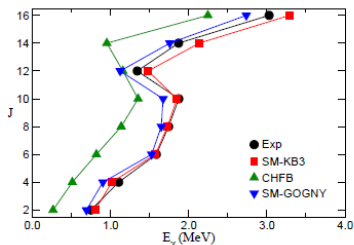
KB3 (semi-empirical interaction
 in pf -shell model space)
 Strasbourg-Madrid

For $J < 10$:

$$E_J \sim J(J+1)$$

$$Q_0 = \frac{(J+1)(2J+3)}{3K^2 - J(J+1)} Q_{\text{spec}}(J), \quad K \neq 1$$

$$B(E2; J \rightarrow J-2) = \frac{5}{16\pi} e^2 |(JK20 | J-2, K)|^2 Q_0^2$$



E. Caurier et al, Rev. Mod. Phys. 77 (2005) 427

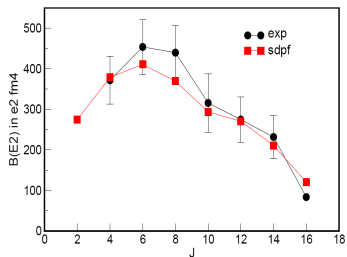
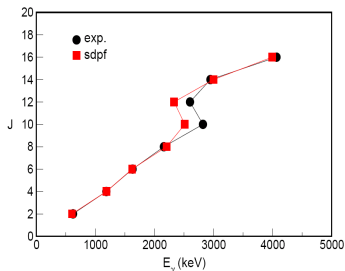
State-of-the-art calculations: superdeformation in ^{36}Ar

Intruder np-nh configurations can lead even to superdeformation !

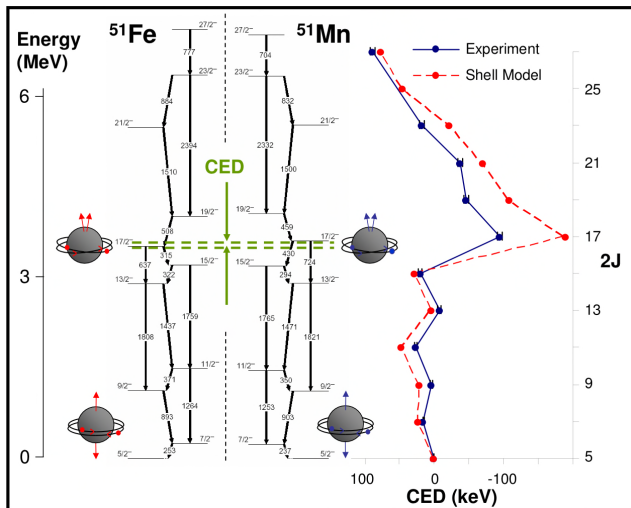
$[\text{sd}]^{16}[\text{pf}]^0$ - 0p0h - spherical configuration

$[\text{sd}]^{12}[\text{pf}]^4$ - 4p4h - deformed configuration

E. Caurier et al, Phys.Rev.Lett. 95, 042502 (2005)



State-of-the-art calculations: mirror bands in $A = 51$



S. Lenzi, A. Zuker, E. Caurier et al

No-Core Shell Model

Non-relativistic Hamiltonian for A nucleons in many $N\hbar\Omega$ harmonic oscillator space

$$\hat{H} = \sum_{i=1}^A \frac{\vec{p}_i^2}{2m} + \sum_{i<j=1}^A W(\vec{r}_i - \vec{r}_j)$$

Problem: excitation of the center-of-mass of the system.

Center-of-mass coordinates

$$\vec{R} = \frac{1}{A} \sum_{i=1}^A \vec{r}_i; \quad \vec{P} = \sum_{i=1}^A \vec{p}_i$$

Translational-invariant Hamiltonian

$$\hat{H} = \sum_{i=1}^A \frac{\vec{p}_i^2}{2m} - \frac{\vec{P}^2}{2mA} + \sum_{i<j=1}^A W(\vec{r}_i - \vec{r}_j)$$

Separation of the harmonic oscillator Hamiltonian into center-of-mass and intrinsic Hamiltonians

$$\begin{aligned} H_{ho} &= \sum_{i=1}^A \left[\frac{\vec{p}_i^2}{2m} + \frac{m\Omega^2 \vec{r}_i^2}{2} \right] \\ &= \underbrace{\frac{1}{2mA} \sum_{i<j=1}^A (\vec{p}_i - \vec{p}_j)^2 + \frac{m\Omega^2}{2A} \sum_{i<j=1}^A (\vec{r}_i - \vec{r}_j)^2}_{\hat{H}_{int}} + \underbrace{\frac{\vec{P}^2}{2mA} + \frac{mA\Omega^2 \vec{R}^2}{2}}_{\hat{H}_{CoM}} \end{aligned}$$

No-Core Shell Model

Translational-invariant Hamiltonian

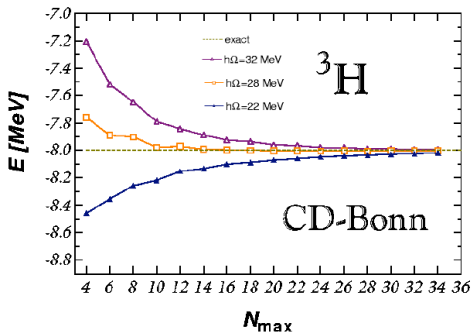
$$\hat{H} = \sum_{i=1}^A \left[\frac{\vec{p}_i^2}{2m} + \frac{m\Omega^2 \vec{r}_i^2}{2} \right] - \frac{\vec{P}^2}{2mA} + \sum_{i<j=1}^A W(\vec{r}_i - \vec{r}_j) - \sum_{i=1}^A \frac{m\Omega^2 \vec{r}_i^2}{2}$$

$$\hat{H} = \sum_{i=1}^A \left[\frac{\vec{p}_i^2}{2m} + \frac{m\Omega^2 \vec{r}_i^2}{2} \right] + \sum_{i<j=1}^A W(\vec{r}_i - \vec{r}_j) - \frac{m\Omega^2}{2A} \sum_{i<j=1}^A (\vec{r}_i - \vec{r}_j)^2 - \hat{H}_{CoM}$$

Separation of the Center-of-Mass motion

$$\hat{H}^\Omega = \hat{H} + \beta \left(\hat{H}_{CoM} - \frac{3}{2} \hbar \Omega \right)$$

^3H with CD-Bonn and N^3LO



Potential	E_{gs} (MeV)
AV18	-7.62
N^3LO	-7.86
CD-Bonn	-8.00
INOY	-8.48
Exp	-8.48

B.R. Barrett, P. Navratil, J. P. Vary, “*Ab-initio* no core shell model”,
Prog. Part. Nucl. Phys. 69 (2013) 131 — and refs. therein.

See lecture of R. Lazauskas for alternative methods.

The Nuclear Shell Model

- The Nuclear Shell Model is a powerful microscopic approach to nuclear structure
- Very good description of energies and transitions at low energies
- High predictive power.
- Important applications :
 - structure of nuclei far from stability (proton rich or neutron-rich nuclei)
 - calculation of weak interaction processes on nuclei for the tests of the Standard Model
 - calculation of the nuclear structure input (masses, half-lives, reaction rates, etc) relevant for astrophysical simulations (*r*-processes, *rp*-process, etc..).

Some references on the Shell model theory

Books

- P. J. Brussaard, P. W. M. Glaudemans, *Shell-Model Applications in Nuclear Spectroscopy* (North-Holland, Amsterdam, 1977).
- K. Heyde, *The Nuclear Shell Model* (Springer-Verlag, Heidelberg, 2004)
- A. DeShalit, I. Talmi, *Nuclear Shell Theory* (Academic Press, New York and London, 1963)
- I. Talmi, *Simple Models of Complex Nuclei, The Shell Model and the Interacting Boson Model* (Harwood Academic Publishers, New York, 1993)
- R. D. Lawson, *Theory of the Nuclear Shell Model* (Clarendon Press, 1980).

Reviews

- E. Caurier et al, *The shell model as a unified view of nuclear structure*, Rev. Mod. Phys. **77** (2005) 427.
- B. R. Barrett, P. Navratil, J. P. Vary, *Ab-initio no core shell model*, Prog. Part. Nucl. Phys. **69** (2013) 131