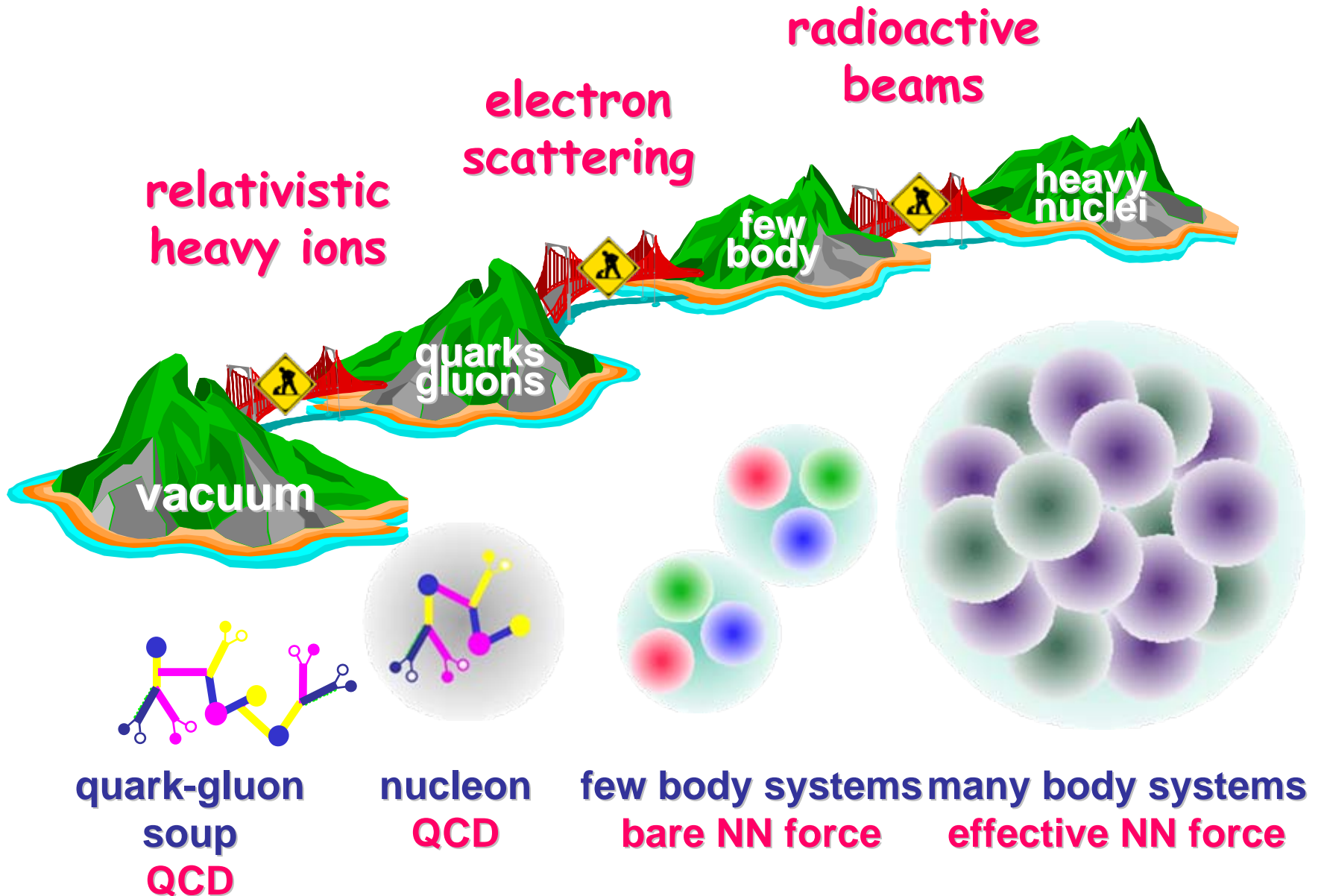
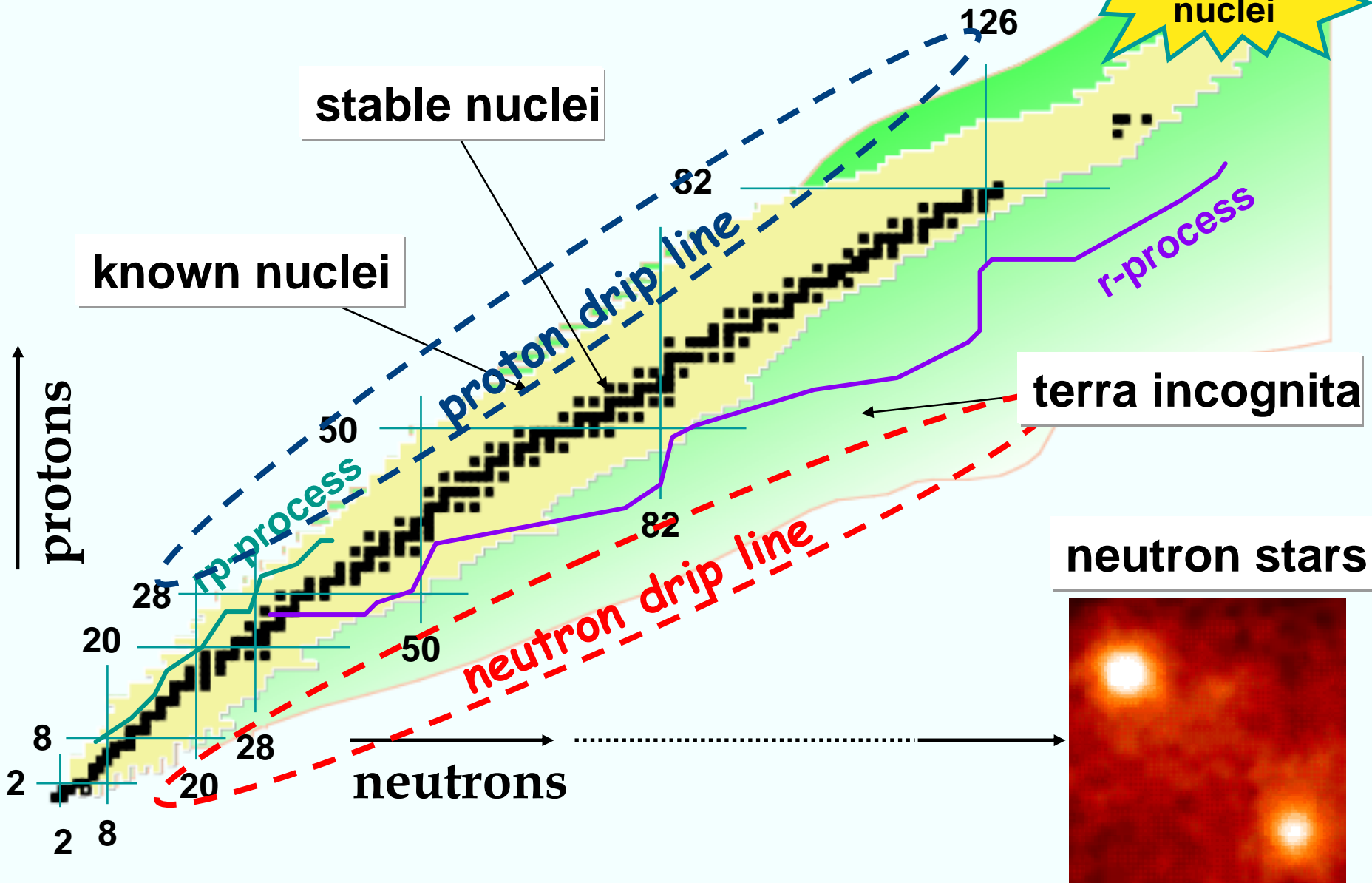


# The Nuclear Many-Body Problem

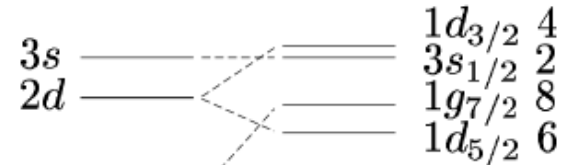


# Nuclear Landscape

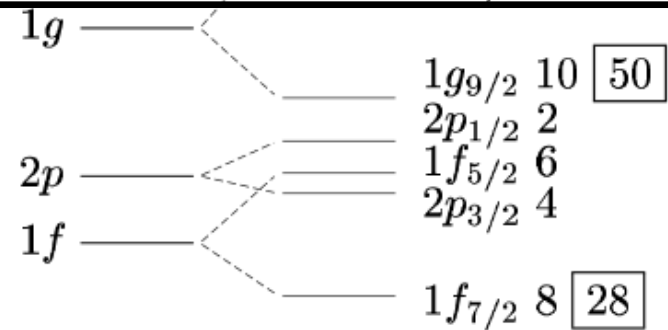


# The shell model: a schematic view

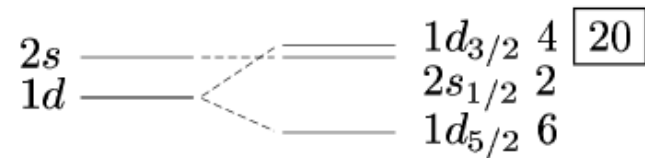
All upper shells neglected



Active space:

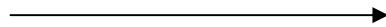


Inactive core  $^{40}\text{Ca}$

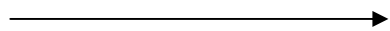


## Active space

_____	$1g_{9/2}$	10	<span style="border: 1px solid black; padding: 2px;">50</span>
_____	$2p_{1/2}$	2	
_____	$1f_{5/2}$	6	
_____	$2p_{3/2}$	4	
_____	$1f_{7/2}$	8	<span style="border: 1px solid black; padding: 2px;">28</span>



**2-body matrix elements  
between all pairs of states**



**Diagonalization of a big matrix in the active space**

**Starting point of the model: nucleons in a potential well.**

**Problem of the model: active space grows very quickly.**

**Respect always symmetries of the 2-body hamiltonian:  
work in the laboratory frame of reference**

# Mean-field methods

====  $1d_{3/2}$  4  
====  $3s_{1/2}$  2  
====  $1g_{7/2}$  8  
====  $1d_{5/2}$  6

====  $1g_{9/2}$  10 50  
====  $2p_{1/2}$  2  
====  $1f_{5/2}$  6  
====  $2p_{3/2}$  4

====  $1f_{7/2}$  8 28

====  $1d_{3/2}$  4 20  
====  $2s_{1/2}$  2  
====  $1d_{5/2}$  6

====  $1p_{1/2}$  2 8  
====  $1p_{3/2}$  4

====  $1s_{1/2}$  2 2

**Construct all the orbitals from the mean**

**2-body interaction of 1 particle with the others**

**All orbitals are active!**

**Caution: not true single particle levels**

**Break symmetries and work in a frame of**

**reference intrinsic to the nucleus!**

The Schrödinger equation is equivalent to a wave function variational principle: minimize  $\langle \Psi | H | \Psi \rangle$  under the constraint  $\langle \Psi | \Psi \rangle = 1$ :

$$\delta \{ \langle \Psi | H | \Psi \rangle / \langle \Psi | \Psi \rangle \} = 0$$

$$H = \sum_i (T_i + U_i) + \sum_{i>j} V(\vec{r}_i - \vec{r}_j)$$

One-body

Two-body

Hartree-Fock method:

the ground state wave function is a Slater determinant.

# The Hartree-Fock method

Wave-function of a many particle system= Slater determinant

$$\Psi = \frac{1}{N!} \sum_P (-)^P \phi_\alpha(\vec{r}_1) \phi_\beta(\vec{r}_2) \dots \phi_\nu(\vec{r}_A)$$

The particles interact through a 2-body interaction  $v(\mathbf{r}_1 - \mathbf{r}_2)$

They are also confined by a central potential  $v_{\text{ext}}(\mathbf{r})$ .

The total energy is:

$$E = \sum_{i,\sigma} \int d\vec{r} \phi_i^*(\vec{r}, \sigma) \left( -\frac{\hbar^2}{2m} \Delta + v_{\text{ext}}(\vec{r}) \right) \phi_i(\vec{r}, \sigma) + \frac{1}{2} \sum_{i,\sigma} \sum_{j,\sigma'} \int \int d\vec{r} d\vec{r}' \phi_i^*(\vec{r}, \sigma) \phi_j^*(\vec{r}', \sigma') v(\vec{r}, \vec{r}') (\phi_i(\vec{r}, \sigma) \phi_j(\vec{r}', \sigma') - \phi_i(\vec{r}', \sigma') \phi_j(\vec{r}, \sigma))$$

direct
exchange

Minimize the energy with a constraint on norm conservation:

$$\left( -\frac{\hbar^2}{2m}\Delta + v_{ext}(\vec{r}) \right) \phi_i(\vec{r}, \sigma) + \sum_{j, \sigma'} \int d\vec{r}' \phi_j^*(\vec{r}', \sigma') v(\vec{r}, \vec{r}') \phi_i(\vec{r}, \sigma) - \phi_i(\vec{r}', \sigma') \phi_j(\vec{r}, \sigma) = \epsilon_i \phi_i(\vec{r}, \sigma)$$

One defines the one-body diagonal and non diagonal densities:

$$\rho(\vec{r}') = \sum_{j, \sigma'} |\phi_j(\vec{r}', \sigma')|^2$$
$$\rho^{(1)}(\vec{r}, \sigma, \vec{r}', \sigma') = \sum_j \phi_j(\vec{r}', \sigma') \phi_j^*(\vec{r}, \sigma)$$



One rewrites the HF equations as a function of these densities:

$$\left( -\frac{\hbar^2}{2m}\Delta + v_{ext}(\vec{r}) \right) \phi_i(\vec{r}, \sigma) + U(\vec{r}) \phi_i(\vec{r}, \sigma) - \sum_{\sigma} \int d\vec{r}' \rho^{(1)}(\vec{r}, \sigma, \vec{r}', \sigma) v(\vec{r}, \vec{r}') \phi_i(\vec{r}', \sigma) = \epsilon_i \phi_i(\vec{r}, \sigma)$$

where 
$$U(\vec{r}) = \int v(\vec{r} - \vec{r}') \rho(\vec{r}') d\vec{r}'$$

The first line is easy: problem in a potential.

The second line is complicate: non local exchange term.

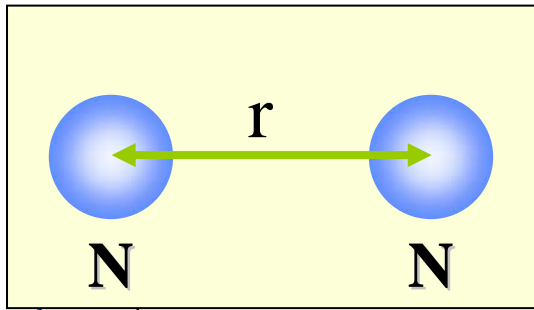
HF single particle energy: 
$$\epsilon_i = t_{ii} + \sum_{j=1}^A (\bar{v}_{ijij})$$
 ↘ 2-body matrix element between i and all other j

Total energy: 
$$E^{HF} = \sum_{i=1}^A \epsilon_i - \frac{1}{2} \sum_{i,j=1}^A \bar{v}_{ij,ij}$$
 ↘ no double counting!

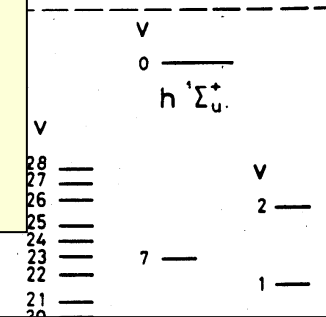
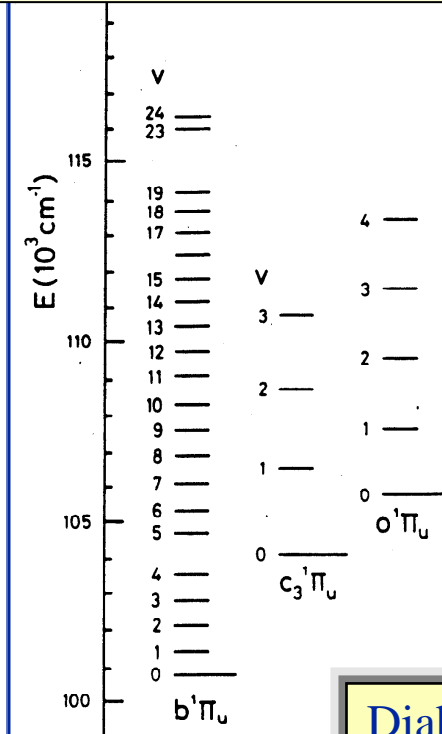
# Mean-field Methods

- Based on an “effective interaction” or a “density functional”  
The (small number of) parameters of the effective interaction are fixed by general considerations (**no local adjustments**)
- Pairing correlations are included at the BCS or better HFB level
- Full self-consistency
- No restrictions to a few shells, mean-field equations are solved as precisely as one wishes.
- Spherical and deformed nuclei are treated on the same footing, no “parametric deformation”

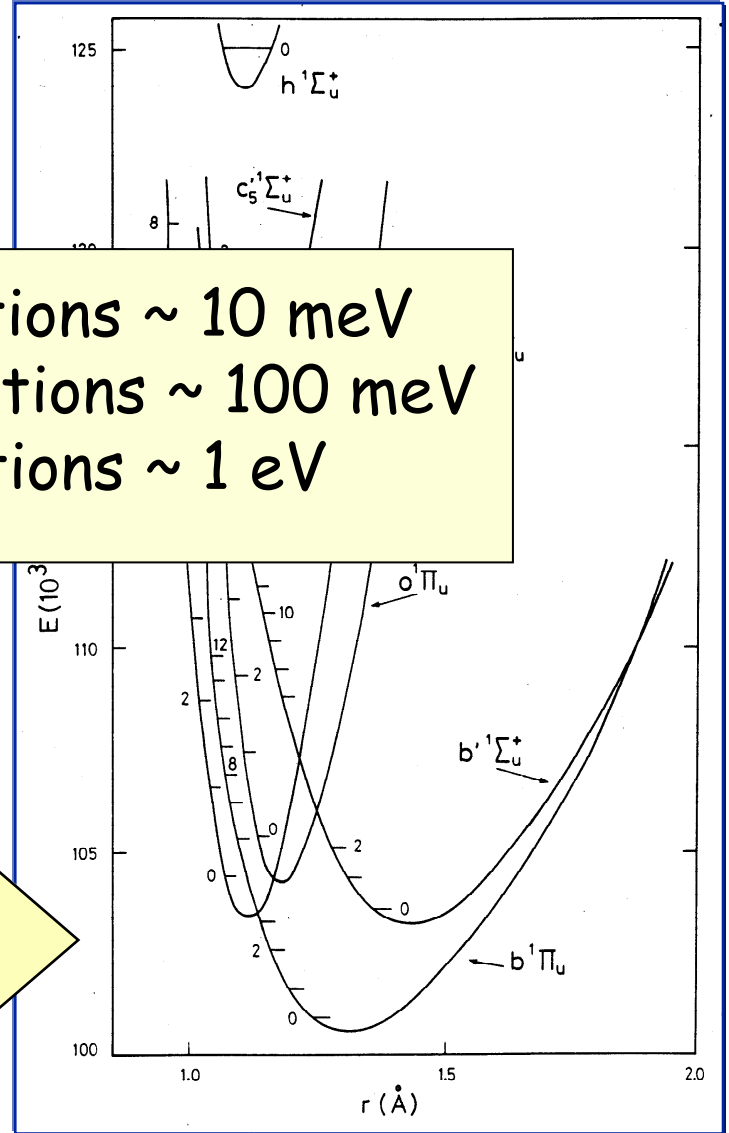
# Excitation spectrum of N<sub>2</sub> molecule



excited  $1^1\Sigma_u^+$  and  $1^1\Pi_u$  states



Rotational Transitions  $\sim 10$  meV  
 Vibrational Transitions  $\sim 100$  meV  
 Electronic Transitions  $\sim 1$  eV



Diabatic potential energy surfaces for excited electronic configurations of N<sub>2</sub>

Deformation of the nucleus introduced by a Lagrange multiplier:

$$H \Rightarrow H - \lambda q$$

by varying  $\lambda$ , one obtains solutions for different deformations

Missing ingredient: pairing correlations (superconductivity)

They can be introduced using the BCS theory:

- single particle states are occupied with a probability  $v^2$  between 1 and 0
- nucleons are grouped in pairs of opposite spin projections

The nuclear density becomes: 
$$\rho(\vec{r}) = \sum_i v_i^2 |\varphi_i(\vec{r})|^2$$

HF equations with modified densities

+ BCS equations to determine the occupations

Total wave function with only the right mean particle number!

# An example of an effective interaction: the Gogny force

It contains:

- A finite range central term:

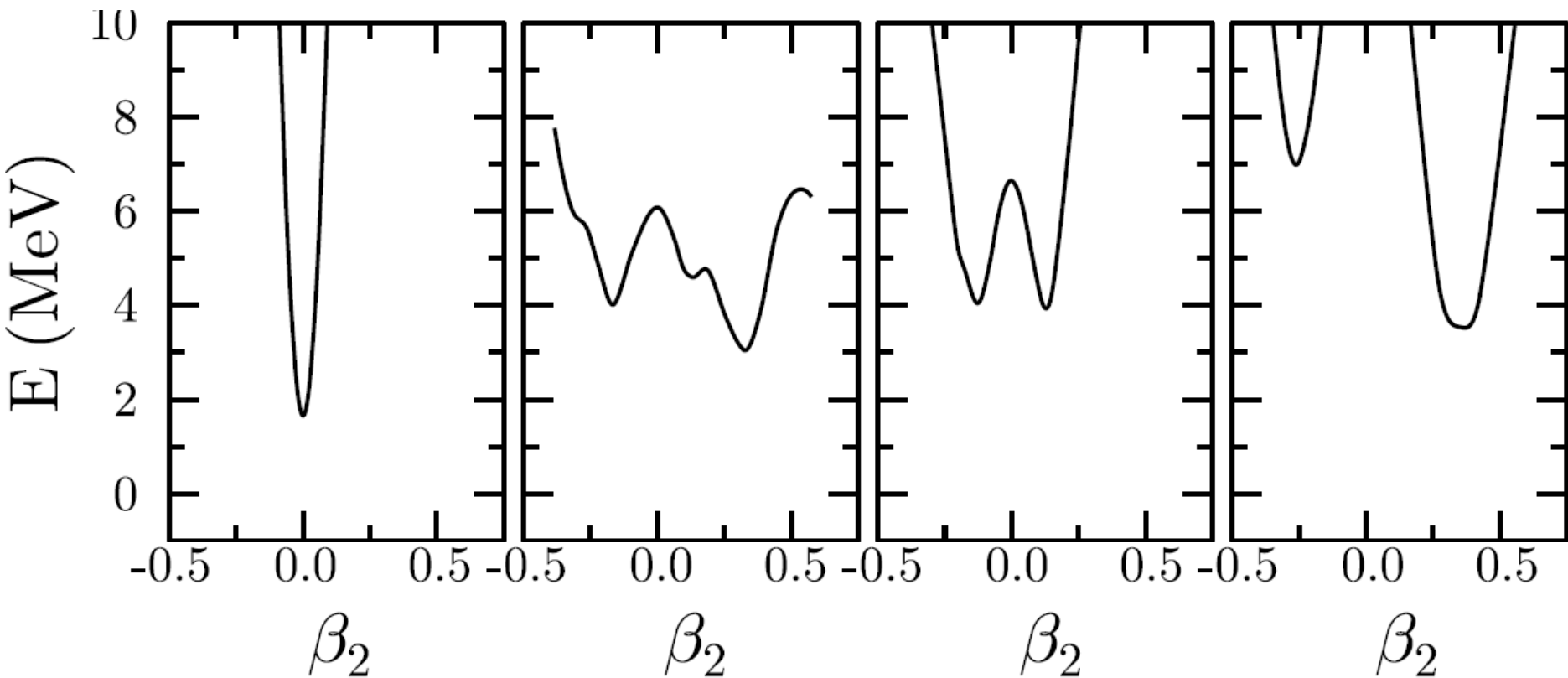
$$10 \quad V_C = \sum_{i=1,2} (V_W^i + V_M^i P^r + V_B^i P^\sigma + V_H^i P^\sigma P^r) \exp(-r^2/b_i^2)$$

- A zero range density dependent term

$$1-3 \quad t_3(1 + x_3 P^\tau) \rho(\vec{r})^\alpha$$

1 - Spin orbit and Coulomb

Parameters are adjusted on nuclear matter properties ( saturation, ...) properties of a few magic nuclei

$^{208}\text{Pb}$  $^{180}\text{Hg}$  $^{202}\text{Rn}$  $^{170}\text{Hf}$ 

Mean-field energy curves ( $\beta_2$  proportional to  $Q$ )

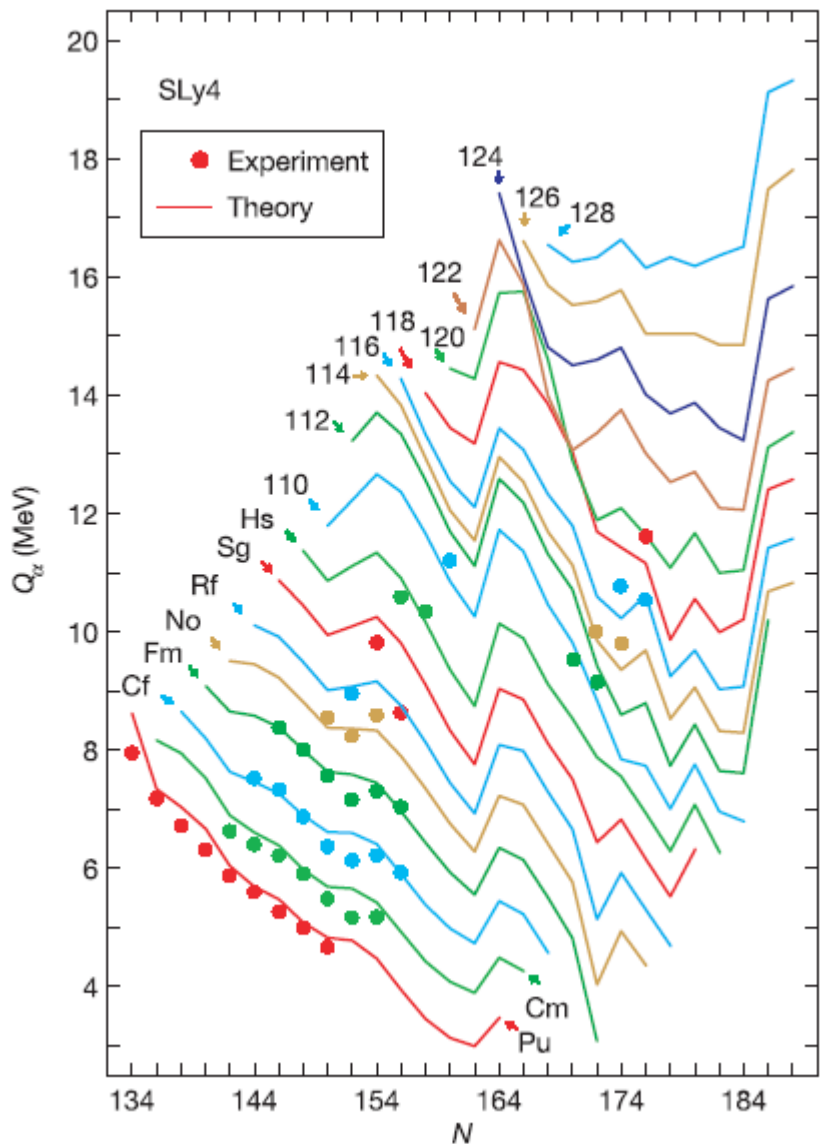
# The main approaches

## Three families:

- **Gogny**: finite range including a density dependence, same interaction for HF and pairing  
(Bruyères le Chatel, Madrid, some Japanese groups)
- **Skyrme**: zero range, specific interaction for pairing, easy  
(France, Poland, Belgium, P. Rheinardt et al., Japanese groups,...)
- **RMF**: relativistic but no exchange, pairing non relativistic  
(Munich-Zagreb, ....)

# Skyrme HFB

$Q_\alpha$  for isotopic chains  
for super heavy elements  
(only even-even)



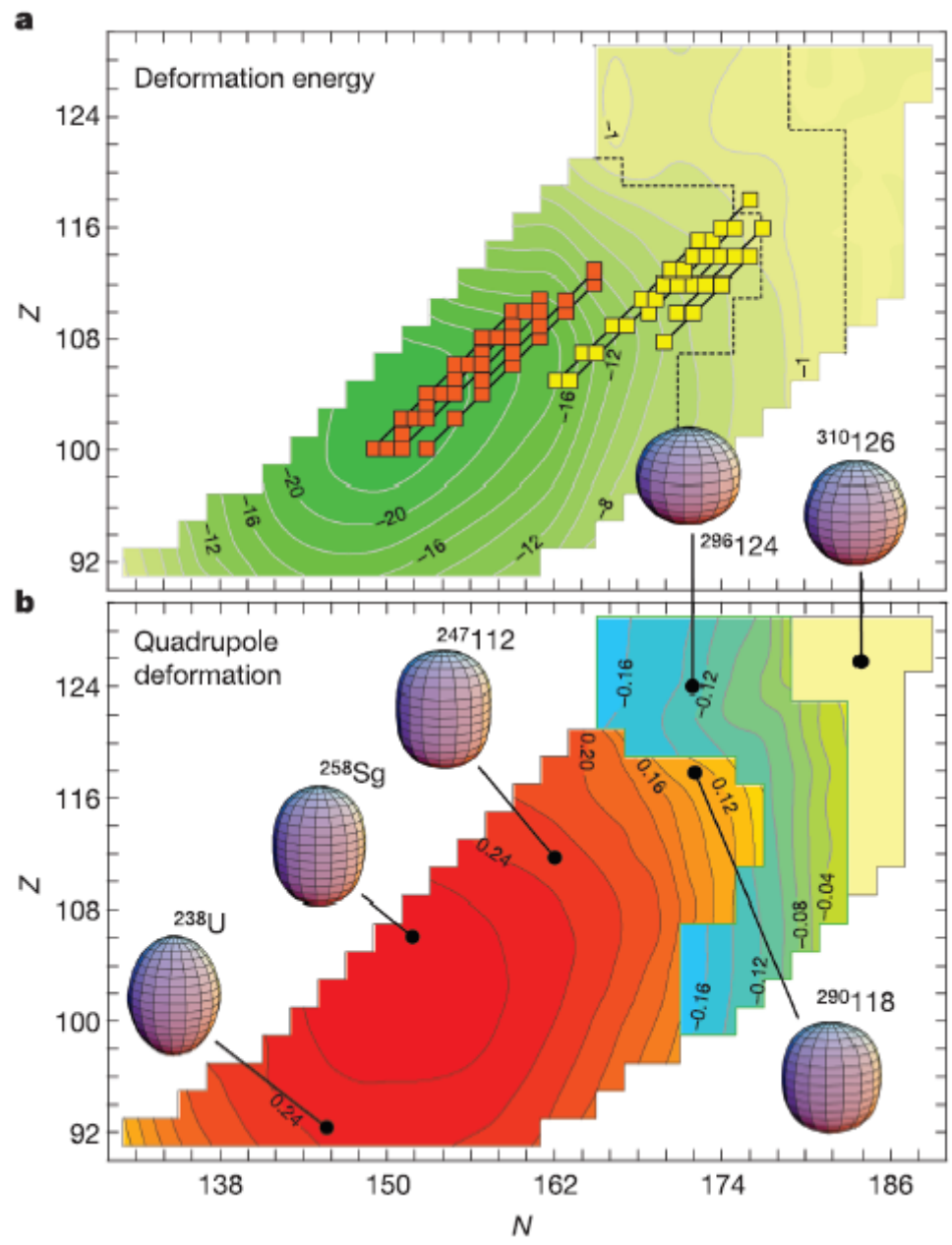
Cwiok, Heenen, Nazarewicz  
Nature 2005





# Skyrme HFB

## Deformation properties of super-heavies



# Beyond ground state properties of even-even nuclei

Breaking of time reversal invariance

by a cranking constraint:

$$H' = H - \omega J_x \quad \text{rotational bands for deformed nuclei}$$

by quasi particle excitations:

Odd nuclei : 1 qp states:

$$\beta_i^\dagger |0\rangle$$

Even nuclei: 2qp states

Still a mean-field method

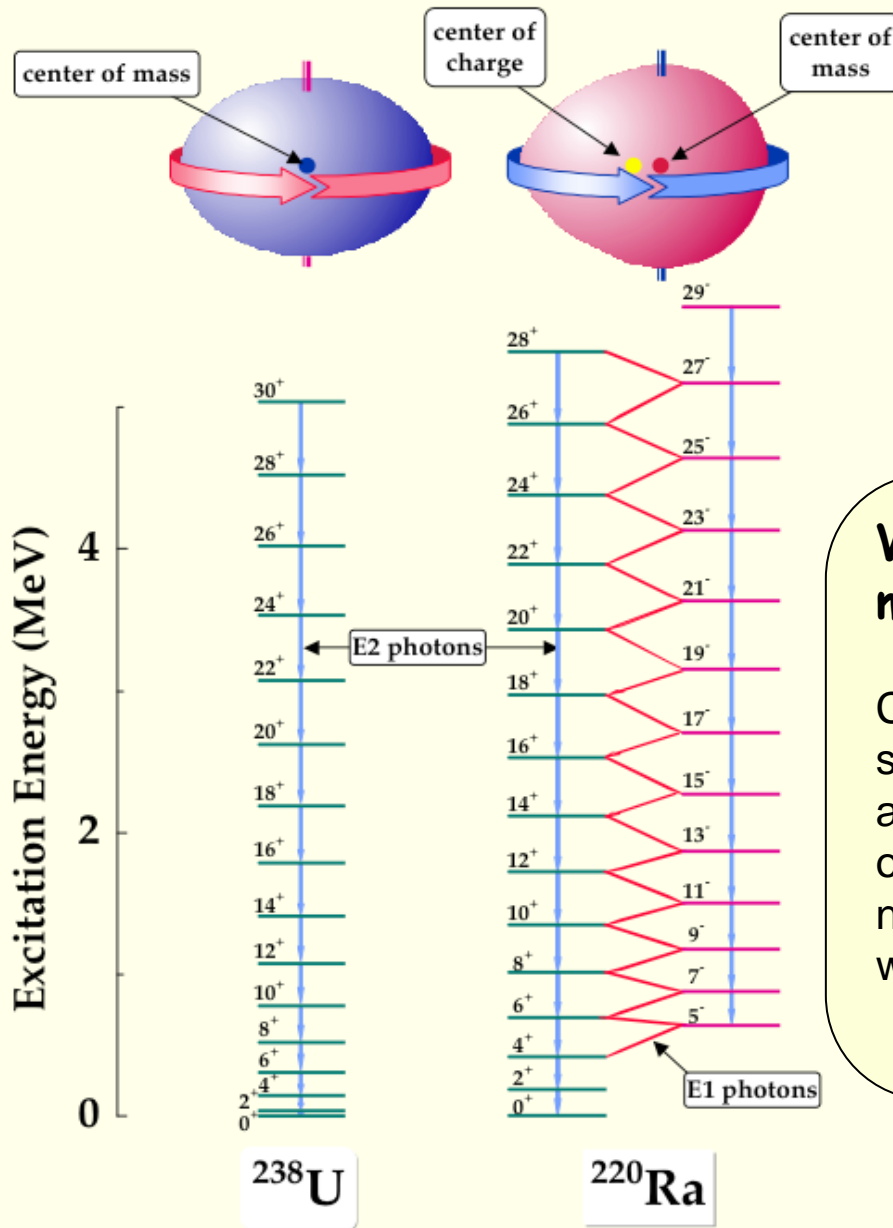
Full self-consistency for mean-field and pairing

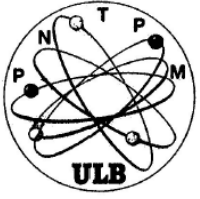
# Nuclear collective motion

Rotational Transitions ~ 0.2-2 MeV  
Vibrational Transitions ~ 0.5-12 MeV  
Nucleonic Transitions ~ 7 MeV

**What is the origin of ordered motion of complex nuclei?**

Complex systems often display astonishing simplicities. Nuclei are no exception. It is astonishing that a heavy nucleus, consisting of hundreds of rapidly moving protons and neutrons can exhibit collective motion, where all particles slowly dance in unison.





# Moments of inertia

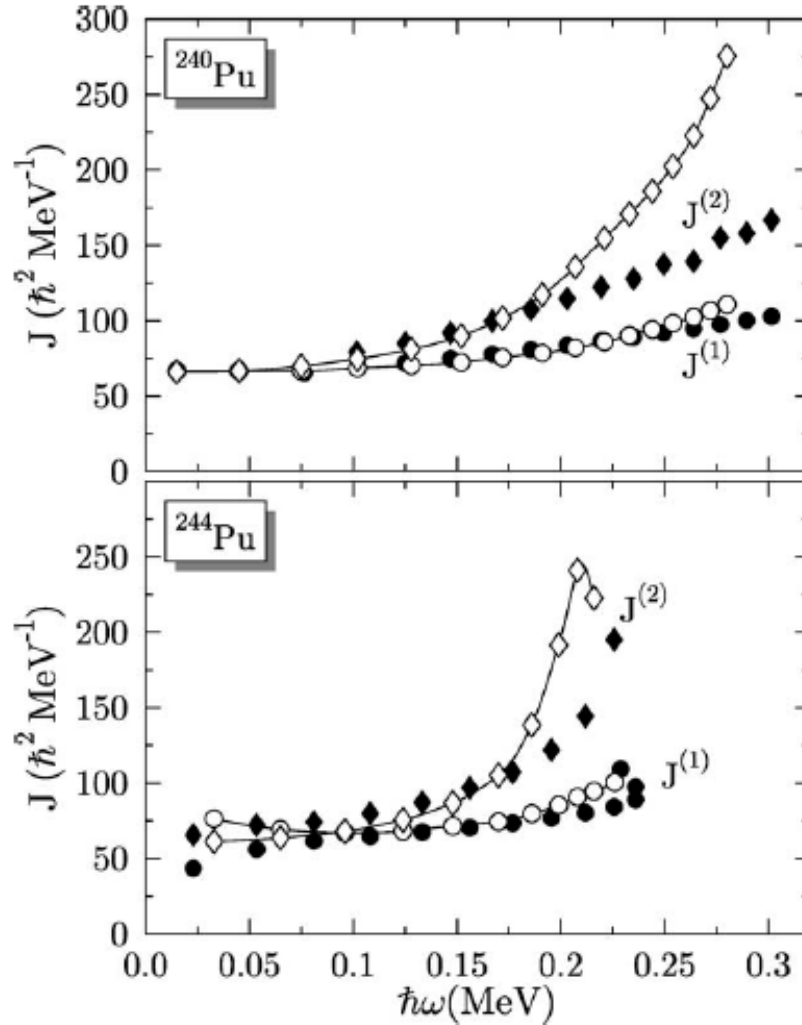
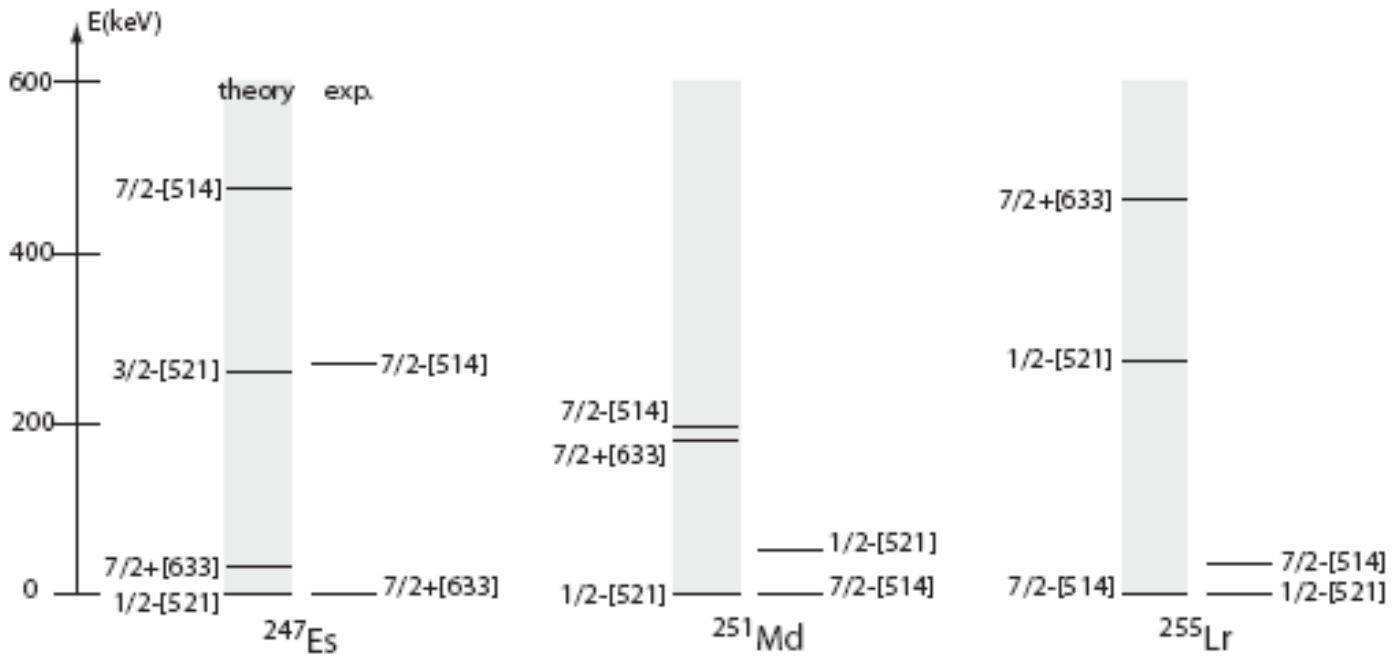


Fig. 3. Kinematical (circles) and dynamical (diamonds) moment of inertia for  $^{240}\text{Pu}$  (top) and  $^{244}\text{Pu}$  (bottom). Open (filled) markers denote calculated (experimental) values.

# Spectra of odd Z nuclei



# Nuclear DFT

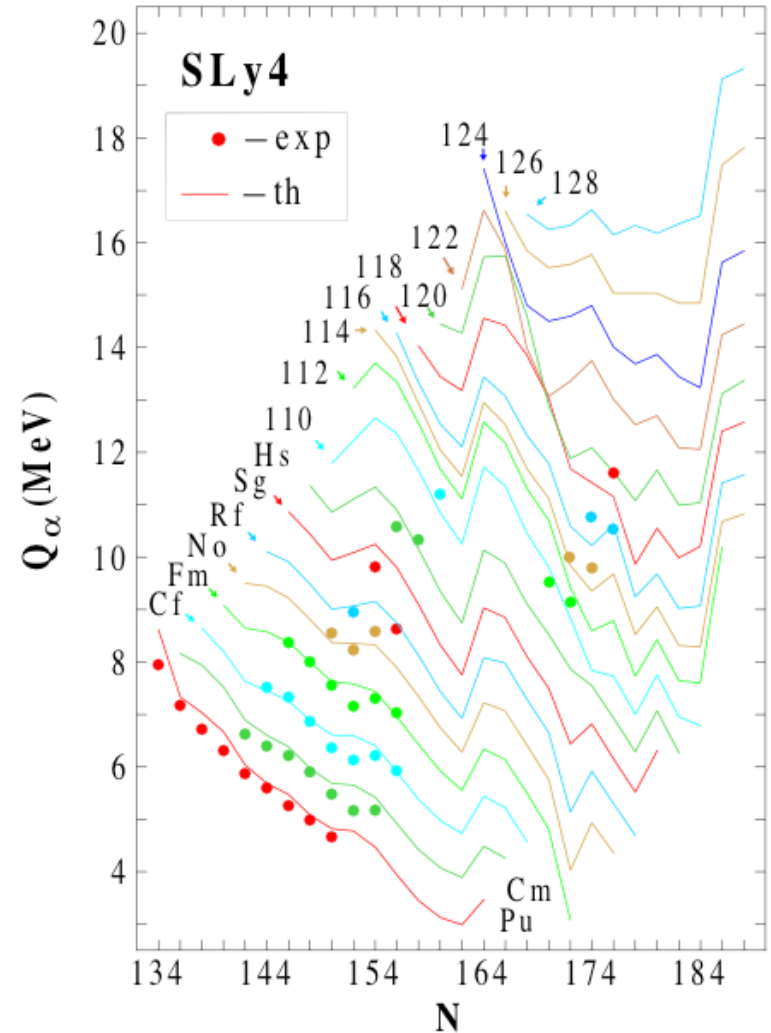
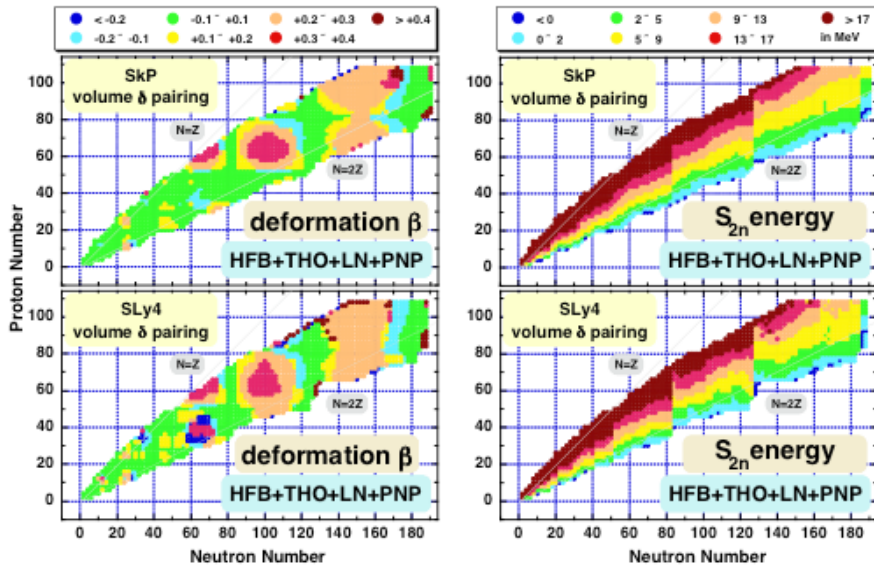
From Qualitative to Quantitative!

S. Cwiok, P.H. Heenen, W. Nazarewicz

## Microscopic Mass Table

M.V. Stoitsov et al., nucl-th/0406075

J. Dobaczewski et al., nucl-th/040407



Deformed Mass Table in one day!

# Towards Nuclear Energy Density Functional

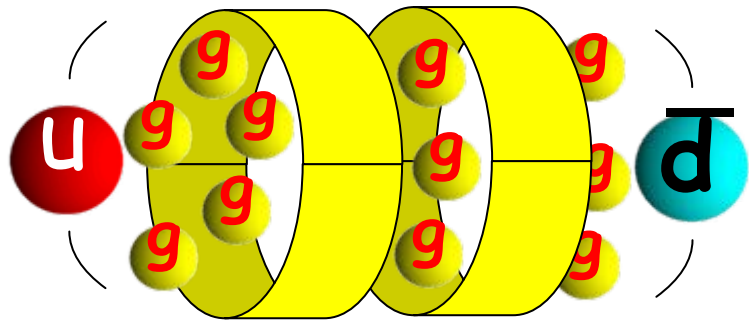
(unified description of nuclei and nuclear matter)

- Self-consistent mean-field theory (HF, HFB, RMF)
- Nuclear density functional theory
- Symmetry breaking crucial
- Symmetry restoration essential (projection techniques, GCM, QRPA)
- Pairing channel extremely important but poorly know

## Challenges:

- better understanding of isovector and density dependence of p-h and p-p interaction
  - how to extrapolate in isospin and mass?
  - time-odd fields
  - spin and isospin pieces
- improved treatment of many-body correlations
  - microscopic treatment
- nuclear matter equation of state at low and high temperatures
  - low density limit and clustering
  - isovector dependence of the symmetry energy

# Energy Scales in Nuclear Physics



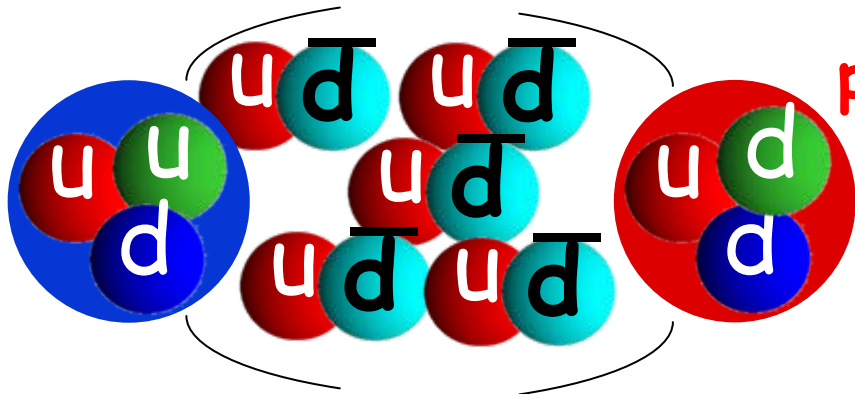
QCD scale



1000 MeV



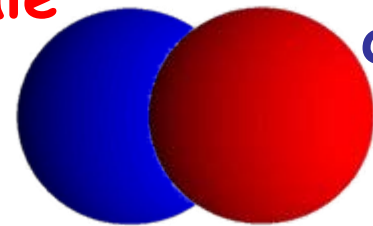
pion  $\pi^+$   
~140 MeV



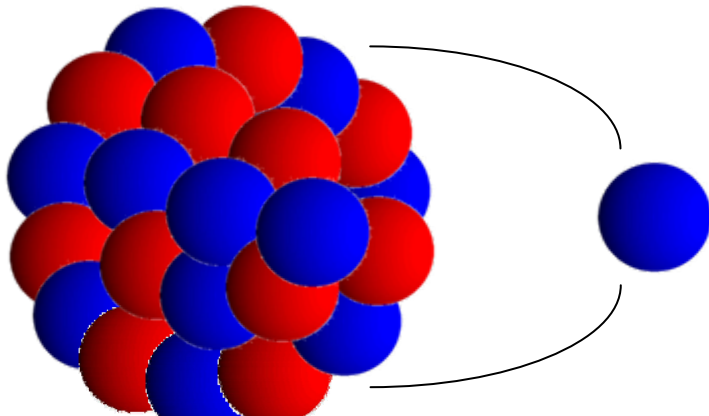
pion-mass scale



100 MeV



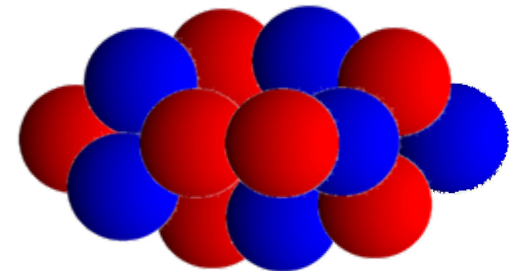
deuteron  
~3 MeV



N-binding scale



10 MeV



collective ~1 MeV