

# RIKEN-Tsukuba Collaboration

Takashi Nakatsukasa (RIKEN Nishina Center)

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# RIKEN-TSUKUBA, RIKEN-Niigata

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## ***Multi-reference EDF***

Projection & Config. Mixing in 3D real space representation

## ***Time-dependent energy-density-functional method (TD-EDF)***

Real-time approaches

Skyrme-TDHF in 3D real space

Canonical-basis TDHFB in 3D real space (Ebata)

Gogny-TDHFB in 3D harmonic oscillator basis (Hashimoto)

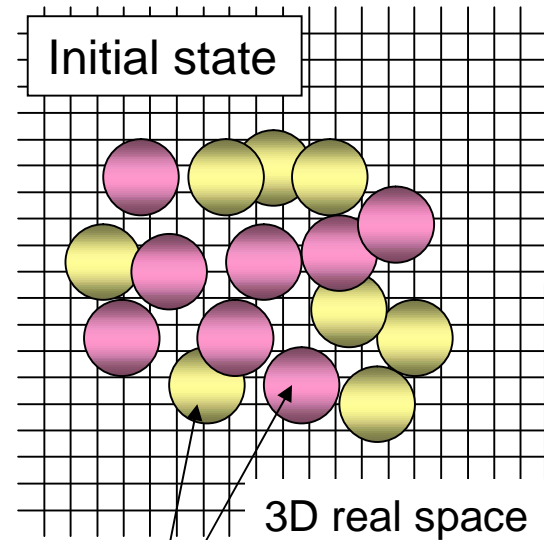
RPA, QRPA

Skyrme QRPA with axially deformed nuclei (Yoshida)

Finite amplitude method for RPA in 3D real space (Inakura)

Finite amplitude method for QRPA (Avogadro)

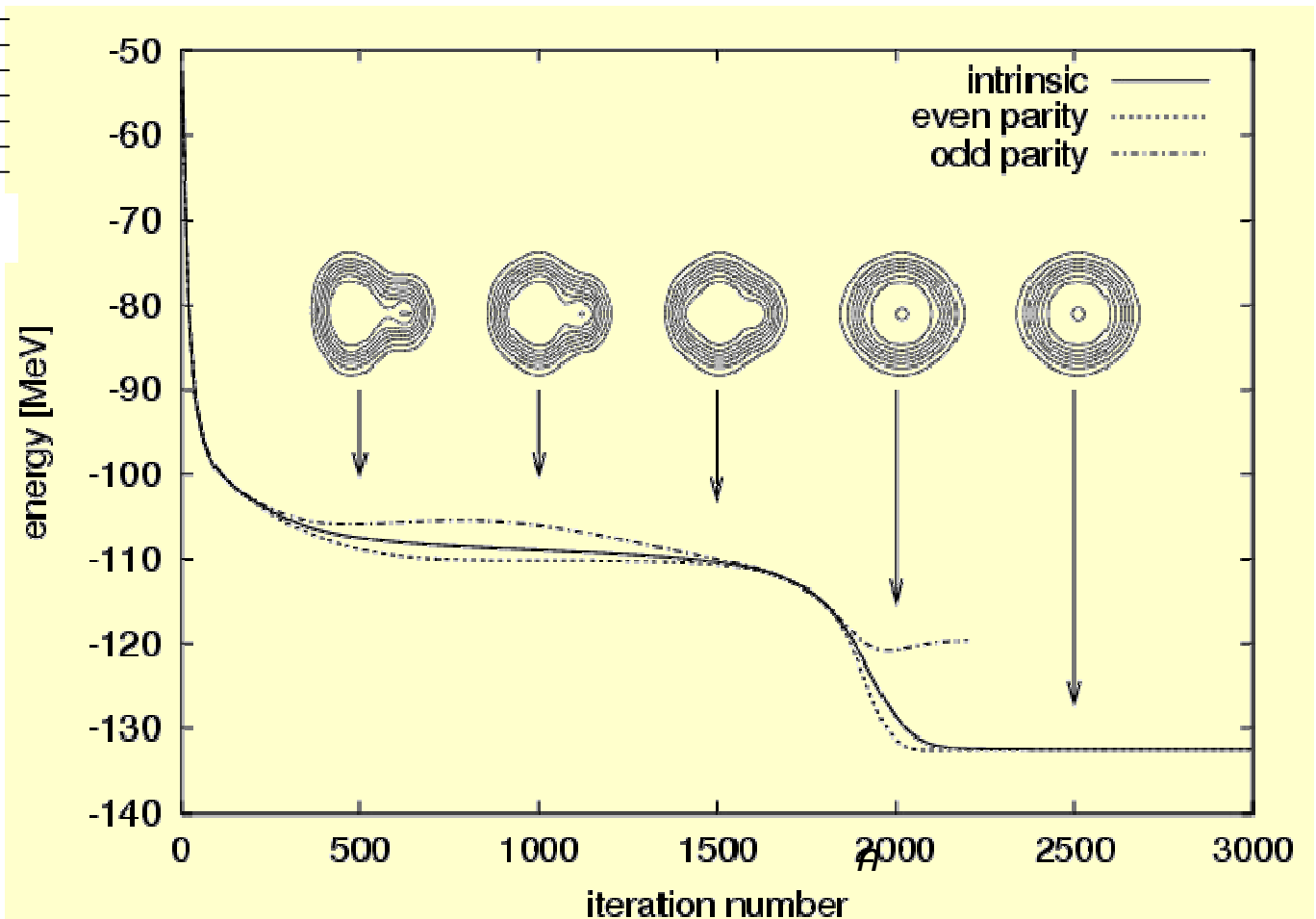
# Generation of many S-det's



$$|\phi_i^{(n+1)}\rangle = e^{-\Delta t h[\rho]} |\phi_i^{(n)}\rangle, \quad i = 1, \dots, A$$

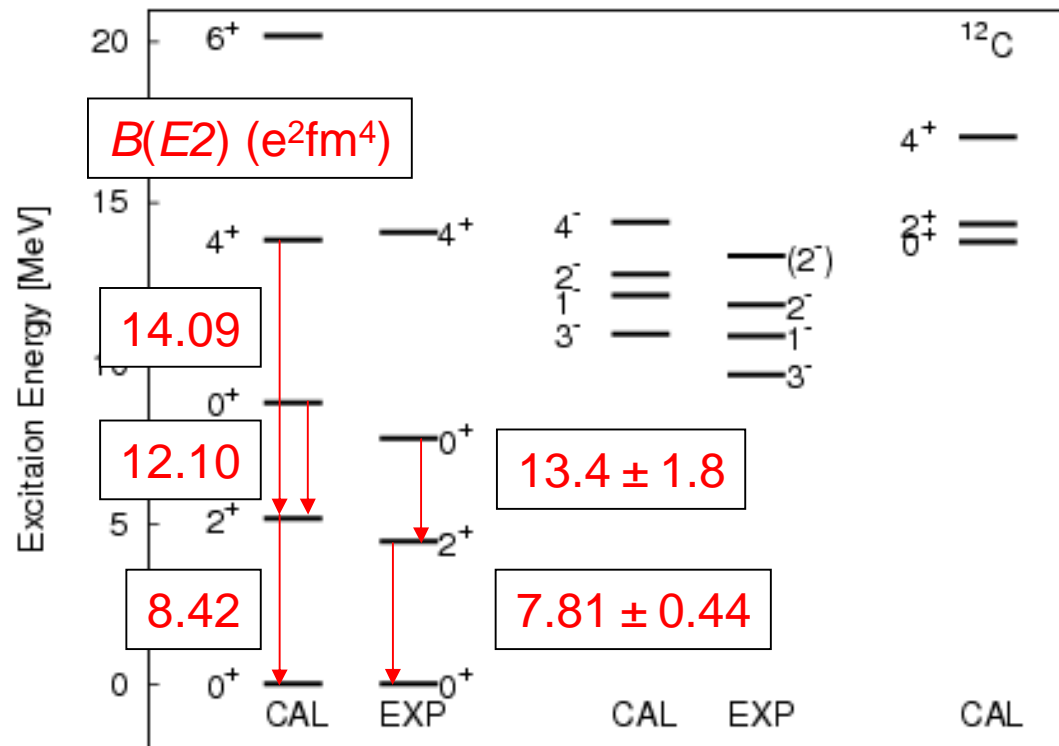
Imaginary-time evolution

Gaussian wave packets (n & p) whose positions are determined by random numbers.



SGII

# Energy spectrum in $^{12}\text{C}$



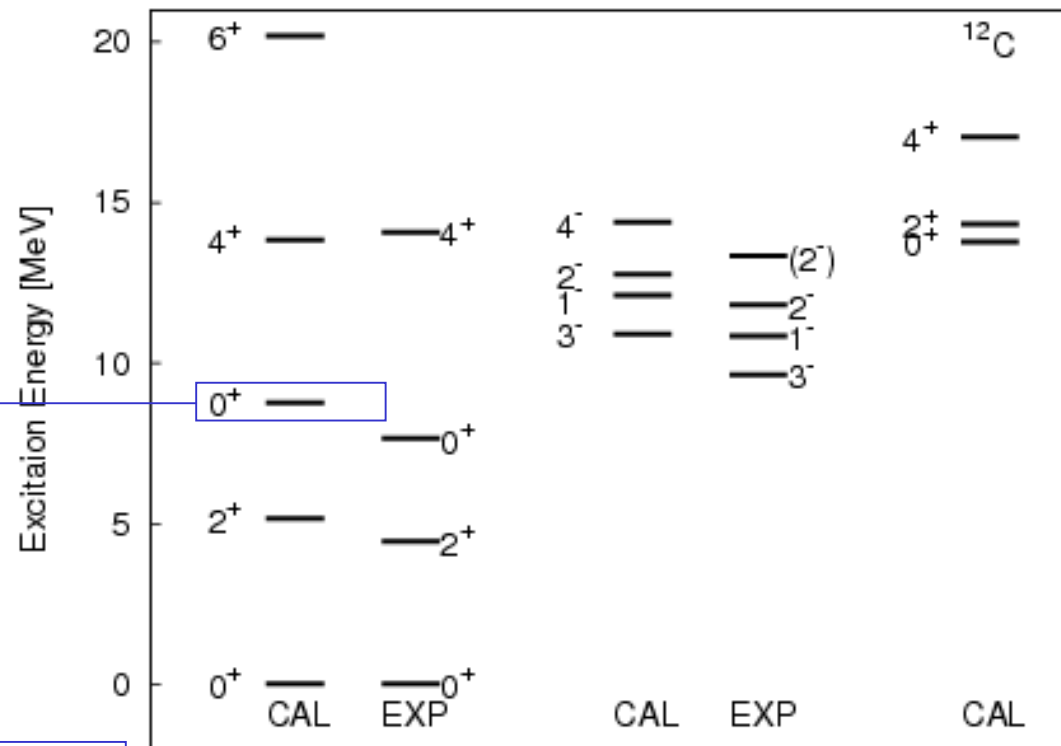
Similar results:

Kanada-En'yo (AMD+VAP), PRL81 (1998) 5292

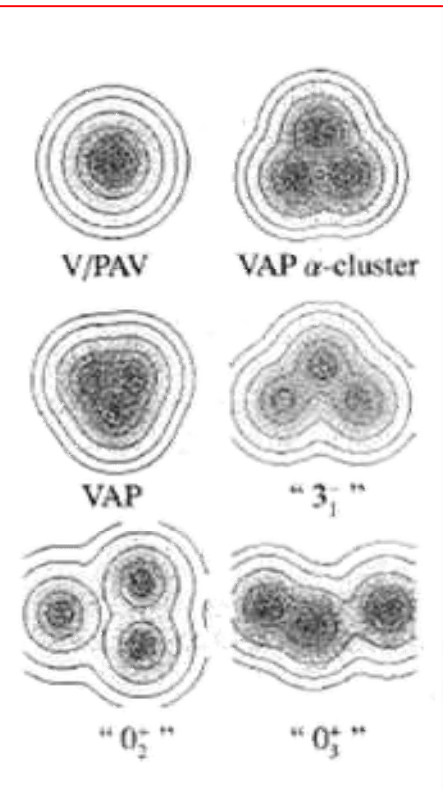
Neff & Feldmeier (FMD+GCM), NPA738 (2004) 357

# Energy spectrum in $^{12}\text{C}$

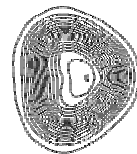
SGII



T. Neff et al. (2004)



second  $0^+$ : 42.4%



ground  $0^+$ : 94.8%



$3^-$ : 78.7%



third  $0^+$ : 62.1%

# Skyrme TDDFT in real space

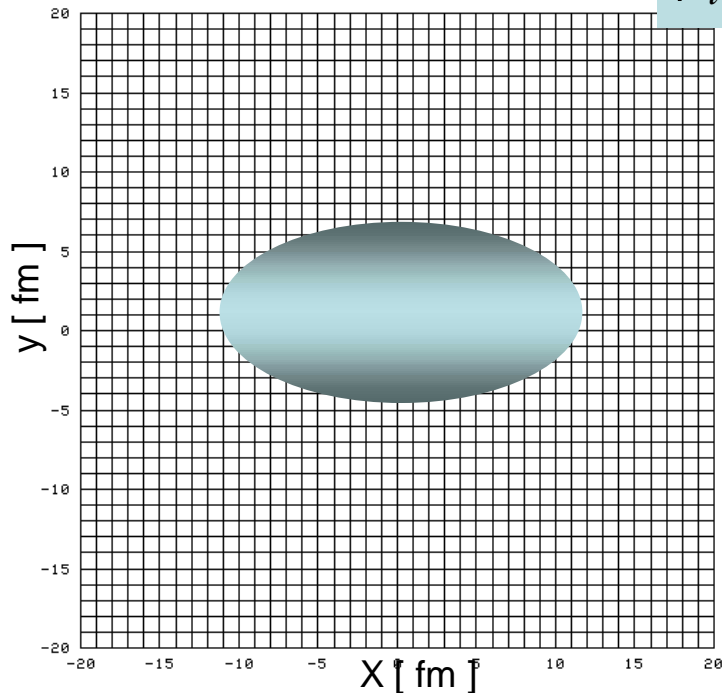
Time-dependent Kohn-Sham equation

$$i \frac{\partial}{\partial t} \psi_i(\mathbf{r} \sigma \tau, t) = \left( h_{\text{KS}}[\rho, \tau, \mathbf{j}, \mathbf{s}, \vec{\mathbf{J}}](t) + V_{\text{ext}}(t) - i \tilde{\eta}(r) \right) \psi_i(\mathbf{r} \sigma \tau, t)$$

3D space is discretized in lattice

Single-particle orbital:

$$\varphi_i(\mathbf{r}, t) = \{ \varphi_i(\mathbf{r}_k, t_n) \}_{k=1, \dots, Mr}^{n=1, \dots, Mt}, \quad i = 1, \dots, N$$



$N$ : Number of particles

$Mr$ : Number of mesh points

$Mt$ : Number of time slices

Spatial mesh size is about 1 fm.

Time step is about 0.2 fm/c

Nakatsukasa, Yabana, Phys. Rev. C71 (2005) 024301

# Inclusion of pairing (superfluidity)

- Simple formulation in Canonical basis
  - Canonical-basis TDHFB (→ Ebata)
  - Simultaneous equations for canonical states and  $(u,v)$ -factors
  - Special pairing functional only
- TDHFB
  - Gogny in 3D harmonic-oscillator basis (→ Hashimoto)
  - Linear-response and non-linear dynamics

# Finite Amplitude Method

T.N., Inakura, Yabana, PRC76 (2007) 024318.

A method to avoid the explicit calculation of the residual fields (interactions)

$$\begin{aligned}\omega |X_i(\omega)\rangle &= (h_0 - \varepsilon_i) |X_i(\omega)\rangle + \hat{Q} \{ \delta h(\omega) + V_{\text{ext}}(\omega) \} |\phi_i\rangle \\ \omega \langle Y_i(\omega) | &= -\langle Y_i(\omega) | (h_0 - \varepsilon_i) - \langle \phi_i | \{ \delta h(\omega) + V_{\text{ext}}(\omega) \} \hat{Q}\end{aligned}\quad (1)$$

Residual fields can be estimated by the finite difference method:

$$\begin{aligned}\delta h(\omega) &= \frac{1}{\eta} (h[\langle \psi' |, |\psi \rangle] - h_0) \\ |\psi_i\rangle &= |\phi_i\rangle + \eta |X_i(\omega)\rangle, \quad \langle \psi'_i | = \langle \phi_i | + \eta \langle Y_i(\omega) | \end{aligned}$$

Starting from initial amplitudes  $X^{(0)}$  and  $Y^{(0)}$ , one can use an iterative method to solve eq. (1).

Programming of the RPA code becomes very much trivial, because we only need calculation of the single-particle potential, with **different bras and kets**.



# Developments in FAM

- Large-scale linear response calculation (→ Inakura)
- Extension to QRPA (→ Avogadro)

# Other participants

- Masaaki Kimura (Hokkaido)
  - Anti-symmetrized Molecular Dynamics + HFB
- Hitoshi Nakada (Chiba)
  - Gogny-HFB, RPA, with gaussian basis
  - M3Y-HFB