



北京大学物理学院  
*School of Physics, Peking University*

# Gamow shell model with realistic nuclear forces

**Furong Xu**

## **I. Model**

**With-Core Gamow Shell Model (CGSM) based on realistic forces**  
**(resonance + continuum)**

## **II. Applications**

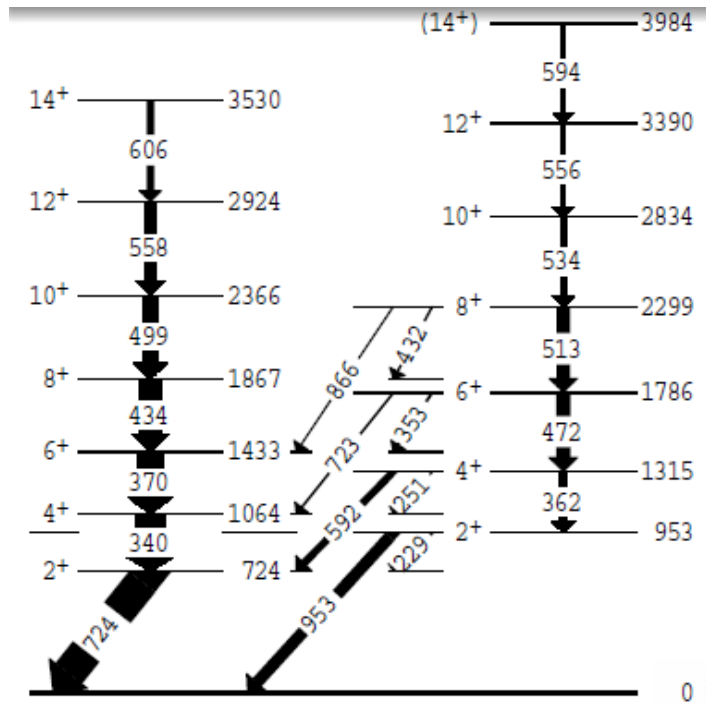
**Neutron-rich oxygen isotopes**

**Excitation spectra**

**Connecting Bound States to the Continuum**  
**Facility for Rare Isotope Beams (FRIB)**  
**June 11-22, 2018**

# $\gamma$ -ray spectra

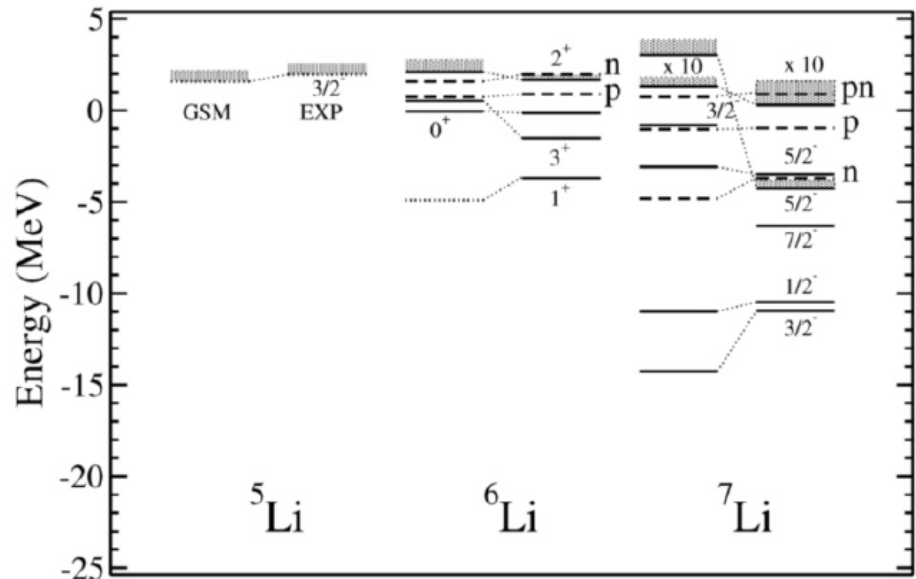
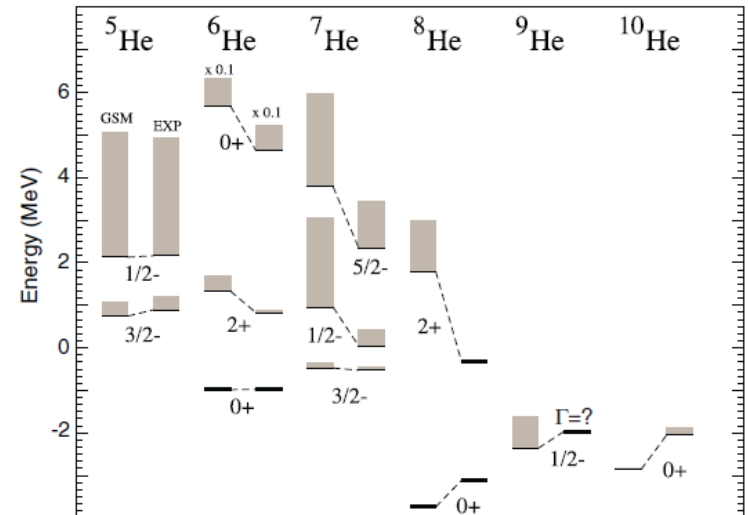
## $^{188}\text{Pb}$ : prolate and oblate bands



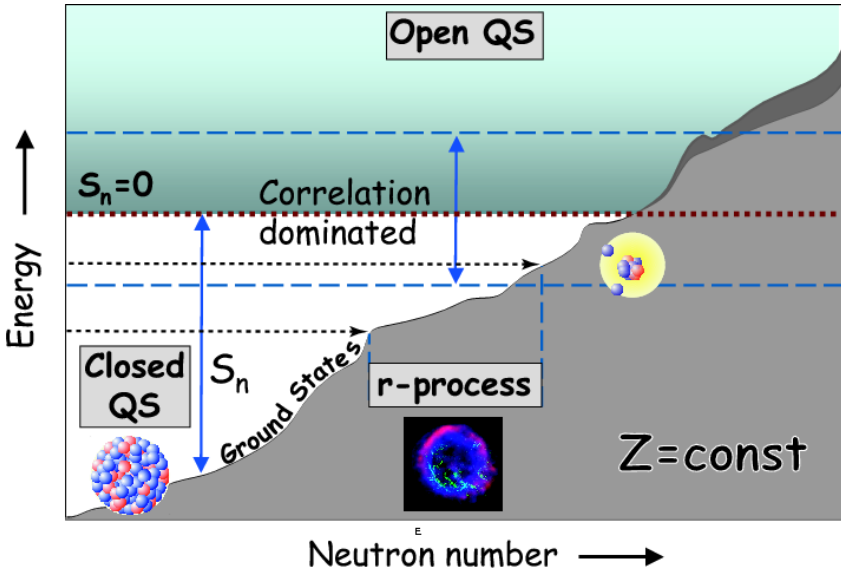
J. Pakarinen et al., PRC 72,  
011304(R) (2005)

# Spectra of resonance states

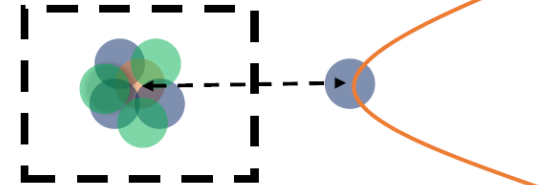
## Energies and resonance widths against particle emissions



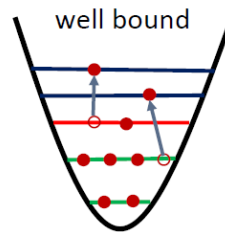
N. Michel, W. Nazarewicz, J. Okolowicz, M. Ploszyczak, Nucl. Phys. A 752, 335c (2005)



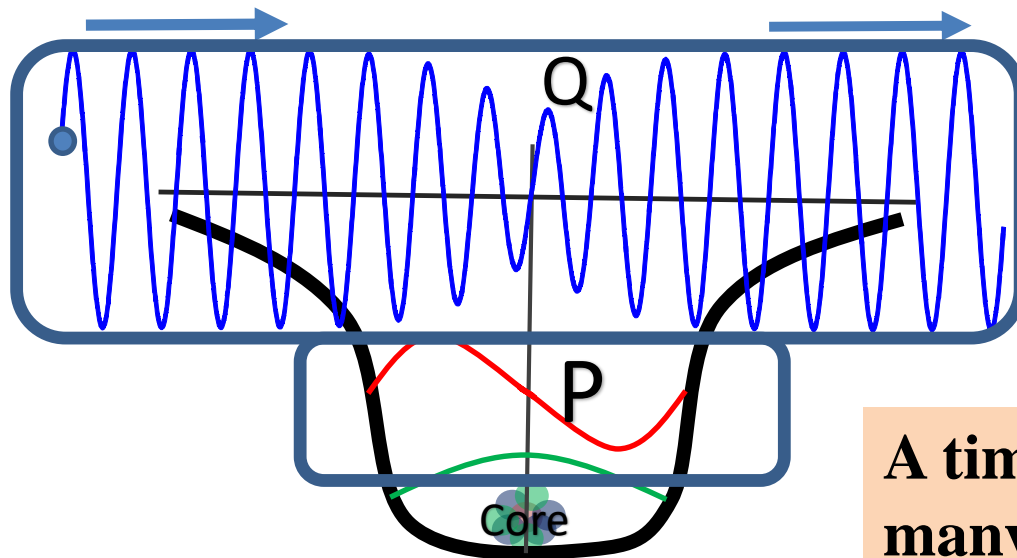
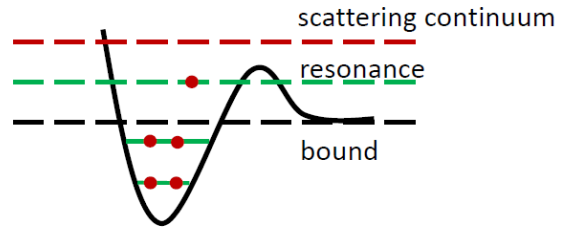
Closed quantum system



Open quantum system



HO basis



**A time-dependent many-body problem**

# Gamow Shell Model

**T. Berggren, Nucl. Phys. A109 (1968) 265**

Single-particle basis in complex- $k$  plane describe bound, resonance and scattering on equal footing.

The radial wave function  $u(r)/r$

$$\frac{d^2 u(k, r)}{dr^2} = \left[ \frac{l(l+1)}{r^2} + \frac{2m}{\hbar^2} U(r) - k^2 \right] u(k, r)$$

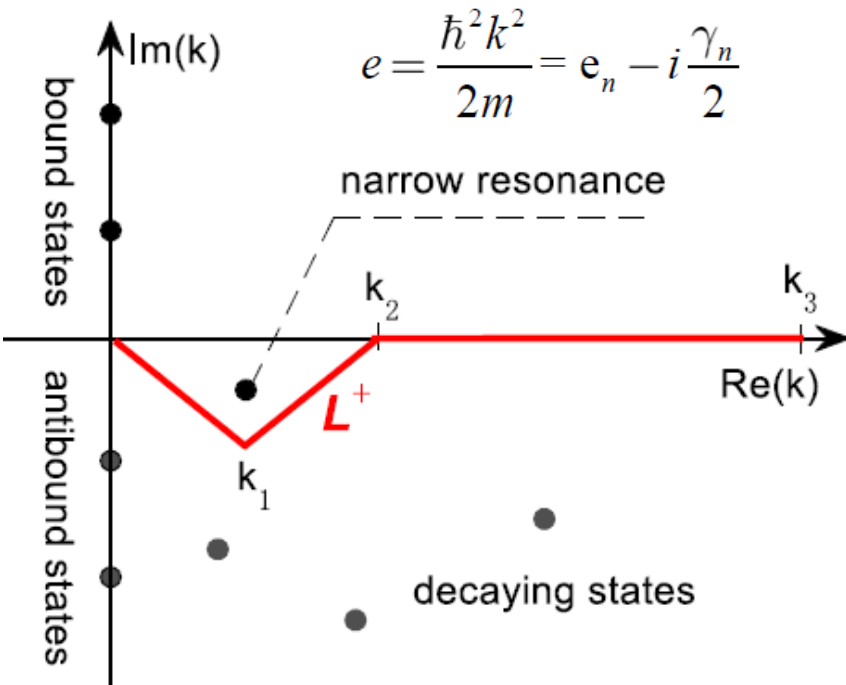
boundary conditions

$$u(0) = 0,$$

$$u(a)O_l'(ka) - u'(a)O_l(ka) = 0$$

$$O_l(kr) \sim e^{i(kr - l\pi/2)}$$

Outgoing solution at large distance



**Orthogonality and Completeness**

$$\delta(r - r') = \sum_n w_n(r, k_n) w_n(r', k_n)$$

$$+ \frac{1}{\pi} \int_{L^+} dq u(r, q) u(r', q)$$

**Discretized**

# Woods-Saxon potential, CD-Bonn, $^{16}\text{O}$ core

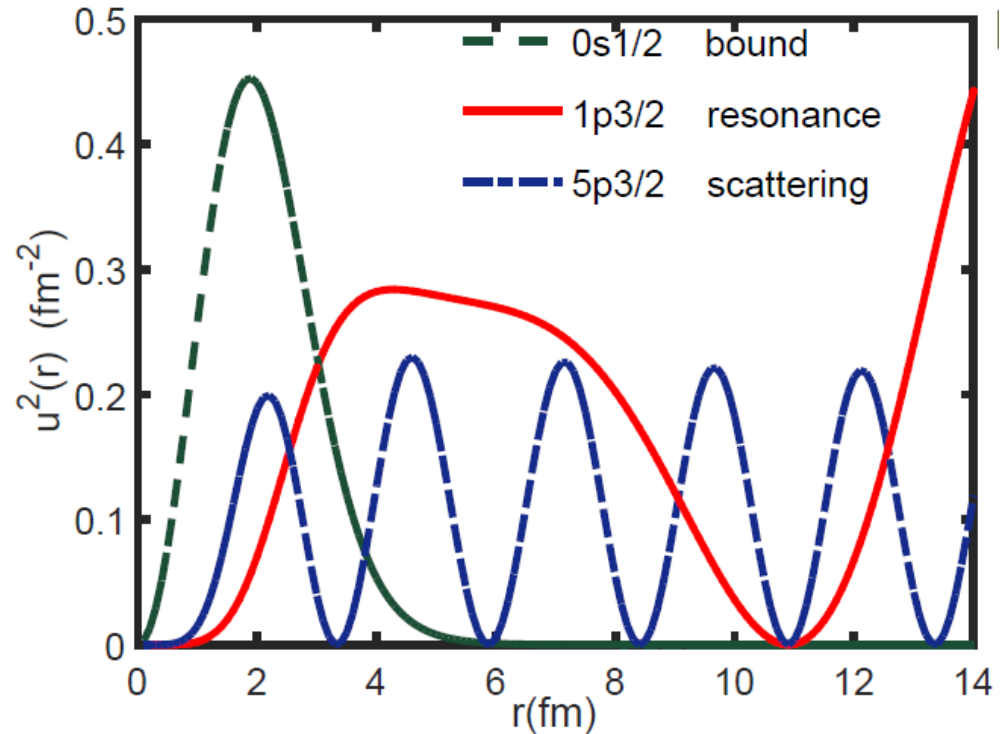
$0d_{3/2}$   $\blacksquare$  1.06-0.089i

-----  $e=0.0$

$1s_{1/2}$   $\blacksquare$  -3.22-0.00i

$0d_{5/2}$   $\blacksquare$  -5.31-0.00i (MeV)

WS SPE's



$^{17}\text{O}$

Details for Berggren basis, see also talks by:

Nazarewicz, Sossez, Ploszajczak, Barrett, Id Betan

$$e = \frac{\hbar^2 k^2}{2m} = e_n - i \frac{\gamma_n}{2}$$

R.J. Liotta *et al.*, PLB 367, 1 (1996)...

used Berggren basis to describe single-particle resonance in nuclei;

later for **two-particle** resonance (Betan *et al.*, PRL 89, 042601 (2002))

using phenomenological potential

## Many-body systems

$$H = \sum_{i=1}^A \frac{p_i^2}{2m} + \sum_{i<j=1}^A v_{ij}^{NN} - \frac{P^2}{2Am} \quad P = \sum_{i=1}^A p_i$$

$$H = \sum_{i=1}^A \frac{p_i^2}{2m} + U + \sum_{i<j=1}^A \left( v_{ij}^{NN} - U - \frac{p_i^2}{2Am} - \frac{p_i p_j}{Am} \right)$$

$$= H_0 + V.$$

$$H_0 = \sum_{i=1}^A \left( \frac{p_i^2}{2m} + U \right)$$

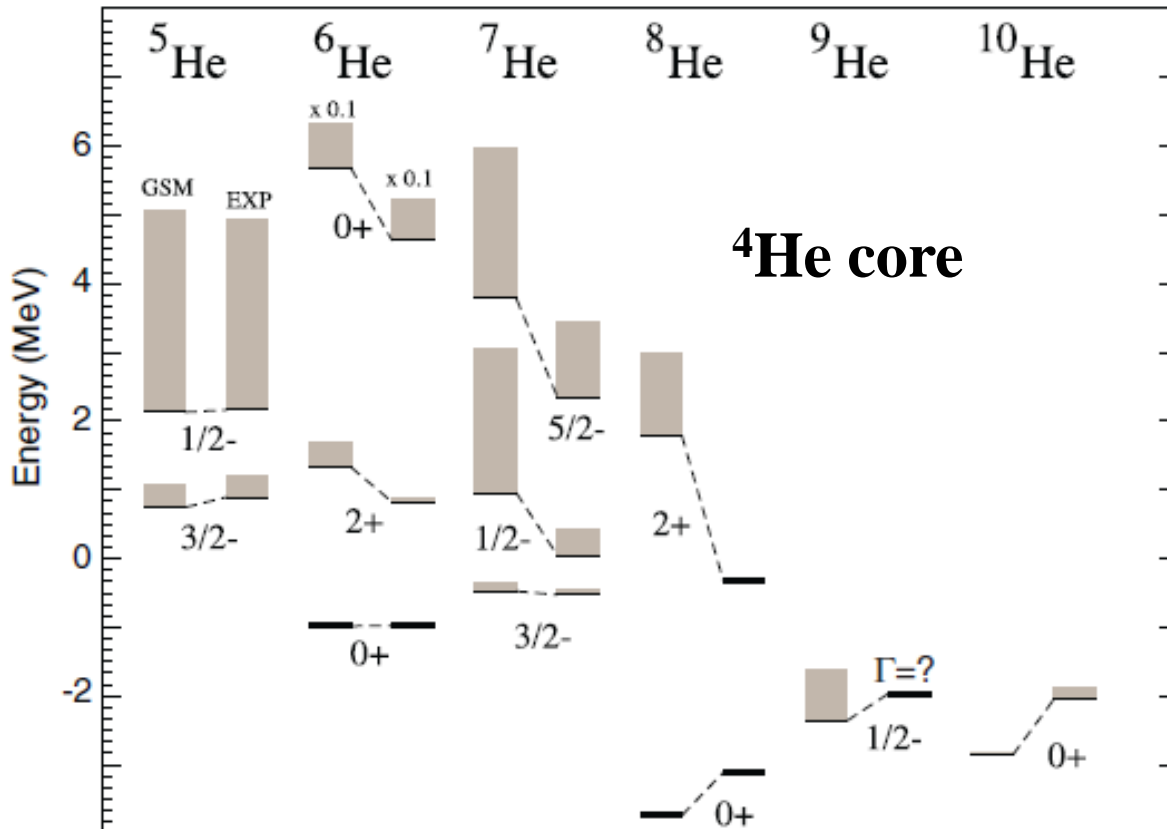
$$E = E_n - i \frac{\Gamma_n}{2}$$

**Michel, Nazarewicz, Płoszajczak, Rotureau *et al.*, 2003--**

$$V = V_{WS} + V_{J,T}(\vec{r}_1, \vec{r}_2)$$

$$V(\mathbf{r}_i, \mathbf{r}_j) = -V_{SGI}^{(J,T)} \exp \left[ - \left( \frac{r_i - r_j}{\mu} \right)^2 \right] \delta(r_i + r_j - 2R_0)$$

$V_{SGI}^{(J)}$  is the strength in the  $JT$  channel

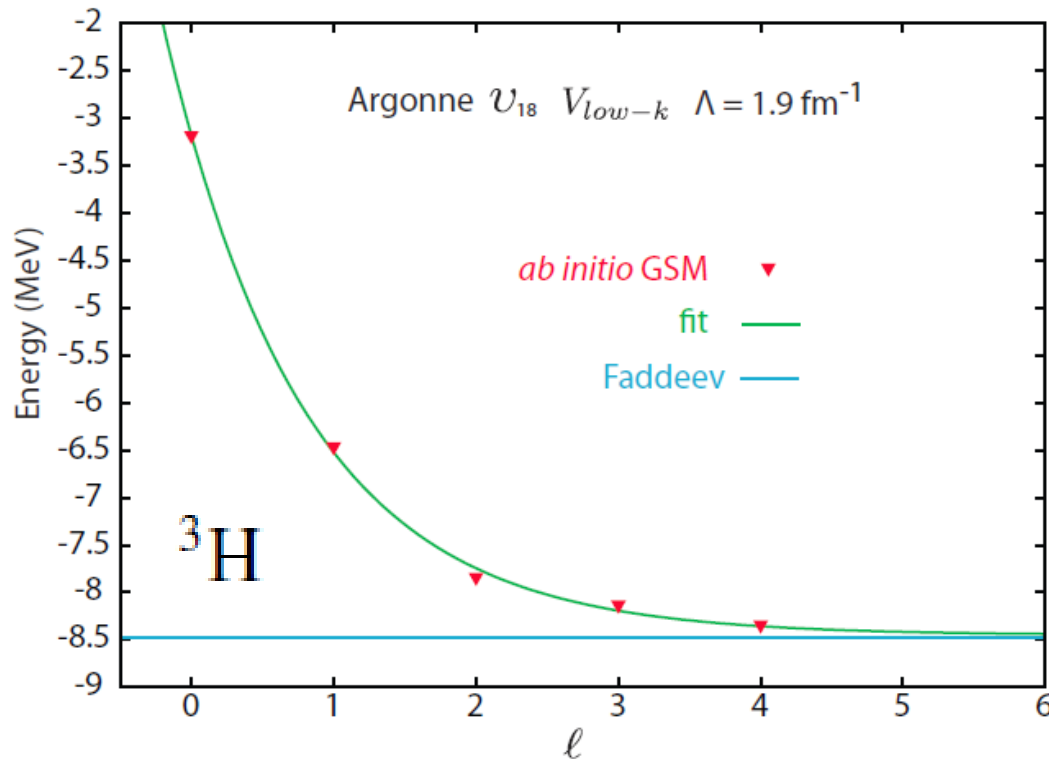


**Michel, Nazarewicz,  
Płoszajczak, Vertse,  
JPG 36 (2009) 013101**

- Hagen, Hjorth-Jensen *et al.*, PRC 73, 064307 (2006): **Core GSM** with **realistic forces**, but neglecting Q-box, applied to two-particle systems (e.g.,  $^{18}\text{O}$ )
- Later, Tsukiyama Hjorth-Jensen, Hagen, PRC 80, 051301 (R) (2009): improving by using Q-box **but no folded-diagrams**.

Papadimitriou *et al.* , Phys. Rev. C 88, 044318 (2013): **realistic forces**

*Ab initio* no-core Gamow shell model for light nuclei





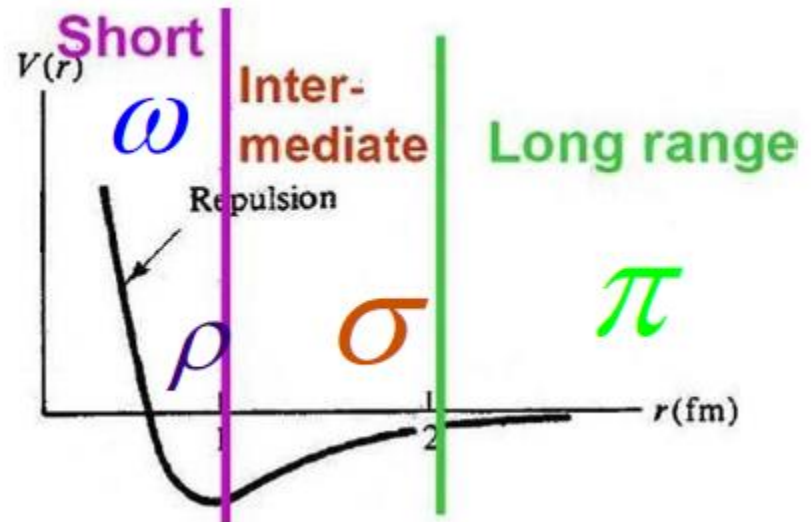
# Gamow shell model with an inert core

1. Start from realistic forces;
2. Take a double magic core

**Q-box + folded diagrams (MBPT)**

3. Calculate resonance spectra

**CD-Bonn**



# CGSM based on realistic nuclear forces

Realistic nuclear forces  $\rightarrow$  Gamow shell model calculations

Taking a doubly closed core

Bare forces:  
Strong repulsion,  
slow convergence

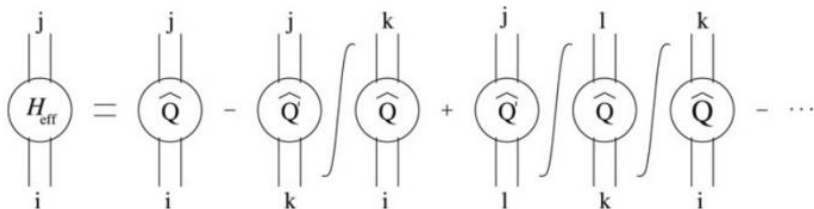
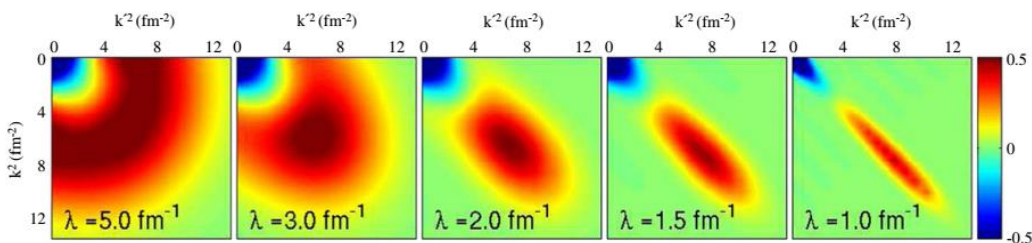
$V_{low k}$  or SRG

To remove hard core,  
but still keep good  
descriptions of NN  
scattering phase shifts

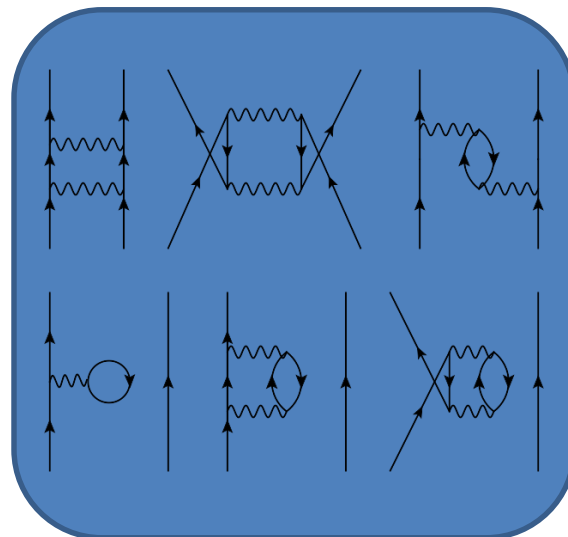
$$\langle \alpha_P | \bar{H}_{\text{eff}} | \alpha_{P'} \rangle = \sum_{\alpha_{P''}} \sum_{\alpha_{P'''}} \sum_{kk'} \sum_{k'k'' \in \mathcal{K}} \langle \alpha_P | \tilde{k}'' \rangle \langle \tilde{k}'' | \alpha_{P''} \rangle \langle \alpha_{P''} | \tilde{k} \rangle E_k \langle \tilde{k} | \alpha_{P'''} \rangle \langle \alpha_{P'''} | \tilde{k}' \rangle \langle \tilde{k}' | \alpha_{P'} \rangle$$

$$\frac{dH_\lambda}{d\lambda} = -\frac{4}{\lambda^5} [[T_{\text{rel}}, H_\lambda], H_\lambda]$$

Q

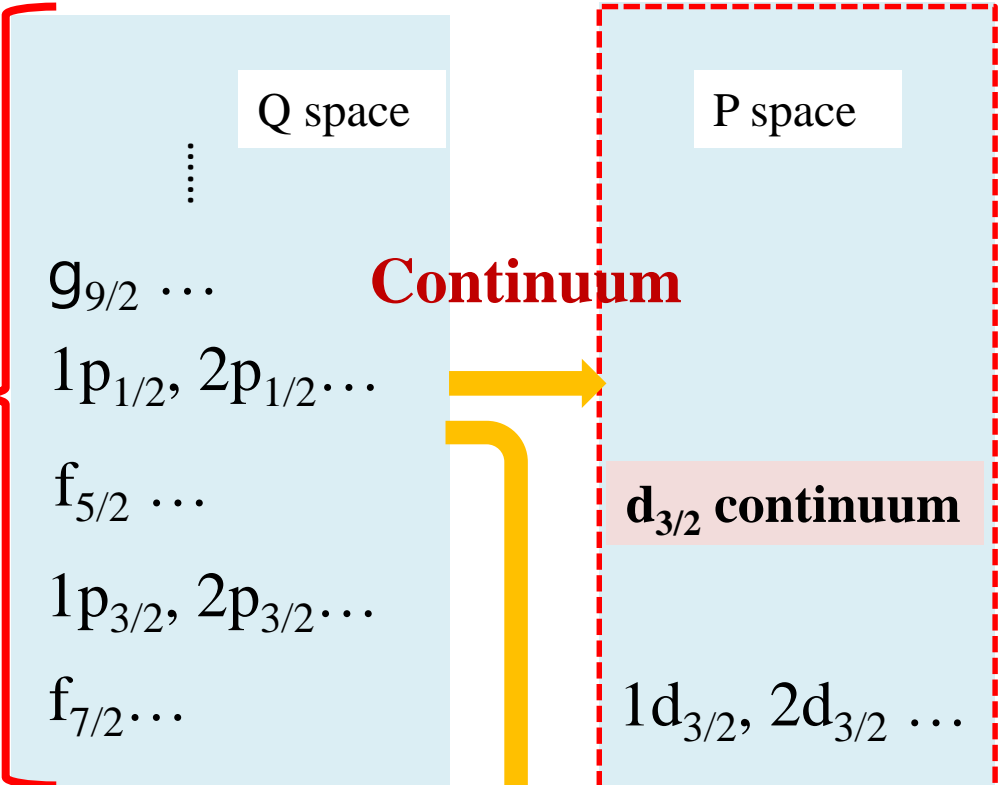


$$\hat{Q}(E) = PVP + PVQ \frac{1}{E - H} QVP;$$

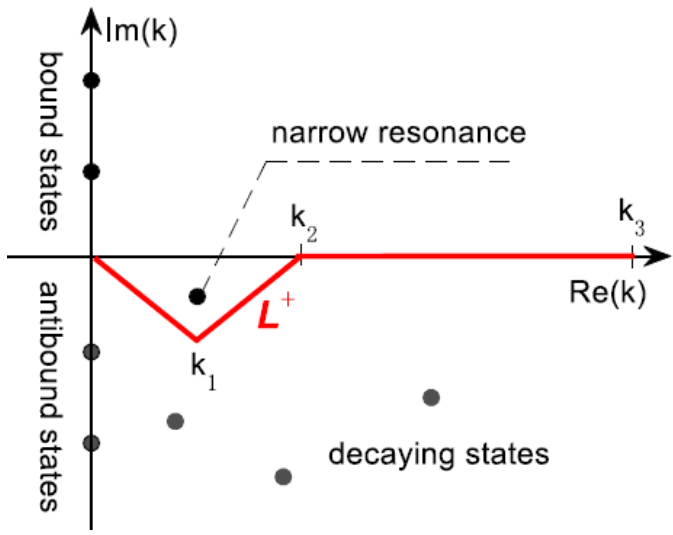


Non-degenerate extended Kuo-Krenciglowa folded-diagram method (EKK) by Takayanagi, NPA 852, 61 (2011);

Continuum



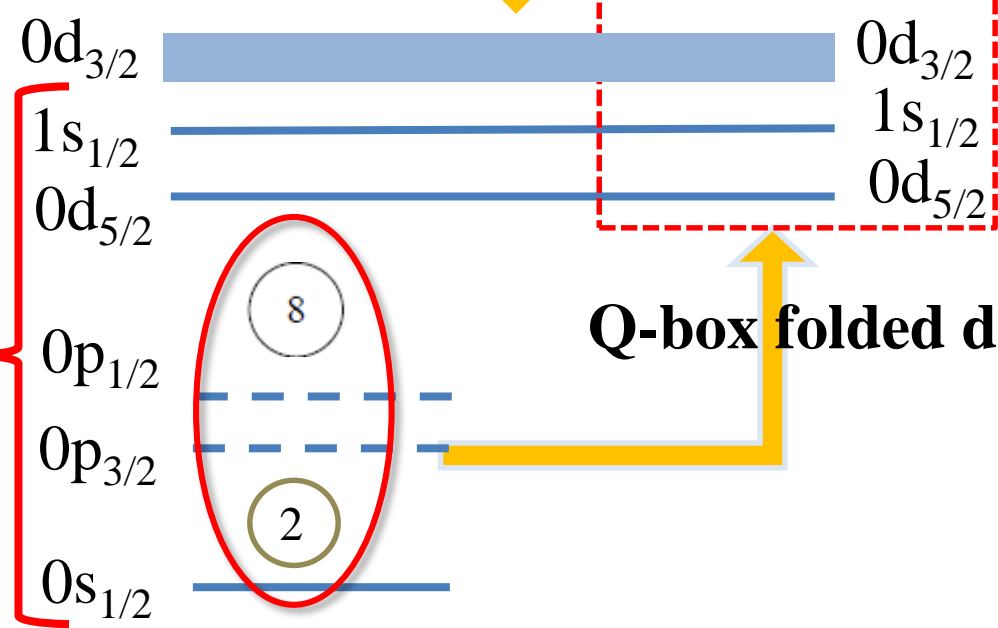
# Model space



**For each given partial wave  
Berggren Completeness Relation**

$$\sum_n |u_n\rangle \langle \tilde{u}_n| + \int_{L^+} |u_k\rangle \langle \tilde{u}_k| dk = 1$$

Bound



**Q-box folded diagrams in complex-k basis**

$$\varepsilon_n = \frac{\hbar^2 k^2}{2m} = e_n - i \frac{\gamma_n}{2}$$

We need to establish the effective Hamiltonian in the model space P, based on realistic forces  
Q-box folded diagram method in complex-energy space

1.  $V_{\text{low-k}}$
2. Using Brody-Mshinsky brackets, NN interaction which is in relative and CoM coordinates is transferred into the laboratory (HO basis)

Truncated by  $N_{\text{shell}} \sim 12$ , an approximate completeness

$$\sum_{\alpha \leq \beta}^{N_{\text{shell}}} |\alpha\beta\rangle\langle\alpha\beta| = \mathbf{1}$$

where  $|\alpha\beta\rangle$  is the two-particle states in HO basis

In HO basis, NN matrix elements:

$$V_{osc} = \sum_{\alpha \leq \beta} \sum_{\gamma \leq \delta} |\alpha\beta\rangle\langle\alpha\beta| V_{\text{low-k}} |\gamma\delta\rangle\langle\gamma\delta|$$

# In Berggren basis

$$\langle ab|V|cd\rangle \approx \sum_{\alpha \leq \beta} \sum_{\gamma \leq \delta} \langle ab|\alpha\beta\rangle \langle \alpha\beta|V_{low-k}|\gamma\delta\rangle \langle \gamma\delta|cd\rangle$$

where  $|ab\rangle$  is two-particle states constructed with Berggren s.p. basis

For identical particles (pp or nn), the expansion coefficients are

$$\langle ab|\alpha\beta\rangle = \frac{\langle a|\alpha\rangle \langle b|\beta\rangle - (-1)^{J-j_\alpha-j_\beta} \langle a|\beta\rangle \langle b|\alpha\rangle}{\sqrt{(1 + \delta_{ab})(1 + \delta_{\alpha\beta})}}$$

For np:  $\langle ab|\alpha\beta\rangle = \langle a|\alpha\rangle \langle b|\beta\rangle$

one-body expansion coefficients are calculated with

$$\langle a|\alpha\rangle = \int dr r^2 u_a(r) R_\alpha \delta_{l_a l_\alpha} \delta_{j_a j_\alpha} \delta_{t_a t_\alpha}$$

$u_a(r)$  – Berggren basis;  $R_\alpha$  - HO basis

$$H = \sum_{i=1}^A \frac{p_i^2}{2m} + U + \sum_{i < j=1} \left( v_{ij}^{NN} - U - \frac{p_i^2}{2Am} - \frac{\mathbf{p}_i \mathbf{p}_j}{Am} \right)$$

$$= H_0 + V.$$

using the exterior complex scaling technique

$$\frac{p_i^2}{2Am}$$

$$\frac{\mathbf{p}_i \mathbf{p}_j}{Am}$$

Q-box

$$\hat{Q}(E) = PVP + PVQ \frac{1}{E - QHQ} QVP$$

$$\hat{Q}(E) = PVP + PV \frac{Q}{E - QH_0Q} VP + PV \frac{Q}{E - QH_0Q} VP \frac{Q}{E - QH_0Q} VP + \dots$$

2<sup>nd</sup> order perturbation                      3<sup>rd</sup> order perturbation

In a degenerate s.p. space, E can be assumed approximately,  $2e_i$ , E is the starting energy

Q-box folded diagrams

$$V_{eff} = \hat{Q}(\epsilon_0) - \hat{Q}'(\epsilon_0) \int \hat{Q}(\epsilon_0) + \hat{Q}'(\epsilon_0) \int \hat{Q}(\epsilon_0) \int \hat{Q}(\epsilon_0) \dots$$

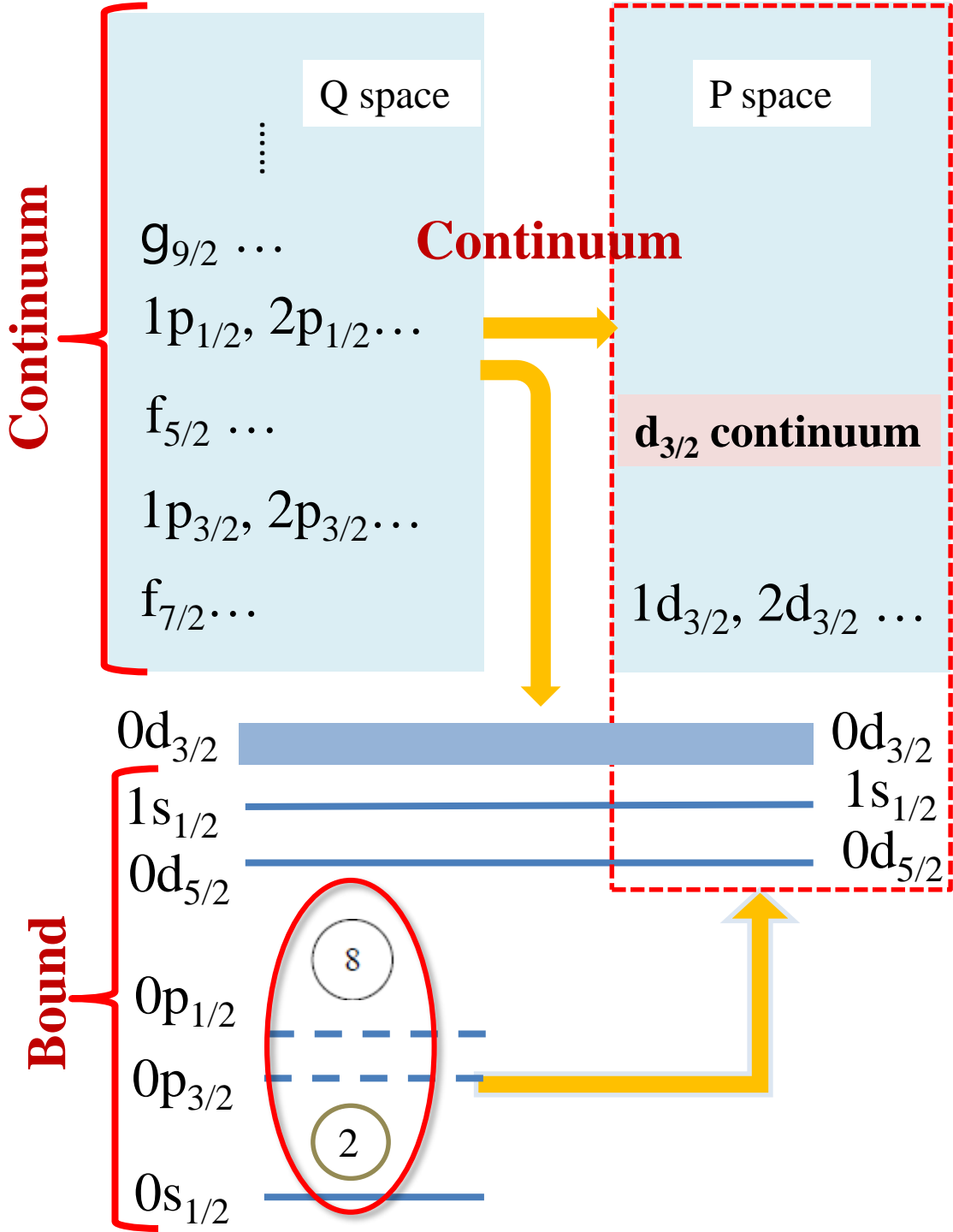
$$V_{eff} = \hat{Q}(\epsilon_0) + \sum_{k=1}^{\infty} \hat{Q}_k(\epsilon_0) [V_{eff}]^k \quad \epsilon_0 = \epsilon_c + \epsilon_d \quad (\text{i.e., the starting energy } E)$$

Q-box derivatives

$$\hat{Q}_k(E) = \frac{1}{k!} \frac{d^k \hat{Q}(E)}{dE^k}$$

$$= (-1)^k PVQ \frac{1}{(E - QHQ)^{k+1}} QVP$$

Kuo-Krenciglowa (KK) method



# Model space

$$P + Q = 1$$

P is the model space

Q is the excluded space  
(including the core)

The Berggren space must be  
nondegenerate

# Q-Box folded diagrams for nondegenerate space: Extended Kuo-Krenciglowa (EKK)

$$H_{eff} = PH_0P + PH_1Q \frac{1}{E - QHQ} QH_1P$$

$$H_{eff} - E_x = PH_0P + V_{eff} - E_x \quad (\text{Meaning: expanded around } E_x)$$

$$\tilde{H}_{eff} = PH_0P - E_x + \hat{Q}(E_x) + \sum_{k=1}^{\infty} \hat{Q}_k(E_x) [\tilde{H}_{eff}]^k$$

$$PH_0P - E_x = e(a) + e(b) - w$$

$\omega$  - starting energy

$$\tilde{H}_{eff}^{[n]} = \tilde{H}_{BH}(E) + \sum_{k=1}^{\infty} \hat{Q}_k(E) \{\tilde{H}_{eff}^{[n-1]}\}^k$$

$$\hat{Q}_k(E) = \frac{1}{k!} \frac{d^k \hat{Q}(E)}{dE^k}$$

$$= (-1)^k PVQ \frac{1}{(E - QHQ)^{k+1}} QVP$$

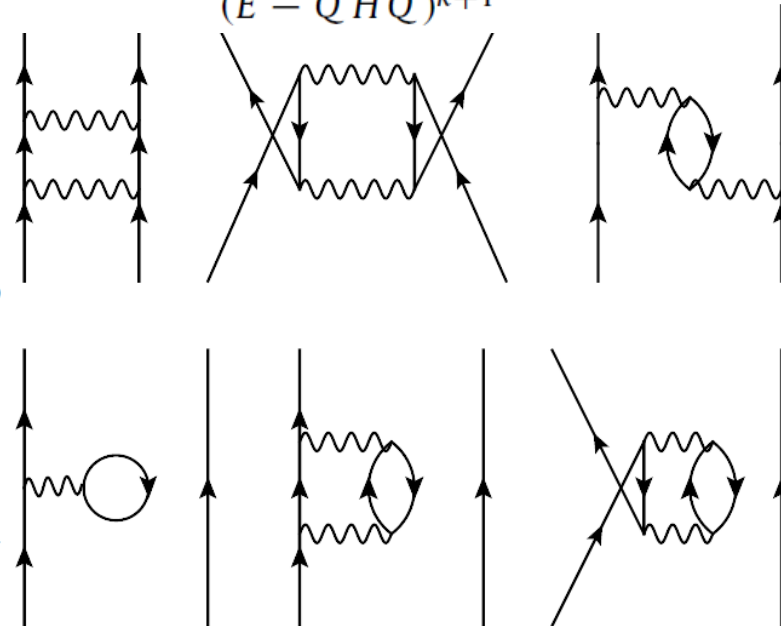
$\tilde{H}_{eff}^{[n]}$  stands for  $\tilde{H}_{eff} = H_{eff} - E$  at the  $n$ -th iteration

$\tilde{H}_{BH} = H_{BH}(E) - E$  is the Block-Horowitz Hamiltonian an energy  $E$ , with

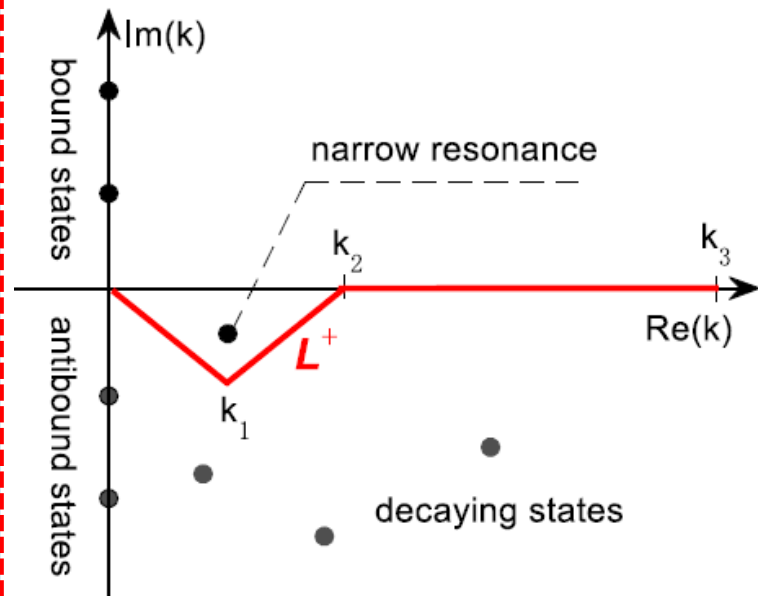
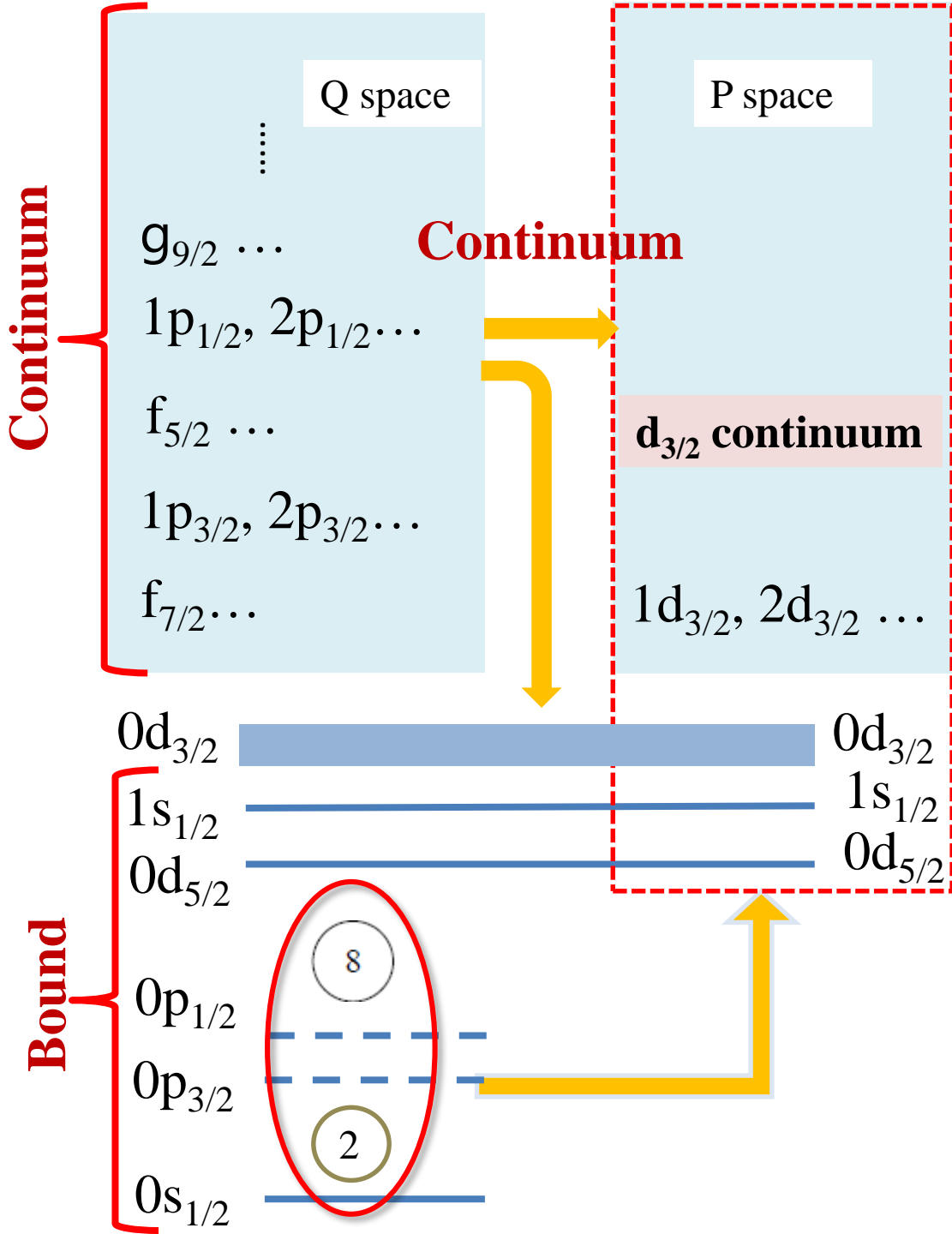
$$H_{BH} = PHP + PHQ \frac{1}{E - QHQ} QHP$$

The effective Hamiltonian is obtained by  $H_{eff} = \tilde{H}_{eff} + E$

effective interaction



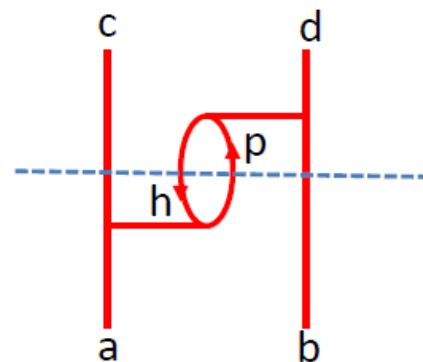




$$H_{eff} = PH_0P + PH_1Q \frac{1}{E - QHQ} QH_1P$$

$$\frac{1}{E - [(e_c + e_b) + (e_p - e_h)]}$$

$$= \frac{1}{[E - (e_c + e_b)] + (e_h - e_p)}$$



## CoM correction

$$H = \sum_{i=1}^A \left( \frac{p_i^2}{2m} + U \right) + \sum_{i < j} \left( v_{ij} - U - \frac{p_i^2}{2Am} - \frac{\mathbf{p}_i \cdot \mathbf{p}_j}{Am} \right)$$

Lawson method is no longer valid

Wave functions?

In cluster orbital coordinates (COSM):  $R, r_i$

Y. Suzuki, K. Ikeda, RC 38, 410 (1988).

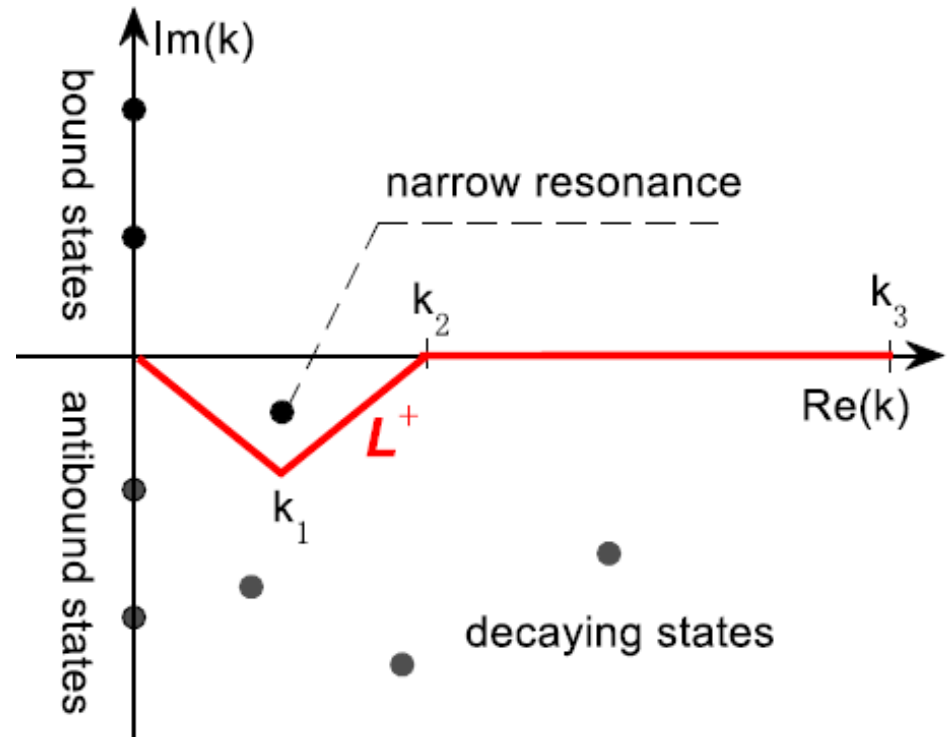
**But with realistic forces:**

$$\langle ab|V|cd \rangle \approx \sum_{\alpha \leq \beta} \sum_{\gamma \leq \delta} \langle ab|\alpha\beta \rangle \langle \alpha\beta|V_{low-k}|\gamma\delta \rangle \langle \gamma\delta|cd \rangle$$

$$\langle ab|\alpha\beta \rangle = \langle a|\alpha \rangle \langle b|\beta \rangle \quad \langle a|\alpha \rangle = \int dr r^2 u_a(r) R_\alpha \delta_{l_a l_\alpha} \delta_{j_a j_\alpha} \delta_{t_a t_\alpha}$$

In our CGSM calculations, for low-lying states we assume small CoM effects due to wave functions expressed in the laboratory coordinates.

Convergence against  
discretization number  $N_L$



$$\tilde{E}_n = E_n - i\Gamma/2$$

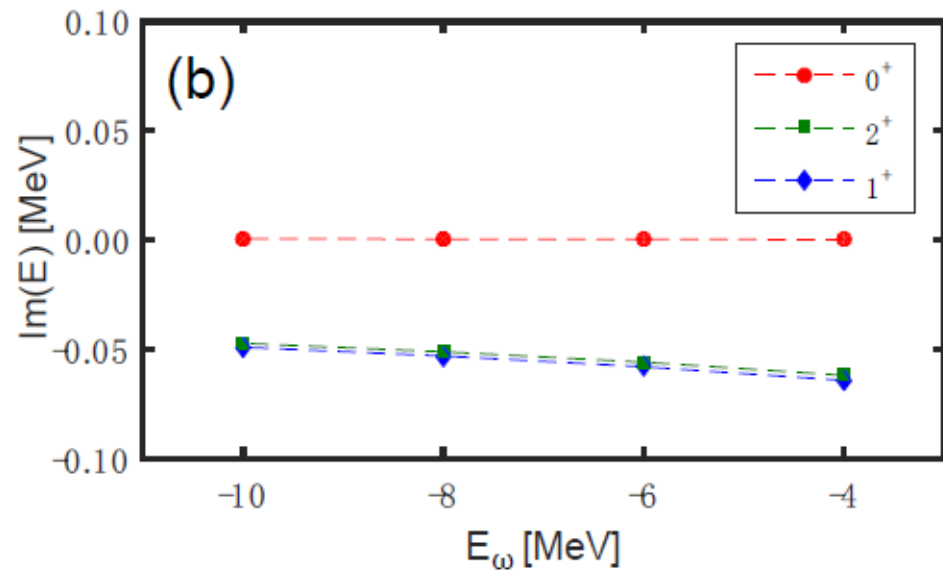
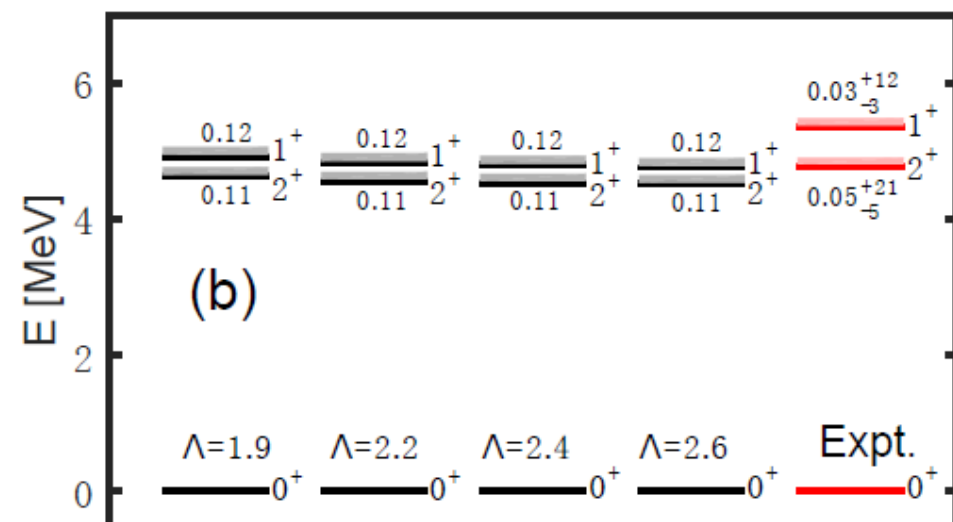
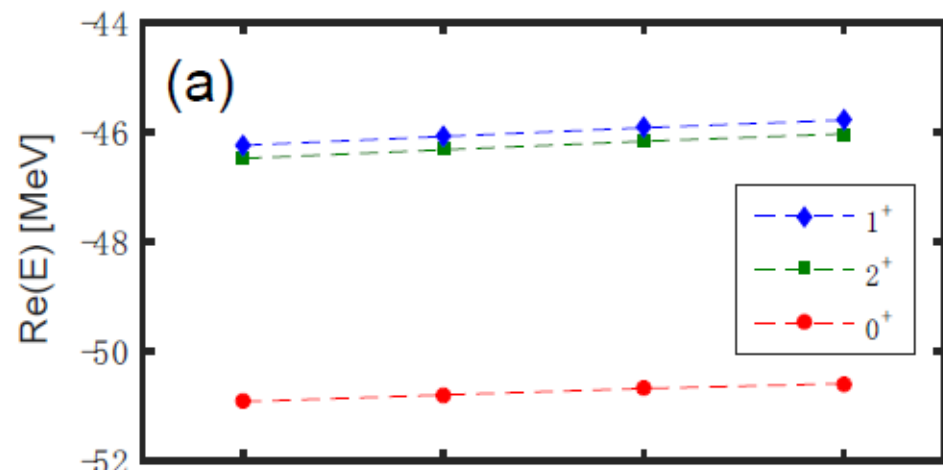
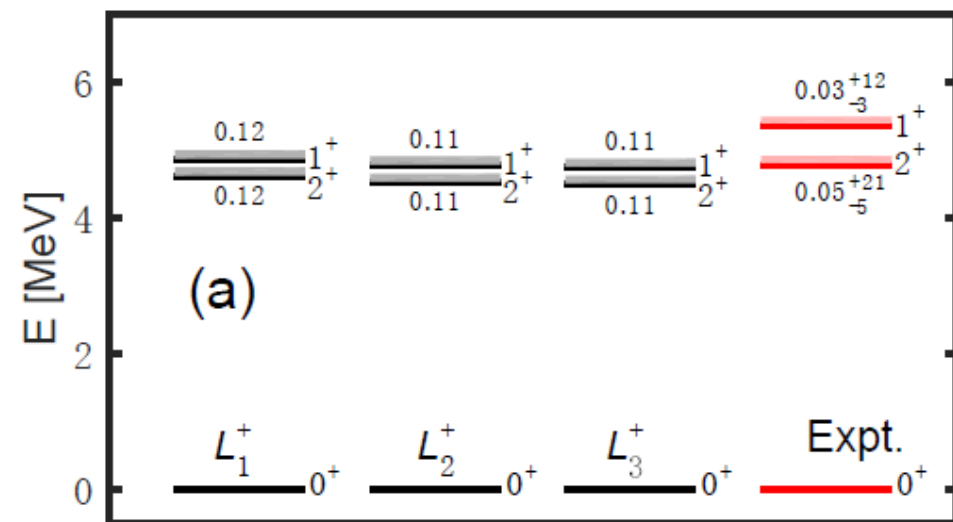
$^{24}\text{O}$

$\Lambda = 2.6 \text{ fm}^{-1}$

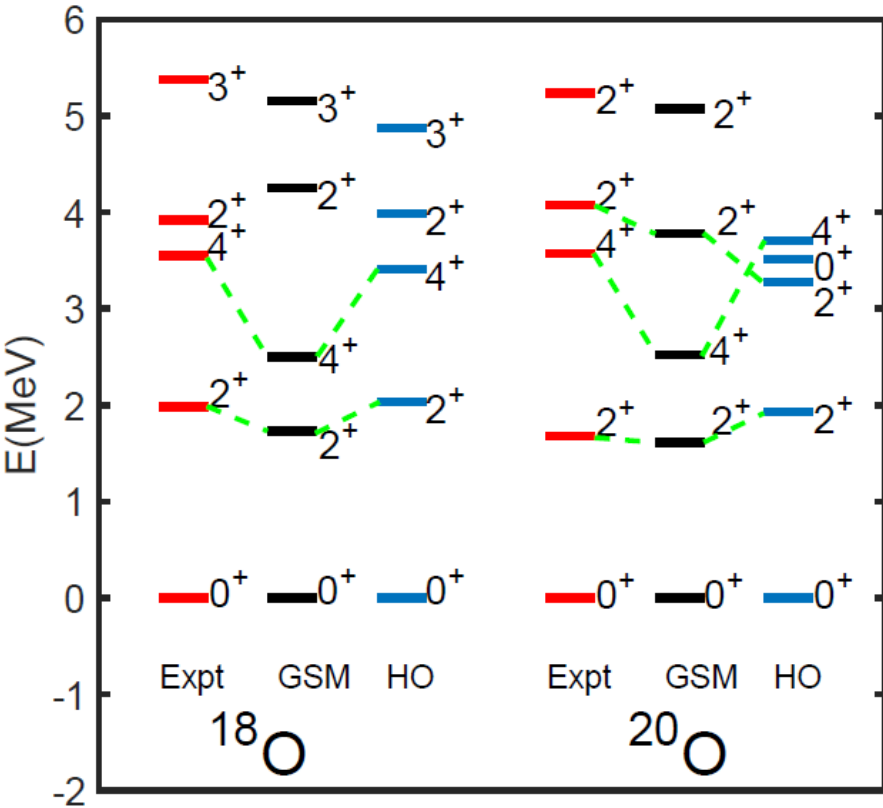
$N_L$	$0^+$	$2^+$	$1^+$
16	$-50.642 + 0.013i$	$-46.172 - 0.004i$	$-45.922 - 0.009i$
18	$-50.716 + 0.002i$	$-46.262 - 0.046i$	$-46.017 - 0.049i$
20	$-50.711 - 0.001i$	$-46.219 - 0.054i$	$-45.976 - 0.056i$
22	$-50.712 + 0.000i$	$-46.218 - 0.053i$	$-45.974 - 0.056i$

# Convergences of spectroscopic calculations

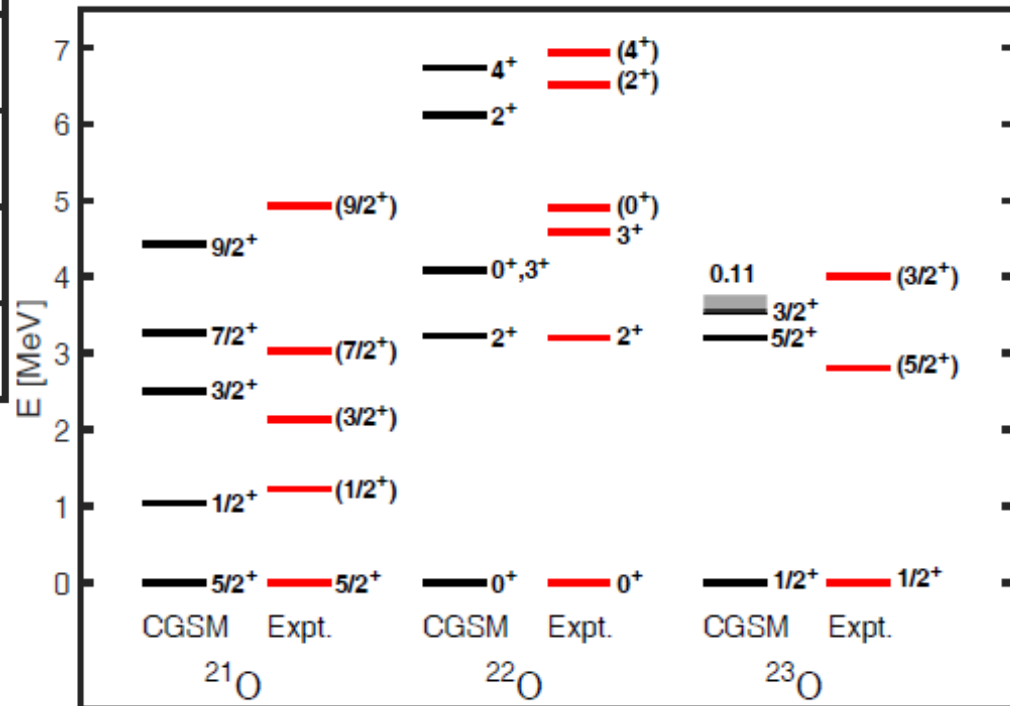
$^{24}\text{O}$

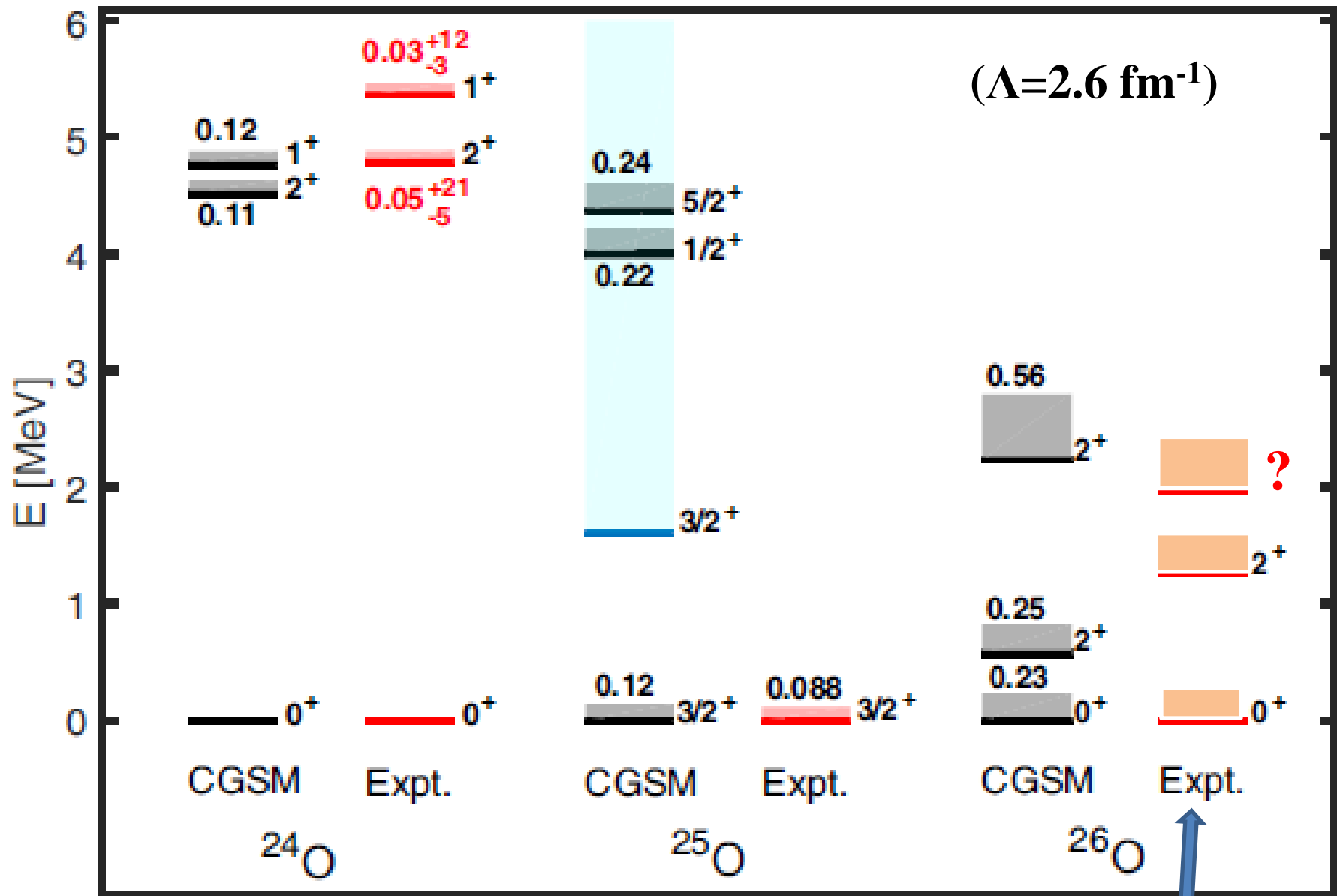


# CD-Bonn CGSM, compared with conventional H.O. SM



Hard cutoff  $\Lambda=2.6 \text{ fm}^{-1}$  to reduce 3NFs





Y. Kondo *et al.*, PRL 116, 102503 (2016)

# Binding energies, one-neutron separation energies

$$\tilde{e}_n = e_n - i\gamma_n/2$$

**0d<sub>3/2</sub>** **1.06-0.089i** **0.94-0.048i**  
 -----  
 -----  
 -----  
**e=0.0**

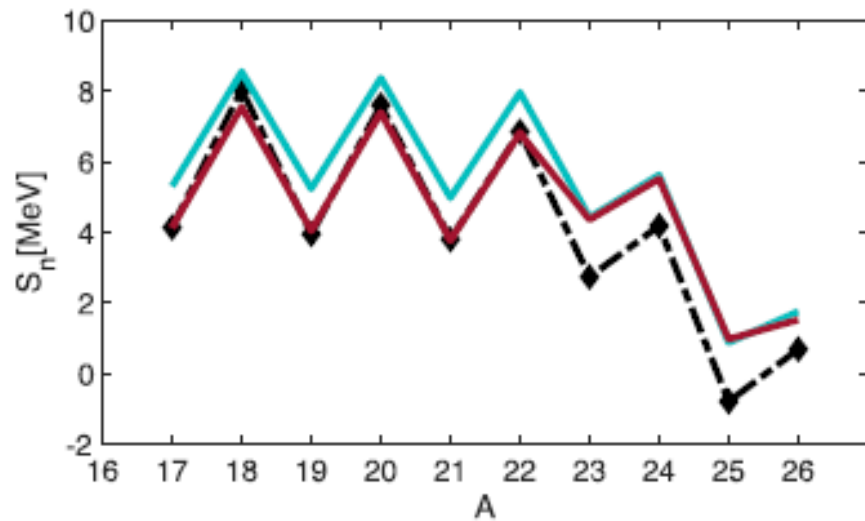
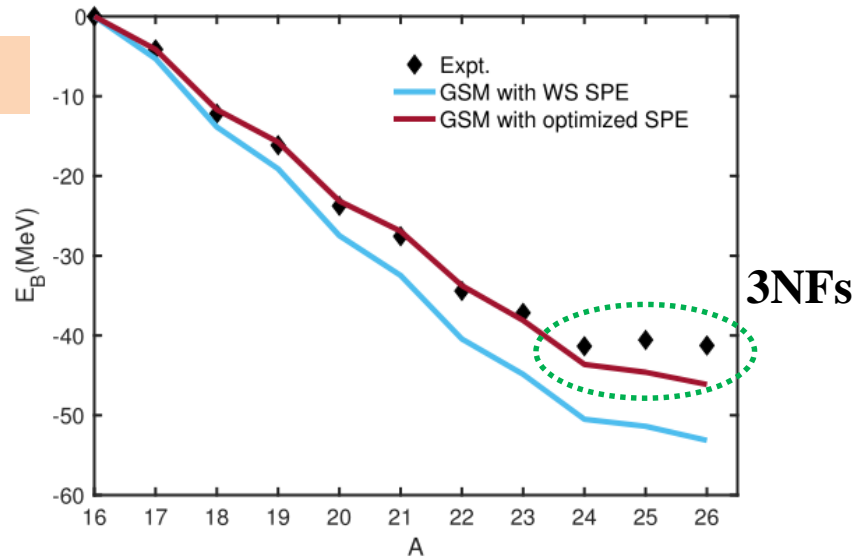
**1s<sub>1/2</sub>** ----- **-3.22-0.00i** ----- **-3.27-0.00i**  
 -----  
 ----- **-4.14-0.00i**

**0d<sub>5/2</sub>** ----- **-5.31-0.00i (MeV)**

WS SPE's

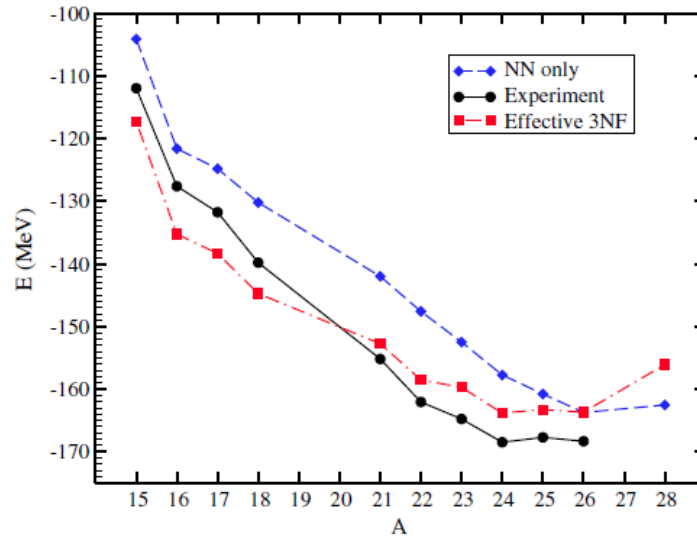
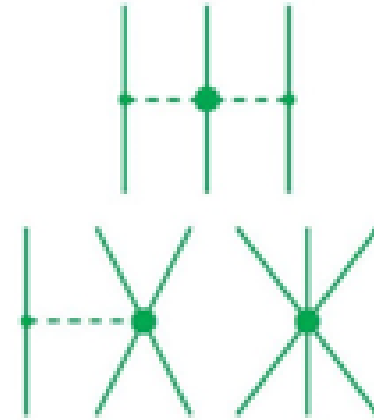
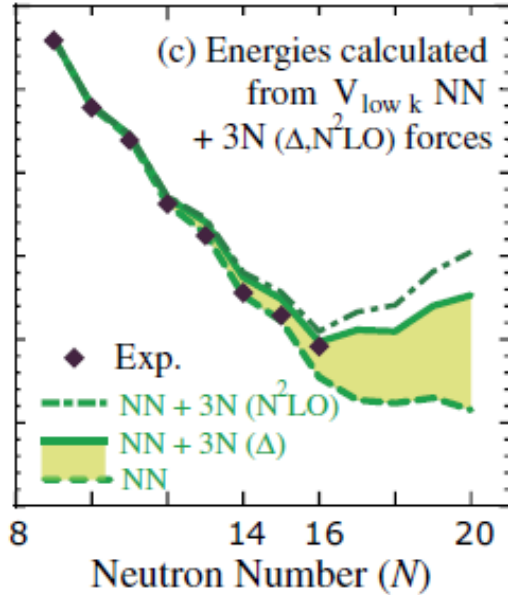
Expt. SPE's

[Extracted from <sup>17</sup>O,  
 by Michel *et al.*, PRC 67, 054311 (2003)]

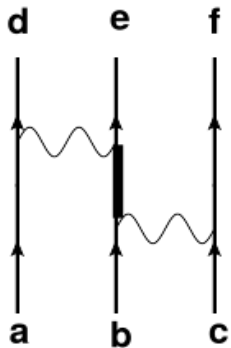
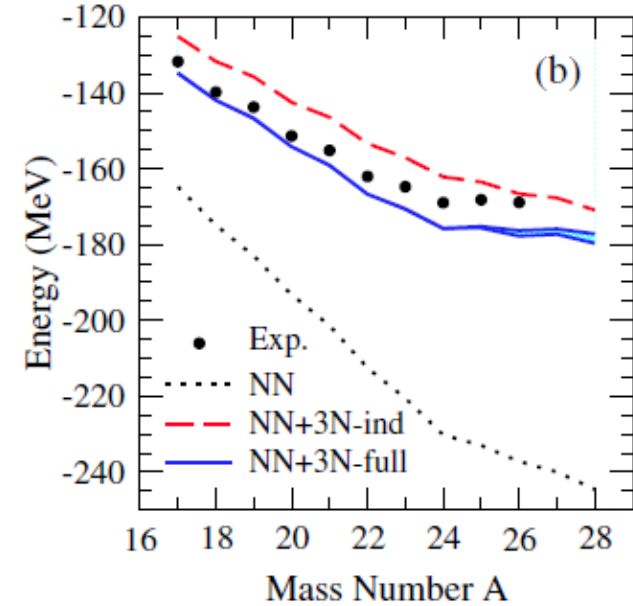


# 3NFs are important for binding energy calculations

Otsuka *et al.*, PRL 105, 032501 (2010): SM,  $N^3LO$ ,  $N^2LO$  3NF



Bogner *et al.*, PRL 113, 142501 (2014): In-medium SRG,  $N^3LO$ , 3NF=induced / initial ( $N^2LO$ )

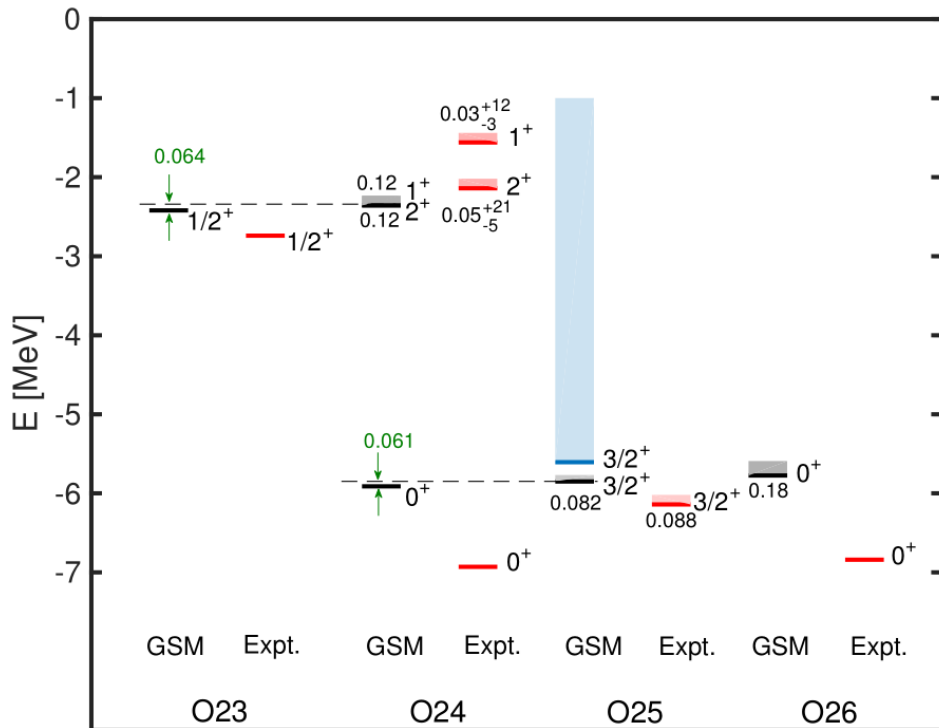


G. Hagen *et al.*, PRL 108, 242501 (2012): Continuum CC,  $N^3LO$ ,  $N^2LO$  3NF

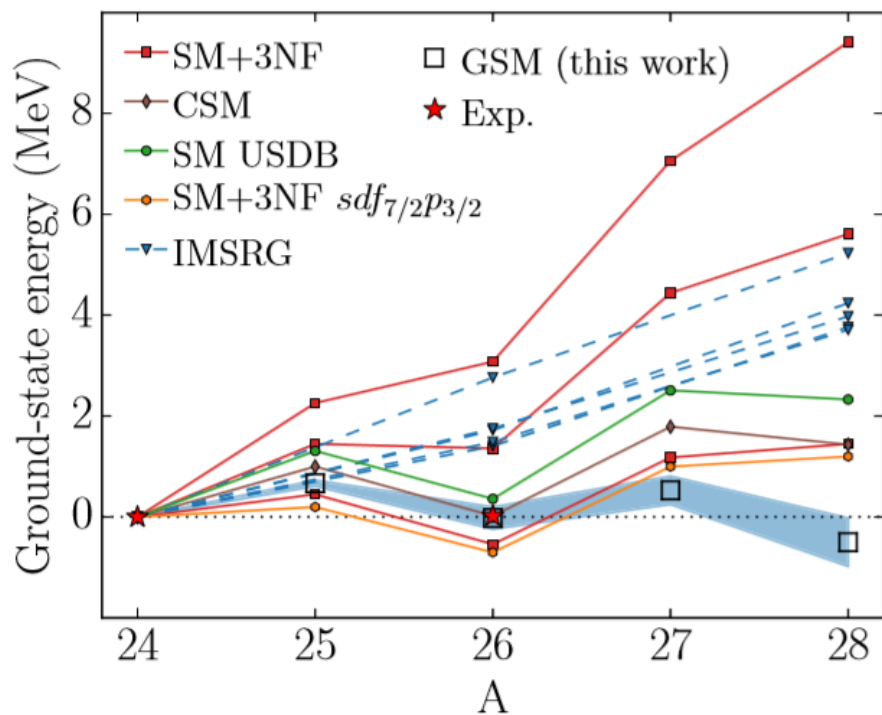


# $^{22}\text{O}$ core (N=14 closed shells)

## 3NF effects



K. Fosse, J. Rotureau, N. Michel, W. Zarewicz, PRC 96, 024308 (2017)



# Summary

**Realistic nuclear forces (CD Bonn)**



**Renormalization by  $V_{\text{low-}k}$**



**Many-body solutions by CGSM**

**Full Q-box folded diagrams in nondegenerate complex- $k$  space, which includes contributions from core polarization and excluded space.**

**✓ Successfully applied to excitation spectra of weakly-bound or unbound oxygen isotopes.**



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# Resonance and continuum Gamow shell model with realistic nuclear forces



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*Thank you for your attention*

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