

Ab initio coupled-cluster computations of nuclei

Thomas Papenbrock



and

OAK RIDGE NATIONAL LABORATORY

G. Hagen

D. J. Dean

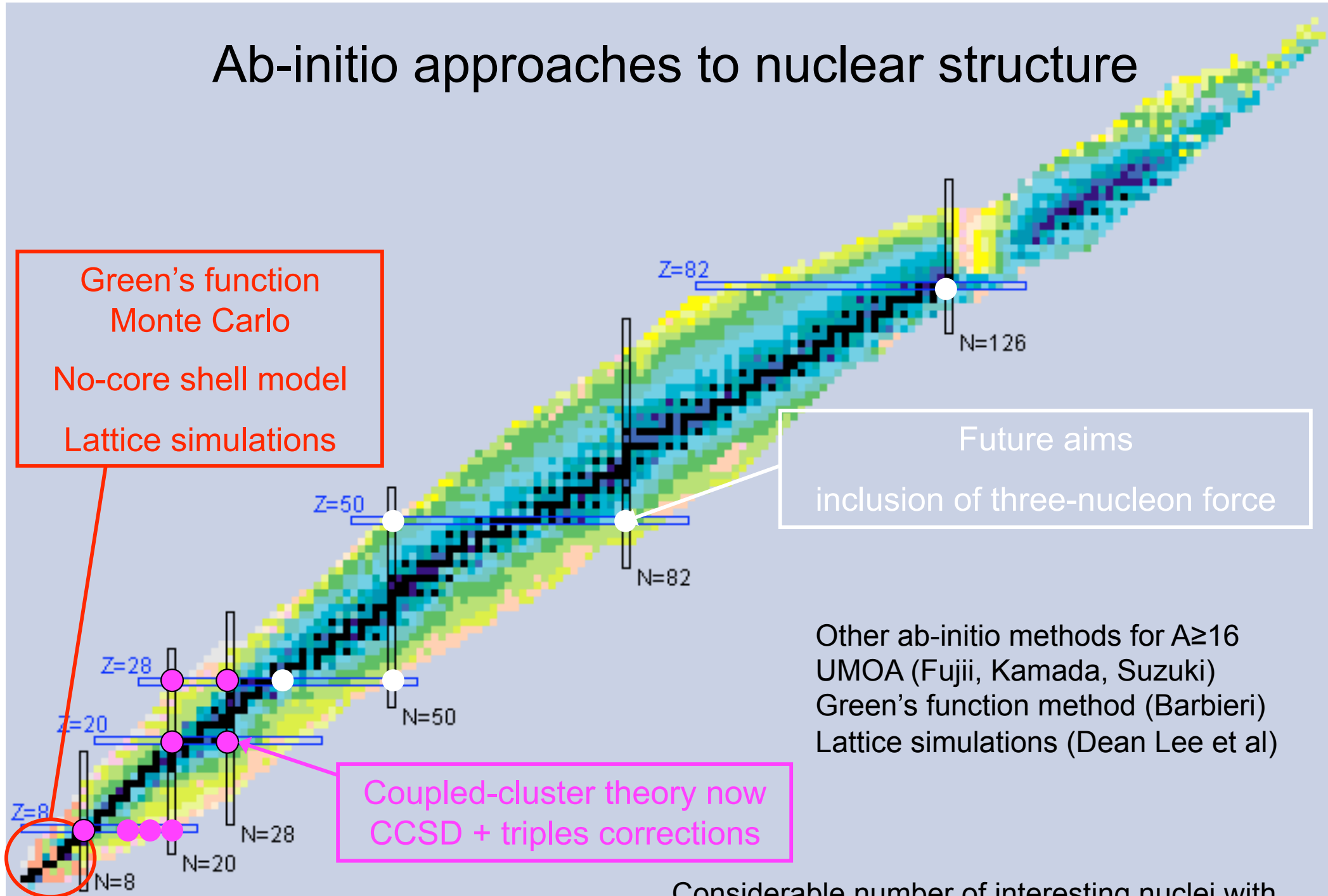
M. Hjorth-Jensen

B. Velamuri Asokan

Overview

1. Introduction
2. Medium-mass nuclei – saturation properties of NN interactions
[Hagen, TP, Dean, Hjorth-Jensen, Phys. Rev. Lett. 101, 092502 (2008)]
3. Practical solution to the center-of-mass problem
[Hagen, TP, Dean, Phys. Rev. Lett. 103, 062503 (2009)]
4. Proton-halo state in ^{17}F
[Hagen, TP, Hjorth-Jensen, arXiv:1003.1995]
5. Does ^{28}O exist?
[Hagen, TP, Dean, Horth-Jensen, Velamur Asokan, Phys. Rev. C 80, 021306(R) (2009)]

Ab-initio approaches to nuclear structure



Coupled-cluster method (CCSD)

Ansatz: $|\Psi\rangle = e^T |\Phi\rangle$

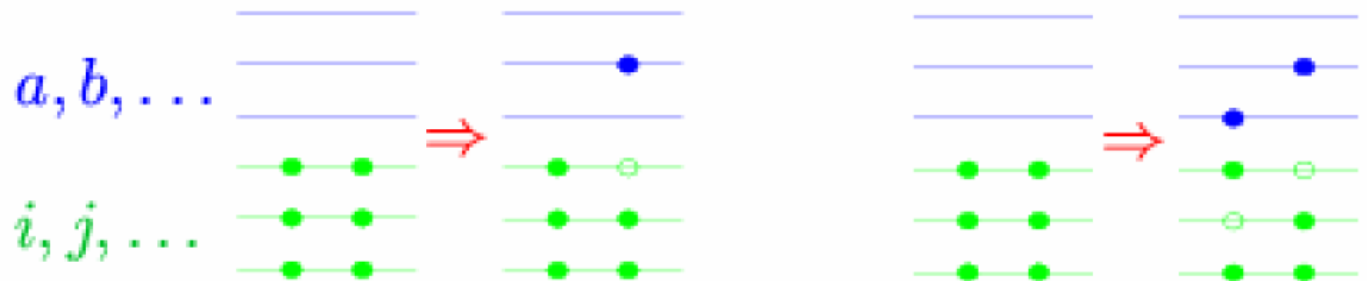
$$T = T_1 + T_2 + \dots$$

$$T_1 = \sum_{ia} t_i^a a_a^\dagger a_i$$

$$T_2 = \sum_{ijab} t_{ij}^{ab} a_a^\dagger a_b^\dagger a_j a_i$$

- ☺ Scales gently (polynomial) with increasing problem size $\mathcal{O}(u^4)$.
- ☺ Truncation is the only approximation.
- ☺ Size extensive (error scales with A)

Correlations are *exponentiated* 1p-1h and 2p-2h excitations. Part of np-nh excitations included!



Coupled cluster equations

$$E = \langle \Phi | \bar{H} | \Phi \rangle$$

$$0 = \langle \Phi_i^a | \bar{H} | \Phi \rangle$$

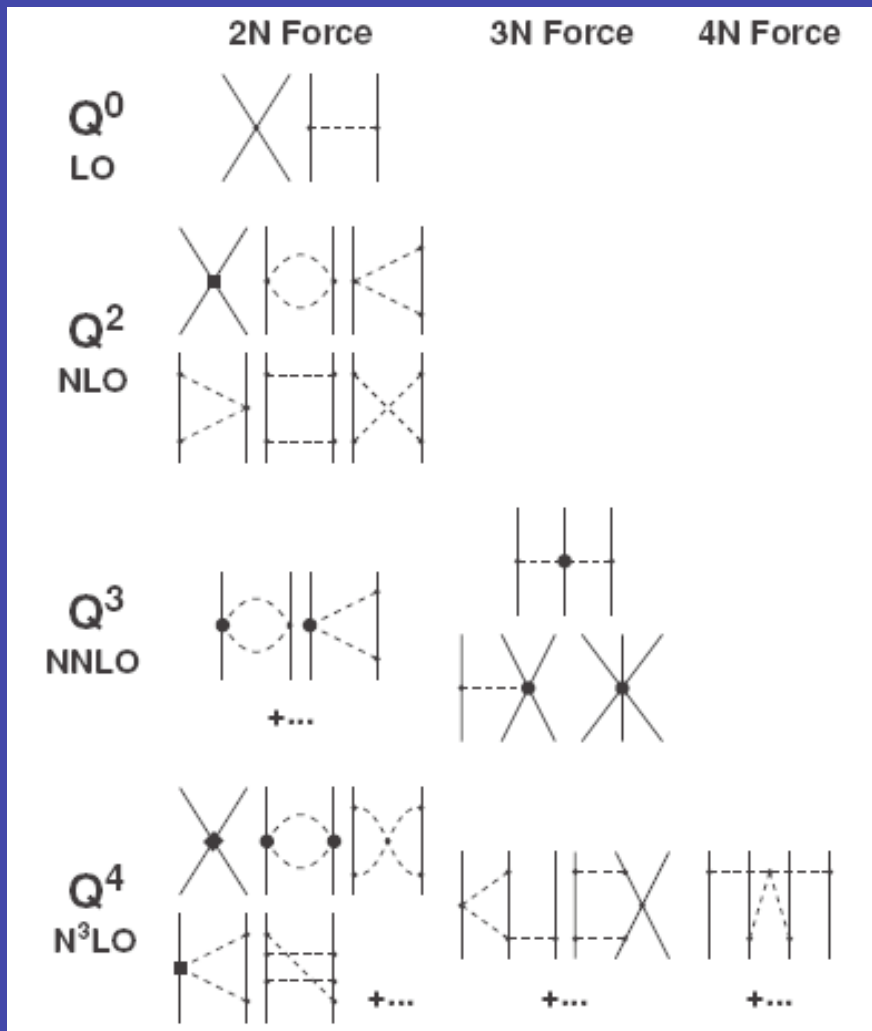
$$0 = \langle \Phi_{ij}^{ab} | \bar{H} | \Phi \rangle$$

Alternative view: CCSD generates similarity transformed Hamiltonian with no 1p-1h and no 2p-2h excitations.

$$\bar{H} \equiv e^{-T} H e^T = (H e^T)_c = \left(H + H T_1 + H T_2 + \frac{1}{2} H T_1^2 + \dots \right)_c$$

Nuclear potential from chiral effective field theory

Diagrams



van Kolck (1994); Epelbaum et al (2002);
Machleidt & Entem (2005);

Ab-initio structure calculations with potentials from chiral EFT

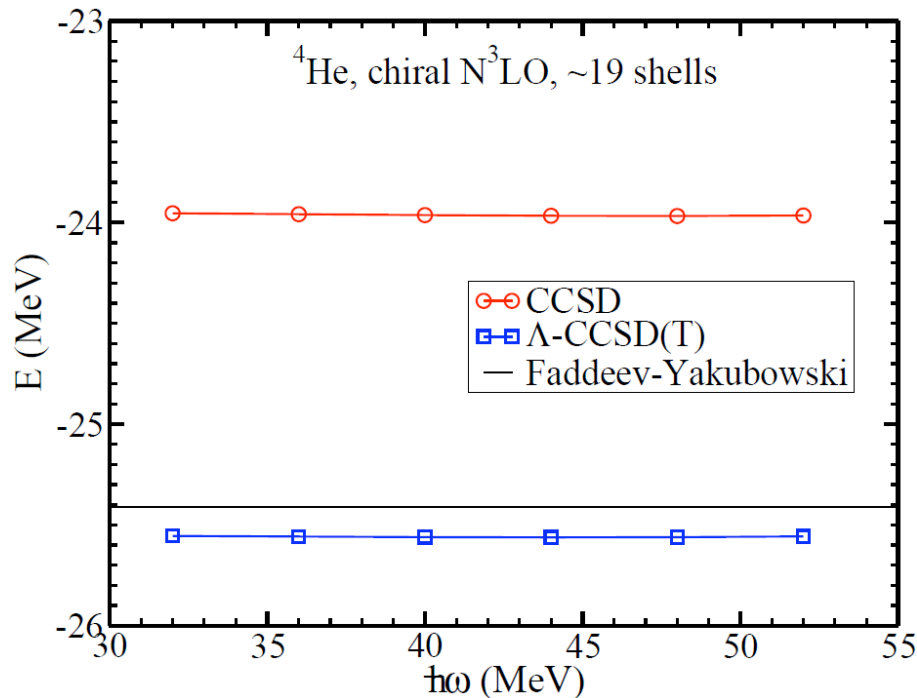
- $A=3, 4$: Faddeev-Yakubowski method
- $A \leq 10$: Hyperspherical Harmonics
- p -shell nuclei: NCSM, GFMC(AV18)
- $^{16,22,24,28}\text{O}$, $^{40,48}\text{Ca}$, ^{48}Ni : Coupled cluster, UMOA, Green's functions (NN so far)
- Lattice simulations
- Nuclear matter

Questions:

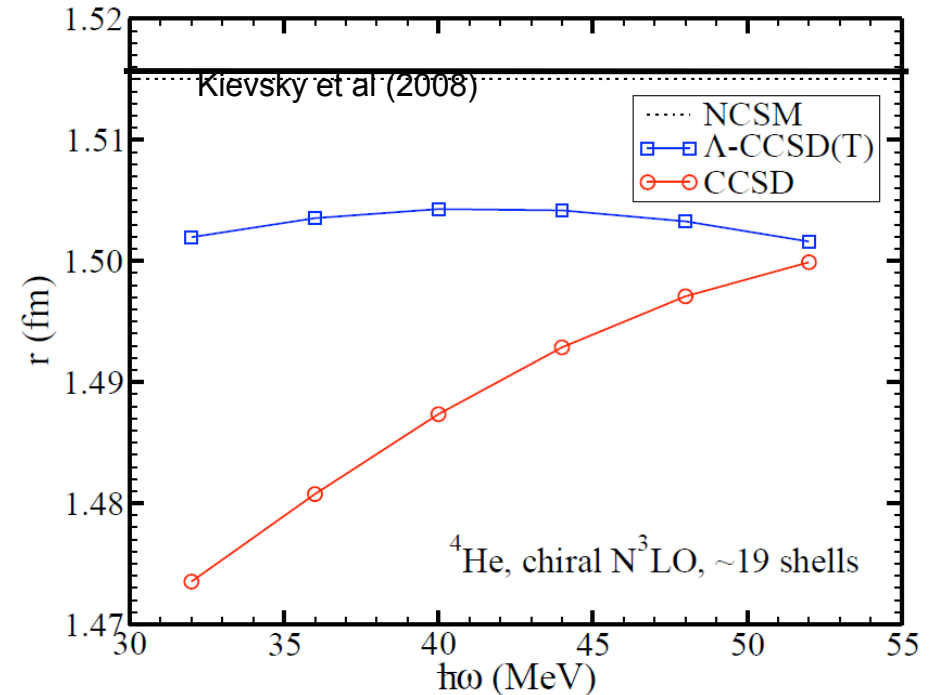
1. Can we compute nuclei from scratch?
2. Role/form of three-nucleon interaction
3. Saturation properties

Precision and accuracy: ^4He , chiral N^3LO [Entem & Machleidt]

Ground-state energy



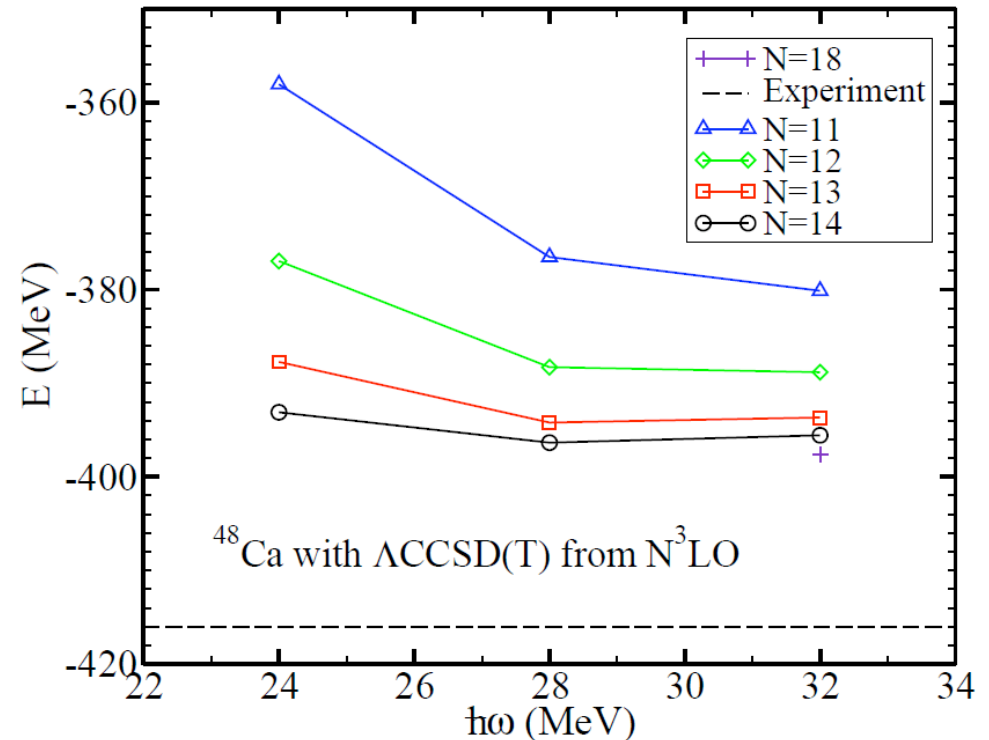
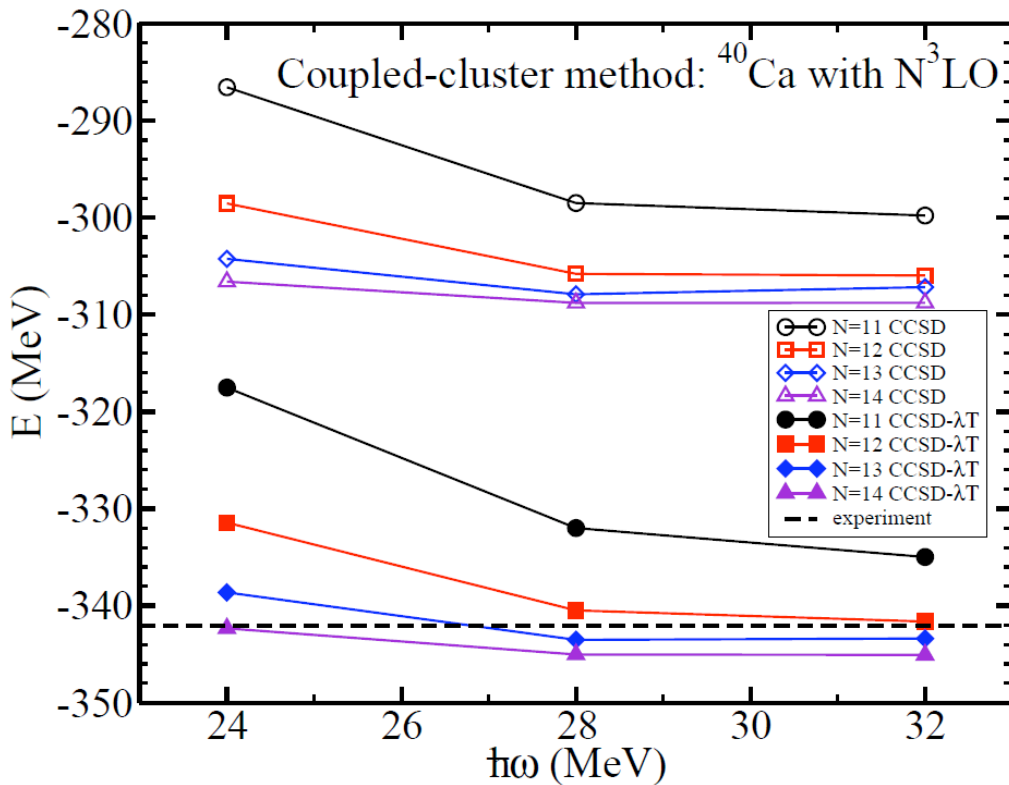
Matter radius



1. Results exhibit very weak dependence on the employed model space.
2. The coupled-cluster method, in its Λ -CCSD(T) approximation, overbinds by 150keV; radius too small by about 0.01fm.
3. Independence of model space of N major oscillator shells with frequency ω :
 - $N\hbar\omega > \hbar^2\Lambda_\chi^2/m$ to resolve momentum cutoff Λ_χ
 - $\hbar\omega < N\hbar^2/(mR^2)$ to resolve nucleus of radius R
4. Number of single-particle states $\sim (R\Lambda_\chi)^3$

Ground-state energies of medium-mass nuclei

CCSD results for chiral N^3LO (NN only)



Binding energy per nucleon

Nucleus	CCSD	Λ -CCSD(T)	Experiment
^4He	5.99	6.39	7.07
^{16}O	6.72	7.56	7.97
^{40}Ca	7.72	8.63	8.56
^{48}Ca	7.40	8.28	8.67

Compare ^{16}O to different approach
Fujii et al., Phys. Rev. Lett. 103,
182501 (2009)

$B/A=6.62$ MeV (2 body clusters)
 $B/A=7.47$ MeV (3 body clusters)

Center-of-mass coordinate

The nuclear Hamiltonian is invariant under rotations and translations

Approach that preserves both symmetries:

☺ Jacobi coordinates

☹ Antisymmetrization very expensive → limited to $A \leq 10$ or so

[Faddeev Yakubowsky; Hyperspherical Harmonics; Manchester group's CCM].

Antisymmetry best dealt within second quantization:

☹ No single-particle basis available that consists of simultaneous eigenstates of the angular momentum operator and the momentum operator.

☺ Within a complete $N\hbar\omega$ oscillator space, the wave function is guaranteed to factorize

$$\psi = \psi_{\text{cm}} \psi_{\text{in}}$$

Intrinsic wave function ψ_{in} invariant under translation

Center-of-mass wave function ψ_{cm} is Gaussian whose width is set by the oscillator length of the employed oscillator basis

The factorization is key. The form of ψ_{cm} is irrelevant.

Center-of-mass coordinate (cont'd)

Intrinsic nuclear Hamiltonian

$$H_{\text{in}} = T - T_{\text{cm}} + V ,$$
$$= \sum_{1 \leq i < j \leq A} \left(\frac{(\vec{p}_i - \vec{p}_j)^2}{2mA} + V(\vec{r}_i - \vec{r}_j) \right)$$

Obviously, H_{in} commutes with any center-of-mass Hamiltonian H_{cm} .

Situation: The Hamiltonian depends on $3(A-1)$ coordinates, and is solved in a model space of $3A$ coordinates. What is the wave function in the center-of-mass coordinate?

Q:How can one demonstrate the factorization of wave function ψ :

A: Find a suitable center-of-mass Hamiltonian H_{cm} whose eigenstate is ψ .

Our approach:

Demonstrate that $\langle H_{\text{cm}} \rangle \approx 0$ for a center-of-mass Hamiltonian with zero-energy ground state.

$$H_{\text{cm}}(\tilde{\omega}) = T_{\text{cm}} + \frac{1}{2}mA\tilde{\omega}^2 R_{\text{cm}}^2 - \frac{3}{2}\hbar\tilde{\omega}$$

Frequency $\tilde{\omega}$ to be determined.

Toy problem

Two particles in one dimension
with intrinsic Hamiltonian

$$H = \frac{p^2}{2m} + V(x)$$

$$V(x) = -V_0 \exp(-(x/l)^2)$$

$$x = (x_1 - x_2) / \sqrt{2}$$

$$p = (p_1 - p_2) / \sqrt{2}$$

Single-particle basis of
oscillator wave functions with
 $m, n = 0, \dots, N$

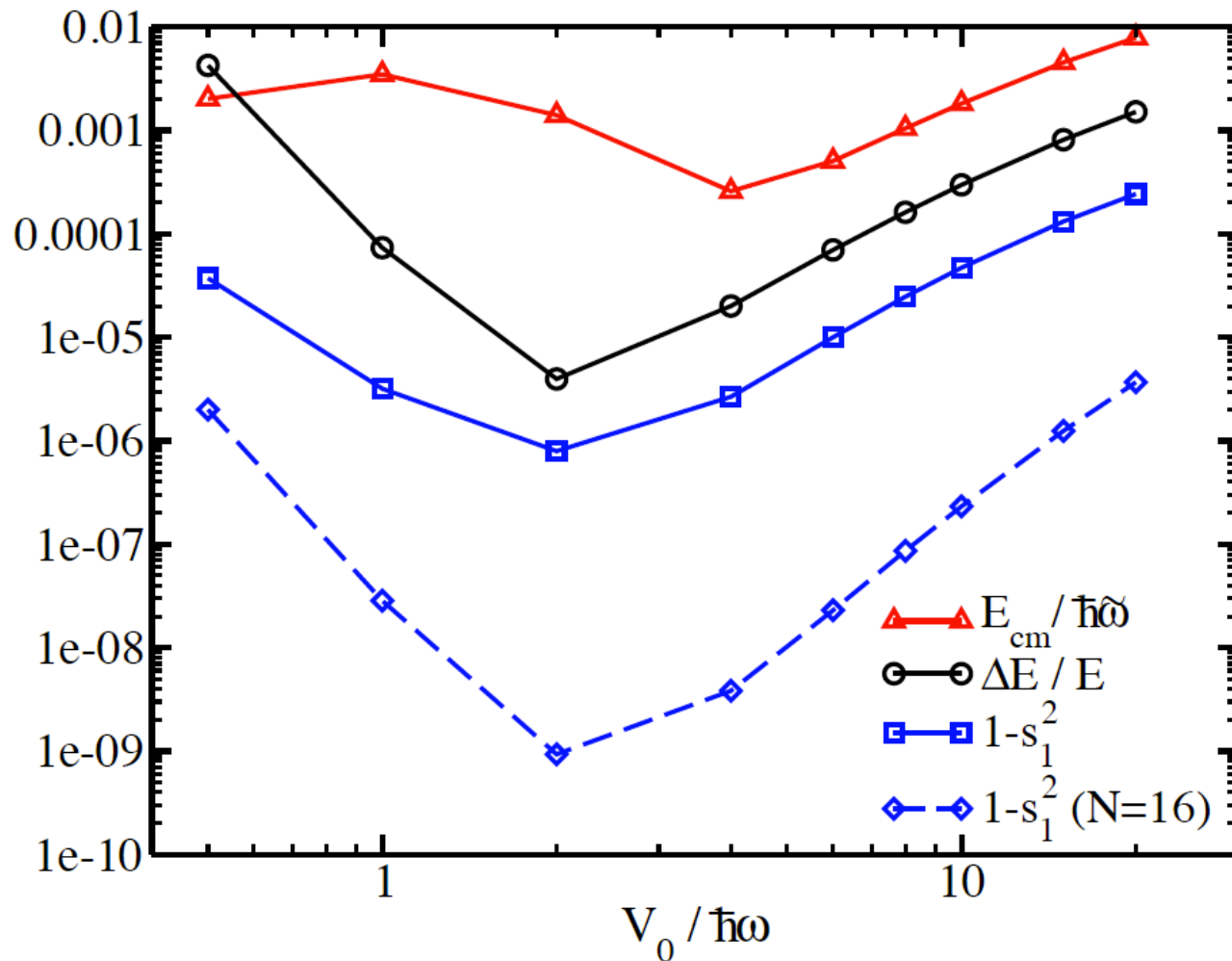
$$\Phi_m(x_1/l) \Phi_n(x_2/l)$$

Results:

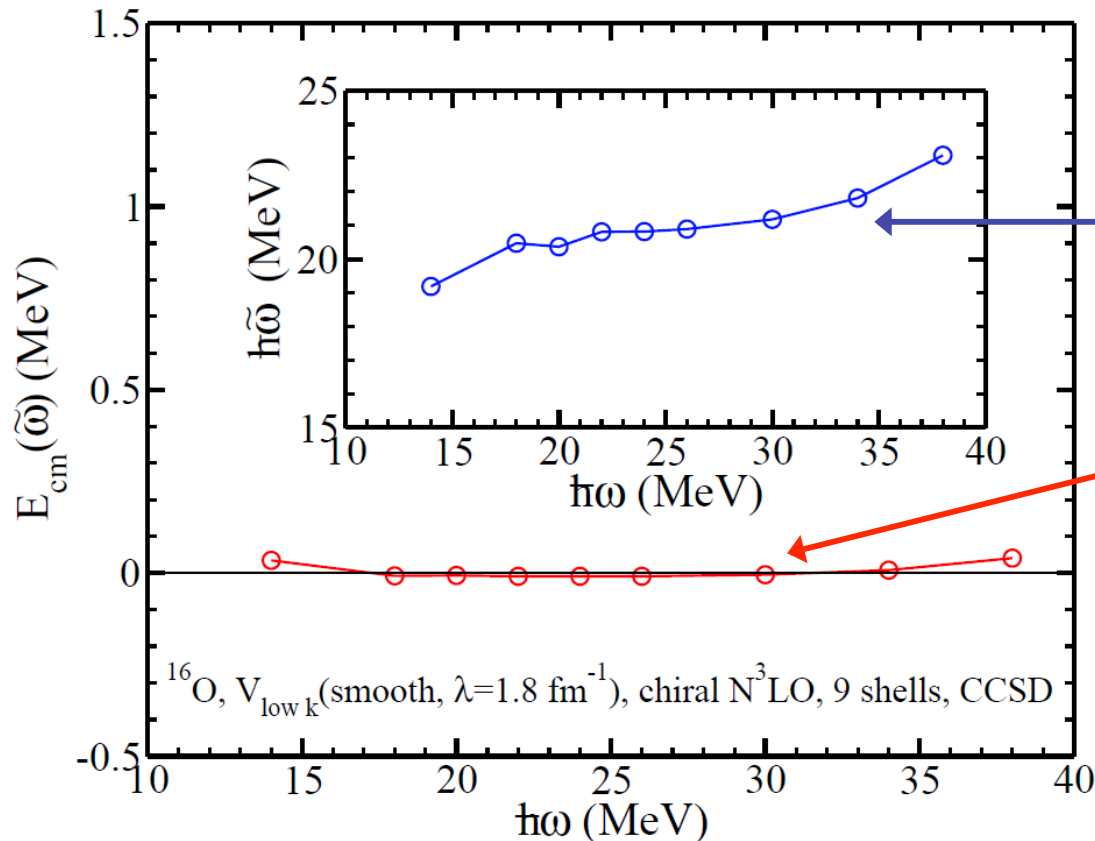
1. Ground-state is factored
with $s_1 \approx 1$

$$\psi_A = \sum_j s_j \psi_{\text{cm}}^{(j)} \psi_{\text{in}}^{(j)}$$

2. CoM wave function is
approximately a Gaussian



Coupled-cluster wave function factorizes to a very good approximation!



Curve becomes practically constant in larger model spaces

E_{cm} is slightly negative (size -0.01 MeV) due to non-variational character of CCSD.

Note: spurious CoM excitations are of order 20 MeV $\ll E_{\text{cm}}$.

Coupled-cluster state is ground state of suitably chosen center-of-mass Hamiltonian.

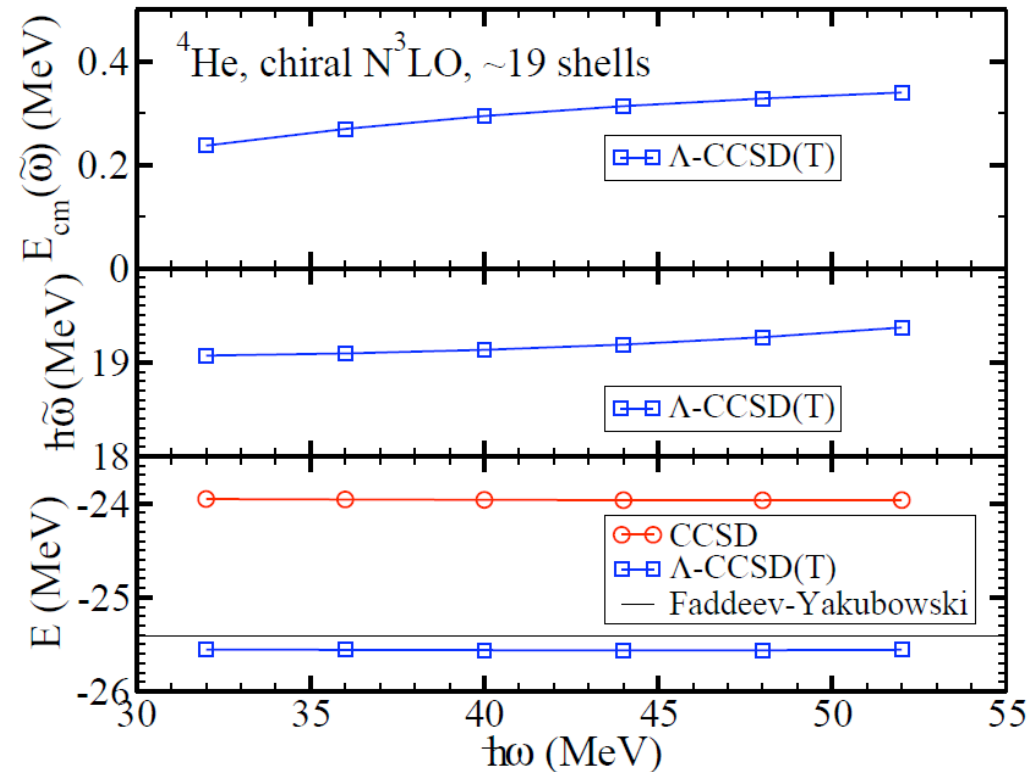
Factorization between intrinsic and center-of-mass coordinate realized within high accuracy.

Note: Both graphs become flatter as the size of the model space is increased.

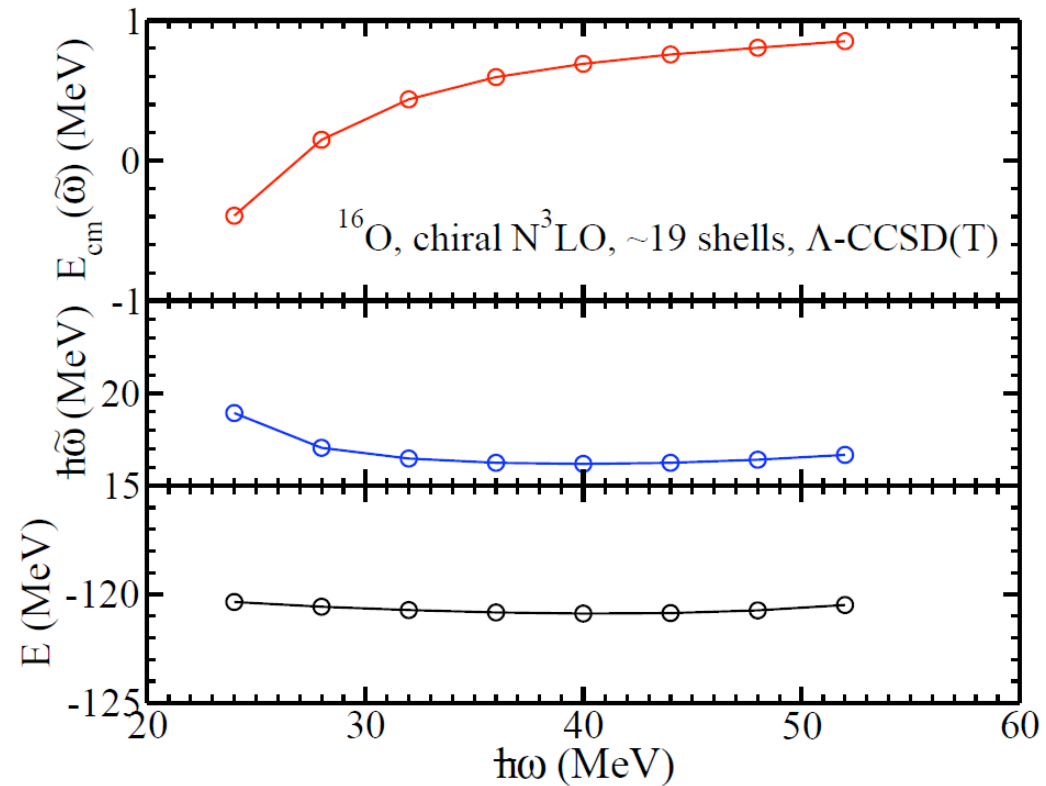
Approximate factorization also for “hard” interactions:

^4He , ^{16}O , and ^{48}Ca from Entem & Machleidt’s chiral N^3LO

^4He



^{16}O



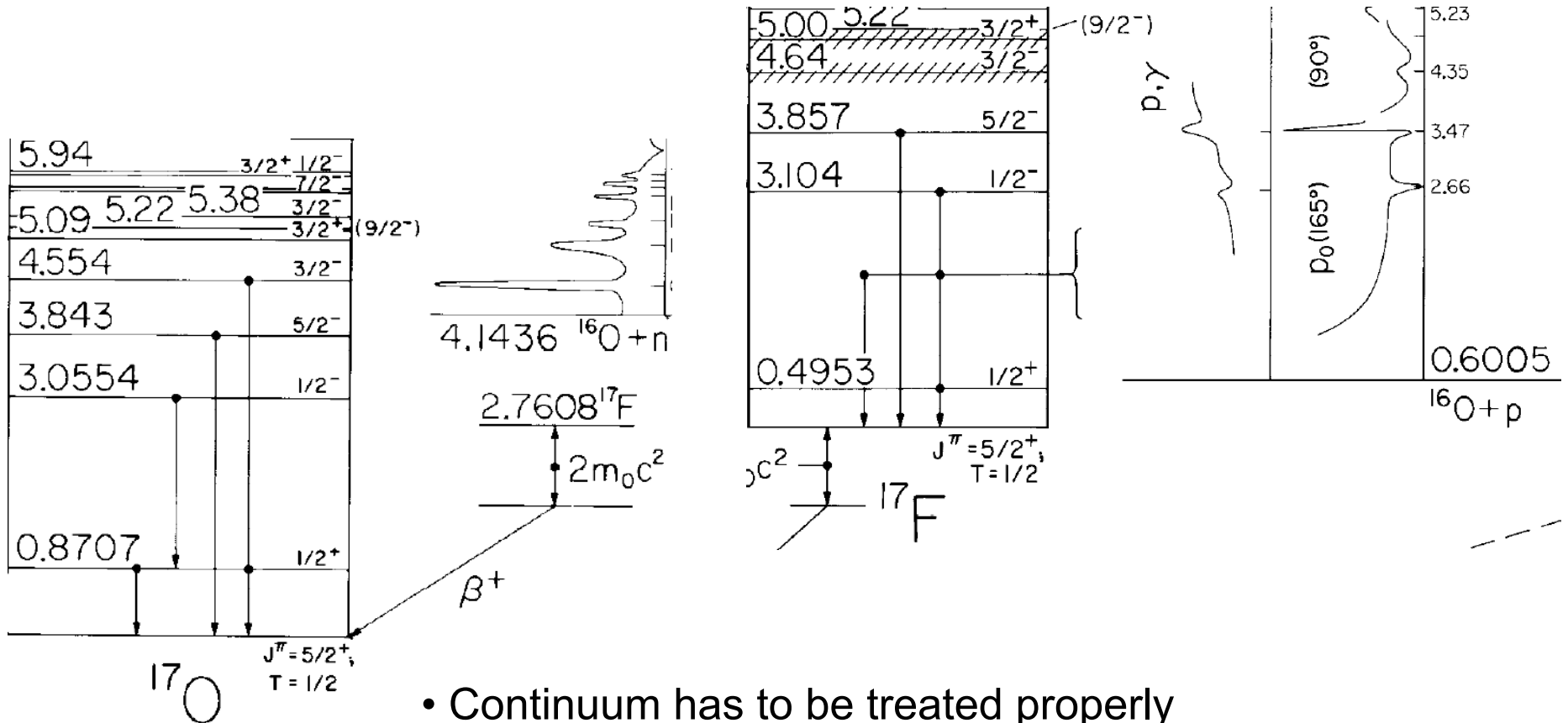
Coupled-cluster wave function factorizes approximately.

Note: spurious states are separated by about $15 - 20 \text{ MeV} \gg E_{\text{cm}}$.

No understanding of Gaussian CoM wave function (yet).

Nucleus	$\hbar\tilde{\omega}$
^4He	19.1 MeV
^{16}O	16.5 MeV
^{48}Ca	14.9 MeV

Low lying states in A=17 nuclei



- Continuum has to be treated properly
- Our focus is on single-particle states
- Previous study: shell model in the continuum with ^{16}O core [Bennaceur et al Phys. Lett. B 488, 75 (2000)]

Bound states and resonances in ^{17}F and ^{17}O

Single-particle basis consists of bound, resonance and scattering states

- Gamow basis for $s_{1/2}$ $d_{5/2}$ and $d_{3/2}$ single-particle states
- Harmonic oscillator states for other partial waves

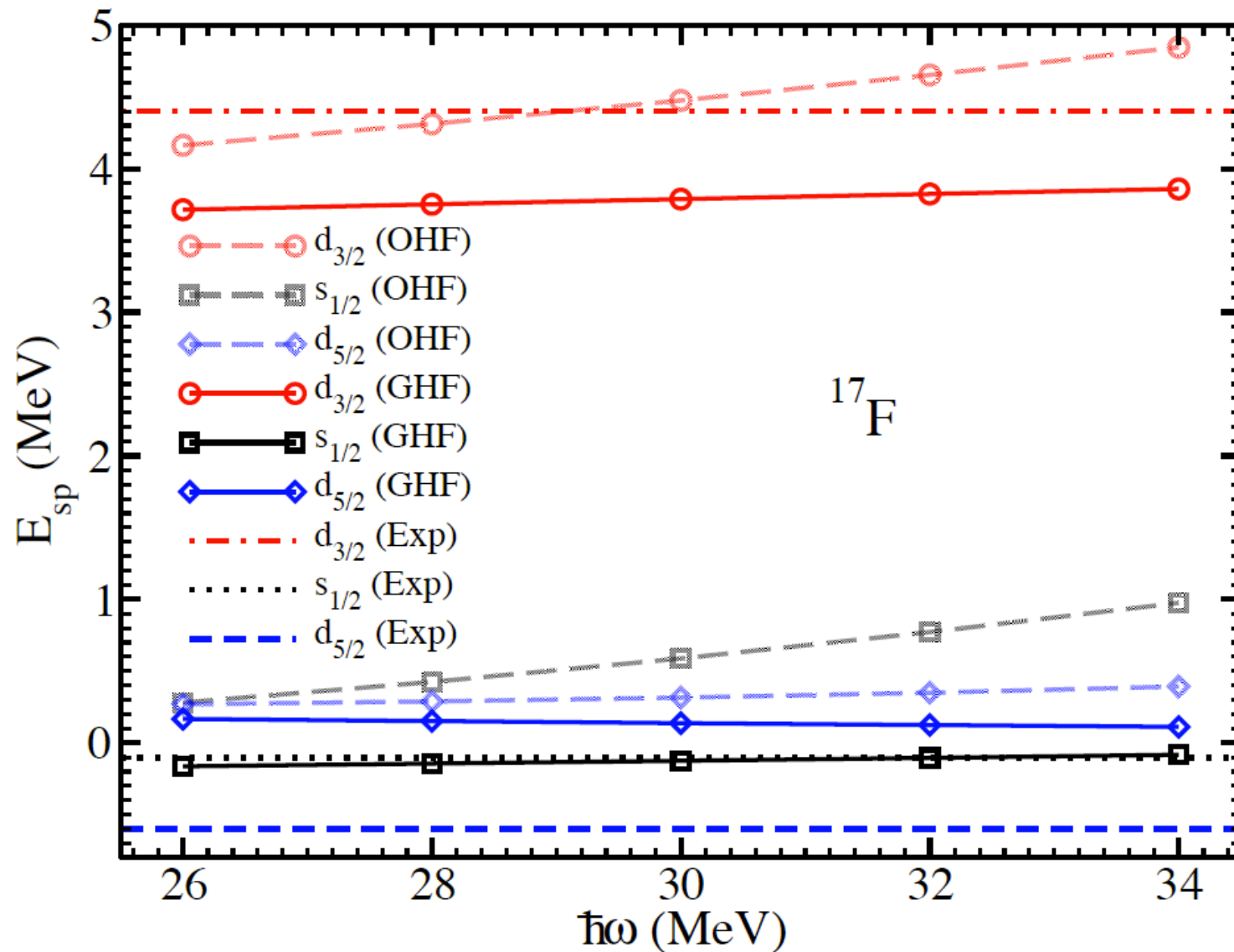
Computation of single-particle states via “Equation-of-motion CCSD”

- Excitation operator acting on closed-shell reference
- Here: superposition of one-particle and 2p-1h excitations

$$R_\mu = r^a a_a^\dagger + \frac{1}{2} r_j^{ab} a_a^\dagger a_b^\dagger a_j$$

$$[\overline{H}, R_\mu] |\phi_0\rangle = \omega_\mu R_\mu |\phi_0\rangle$$

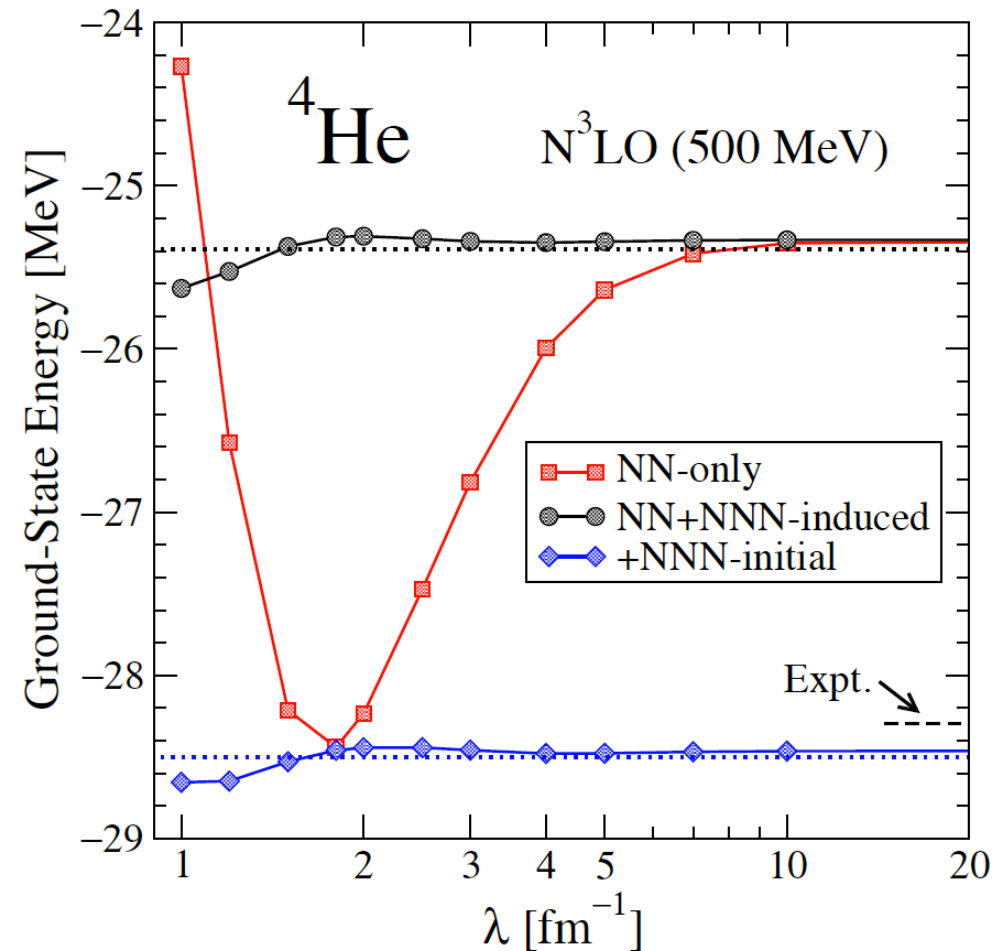
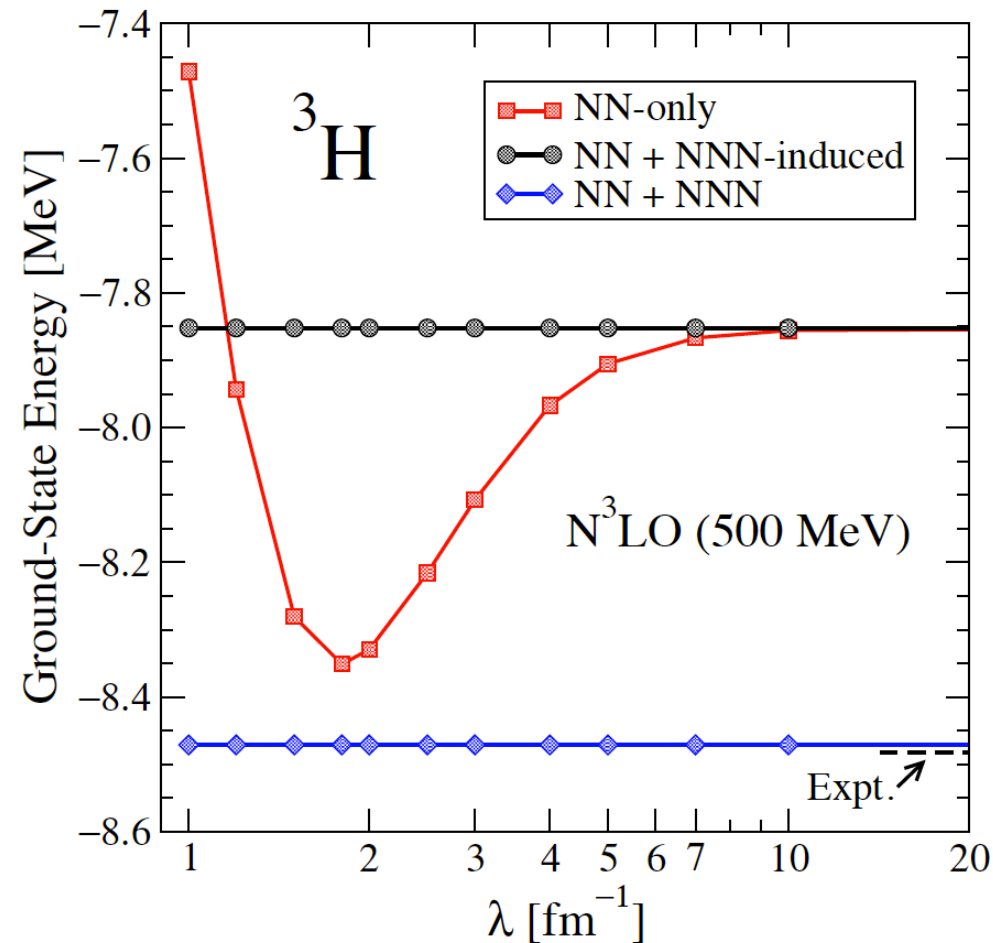
Bound states and resonances in ^{17}F



- Gamow basis weakly dependent on oscillator frequency
- $d_{5/2}$ not bound; spin-orbit splitting too small
- $s_{1/2}$ proton halo state close to experiment

Insights from cutoff variation

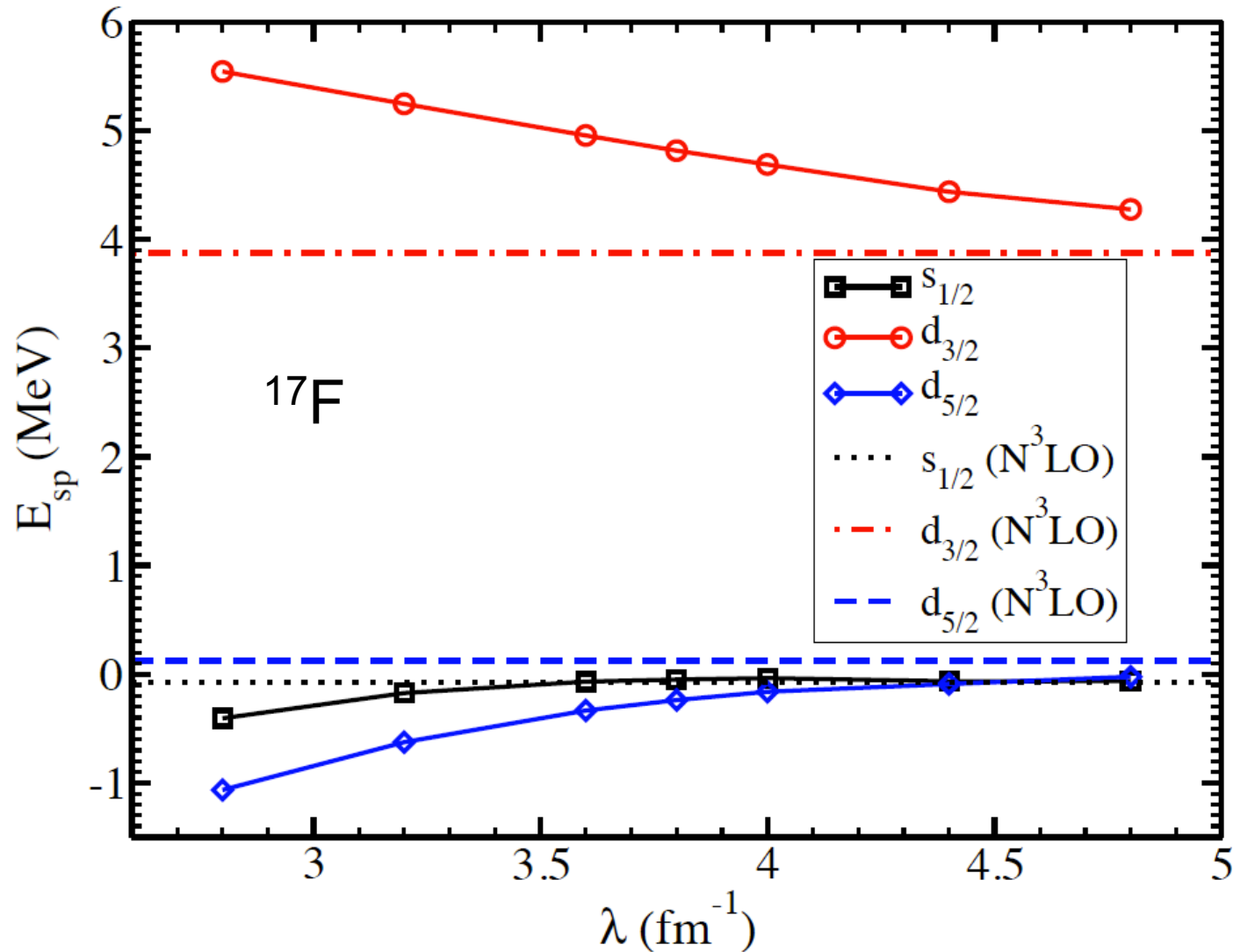
^3H and ^4He with induced and initial 3NF



Jurgenson, Navratil & Furnstahl, Phys. Rev. Lett. 103, 082501 (2009)

Cutoff-dependence hints at missing physics, specifically short-ranged many-body forces.

Variation of cutoff probes omitted short-range forces



- Proton-halo state ($s_{1/2}$) very weakly sensitive to variation of cutoff
- Spin-orbit splitting increases with decreasing cutoff

Results for single-particle energies and decay widths

	^{17}O			^{17}F		
	$1/2^+$	$5/2^+$	E_{so}	$1/2^+$	$5/2^+$	E_{so}
GHF	-2.8	-3.2	4.3	-0.082	0.11	3.7
Exp.	-3.272	-4.143	5.084	-0.105	-0.600	5.000

- Level ordering correctly reproduced in ^{17}O
- Spin-orbit splitting too small

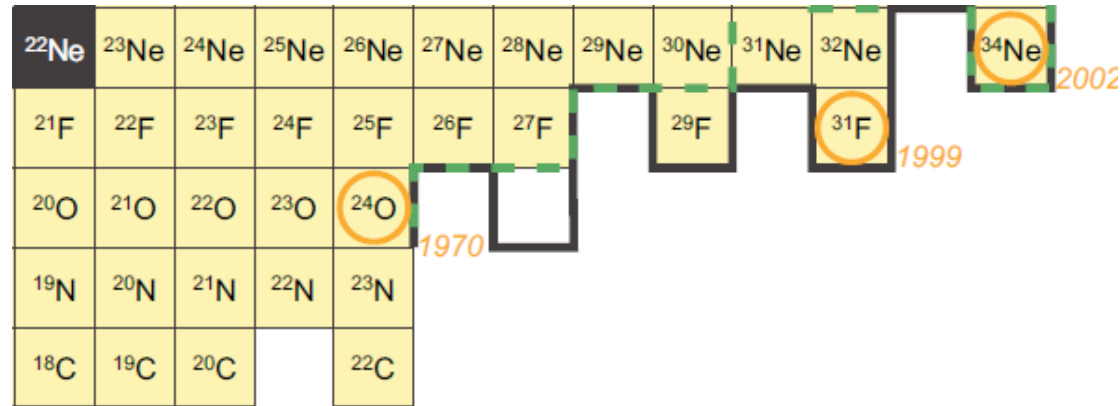
Life times of resonant states

	$^{17}\text{O } 3/2^+$		$^{17}\text{F } 3/2^+$	
	E_{sp}	Γ	E_{sp}	Γ
This work	1.1	0.014	3.9	1.0
Experiment	0.942	0.096	4.399	1.530

Neutron drip line in oxygen isotopes

Experimental situation

- “Last” stable oxygen isotope ^{24}O
- ^{25}O unstable (Hoffman et al 2008)
- $^{26,28}\text{O}$ not seen in experiments
- ^{31}F exists (adding on proton shifts drip line by 6 neutrons!?)



Theoretical situation

- USD interaction predicts stable $^{26,28}\text{O}$ (Brown)
- sd-pf shell calculation can reproduce data after adjusting TBME (Otsuka et al.)
- Shell model w/ continuum couplings employs two different interactions for oxygen isotopes near and far away from β -stability to reproduce data (Volya & Zelevinsky)
- Shell model with 3NF: ^{24}O is last bound isotope (Otsuka, Suzuki, Holt, Schwenk, Akaishi).

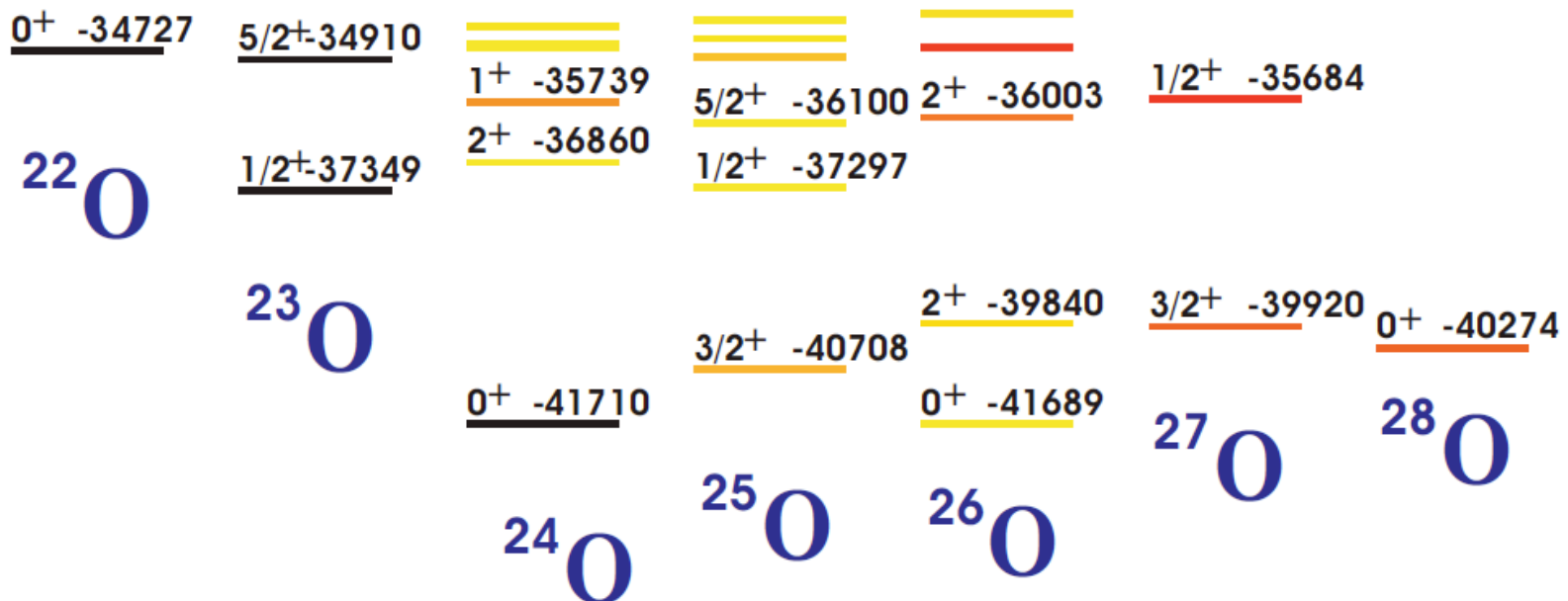
Most theoretical papers rule out a stable ^{28}O .

No approach flawless, i.e. no approach includes everything (continuum effects, 3NFs, no adjustments of interaction)

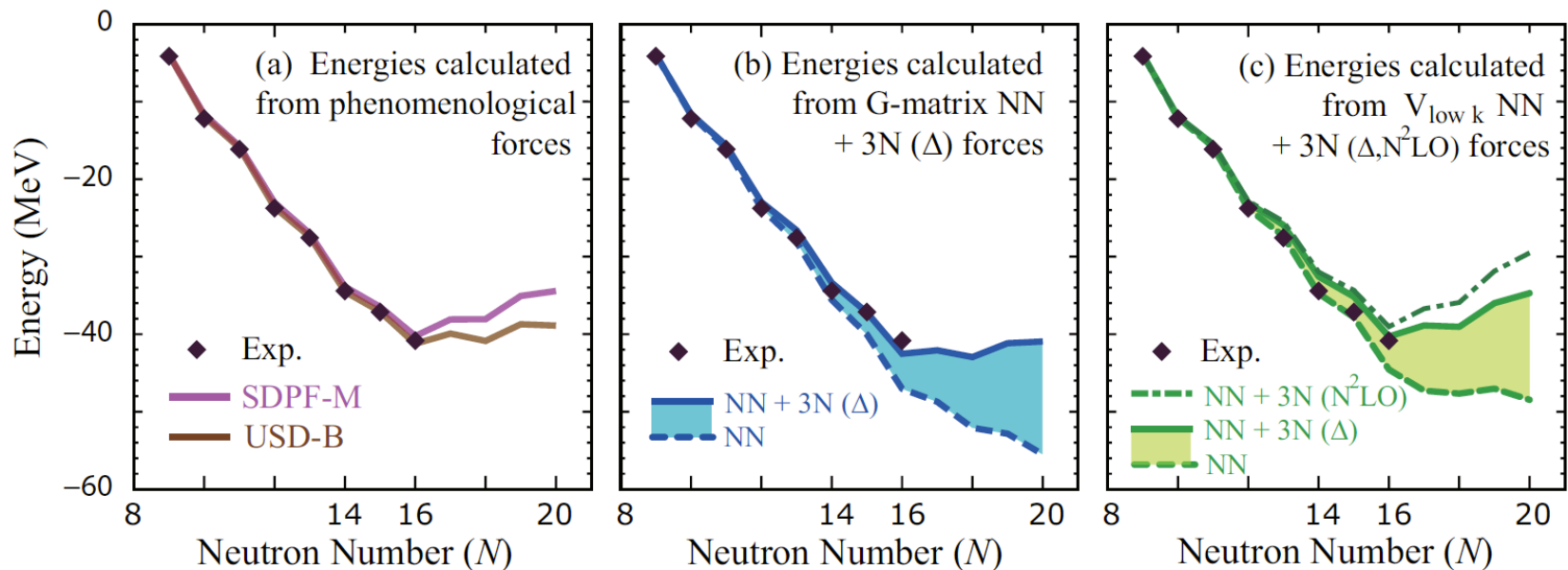
Theoretical difficulties: uncertainties in the effective interaction, quantify the resulting errors.

→ ab-initio calculations: coupled-cluster can address closed sub-shell nuclei $^{22,24,28}\text{O}$ with chiral interactions; study cutoff dependence

Examples of theoretical calculations

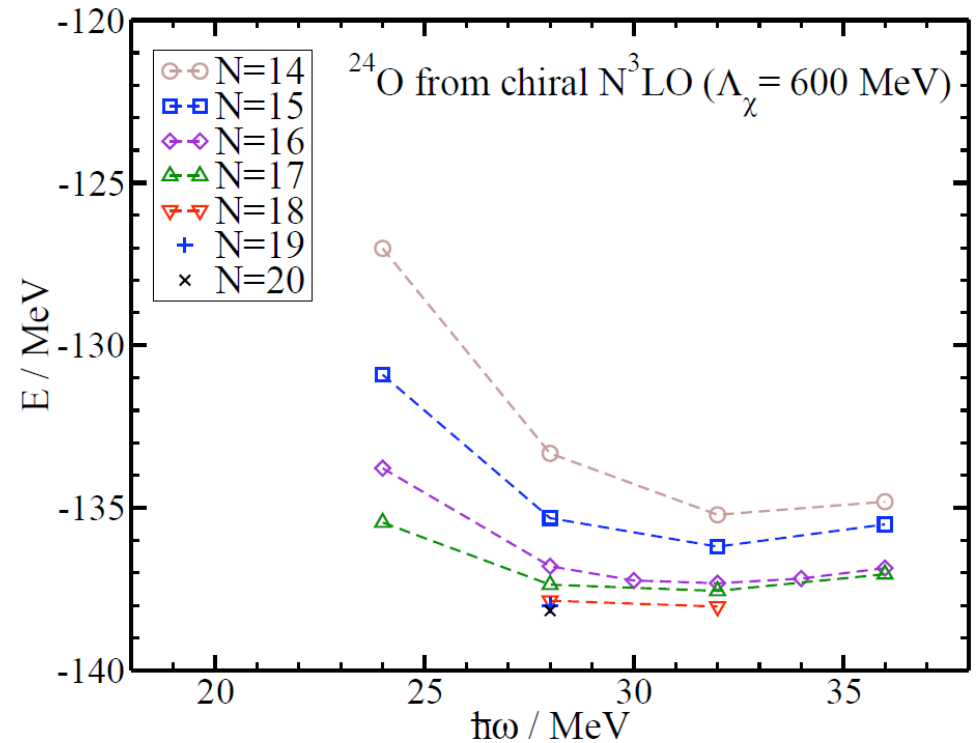
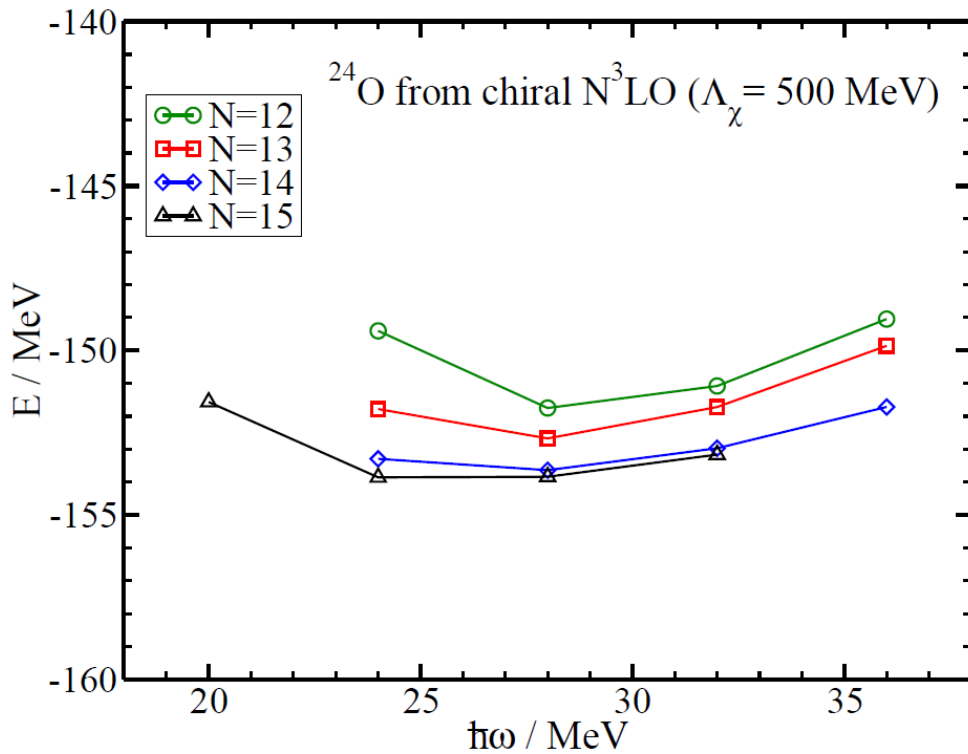


Volya & Zelevinsky, Phys. Rev. Lett. 94 (2005) 052501: Continuum + empirical interaction



Otsuka, Suzuki, Holt, Schwenk, Akaishi, arXiv:0908.2607: 3NF (T=1) in small model space

Neutron-rich oxygen isotopes



$\Lambda_\chi = 500$ MeV potential converges in about 15 major oscillator shells

$\Lambda_\chi = 600$ MeV potential converges in about 20 shells

Summary of results

Energies	^{16}O	^{22}O	^{24}O	^{28}O
$(\Lambda_\chi = 500 \text{ MeV})$				
E_0	24.11	50.37	56.19	71.58
ΔE_{CCSD}	-144.77	-175.79	-190.39	-207.67
ΔE_3	-13.31	-19.22	-19.64	-19.85
E	-120.66	-144.64	-153.84	-155.94
$(\Lambda_\chi = 600 \text{ MeV})$				
E_0	22.08	46.33	52.94	68.57
ΔE_{CCSD}	-119.04	-156.51	-168.49	-182.42
ΔE_3	-14.95	-20.71	-22.49	-22.86
E	-111.91	-130.89	-138.04	-136.71
Experiment	-127.62	-162.03	-168.38	

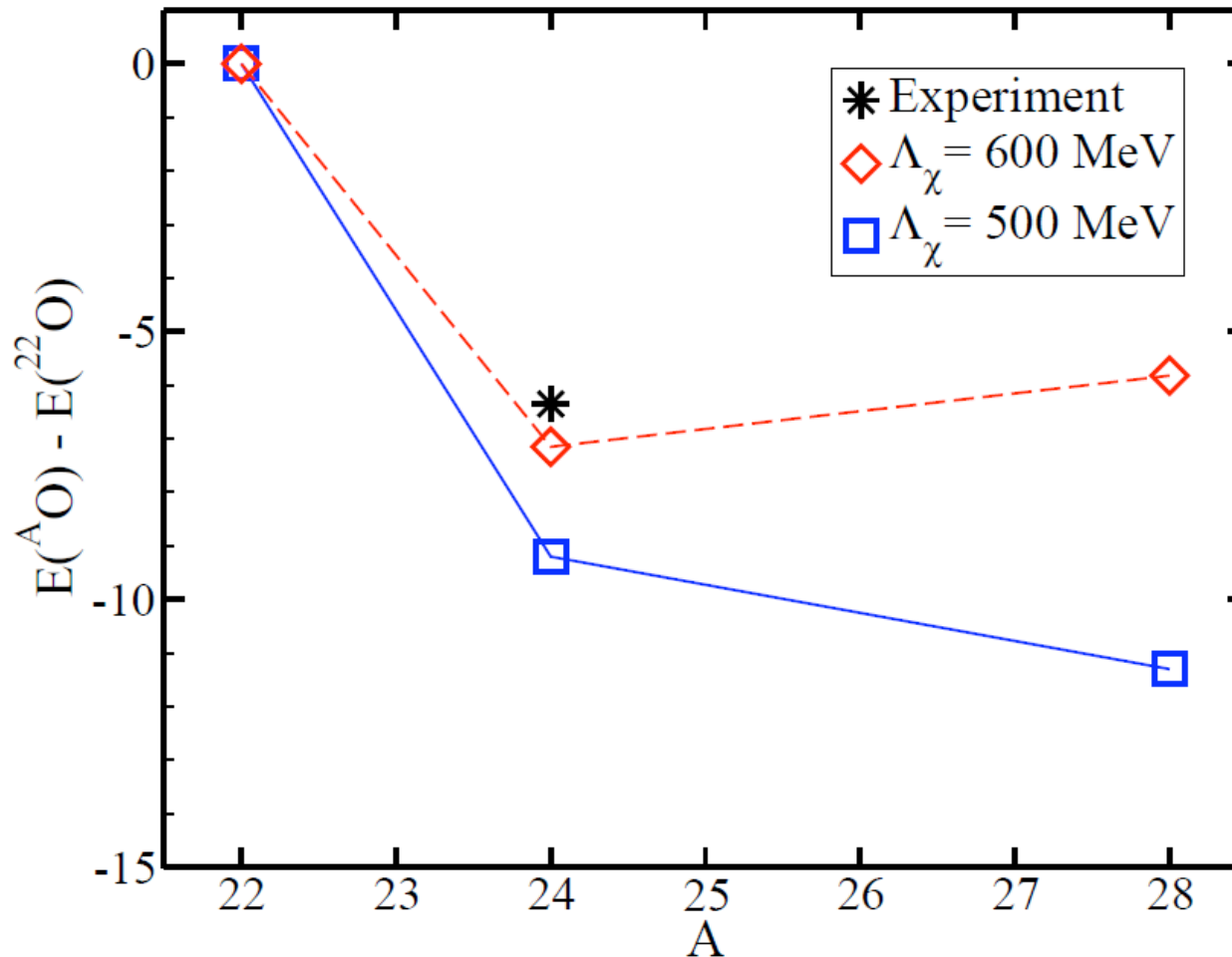
← ~90% of correlation energy

← ~10% of correlation energy

Estimate of theoretical uncertainties:

1. Finite model space ~2MeV
2. Truncation at triples clusters ~2MeV (educated guess)
3. Omission of three-nucleon forces (cutoff dependence) ~15MeV

Is ^{28}O bound relative to ^{24}O ?



Too close to call. Theoretical uncertainties \gg differences in binding energies.

Chiral potentials by Entem & Machleidt's different from G-matrix-based interactions.

Ab-initio theory cannot rule out a stable ^{28}O .

Three-body forces largest potential contribution that decides this question.

Summary

Saturation properties of medium-mass nuclei:

- “Bare” interactions from chiral effective field theory can be converged in large model spaces
- Chiral NN potentials miss ~ 0.4 MeV per nucleon in binding energy in medium-mass nuclei

Practical solution to the center-of-mass problem:

- Demonstration that coupled-cluster wave function factorizes into product of intrinsic and center-of-mass state
- Center-of-mass wave function is Gaussian
- Factorization very pure for “soft” interactions and approximate for “hard” interaction

A=17 nuclei:

- Equation-of-motion CCSD combined with a Gamow basis
- Accurate computation of proton-halo state in ^{17}F ; halo weakly dependent on cutoff

Neutron-rich oxygen isotopes:

- Ab-initio theory cannot rule out a stable ^{28}O
- Greatest uncertainty from omitted three-nucleon forces

Outlook

Towards heavier masses (Ca, Ni, Sn, Pb isotopes)

Inclusion of three-nucleon forces

α -particle excitations (low-lying 0^+ states in doubly magic nuclei)