

# Importance Truncated No-Core Shell Model for Ab Initio Nuclear Structure

Robert Roth  
Institut für Kernphysik



TECHNISCHE  
UNIVERSITÄT  
DARMSTADT

# Overview

- Motivation
- Unitarily Transformed Interactions
  - Unitary Correlation Operator Method
  - Similarity Renormalization Group
- Computational Many-Body Methods
  - No-Core Shell Model
  - Importance Truncated NCSM
  - Center-of-Mass Diagnostics

# From QCD to Nuclear Structure

**Nuclear Structure**

**Low-Energy QCD**

# From QCD to Nuclear Structure

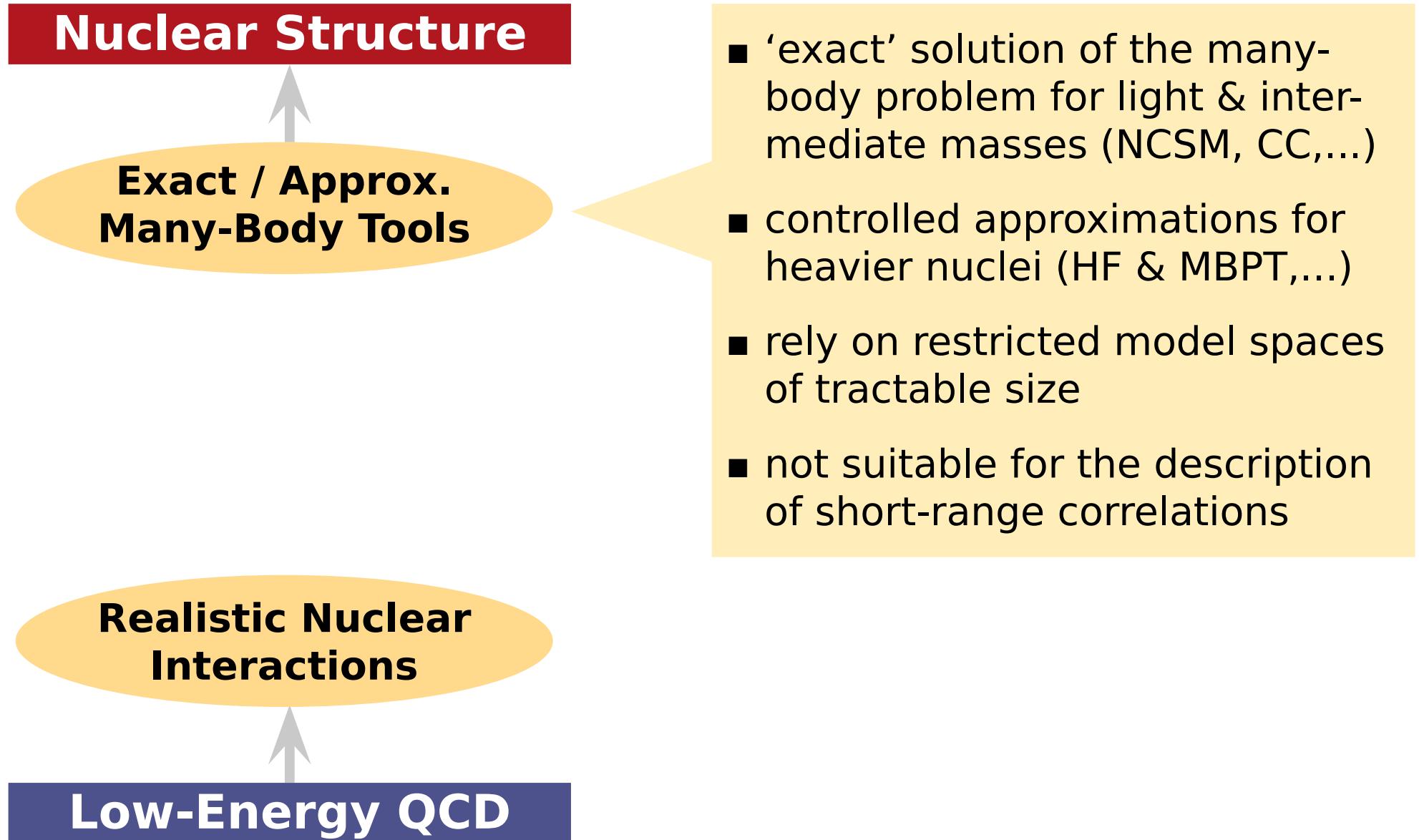
## Nuclear Structure

**Realistic Nuclear  
Interactions**

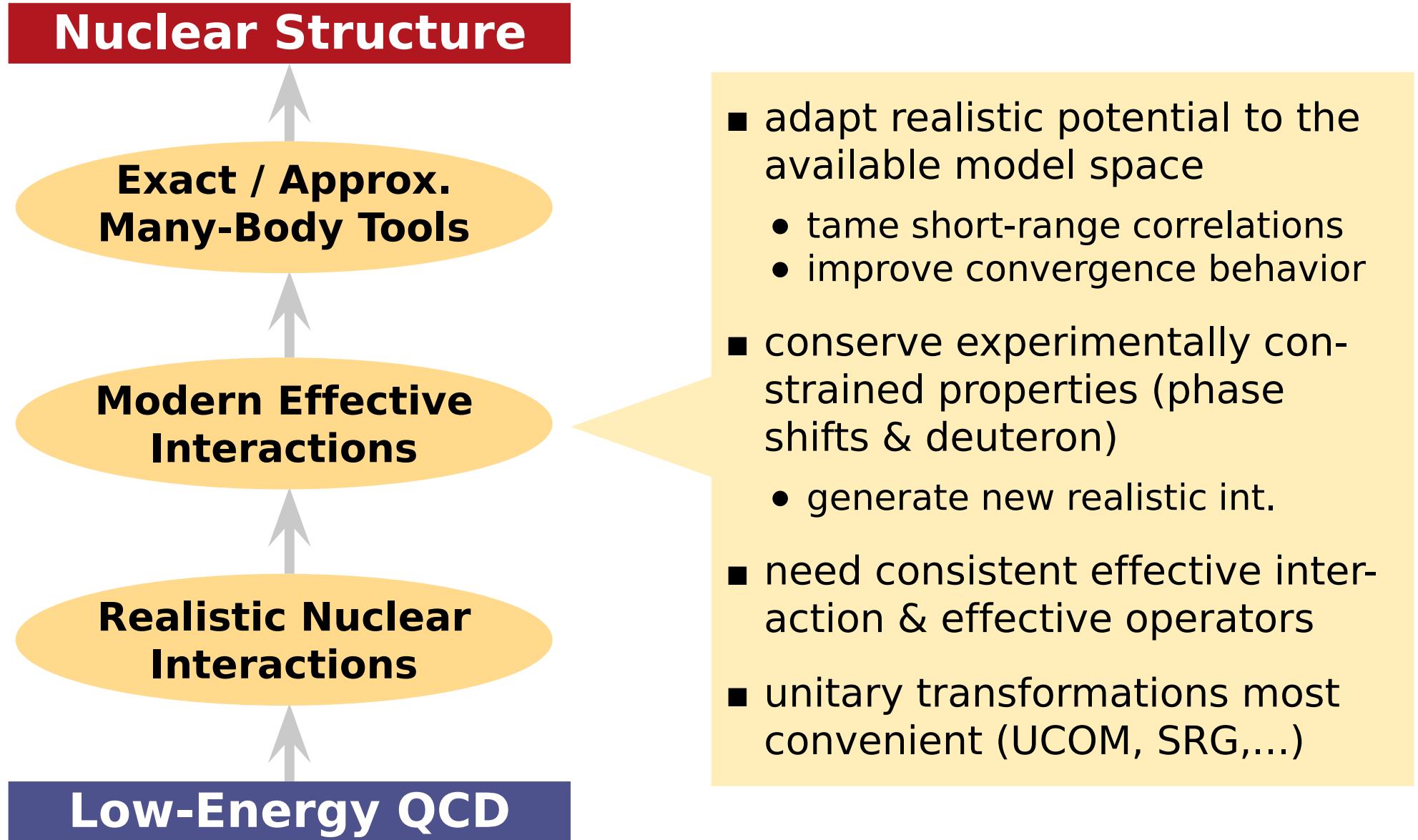
**Low-Energy QCD**

- chiral EFT interactions: consistent NN & 3N interaction derived within  $\chi$ EFT
- traditional NN-interactions: Argonne V18, CD Bonn,...
- reproduce experimental two-body data with high precision
- induce strong short-range central & tensor correlations

# From QCD to Nuclear Structure



# From QCD to Nuclear Structure



Unitarily Transformed Interactions

# Unitary Correlation Operator Method (UCOM)

H. Feldmeier et al. — Nucl. Phys. A 632 (1998) 61

T. Neff et al. — Nucl. Phys. A713 (2003) 311

R. Roth et al. — Nucl. Phys. A 745 (2004) 3

R. Roth et al. — Phys. Rev. C 72, 034002 (2005)

# Unitary Correlation Operator Method

explicit ansatz for unitary transformation operator **motivated by the physics of short-range correlations**

## Central Correlator $C_r$

- radial distance-dependent shift in the relative coordinate of a nucleon pair

$$g_r = \frac{1}{2} [s(r) q_r + q_r s(r)]$$

$$q_r = \frac{1}{2} [\vec{r} \cdot \vec{q} + \vec{q} \cdot \vec{r}]$$

## Tensor Correlator $C_\Omega$

- angular shift depending on the orientation of spin and relative coordinate of a nucleon pair

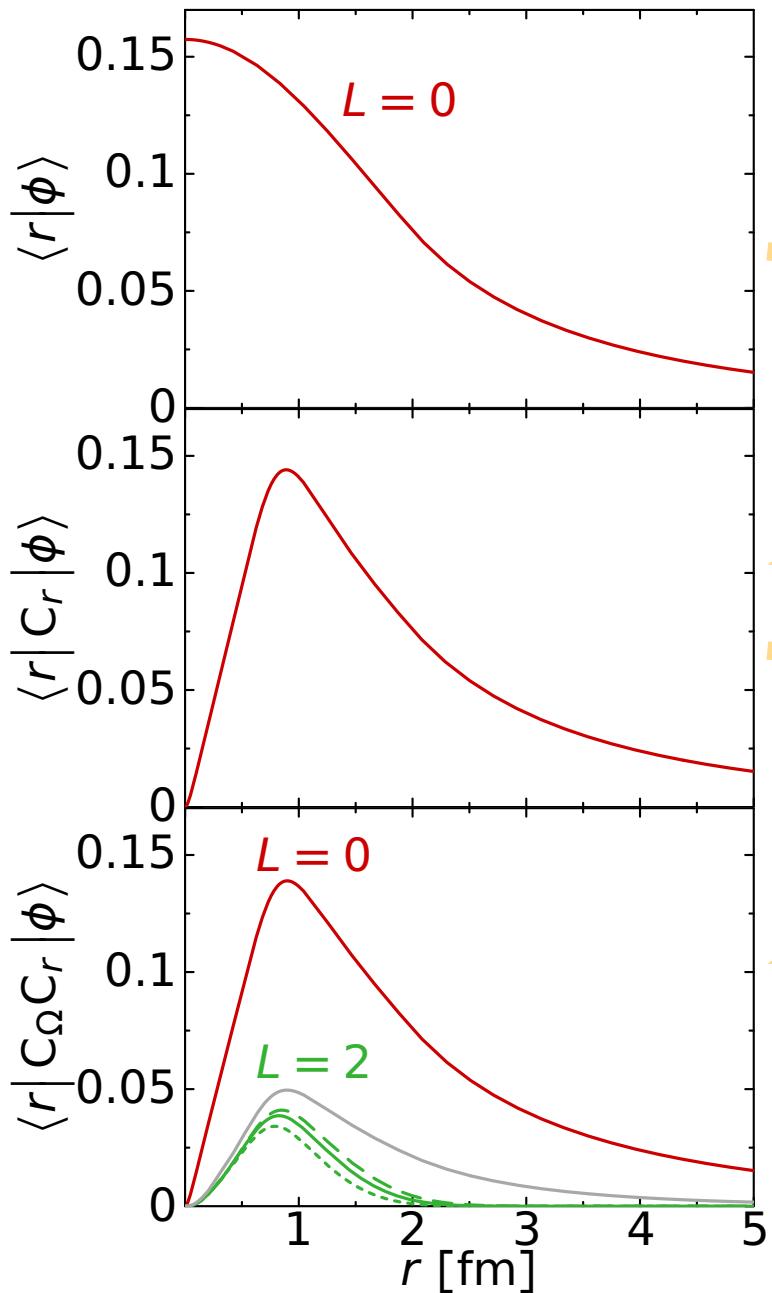
$$g_\Omega = \frac{3}{2} \vartheta(r) [(\vec{\sigma}_1 \cdot \vec{q}_\Omega)(\vec{\sigma}_2 \cdot \vec{r}) + (\vec{r} \leftrightarrow \vec{q}_\Omega)]$$

$$\vec{q}_\Omega = \vec{q} - \frac{\vec{r}}{r} q_r$$

$$C = C_\Omega C_r = \exp\left(-i \sum_{i < j} g_{\Omega,ij}\right) \exp\left(-i \sum_{i < j} g_{r,ij}\right)$$

- $s(r)$  and  $\vartheta(r)$  depend on & are optimized for initial potential

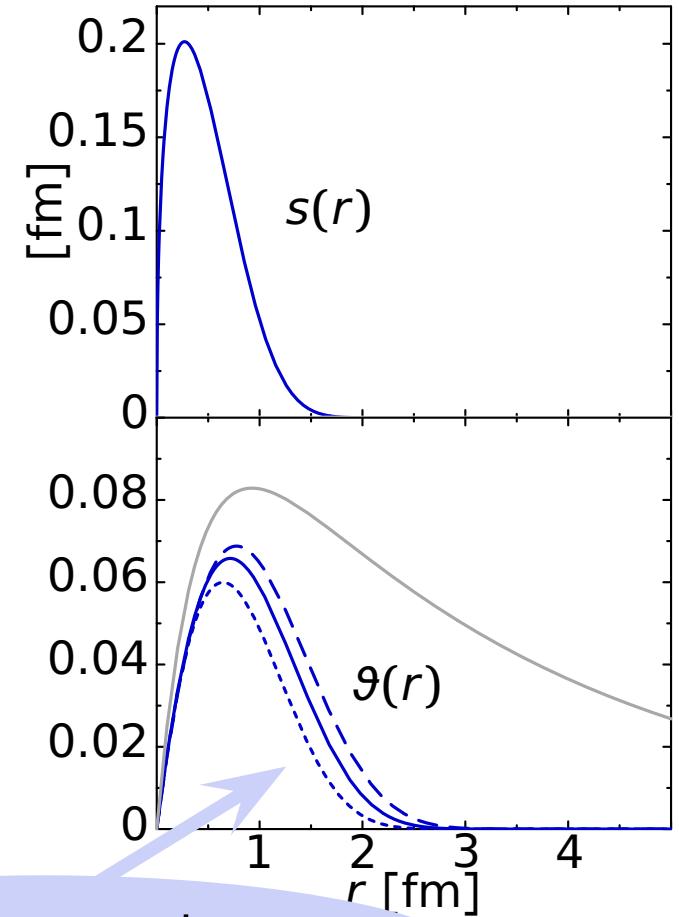
# Correlated States: The Deuteron



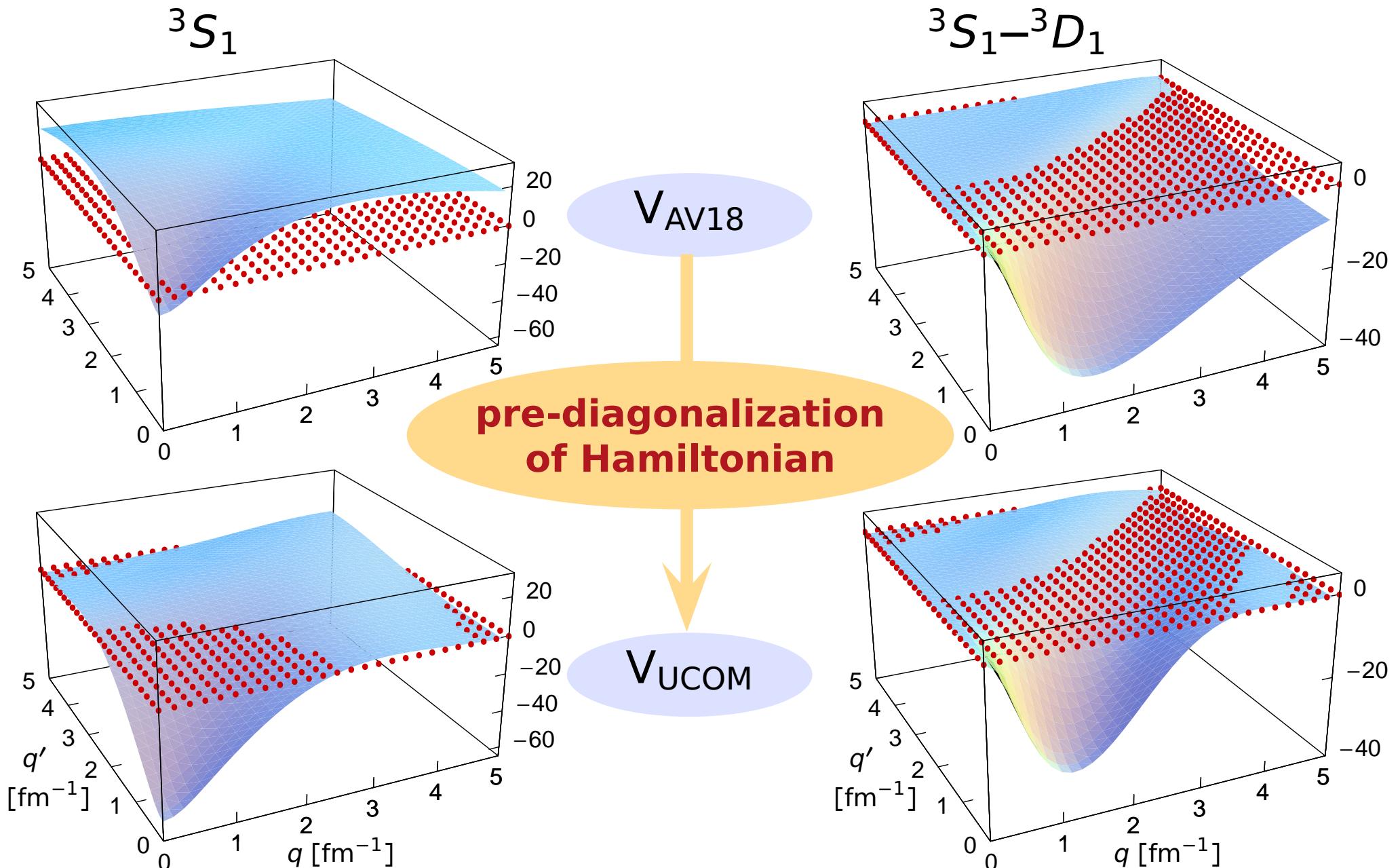
central correlations

tensor correlations

only short-range tensor correlations treated by  $C_\Omega$



# Correlated Interaction: $V_{\text{UCOM}}$



Unitarily Transformed Interactions

# Similarity Renormalization Group (SRG)

Hergert & Roth — Phys. Rev. C 75, 051001(R) (2007)

Bogner et al. — Phys. Rev. C 75, 061001(R) (2007)

Roth, Reinhardt, Hergert — Phys. Rev. C 77, 064033 (2008)

# Similarity Renormalization Group

flow evolution of the **Hamiltonian to band-diagonal form** with respect to uncorrelated many-body basis

## Flow Equation for Hamiltonian

- evolution equation for Hamiltonian

$$\tilde{H}(\alpha) = C^\dagger(\alpha) H C(\alpha) \quad \rightarrow \quad \frac{d}{d\alpha} \tilde{H}(\alpha) = [\eta(\alpha), \tilde{H}(\alpha)]$$

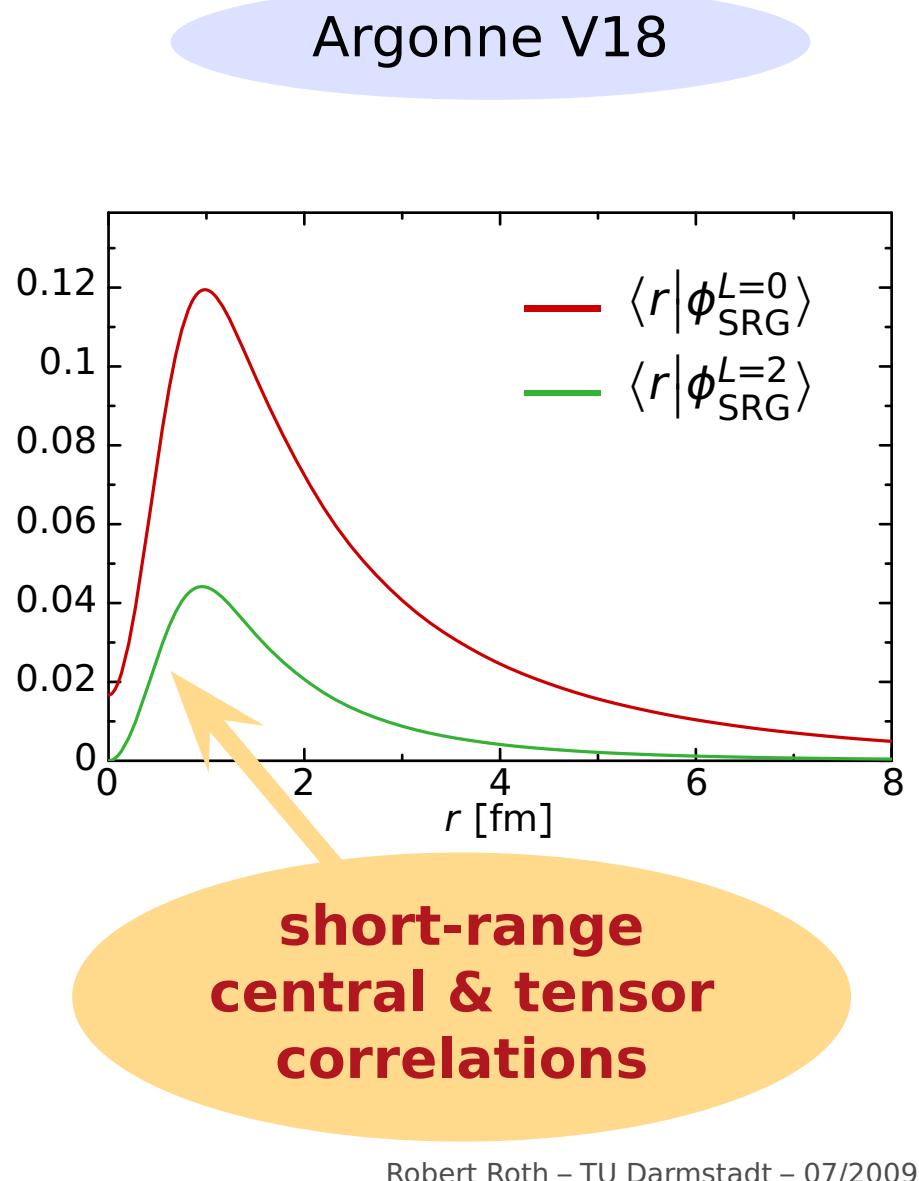
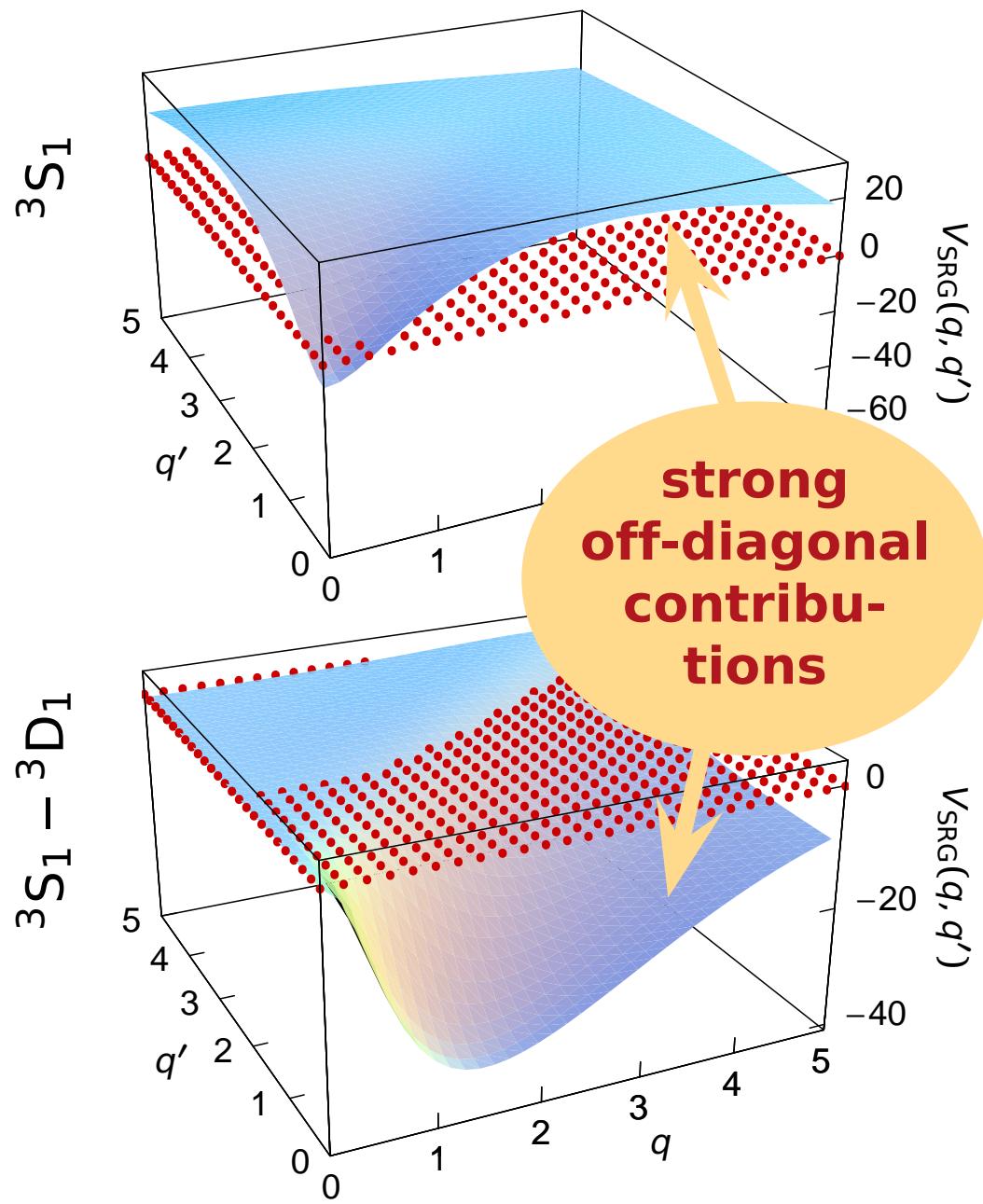
- dynamical generator defined as commutator with the operator in whose eigenbasis  $H$  shall be diagonalized

$$\eta(\alpha) \stackrel{2B}{=} \frac{1}{2\mu} [\vec{q}^2, \tilde{H}(\alpha)]$$

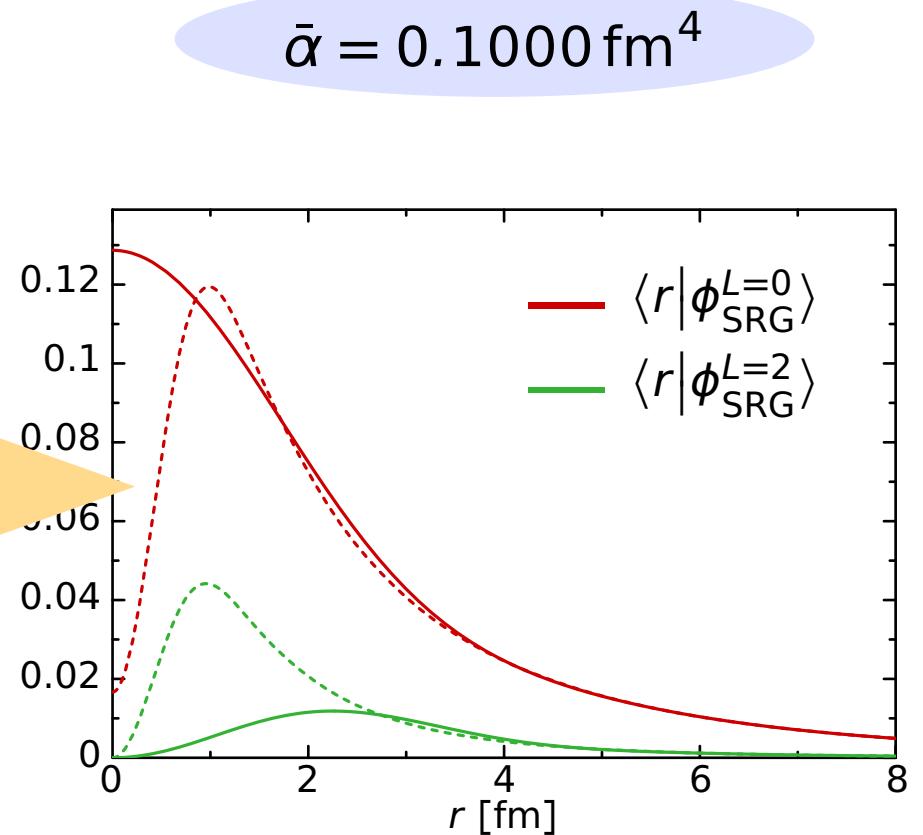
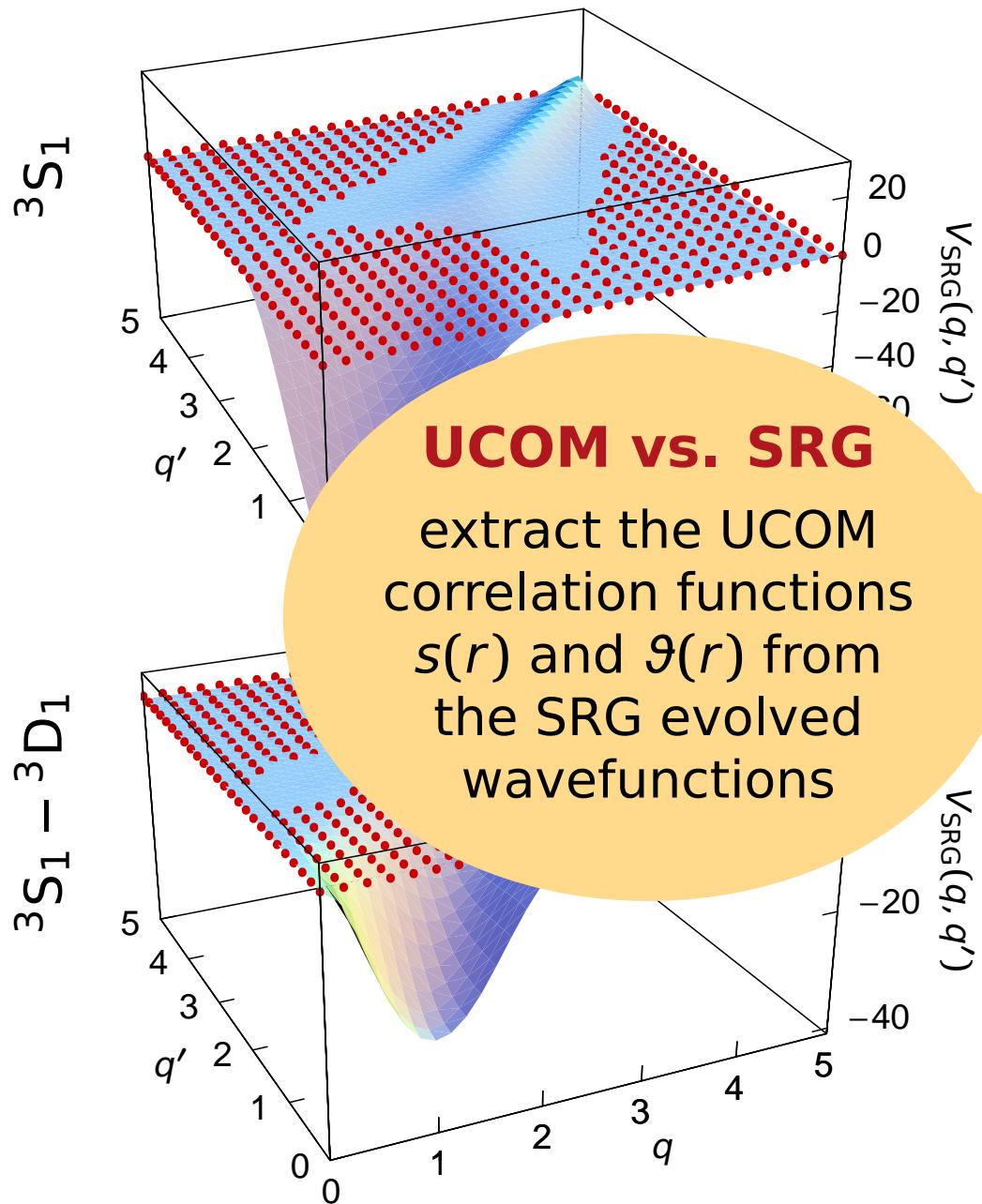
## UCOM vs. SRG

$\eta(0)$  has the same structure as UCOM generators  $g_r$  &  $g_\Omega$

# SRG Evolution: The Deuteron



# SRG Evolution: The Deuteron



# Computational Many-Body Methods

# No-Core Shell Model

Roth et al. — Phys. Rev. C 72, 034002 (2005)

Roth & Navrátil — in preparation

# No-Core Shell Model: Basics

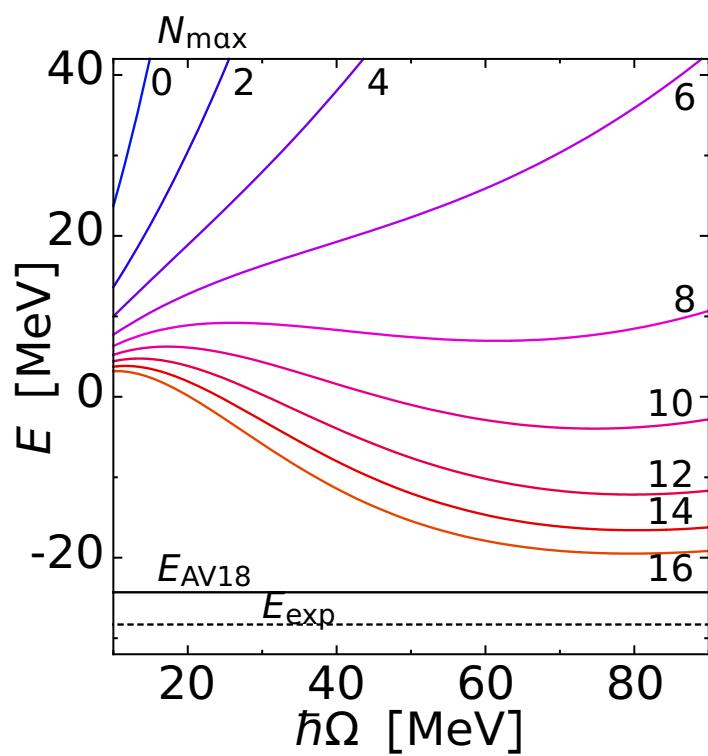
- special case of a **full configuration interaction (CI)** scheme
- **many-body basis**: Slater determinants  $|\Phi_\nu\rangle$  composed of harmonic oscillator single-particle states

$$|\Psi\rangle = \sum_\nu C_\nu |\Phi_\nu\rangle$$

- **model space**: spanned by basis states  $|\Phi_\nu\rangle$  with unperturbed excitation energies of up to  $N_{\max}\hbar\Omega$ 
  - ▶ **exact factorization** of intrinsic and CM component is possible
- numerical solution of **eigenvalue problem** for  $H_{\text{int}}$  within  $N_{\max}\hbar\Omega$  model space via Lanczos methods
  - ▶ model spaces of **up to  $10^9$  basis states** are used routinely
- increase  $N_{\max}$  until **convergence** is observed

# $^4\text{He}$ : NCSM Convergence

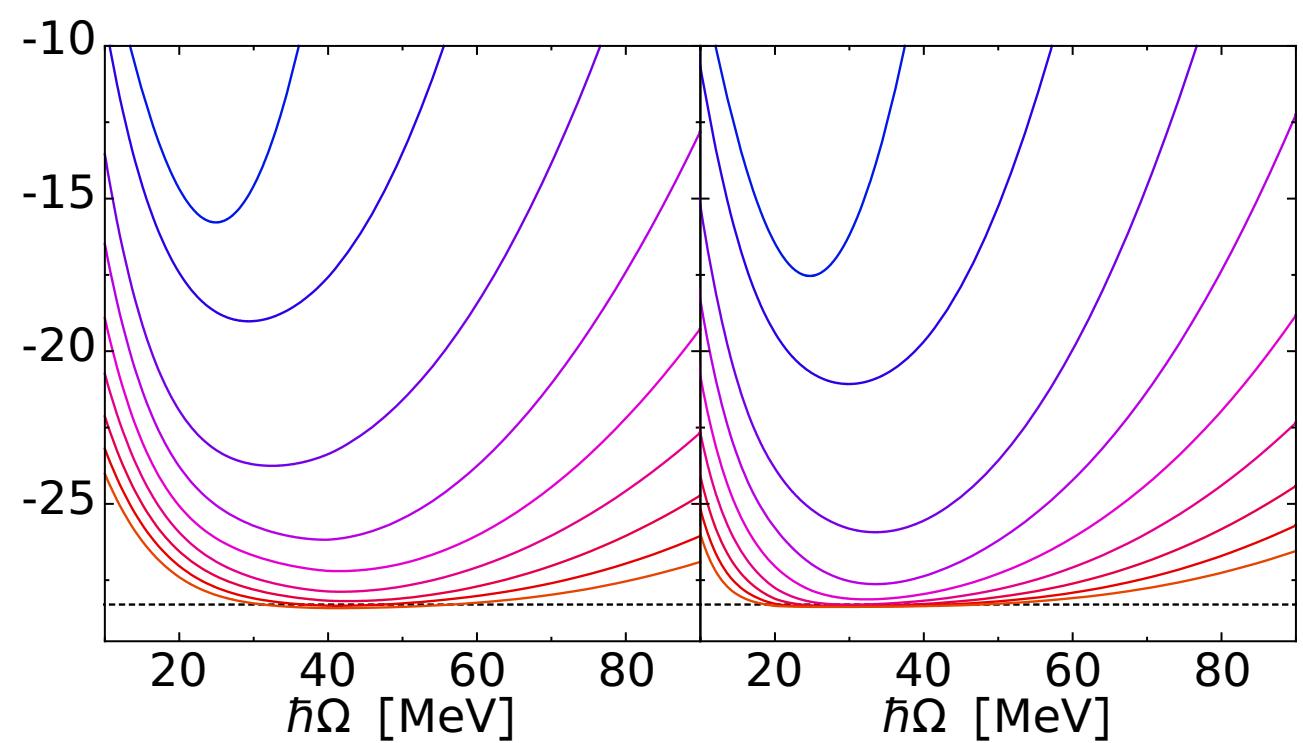
$V_{\text{AV18}}$



$V_{\text{UCOM}}$

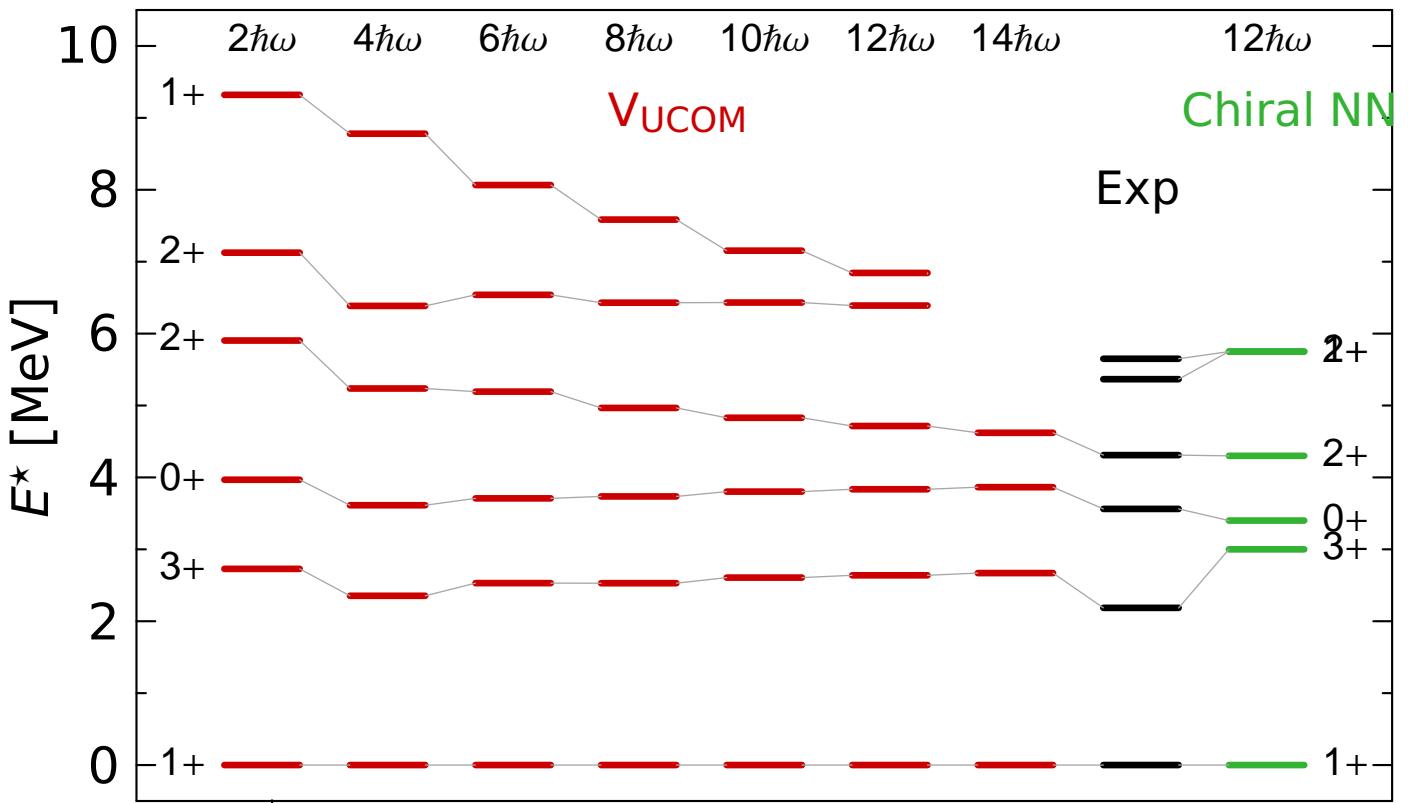
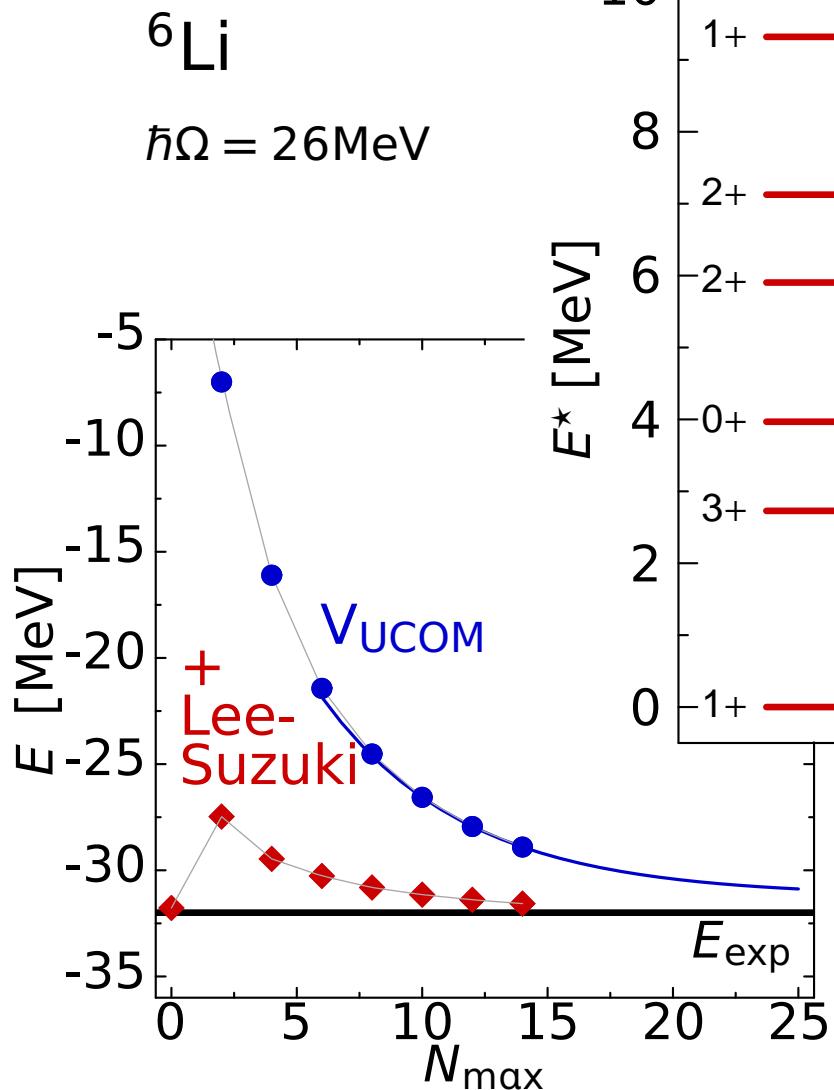
MIN,  $I_9 = 0.09 \text{ fm}^3$

$V_{\text{SRG}}$   
 $\bar{\alpha} = 0.03 \text{ fm}^4$



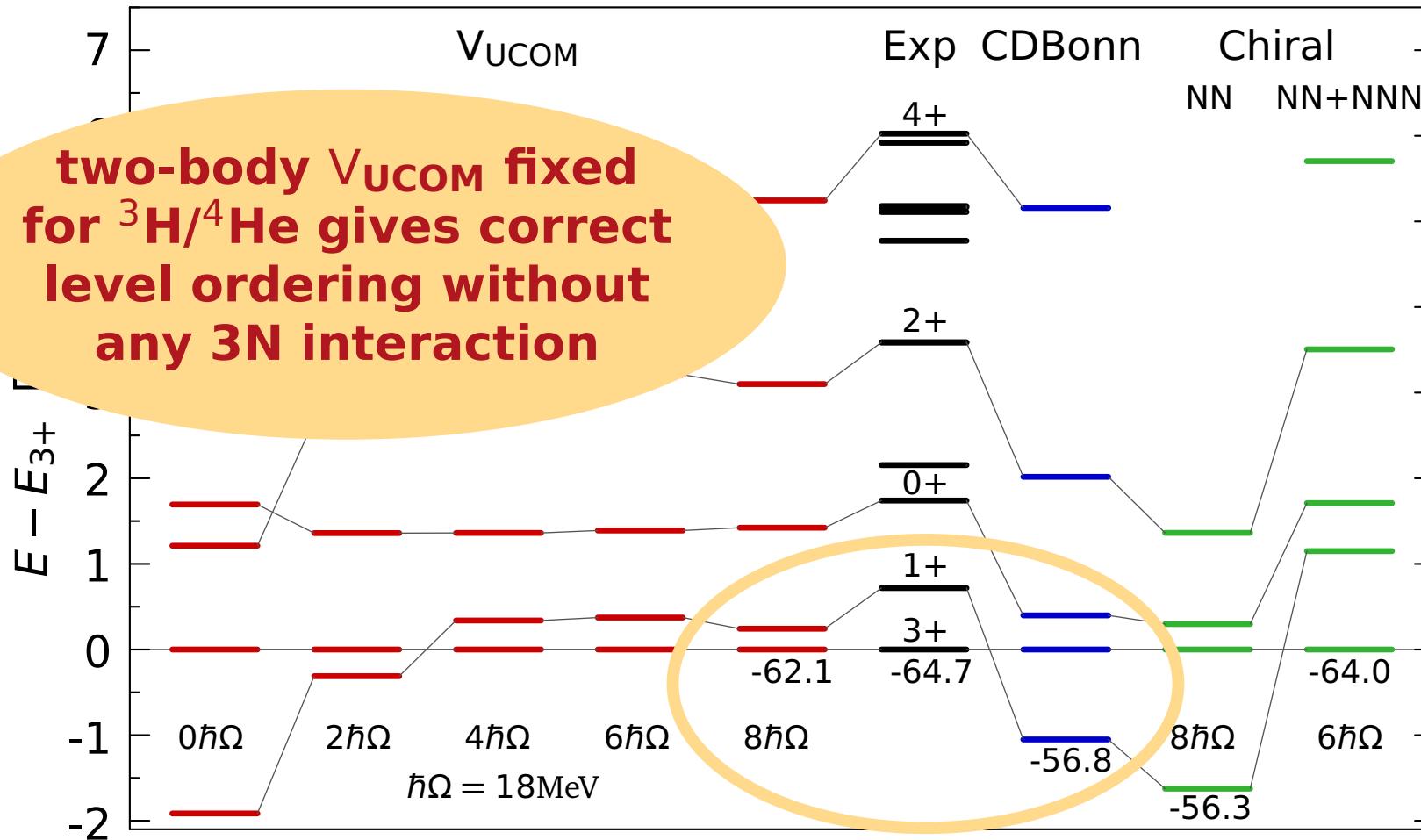
- $I_9$  or  $\bar{\alpha}$  adjusted such that  $^4\text{He}$  binding energy is reproduced

# $^6\text{Li}$ : NCSM throughout the p-Shell



systematic NCSM studies  
for p-shell nuclei up to  
 $A \lesssim 13$  possible

# $^{10}\text{B}$ : Hallmark of a 3N Interaction?



Computational Many-Body Methods

# Importance Truncated No-Core Shell Model

Roth — Phys. Rev. C 79, 064324 (2009)

Roth, Gour & Piecuch — Phys. Rev. C 79, 054325 (2009)

Roth & Navrátil — Phys. Rev. Lett. 99, 092501 (2007)

# Importance Truncated NCSM

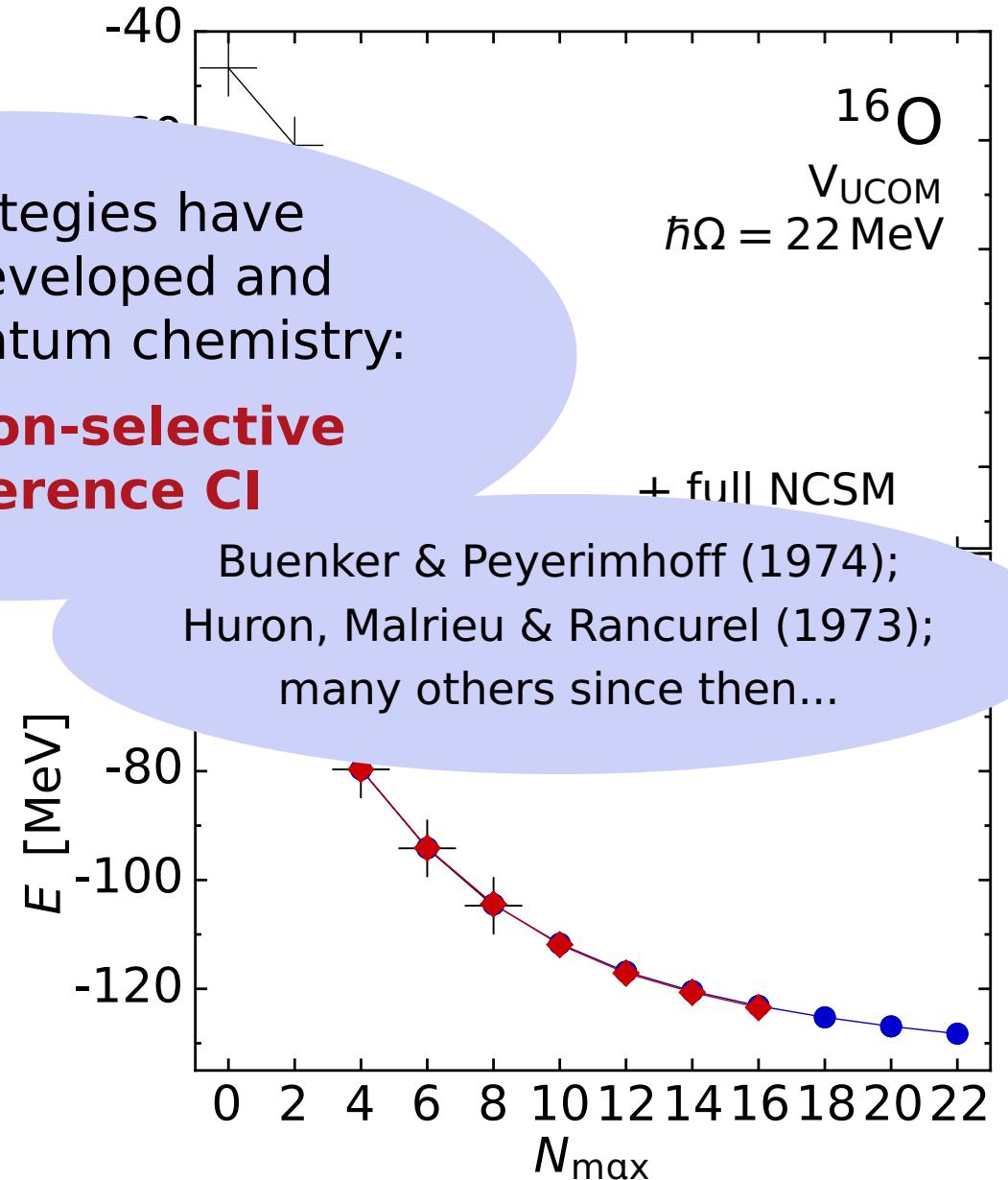
- converged NCSM calculations are essentially restricted to p-shell nuclei
- full 10 orders of truncation for  $^{16}\text{O}$  required (basis dimension ~ 10<sup>10</sup>)

## Importance Truncation

reduce NCSM space to the relevant basis states using an **a priori importance measure** derived from MBPT

similar strategies have first been developed and applied in quantum chemistry:  
**configuration-selective multireference CI**

Buenker & Peyerimhoff (1974);  
Huron, Malrieu & Rancurel (1973);  
many others since then...



# Importance Truncation: General Idea

- given an initial approximation  $|\Psi_{\text{ref}}\rangle$  for the **target state** within a limited **reference space**  $\mathcal{M}_{\text{ref}}$

$$|\Psi_{\text{ref}}\rangle = \sum_{\nu \in \mathcal{M}_{\text{ref}}} C_{\nu}^{(\text{ref})} |\Phi_{\nu}\rangle$$

- **measure the importance** of individual basis state  $|\Phi_{\nu}\rangle \notin \mathcal{M}_{\text{ref}}$  via first-order multiconfigurational perturbation theory

$$\kappa_{\nu} = -\frac{\langle \Phi_{\nu} | H_{\text{int}} | \Psi_{\text{ref}} \rangle}{\epsilon_{\nu} - \epsilon_{\text{ref}}}$$

- construct **importance-truncated space**  $\mathcal{M}(\kappa_{\min})$  spanned by basis states with  $|\kappa_{\nu}| \geq \kappa_{\min}$
- **solve eigenvalue problem** in importance truncated space  $\mathcal{M}(\kappa_{\min})$  and obtain improved approximation of target state

# Importance Truncation: Iterative Scheme

- non-zero importance measure  $\kappa_\nu$  only for states which **differ from  $|\Psi_{\text{ref}}\rangle$  by 2p2h excitation** at most

## IT-NCSM[i] or IT-CI[i]

- simple iterative scheme for arbitrary many-body model spaces

★ start with  $|\Psi_{\text{ref}}\rangle = |\Phi_0\rangle$

① construct importance truncated space containing up to 2p2h on top of  $|\Psi_{\text{ref}}\rangle$

② solve eigenvalue problem

③ use dominant components of eigenstate ( $|C_\nu| \geq C_{\min}$ ) as new  $|\Psi_{\text{ref}}\rangle$ , goto ①

## IT-NCSM(seq)

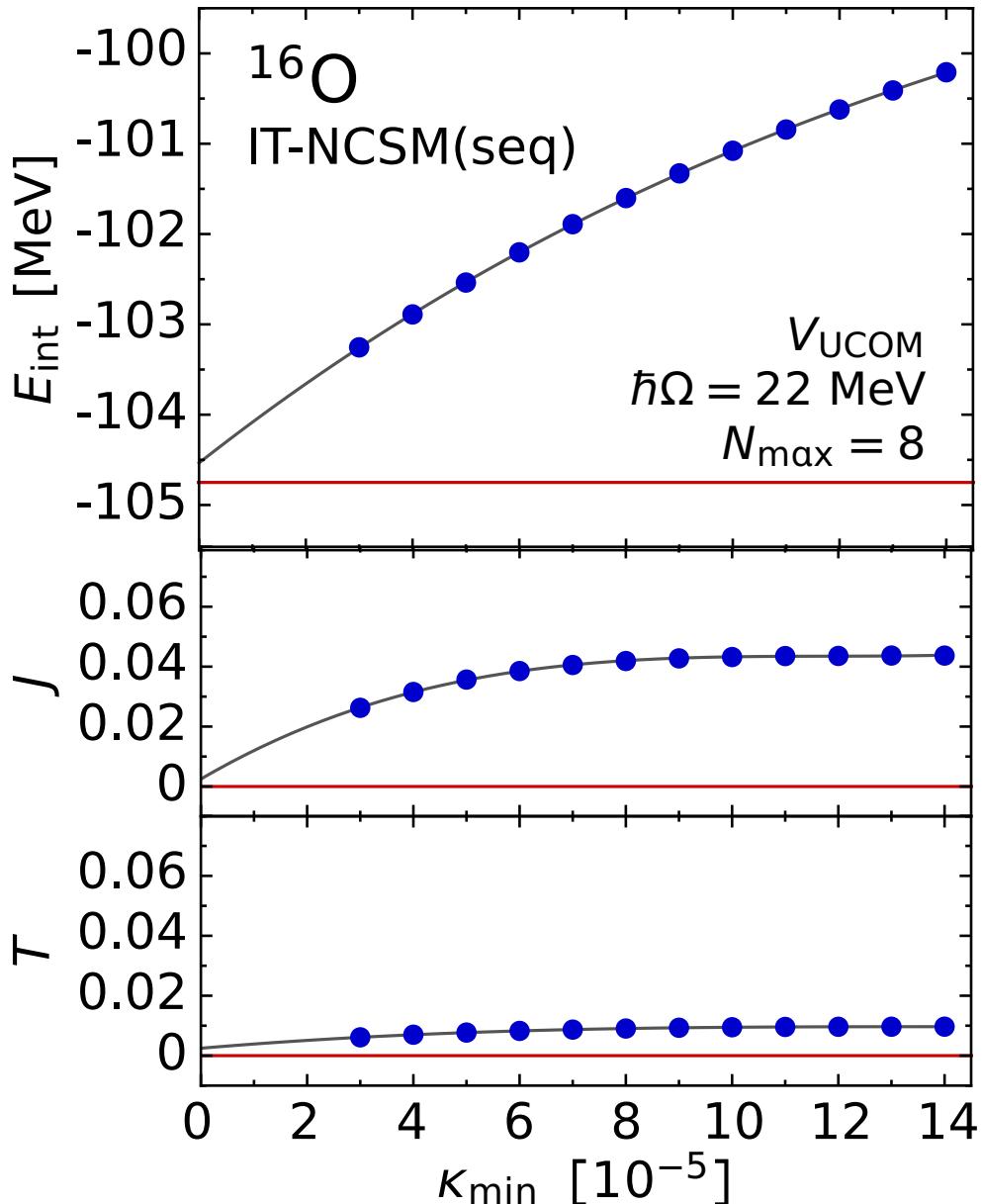
- sequential update scheme for a set of  $N_{\max}\hbar\Omega$  spaces

★ start with  $N_{\max} = 2$  eigenstate from full NCSM as initial  $|\Psi_{\text{ref}}\rangle$

① construct importance truncated space for  $N_{\max} + 2$

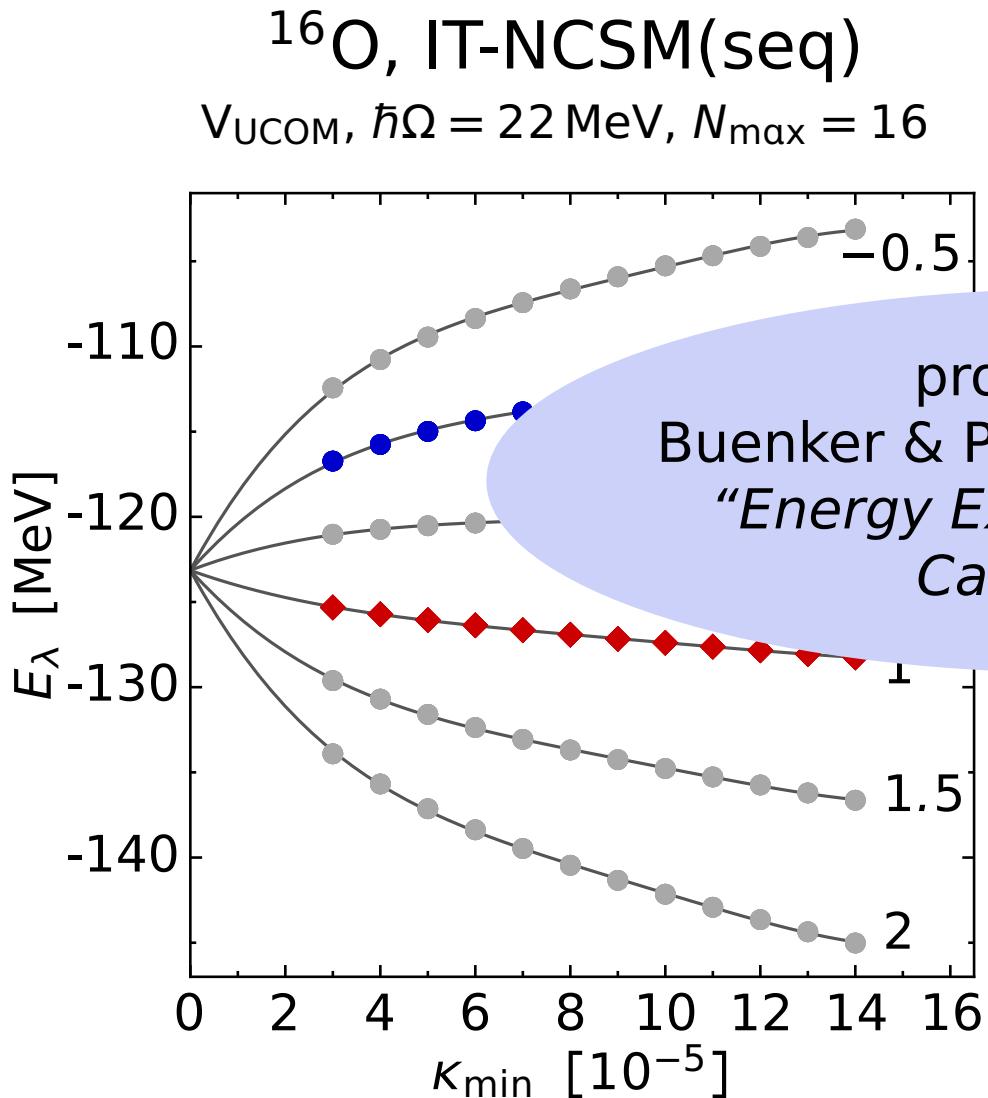
**full NCSM model space is recovered in the limit  $(K_{\min}, C_{\min}) \rightarrow 0$  in IT-NCSM(seq) and IT-NCSM[i<sub>conv</sub>]**

# Threshold Dependence



- do calculations for a **sequence of importance thresholds**  $K_{\min}$
- observables show smooth threshold dependence
- systematic approach to the **full NCSM limit**
- use **a posteriori extrapolation**  $K_{\min} \rightarrow 0$  of observables to account for effect of excluded configurations

# Constrained Threshold Extrapolation

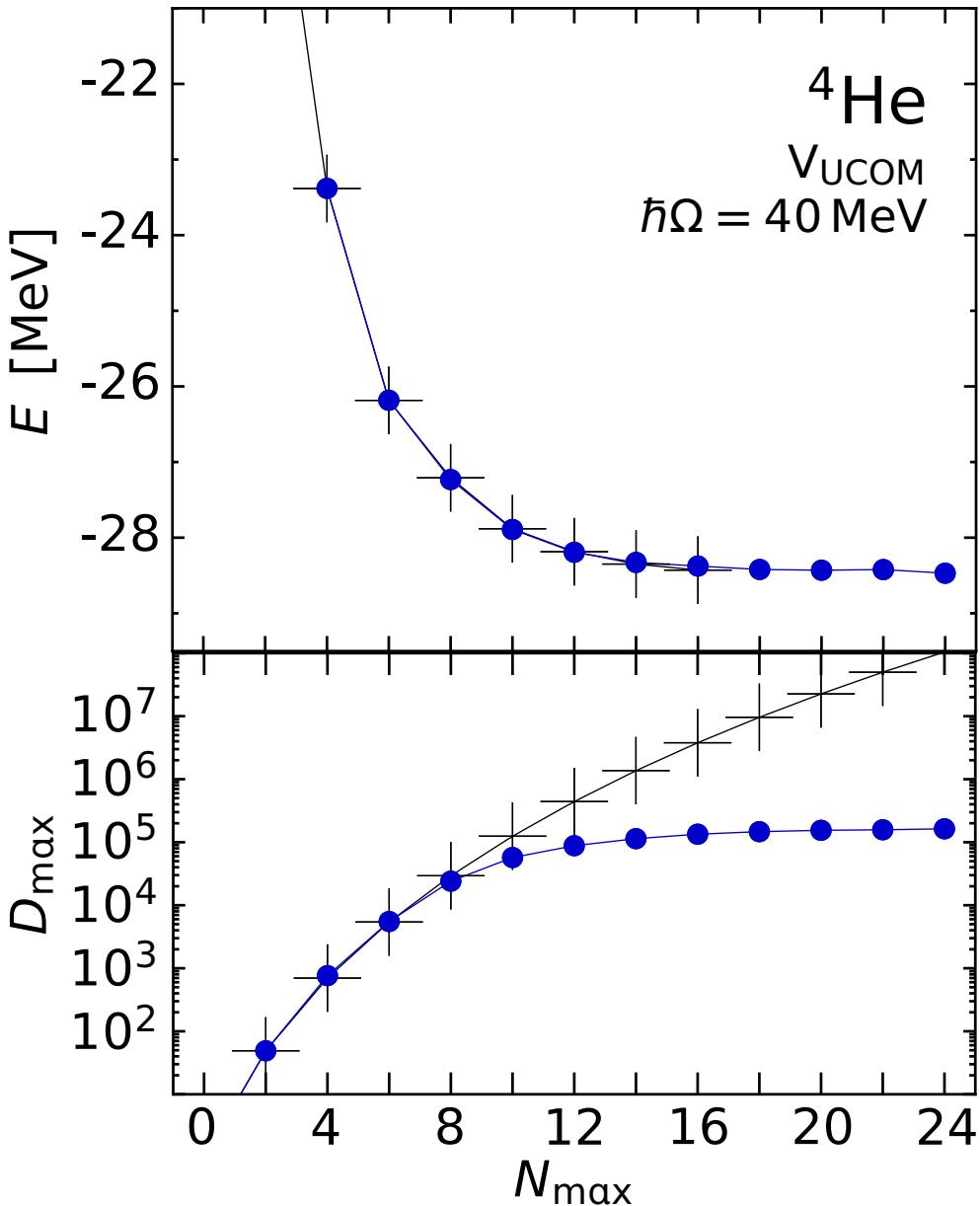


- estimate energy contribution of **excluded states** perturbatively  $\rightarrow \Delta_{\text{excl}}(\kappa_{\min})$

**Hogeneous fit** of combined  $\Delta_{\text{excl}}(\kappa_{\min})$   
for set of  $\lambda$ -values with the constraint  $E_\lambda(0) = E_{\text{extrap}}$

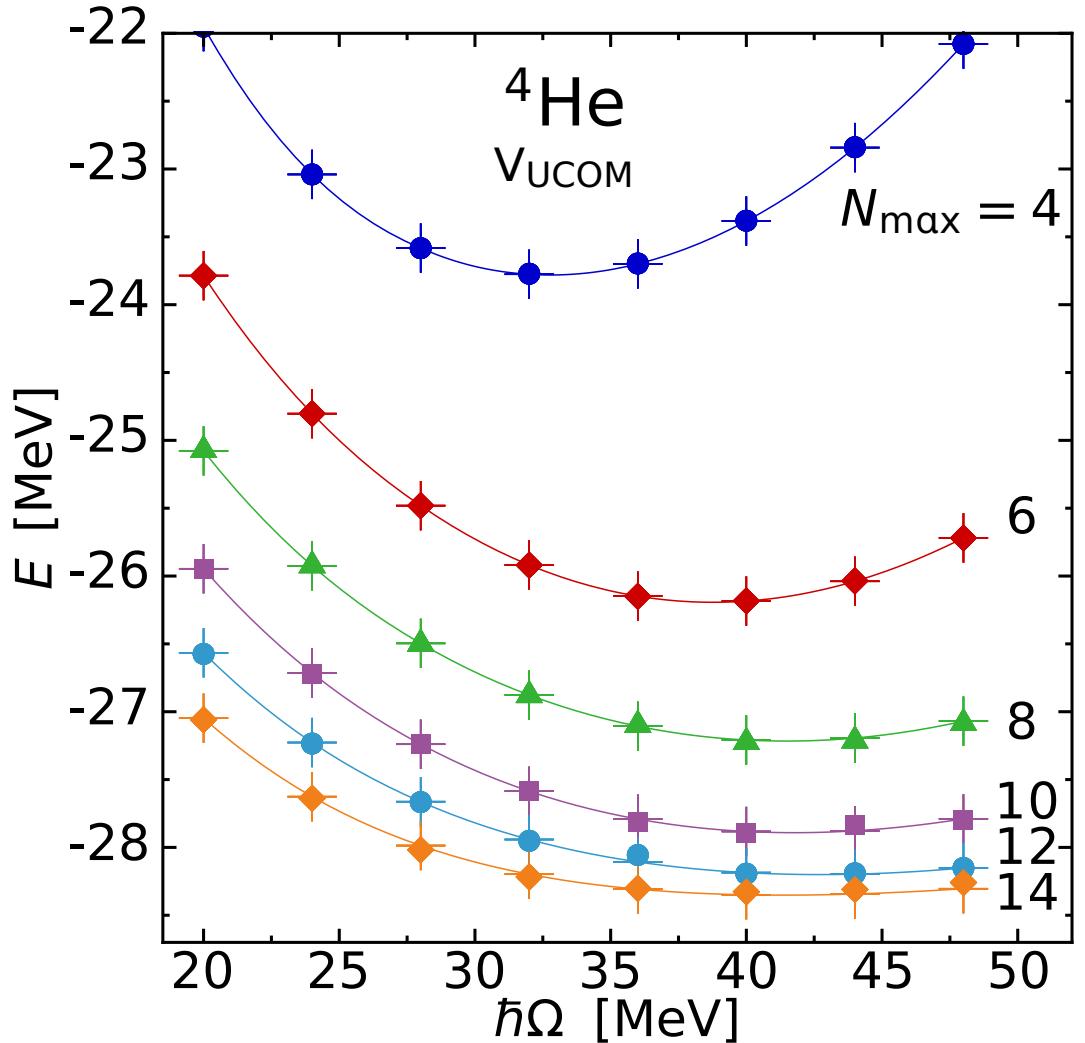
- **robust threshold extrapolation** with error bars determined by variation of the  $\lambda$  set

# $^4\text{He}$ : Importance-Truncated NCSM



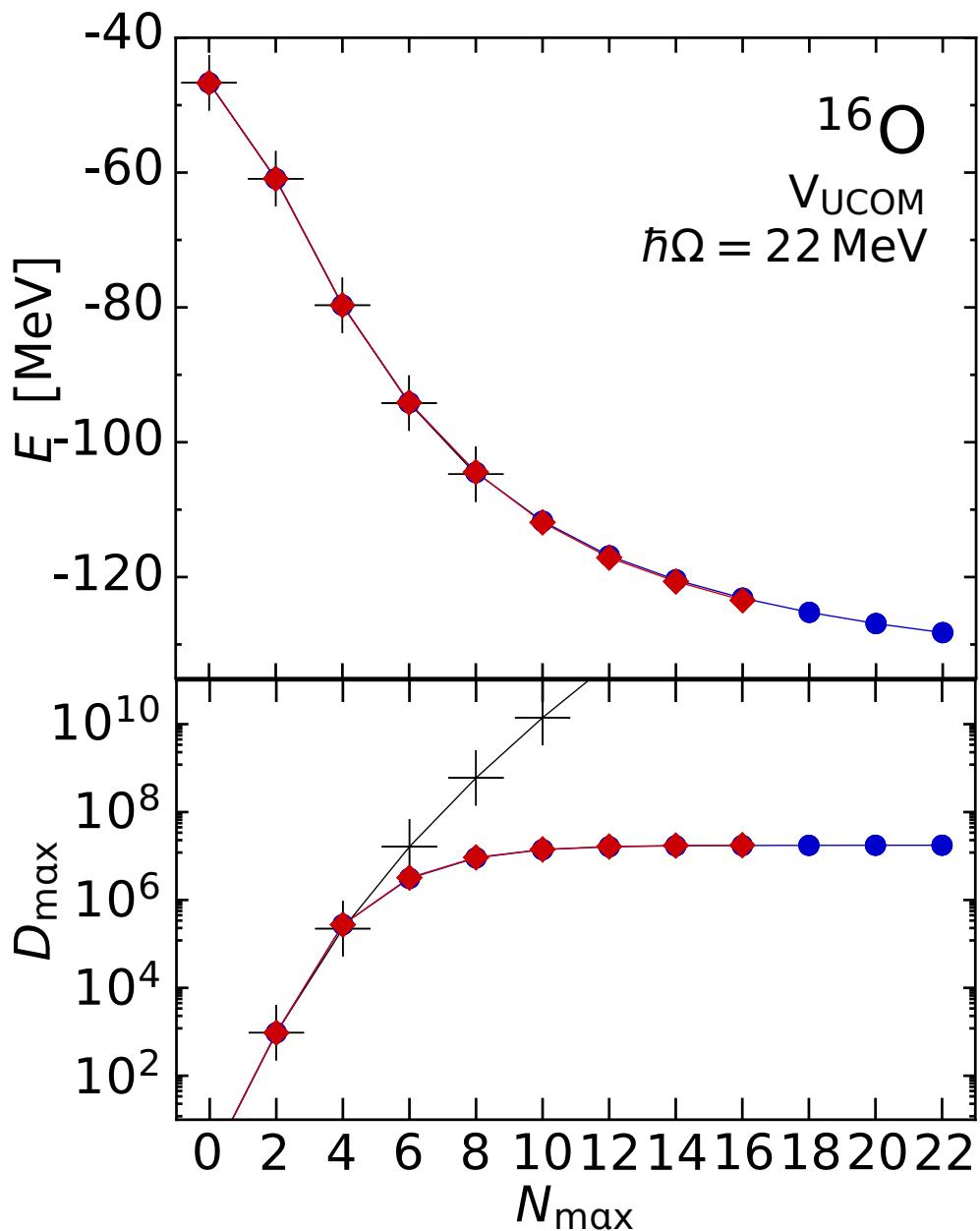
- **sequential IT-NCSM(seq):** single importance update using  $(N_{\max} - 2)\hbar\Omega$  eigenstate as reference
  - **reproduces exact NCSM result** for all  $N_{\max}$
  - reduction of basis by more than two orders of magnitude w/o loss of precision
- + full NCSM  
● IT-NCSM(seq)

# $^4\text{He}$ : Importance-Truncated NCSM



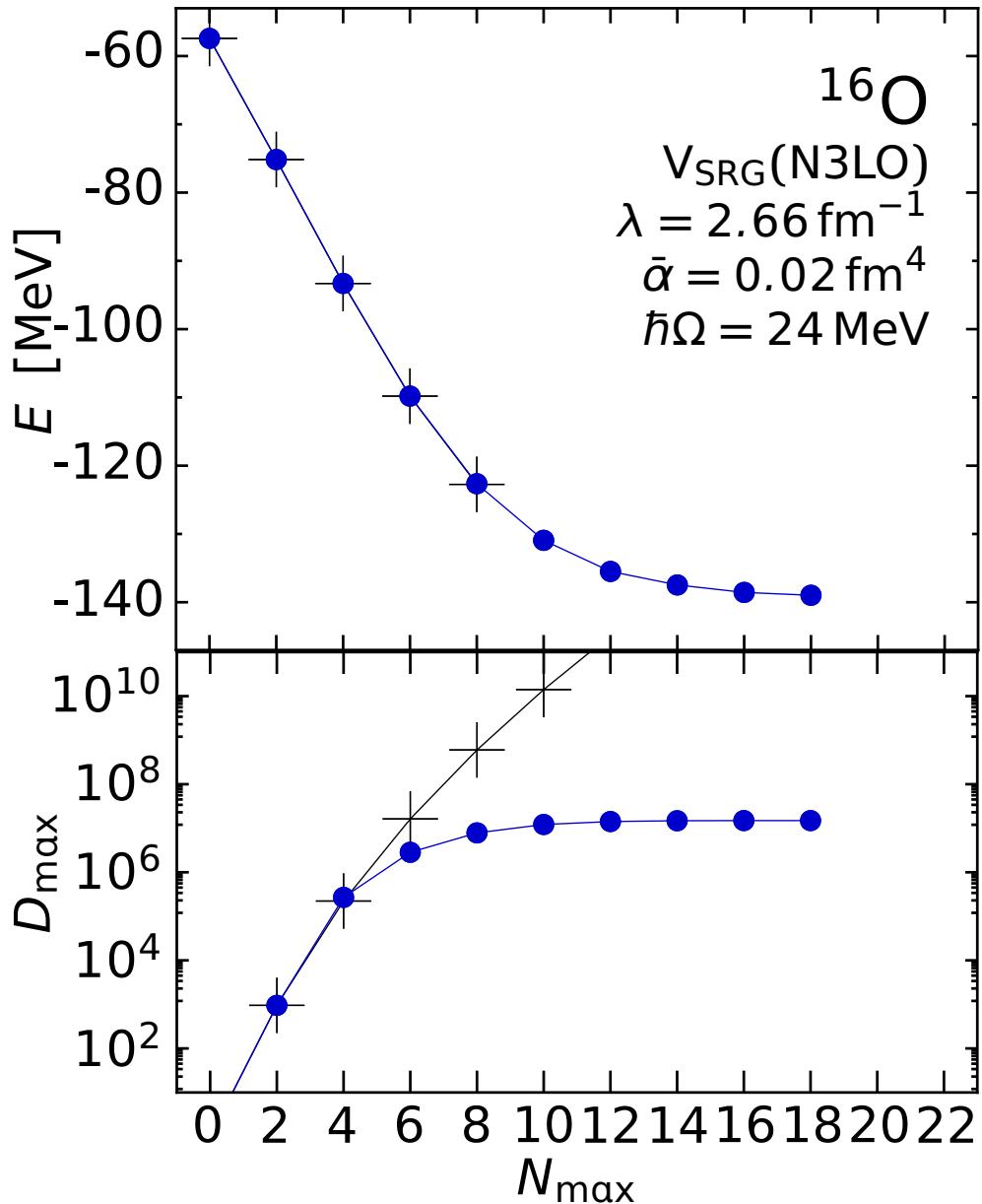
- **reproduces exact NCSM result** for all  $\hbar\Omega$  and  $N_{\max}$
  - importance truncation & threshold extrapolation is robust
  - **no center-of-mass contamination** for any  $N_{\max}$  and  $\hbar\Omega$
- + full NCSM  
● IT-NCSM(seq)

# $^{16}\text{O}$ : Importance-Truncated NCSM



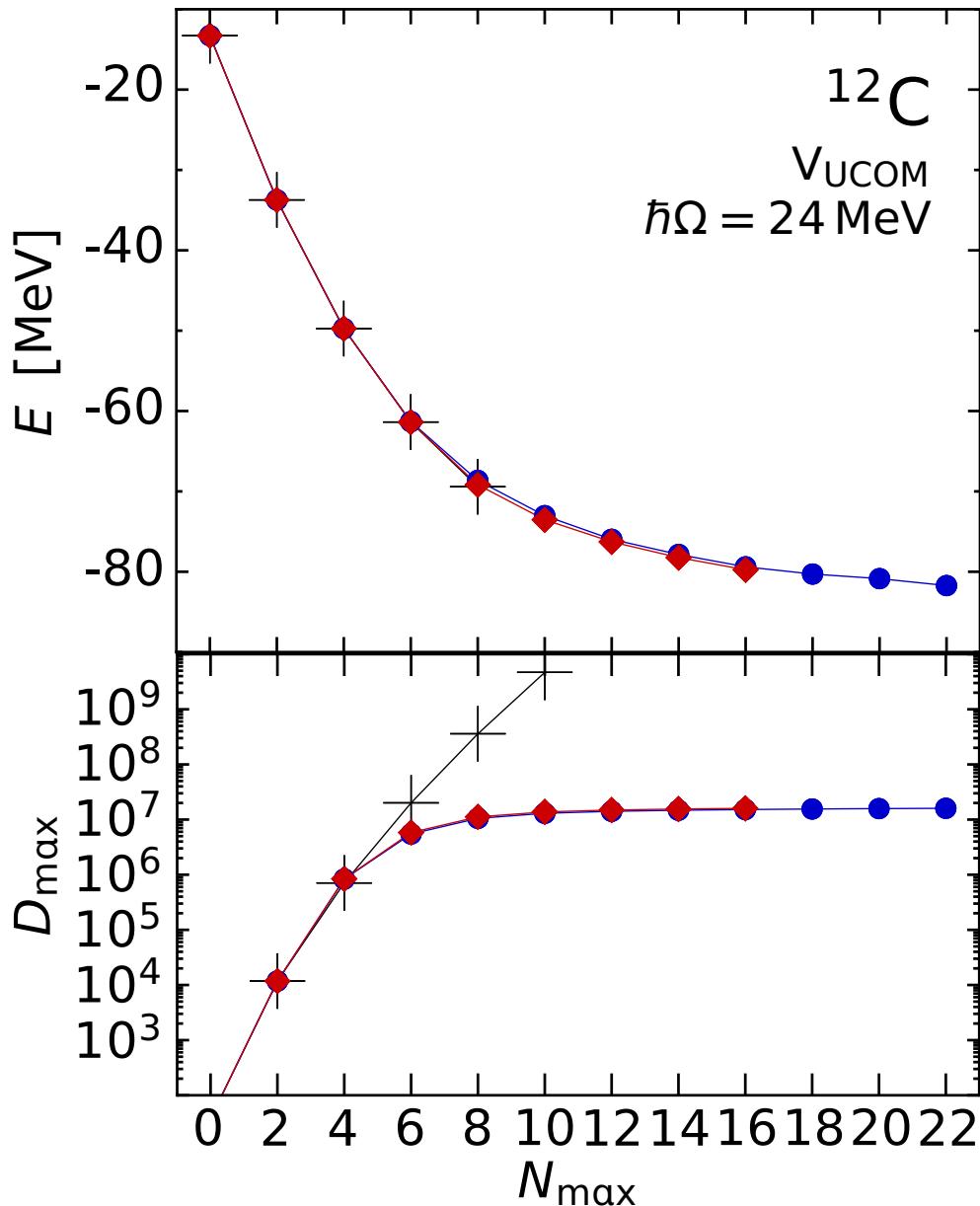
- IT-NCSM(seq) provides **excellent agreement with full NCSM** calculation
- dimension reduced by **several orders of magnitude**
- possibility to go **way beyond** the domain of the full NCSM

# $^{16}\text{O}$ : Importance-Truncated NCSM



- **SRG-evolved N3LO potential** provides a much better convergence behavior
- nevertheless,  $N_{\max} \leq 8$  calculations are not sufficient
- non-exponential behavior observed with  $V_{\text{UCOM}}$  is really due to interaction

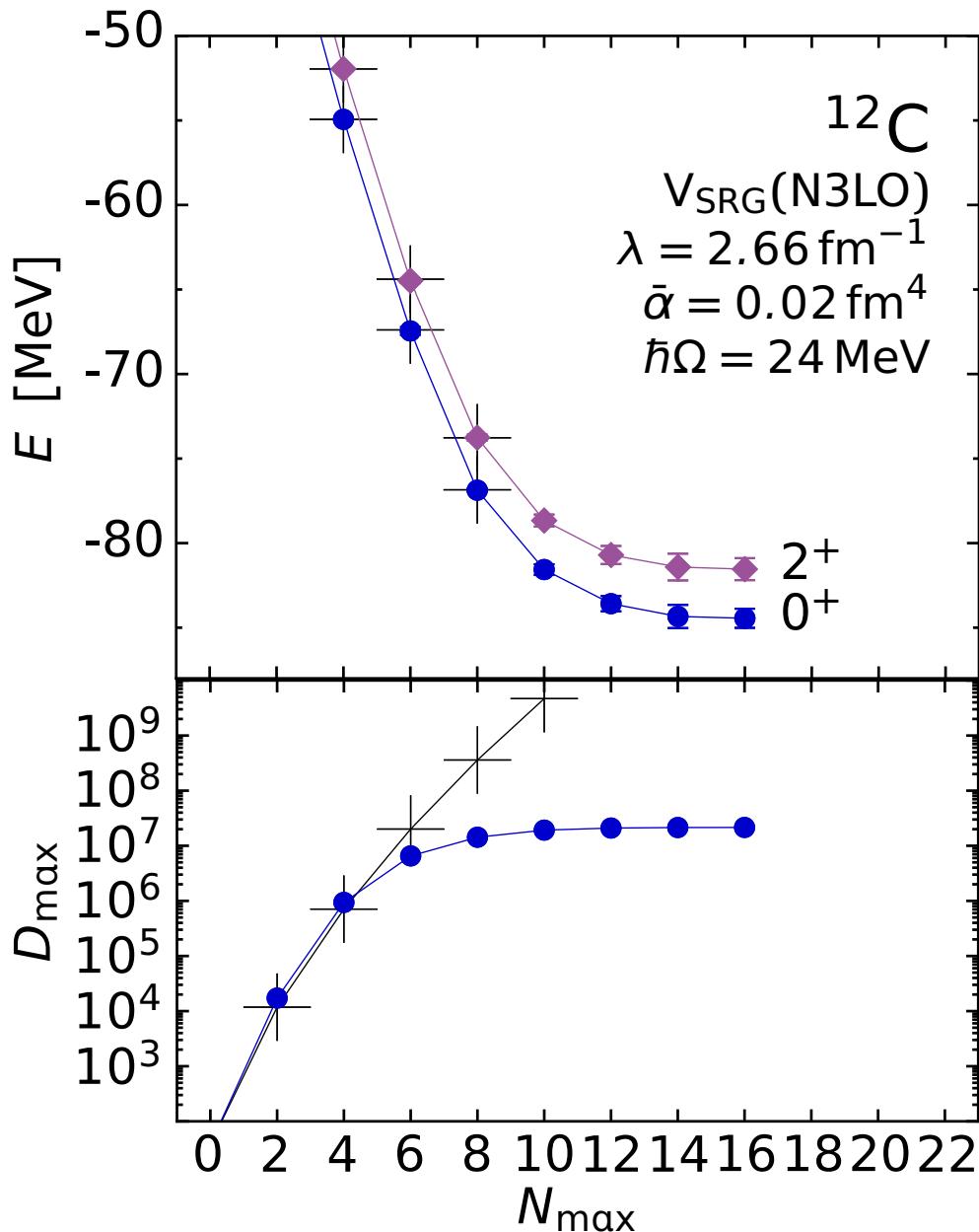
# $^{12}\text{C}$ : IT-NCSM for Open-Shell Nuclei



- excellent agreement with full NCSM calculations
- IT-NCSM(seq) works just as well for non-magic / open-shell nuclei
- all calculations limited by available two-body matrix elements & CPU time only

+ full NCSM  
● IT-NCSM(seq),  $C_{\min} = 0.0005$   
◆ IT-NCSM(seq),  $C_{\min} = 0.0003$

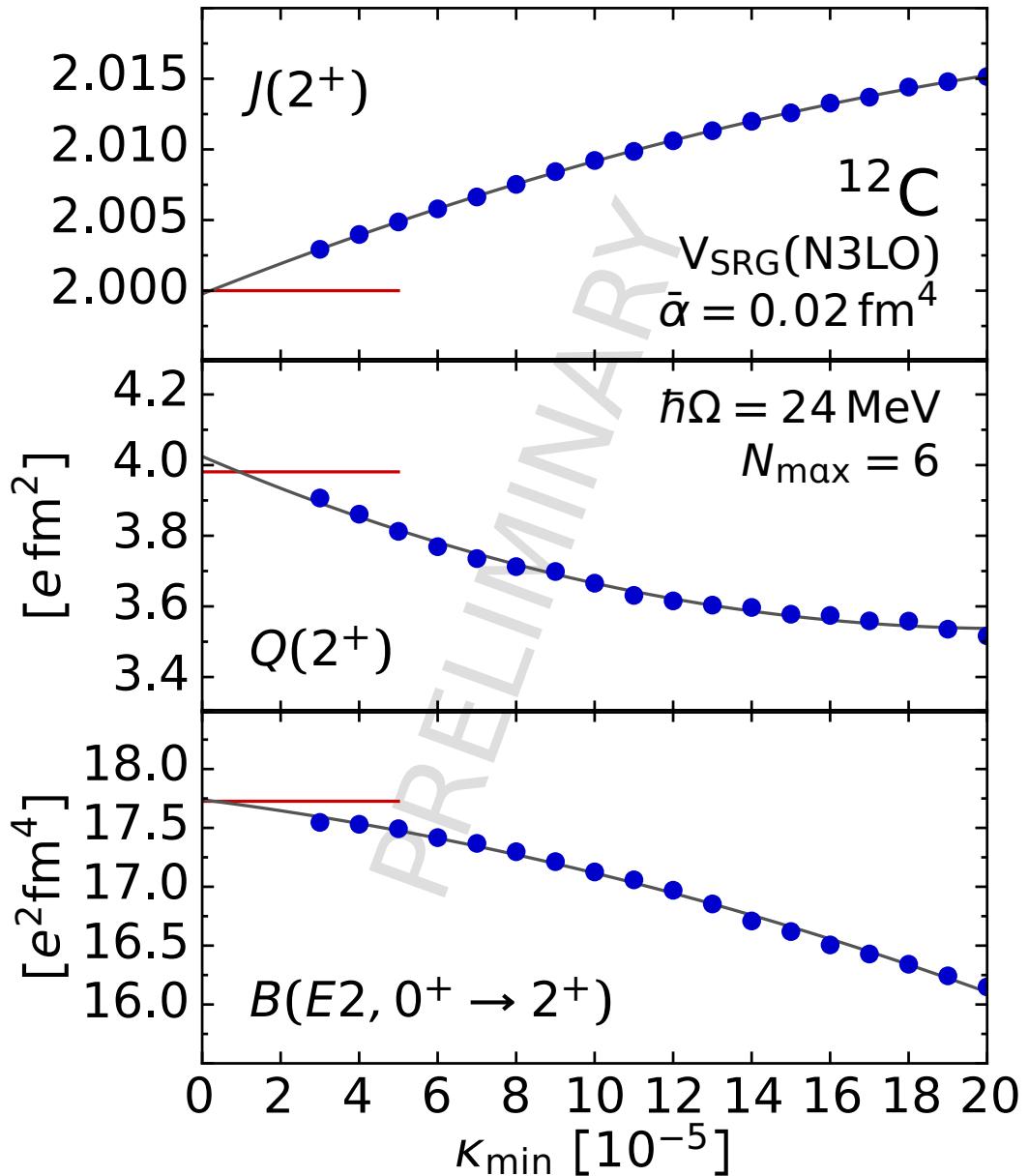
# $^{12}\text{C}$ : IT-NCSM for Excited States



- target ground & excited states simultaneously
  - separate importance measure  $\kappa_\nu^{(n)}$  for each target state
  - basis state is included if  $|\kappa_\nu^{(n)}| \geq \kappa_{\min}$  for any  $n$
- dimension of importance truncated space **grows linearly** with # of target states

+ full NCSM  
● IT-NCSM(seq),  $C_{\min} = 0.0005$

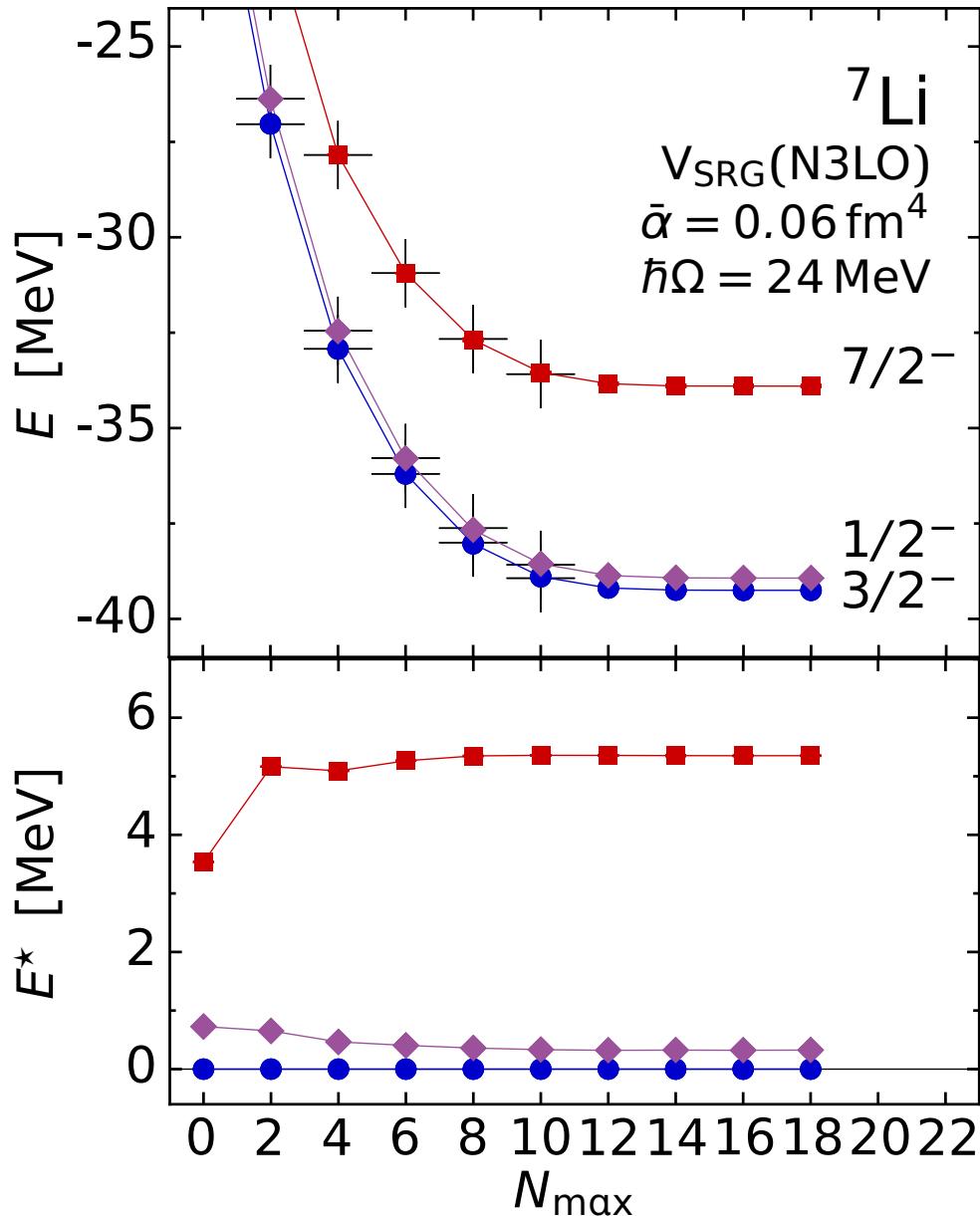
# $^{12}\text{C}$ : IT-NCSM for Spectroscopy



- access to **spectroscopic observables** via eigenstates
- multipole moments, transition strengths, transition form-factors, densities,...
- simple threshold extrapolation essentially **reproduces full NCSM results**

**systematic spectroscopy in p- and sd-shell** with large  $N_{\max}\hbar\Omega$  spaces

# $^7\text{Li}$ : IT-NCSM for Odd Nuclei



- IT-NCSM(seq) treats a ground state & low-lying excited states for open- and closed-shell nuclei **on the same footing**

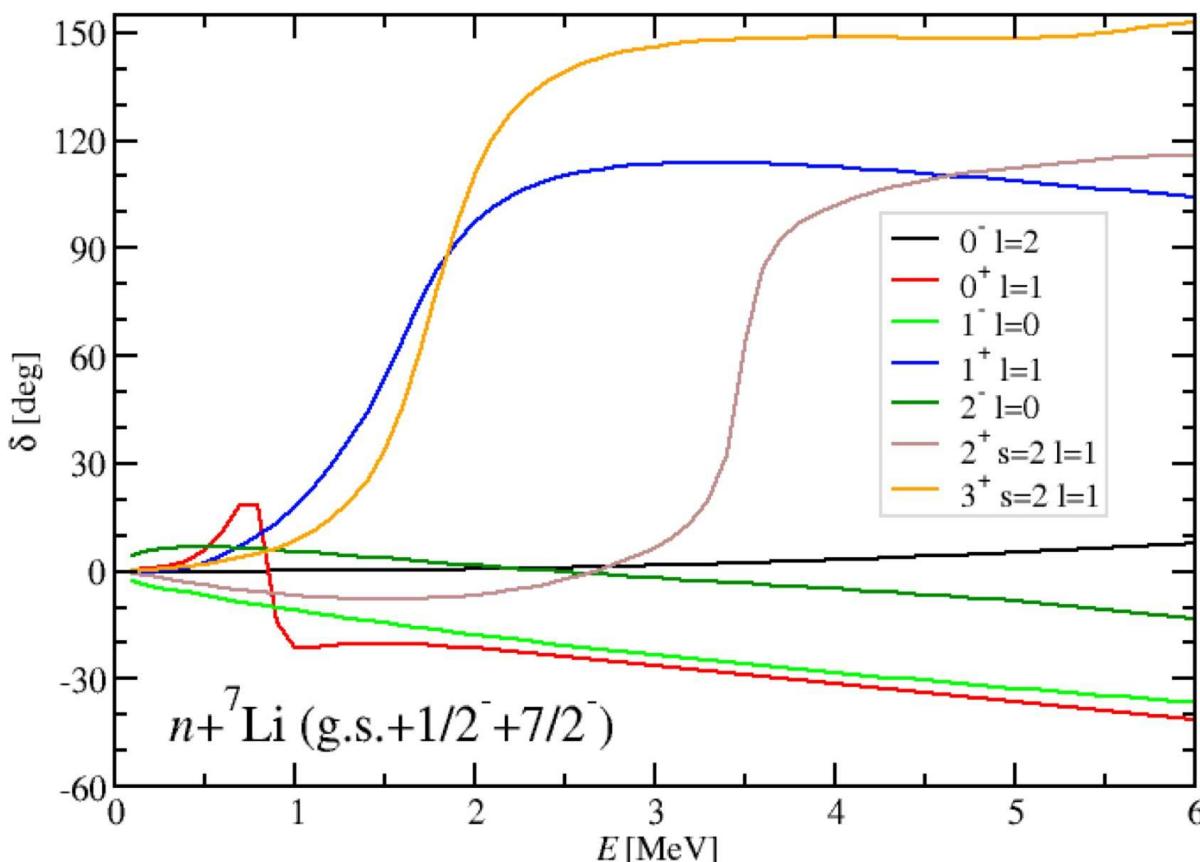
- **excellent agreement with full NCSM** calculations in all cases

+ full NCSM  
●♦■ IT-NCSM(seq),  $C_{\min} = 0.0002$

# RGM & IT-NCSM: Ab Initio Reactions

with Petr Navrátil & Sofia Quaglioni (LLNL)

- **IT-NCSM wave function as input for RGM** (Resonating Group Method) calculations of low-energy nucleon-nucleus scattering



- using 3 lowest  ${}^7\text{Li}$  states
- so-far up to  $N_{\max} = 14$ , here  $N_{\max} = 8$
- phase-shifts with full NCSM and IT-NCSM input agree
- 2 bound states for  ${}^8\text{Li}$
- 4 resonances:  $3^+$  and  $1^+$  are known,  $0^+$  and  $2^+$  resonances are predictions

# IT-NCSM: Pros and Cons

- ✓ **fulfills variational principle** & Hylleraas-Undheim theorem
- ✓ **no center-of-mass contamination** induced by importance truncation in  $N_{\max}\hbar\Omega$  space
- ✓ constrained **threshold extrapolation**  $K_{\min} \rightarrow 0$  recovers contribution of excluded configurations efficiently and accurately
- ✓ **open and closed-shell nuclei** with **ground and excited states** can be treated on the same footing
- ✓ **compatible with shell model**: compute any observable from wave functions in SM representation
- **approximate size-extensivity** after threshold extrapolation in IT-NCSM(seq) or IT-NCSM[ $i_{\text{conv}}$ ] – **no explicit  $n p n h$  truncation**
- ✗ computationally still demanding

Computational Many-Body Methods

# Center-of-Mass Diagnostics

Roth, Gour & Piecuch — arXiv:0906.4276

Roth, Gour & Piecuch — Phys. Rev. C 79, 054325 (2009)

# CM Problem: Bane of Nuclear Structure

- **nucleus is a self-bound system**: intrinsic and CM component of the many-body state have to **factorize**

$$|\Psi\rangle = |\psi_{\text{int}}\rangle \otimes |\psi_{\text{cm}}\rangle$$

- factorization is manifest in Jacobi-coordinate methods
- Slater-determinant methods: only the  $N_{\max} \hbar \Omega$  **space** build from **harmonic oscillator basis** allows for exact factorization
- for any other truncation or single-particle basis one has to ask:
  - Is there a **coupling** between intrinsic and CM component?
  - How strong is the **effect on observables** of interest?
- **CM diagnostics**: perturb CM part and check for effect on the intrinsic part via expectation values of intrinsic observables

# CM Diagnostics

- consider model space built from **HO single-particle basis**
- solve many-body problem with **modified Hamiltonian** (following Palumbo, Gloeckner & Lawson)

$$H_\beta = H_{\text{int}} + \beta H_{\text{cm}}$$

including additional HO Hamiltonian w.r.t. the CM

$$H_{\text{cm}} = \frac{1}{2mA} \vec{P}_{\text{cm}}^2 + \frac{mA\Omega^2}{2} \vec{X}_{\text{cm}}^2 - \frac{3}{2}\hbar\Omega .$$

- Why this particular  $H_{\text{cm}}$  operator?
  - ▶ the exact ground state of  $H_{\text{cm}}$  can be represented in model space (embedded  $0\hbar\Omega$  space)
  - ▶ if there is factorization, then  $H_{\text{cm}}$  will not induce a coupling

# CM Diagnostics

- ① **analyze  $\beta$ -dependence of expectation values of  $H_{\text{int}}$**  computed with eigenstates of  $H_\beta$

$$\langle H_{\text{int}} \rangle_\beta = \langle \Psi_\beta | H_{\text{int}} | \Psi_\beta \rangle$$

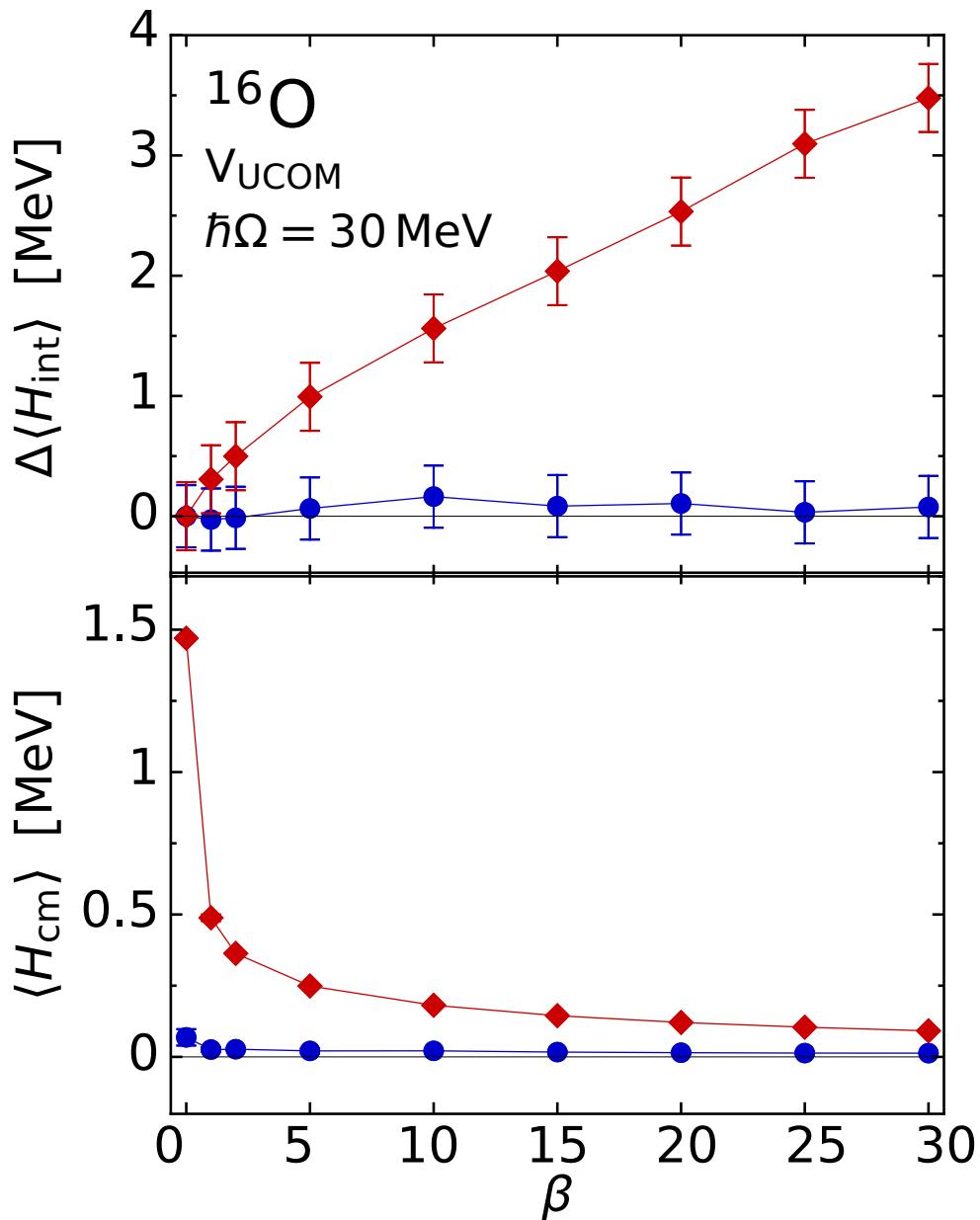
→ any dependence of  $\langle H_{\text{int}} \rangle_\beta$  on  $\beta$  indicates an unphysical coupling between intrinsic and CM motion

- ② **analyze  $\beta$ -dependence of expectation values of  $H_{\text{cm}}$**  computed with eigenstates of  $H_\beta$

$$\langle H_{\text{cm}} \rangle_\beta = \langle \Psi_\beta | H_{\text{cm}} | \Psi_\beta \rangle$$

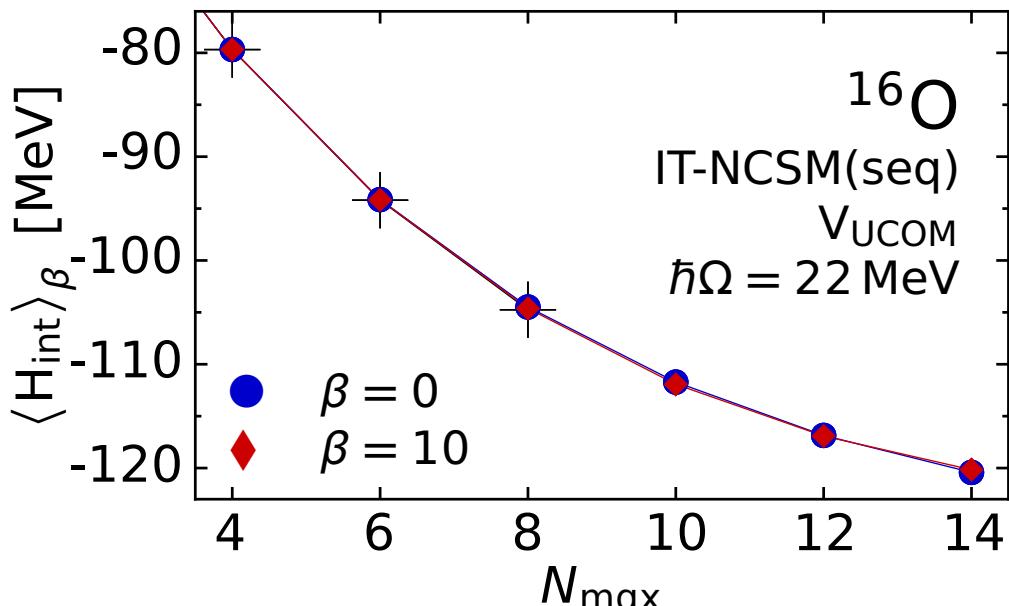
→ any non-zero value of  $\langle H_{\text{cm}} \rangle_\beta$  for  $\beta > 0$  indicates an unphysical coupling between intrinsic and CM motion

# CM Diagnostics: IT-NCSM vs. IT-CI

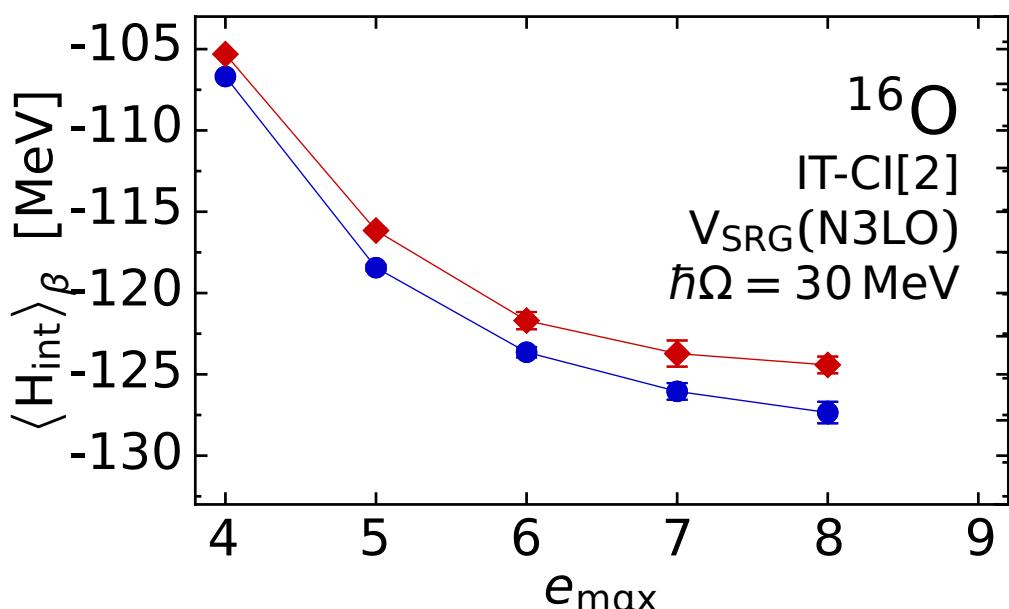


- IT-NCSM(seq) –  $N_{\text{max}}\hbar\Omega$  trunc.
    - $\Delta\langle H_{\text{int}} \rangle_\beta$  and  $\langle H_{\text{cm}} \rangle_\beta$  are practically zero for all  $\beta > 0$
    - **CM is decoupled** to a very good approximation
  
  - IT-CI[2] – single-particle trunc.
    - sizable dependence of  $\Delta\langle H_{\text{int}} \rangle_\beta$  on  $\beta$  and thus **sizable CM contamination**
- IT-NCSM(seq),  $N_{\text{max}} = 8$   
◆ IT-CI[2],  $e_{\text{max}} = 5$

# CM Diagnostics: IT-NCSM vs. IT-CI



- IT-NCSM shows excellent decoupling for all model spaces
  - ▶ importance truncation preserves the **translational invariance** of the  $N_{\text{max}}\hbar\Omega$  space



- IT-CI exhibits **sizable coupling** which does not improve with increasing  $e_{\text{max}}$

# IT-NCSM: Perspectives

importance truncation extends the range of applicability of the NCSM to larger  $N_{\max}$  and  $A$  while preserving most of its advantages

- full **ab-initio spectroscopy** for low-lying states **in p- and sd-shell** ( $A \lesssim 40$ )
- use eigenstates as input for secondary calculation: **RGM for nucleon-nucleus phase shifts**
- include **three-body interactions**, at least approximately
- **algorithmic and conceptual improvements** to extend the mass range

# Epilogue

## ■ thanks to my group & my collaborators

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Lawrence Livermore National Laboratory, USA
- P. Piecuch, J. Gour  
Michigan State University, USA
- H. Feldmeier, T. Neff,...  
Gesellschaft für Schwerionenforschung (GSI)

Deutsche  
Forschungsgemeinschaft  
**DFG**



 LOEWE – Landes-Offensive  
zur Entwicklung Wissenschaftlich-  
ökonomischer Exzellenz