

Nuclear Structure in the UCOM Framework: Few-Body Systems

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Overview

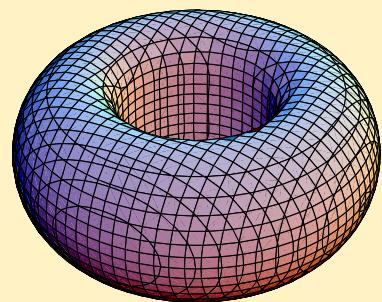
- The Unitary Correlation Operator Method (UCOM)
- No-Core Shell Model (NCSM) Calculations
 - Short- & Long-Range Correlations
 - Tjon-Line
 - p -Shell: ${}^6\text{Li}$ and ${}^{10}\text{B}$
- Summary & Outlook

Unitary Correlation Operator Method (UCOM)

Motivation

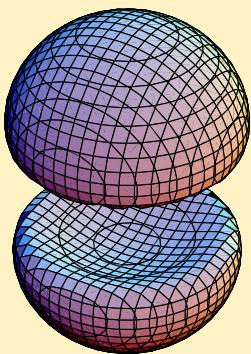
Argonne V18 Deuteron Solution

$$M_S = 0 \\ \frac{1}{\sqrt{2}}(|\uparrow\downarrow\rangle + |\downarrow\uparrow\rangle)$$



$$\rho_{1,M_S}^{(2)}(\vec{r})$$

$$M_S = \pm 1 \\ |\uparrow\uparrow\rangle, |\downarrow\downarrow\rangle$$



- **central correlations:** two-body density is suppressed at low distances

- **tensor correlations:** angular distribution depends on the relative spin alignments

use very large many-body Hilbert spaces
⇒ **high computational effort**

or

use numerically affordable Hilbert spaces and
treat strong correlations explicitly

Central and Tensor Correlators

Central Correlator C_r

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

$$C_r = \exp(-i \sum_{i,j}^A g_{r,ij})$$

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

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

Tensor Correlator C_Ω

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

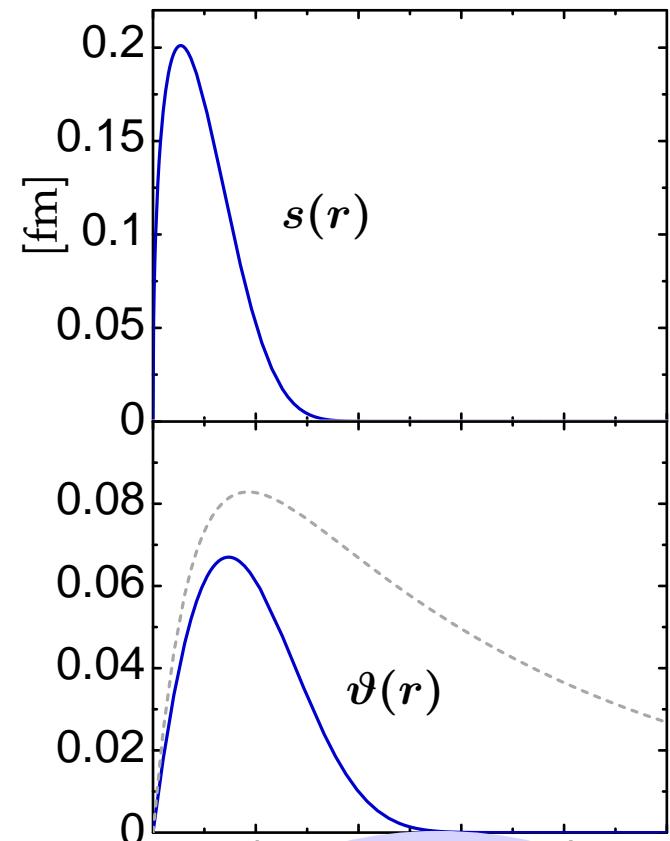
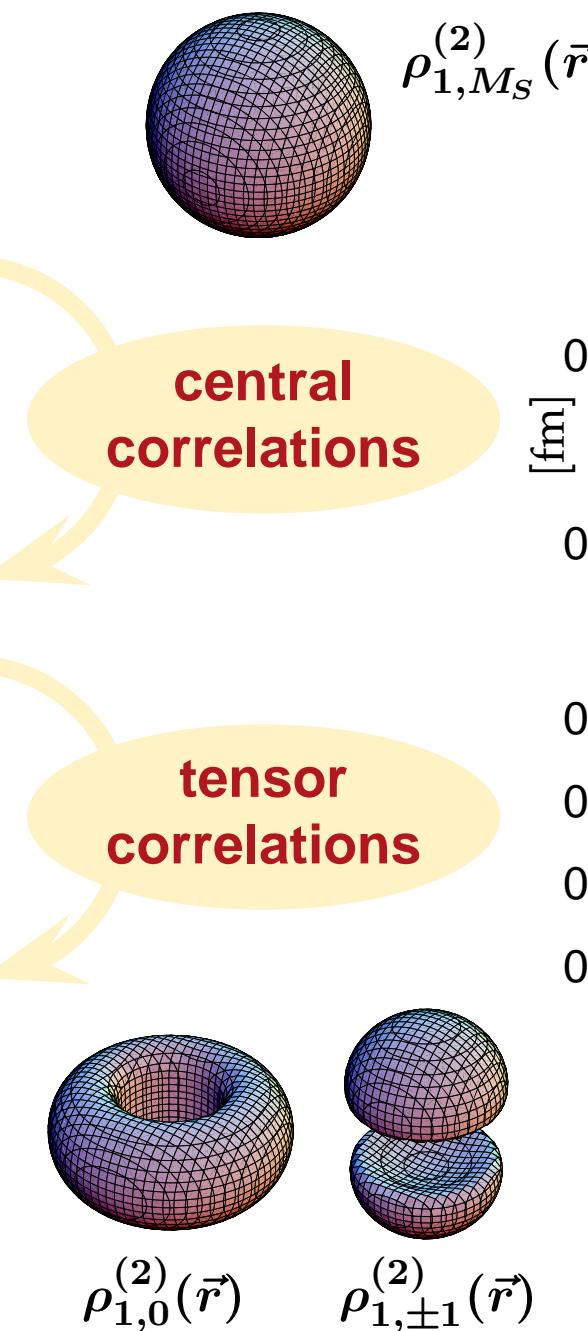
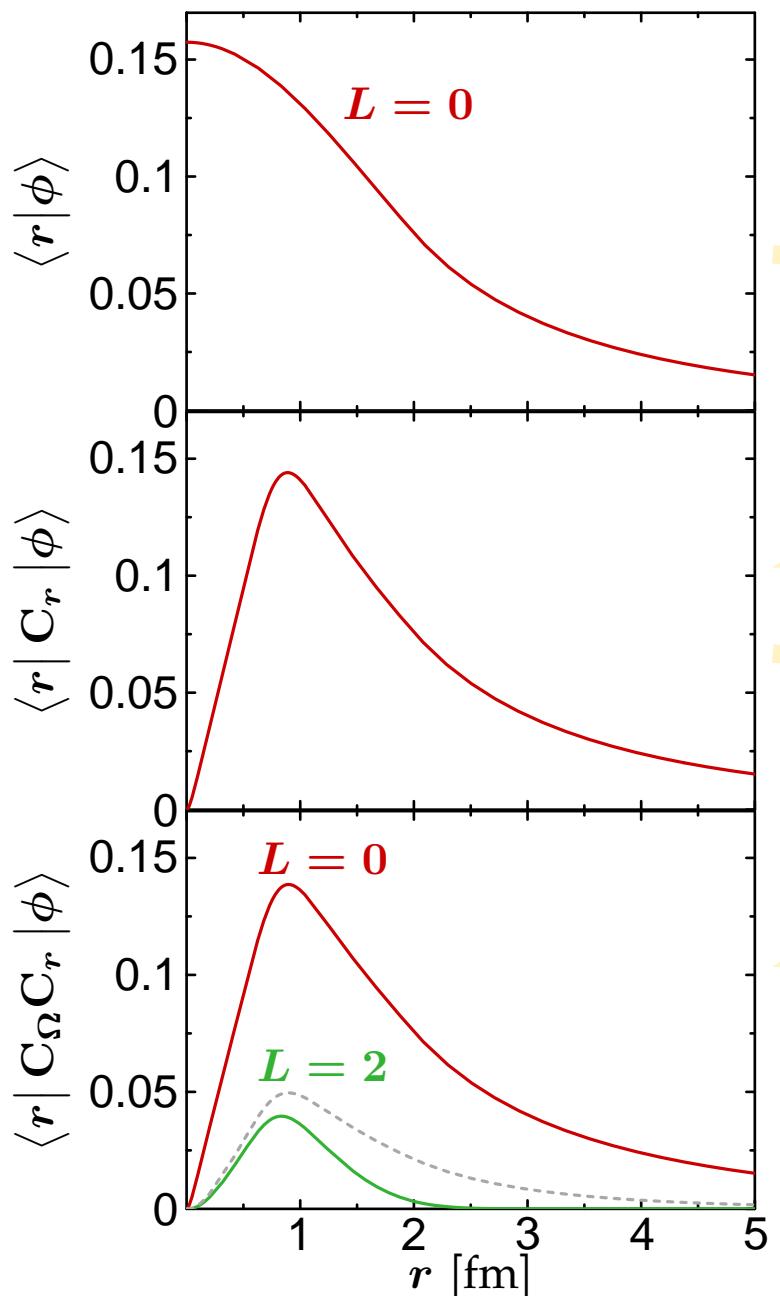
$$C_\Omega = \exp(-i \sum_{i,j}^A g_{\Omega,ij})$$

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

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

$s(r)$ and $\vartheta(r)$
encapsulate the physics of
short-range correlations.

Correlated States



A purple speech bubble at the bottom right states: "tensor corr. range becomes a parameter".

Correlated Interaction

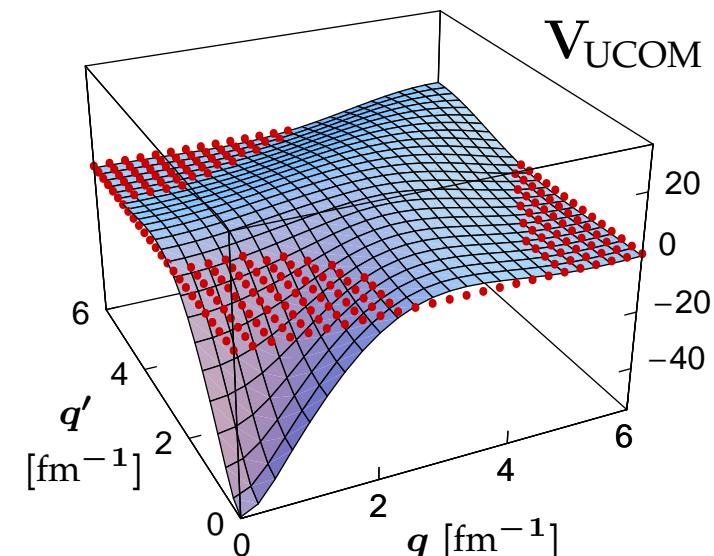
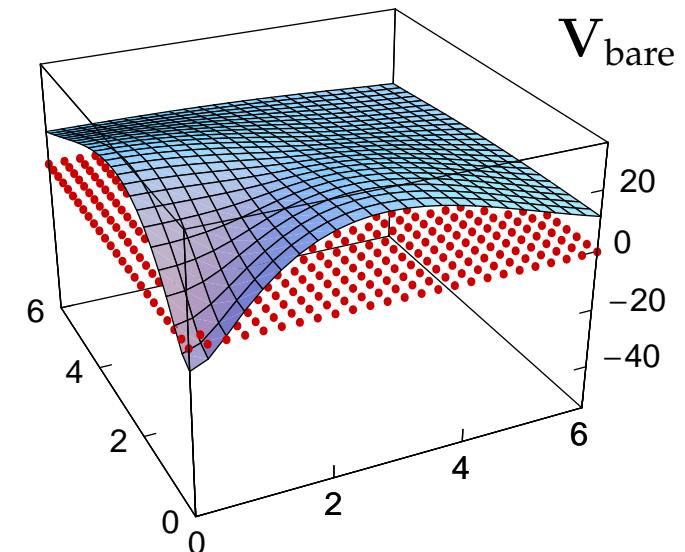
Correlated Hamiltonian

$$\tilde{H} = T^{[1]} + V_{\text{UCOM}} + V_{\text{UCOM}}^{[3]} + \dots$$

- **closed operator representation** of V_{UCOM} in two-body approximation
⇒ usable with **arbitrary many-body basis**
- V_{UCOM} is **phase-shift equivalent** to the underlying bare nucleon-nucleon interaction
- V_{UCOM} is pre-diagonalized in momentum space, i. e. **high-momentum components are decoupled** (similar to $V_{\text{low-}k}$)

AV18

3S_1



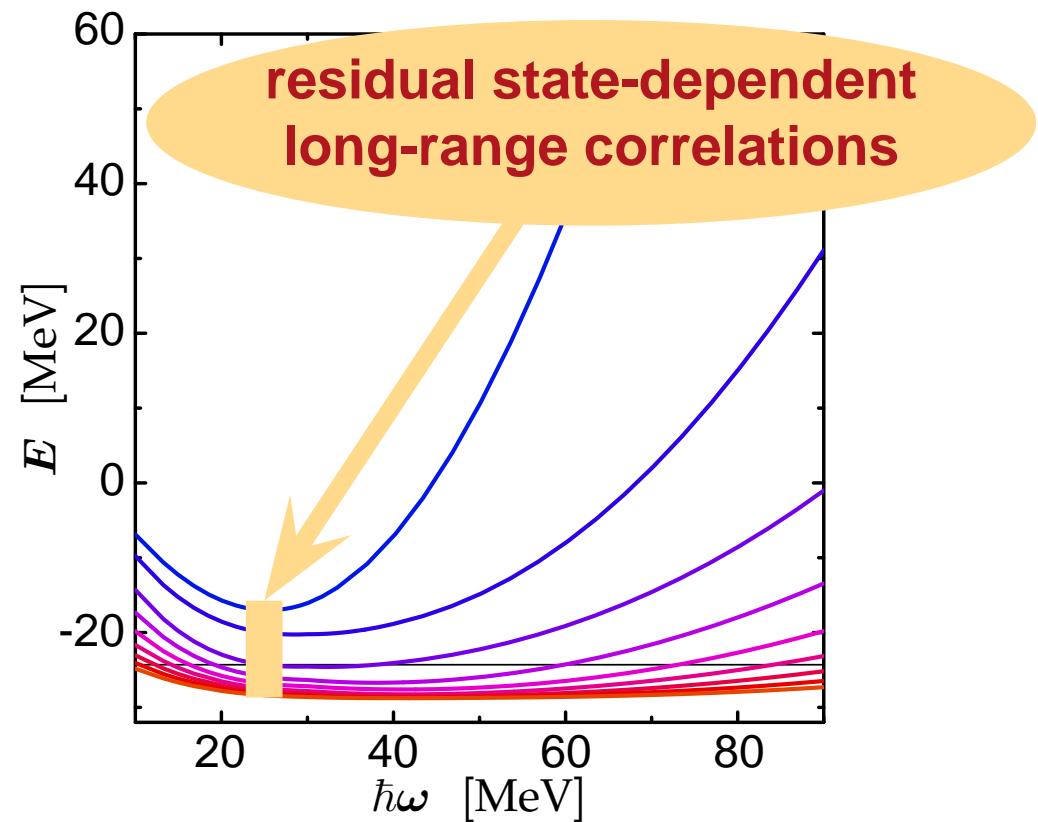
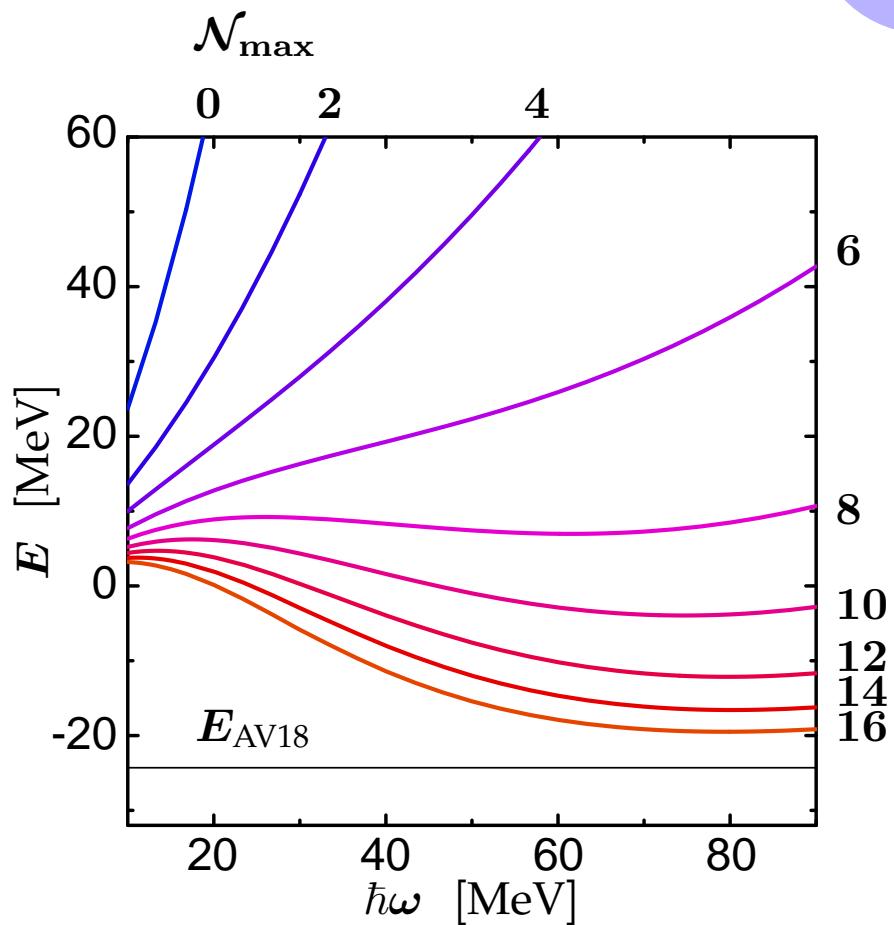
No-Core Shell Model (NCSM) Calculations

^4He : Convergence

AV18

^4He

VUCOM



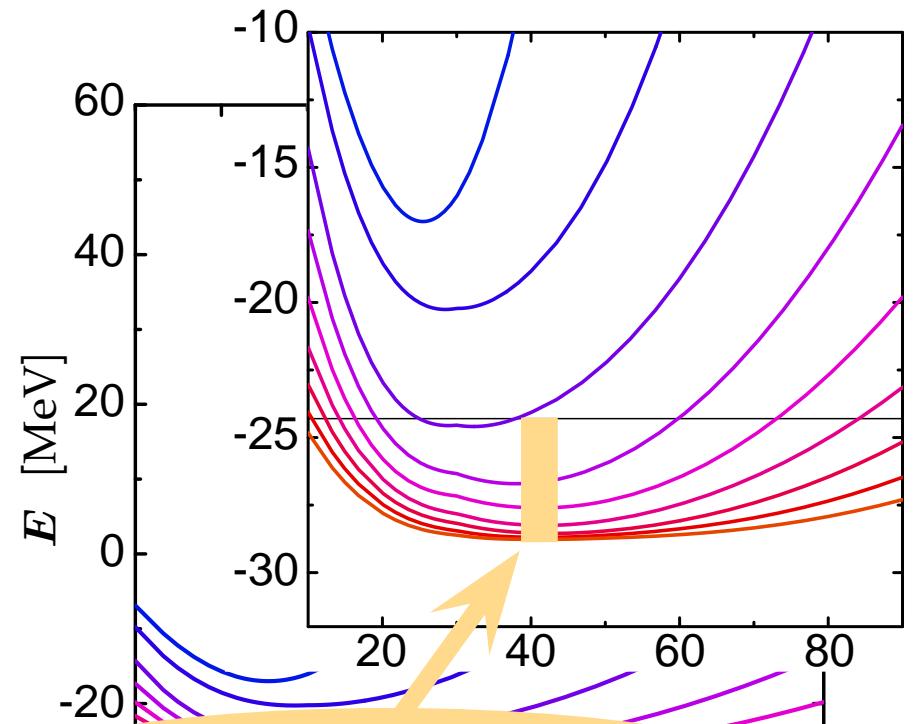
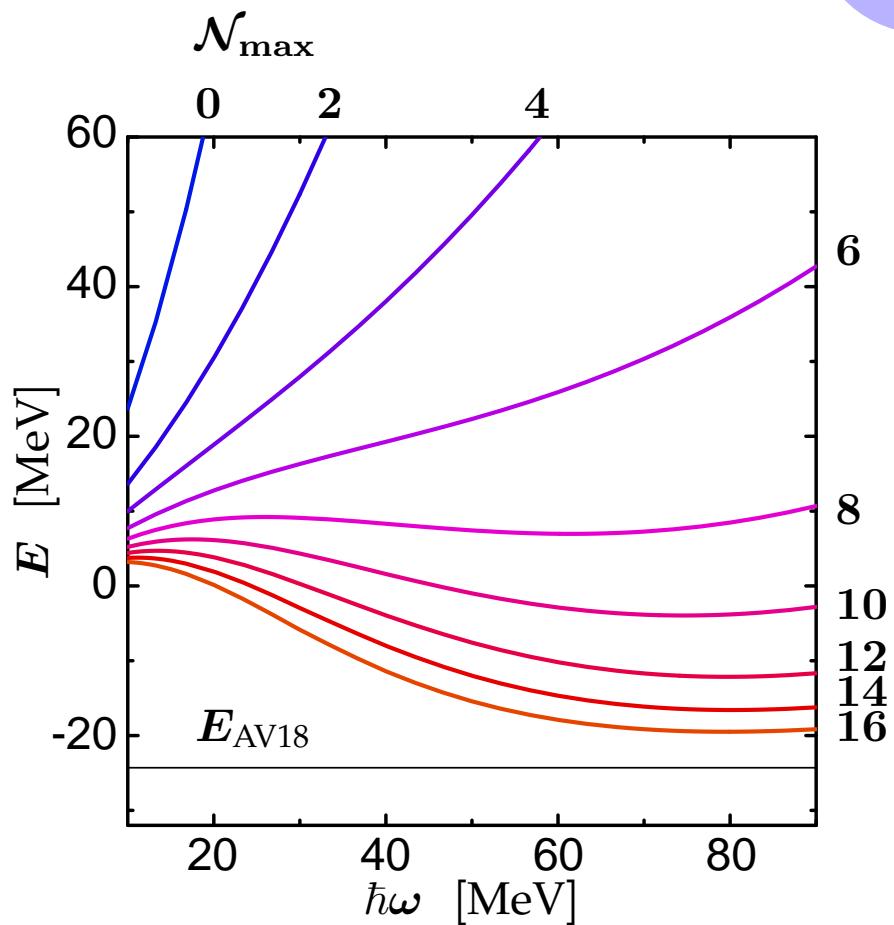
NCSM code by P. Navrátil [PRC 61, 044001 (2000)]

^4He : Convergence

AV18

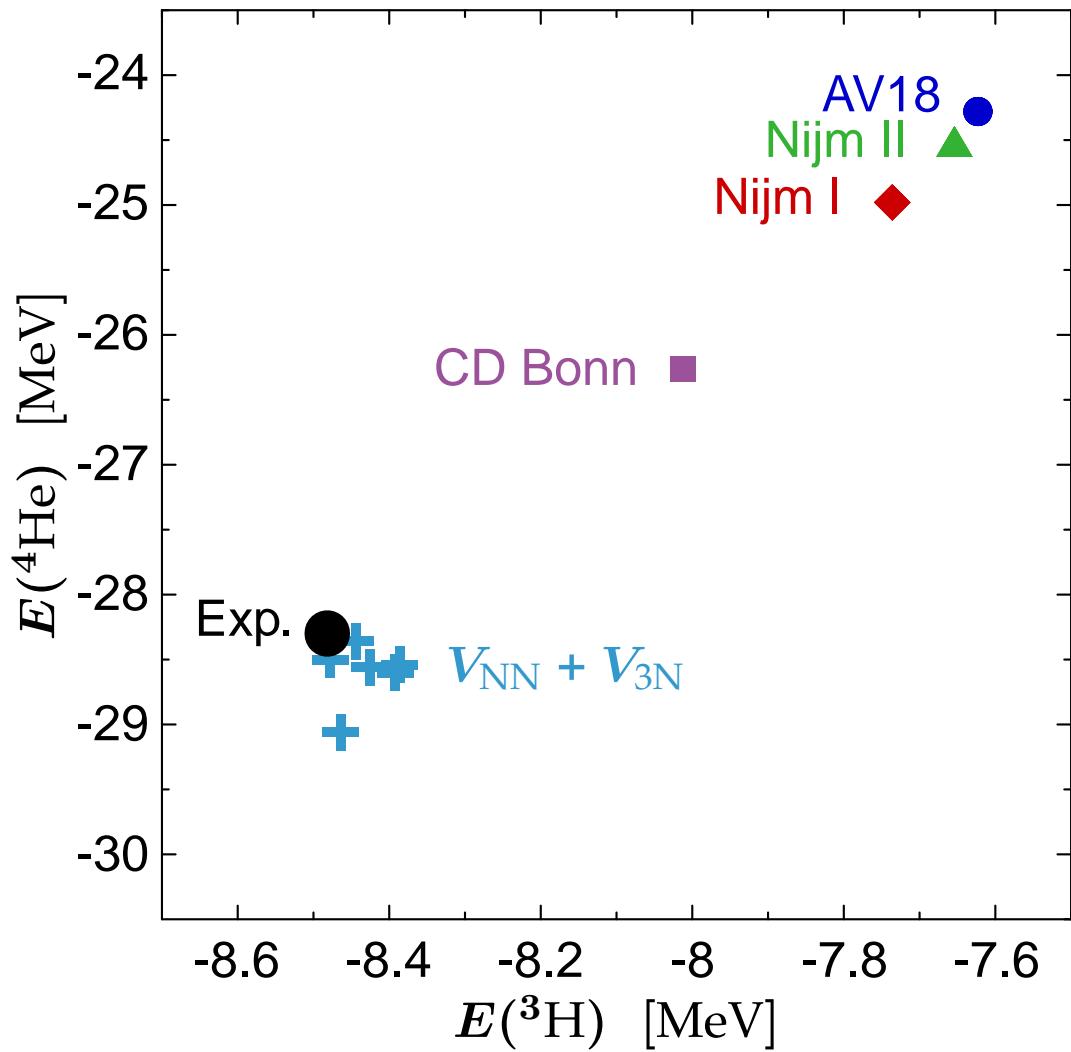
^4He

VUCOM



omitted higher-order
cluster contributions

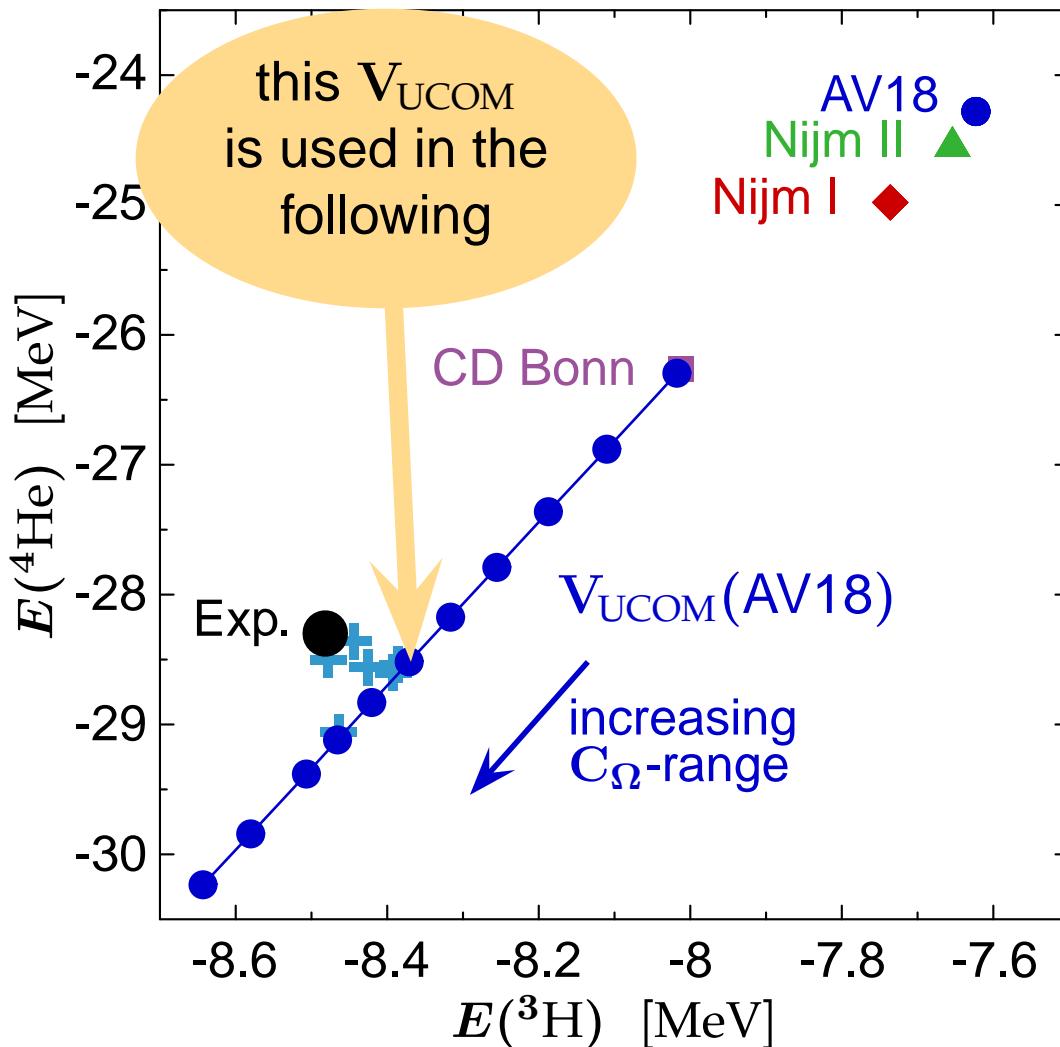
Tjon-Line and Correlator Range



- **Tjon-line:** $E(^4\text{He})$ vs. $E(^3\text{H})$ for phase-shift equivalent NN-interactions

Data points: A. Nogga et al., Phys. Rev. Lett. **85**, 944 (2000)

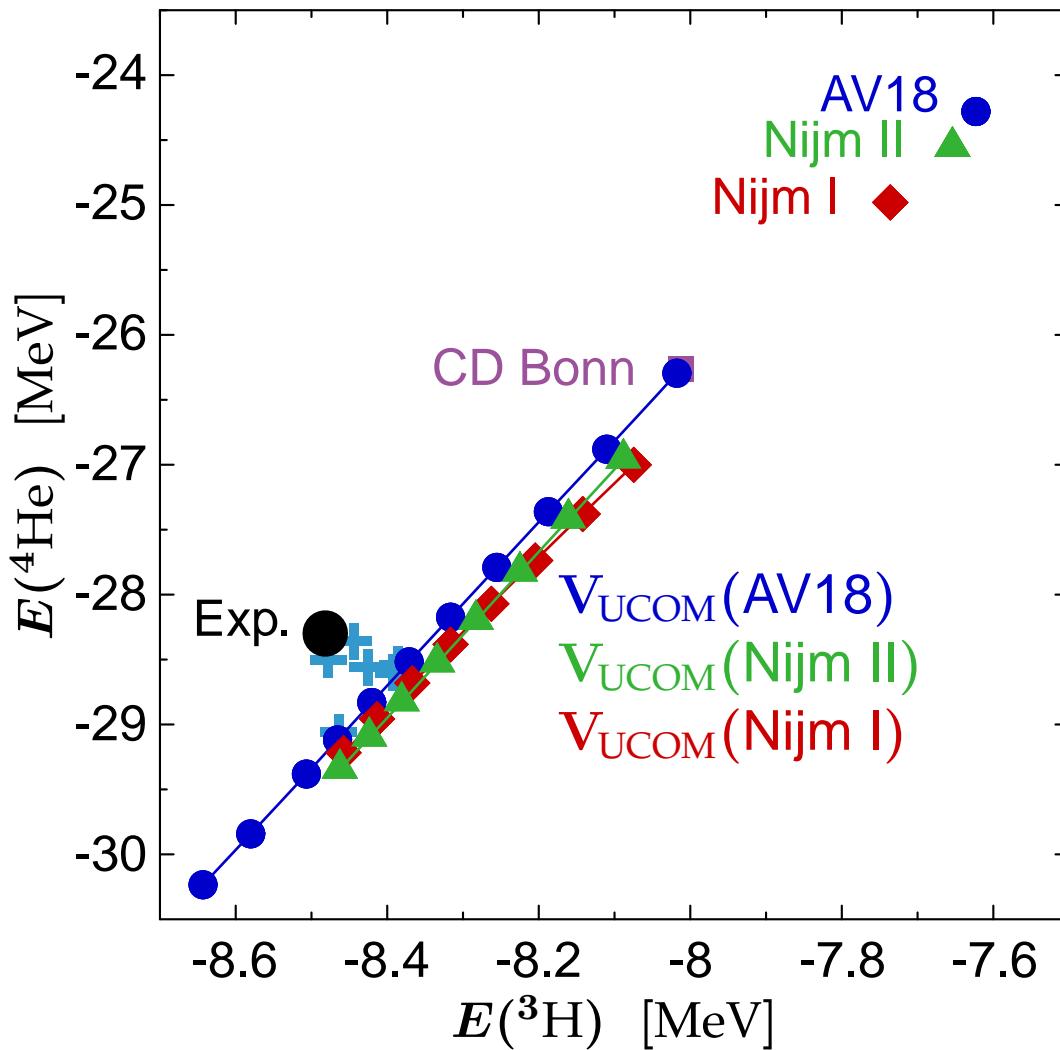
Tjon-Line and Correlator Range



- **Tjon-line:** $E(^4\text{He})$ vs. $E(^3\text{H})$ for phase-shift equivalent NN-interactions
- change of C_Ω -correlator range results in shift along Tjon-line

minimize net three-body force
by choosing correlator with energies close to experimental value

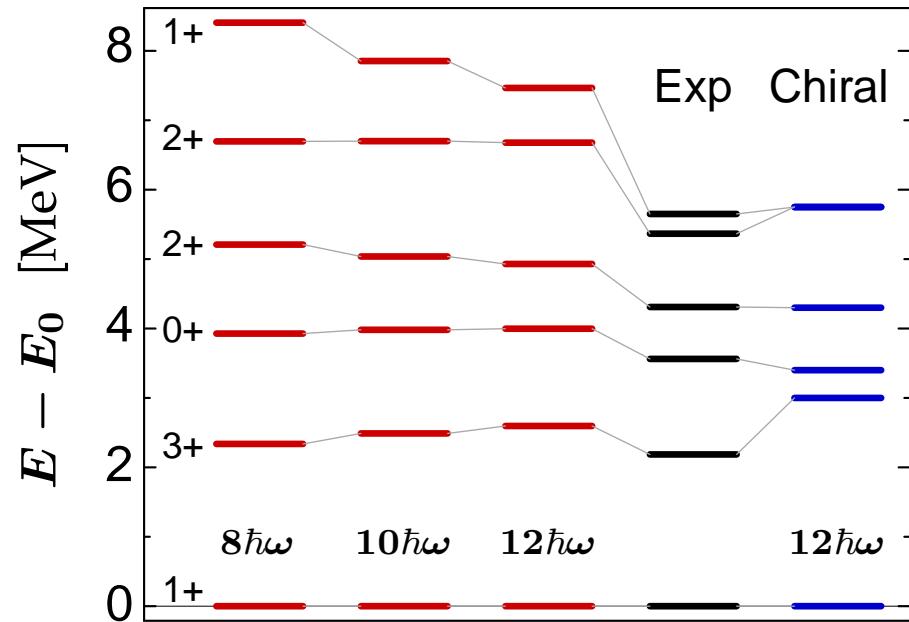
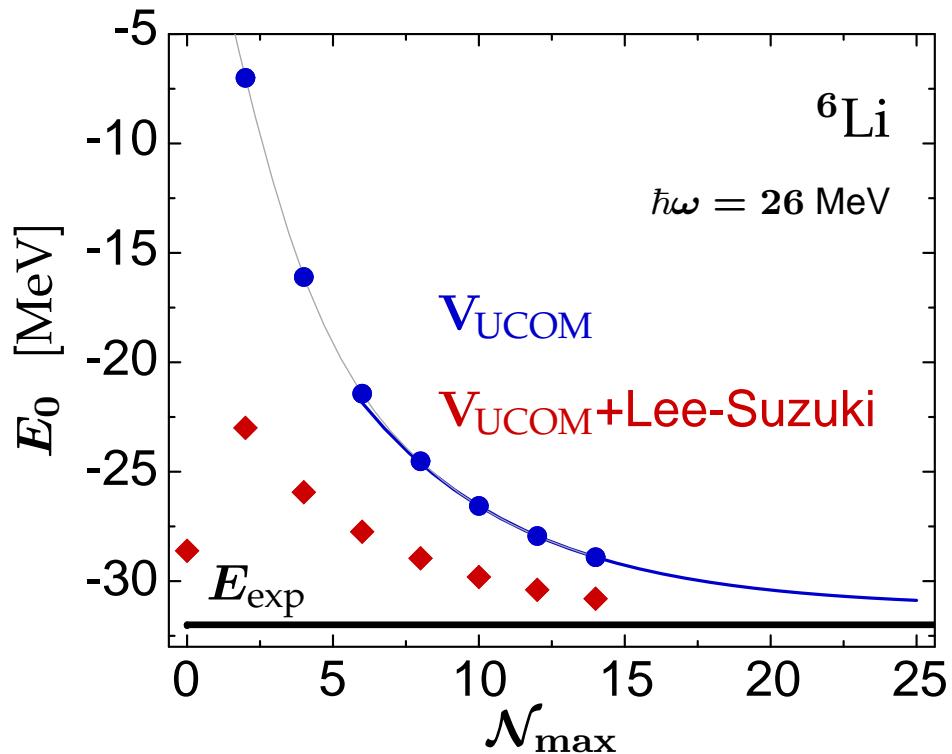
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**minimize net
three-body force**
by choosing correlator
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^6Li : NCSM + UCOM for p -Shell Nuclei

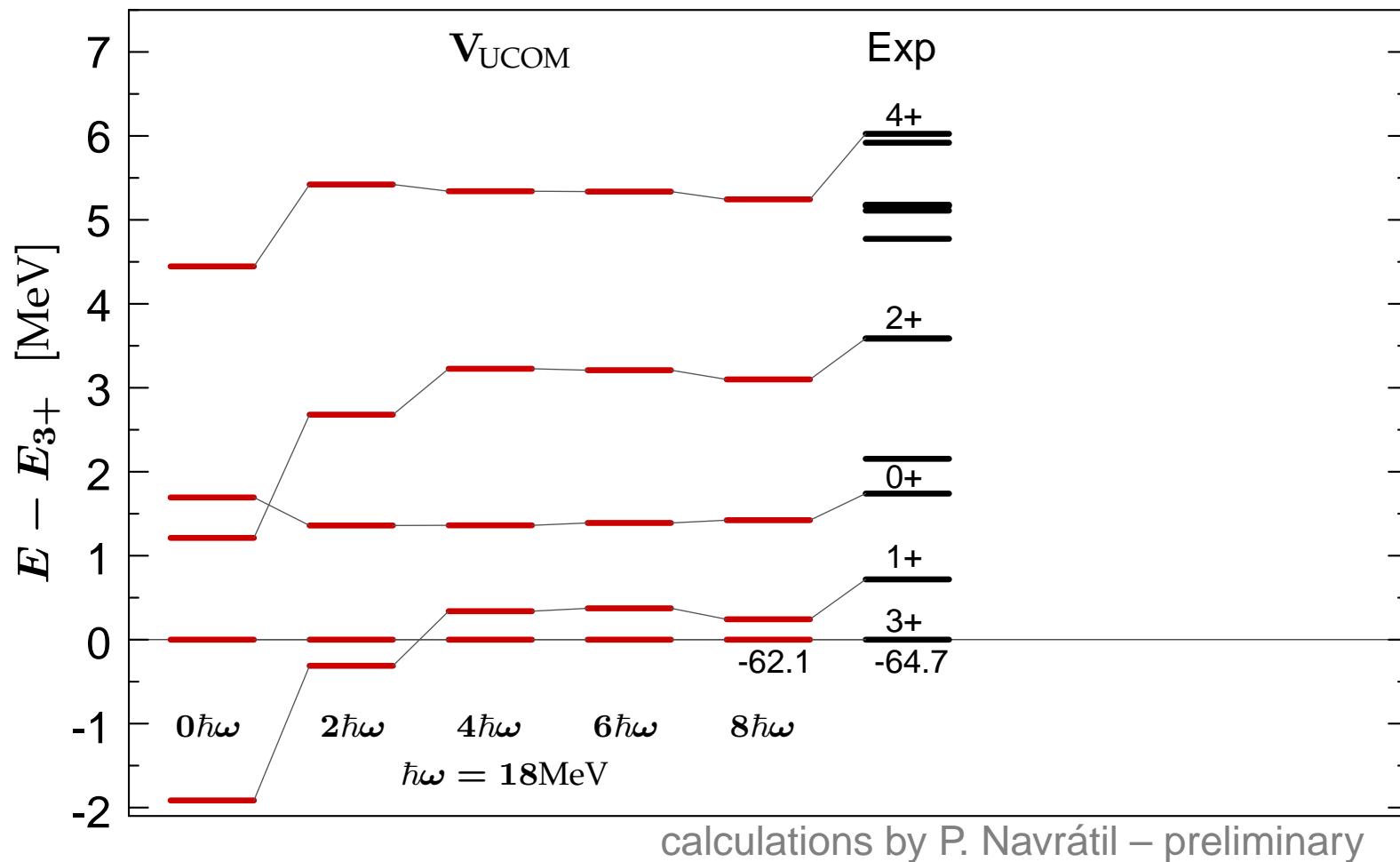


E_0 [MeV] (extrapolation)	-31.23
E_0 [MeV] (experiment)	-32.00

NCSM calculations by P. Navrátil (LLNL)

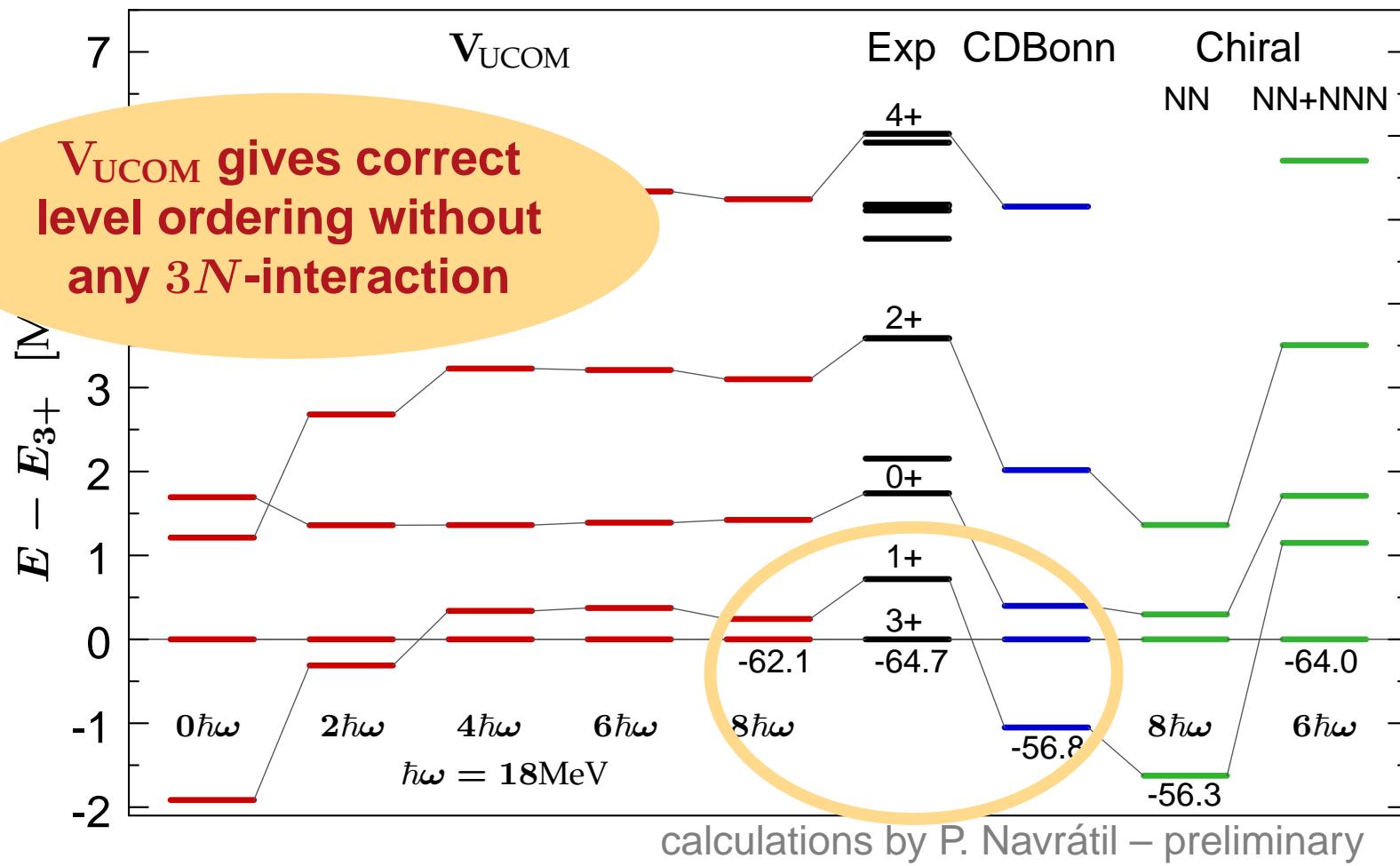
Without
 $3N$ -Interaction!

^{10}B : Benchmarking V_{UCOM}



- stable spectrum from $\sim 4\hbar\omega$
- E_0 in good agreement with experiment

^{10}B : Benchmarking V_{UCOM}



- stable spectrum from $\sim 4\hbar\omega$
- E_0 in good agreement with experiment

Summary & Outlook

Summary

- UCOM treats **short-range central** and **tensor correlations explicitly**
 - ⇒ **improved convergence** through pre-diagonalization
 - ⇒ **minimize net three-body forces**
- **few-body calculations** (NCSM) can be used to **constrain and optimize** V_{UCOM}
- V_{UCOM} is suited for a wide range of many-body methods:
NCSM, FMD, HF, MBPT, RPA, ...
 - ⇒ “**Universal**” phase-shift equivalent NN -interaction!

Outlook

- systematic NCSM study of *p*-shell Nuclei
- construction & inclusion of simple **three-nucleon forces**
(encouraging early results)
- **reactions & resonances** in Fermionic Molecular Dynamics
(UCOM-FMD)
- further & extended applications: **pairing** (HFB), etc.

(In)Famous Last Words...

My Collaborators

- R. Roth, N. Paar, P. Papakonstantinou, A. Zapp

Institut für Kernphysik, TU Darmstadt

- T. Neff

NSCL, Michigan State University

- H. Feldmeier

Gesellschaft für Schwerionenforschung (GSI)

Recent References

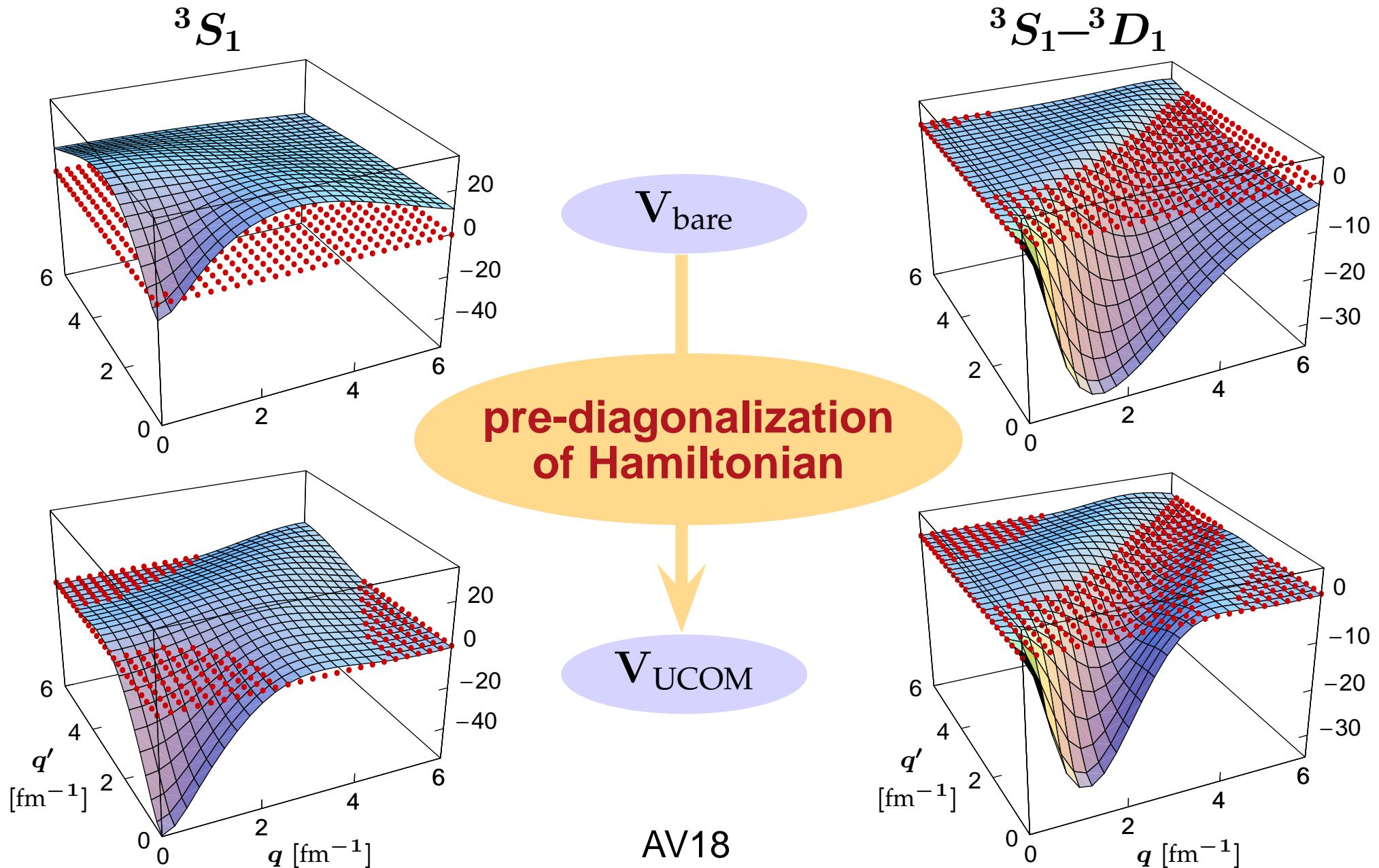
- R. Roth, H. Hergert, P. Papakonstantinou, T. Neff, and H. Feldmeier, Phys. Rev. C**72**, 034002 (2005)

- R. Roth, P. Papakonstantinou, N. Paar, H. Hergert, T. Neff, and H. Feldmeier, nucl-th/0510036, accepted for publication in Phys. Rev. C

- <http://crunch.ikp.physik.tu-darmstadt.de/tnp/>

Optional / Appendices

Momentum-Space Matrix Elements



UCOM-FMD Approach

Gaussian Single-Particle States

$$|q\rangle = \sum_{\nu=1}^n \mathbf{c}_{\nu} \ |a_{\nu}, \vec{b}_{\nu}\rangle \otimes |\chi_{\nu}\rangle \otimes |m_t\rangle$$

$$\langle \vec{x} | a_{\nu}, \vec{b}_{\nu} \rangle = \exp \left[- \frac{(\vec{x} - \vec{b}_{\nu})^2}{2 a_{\nu}} \right]$$

a_{ν} : complex width

χ_{ν} : spin orientation

\vec{b}_{ν} : mean position & momentum

Slater Determinant

$$|Q\rangle = \mathcal{A} (|q_1\rangle \otimes |q_2\rangle \otimes \cdots \otimes |q_A\rangle)$$

Correlated Hamiltonian

$$\tilde{\mathbf{H}} = \mathbf{T} + \mathbf{V}_{\text{UCOM}} + \delta V_{c+p+ls}$$

Variation

$$\frac{\langle Q | \tilde{\mathbf{H}} - \mathbf{T}_{\text{cm}} | Q \rangle}{\langle Q | Q \rangle} \rightarrow \min$$

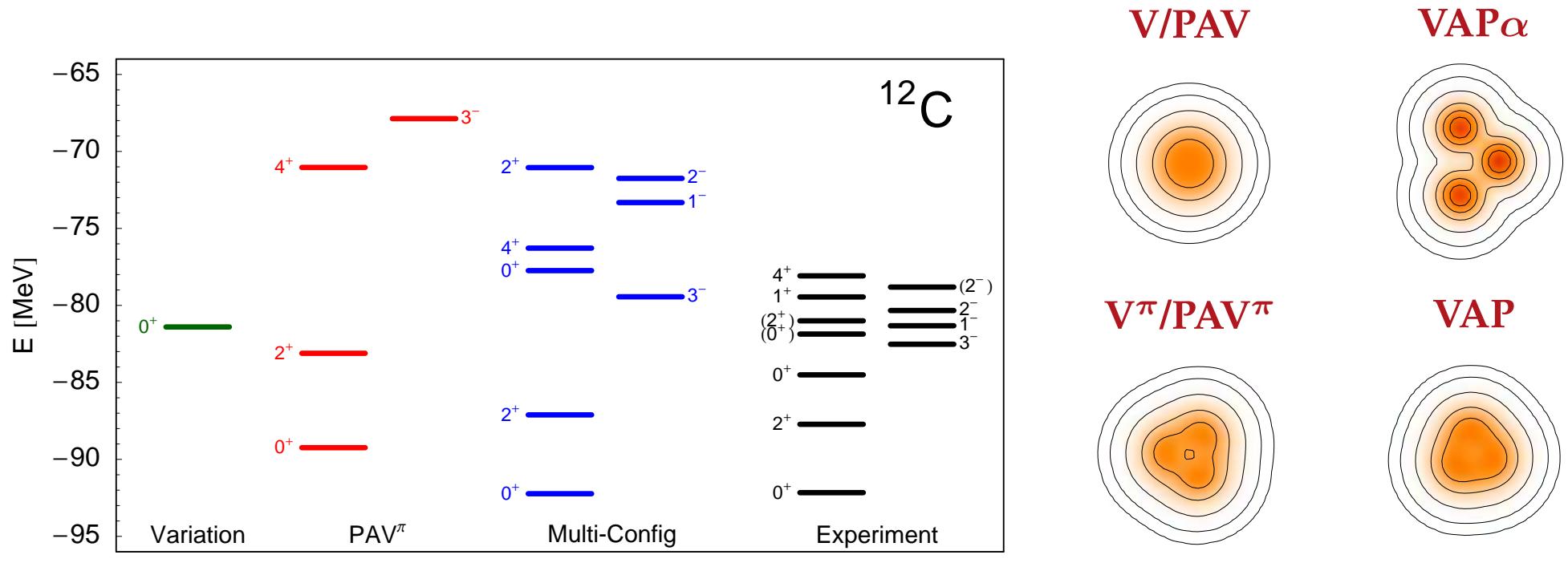
Projection

restoration of parity and
rotational symmetry
PAV / VAP

Multi- Configuration

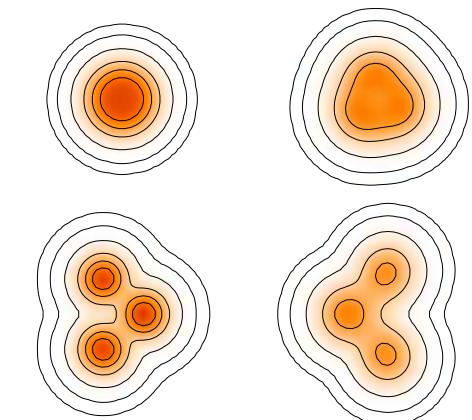
mixing of several
intrinsic configurations
GCM

Structure of ^{12}C

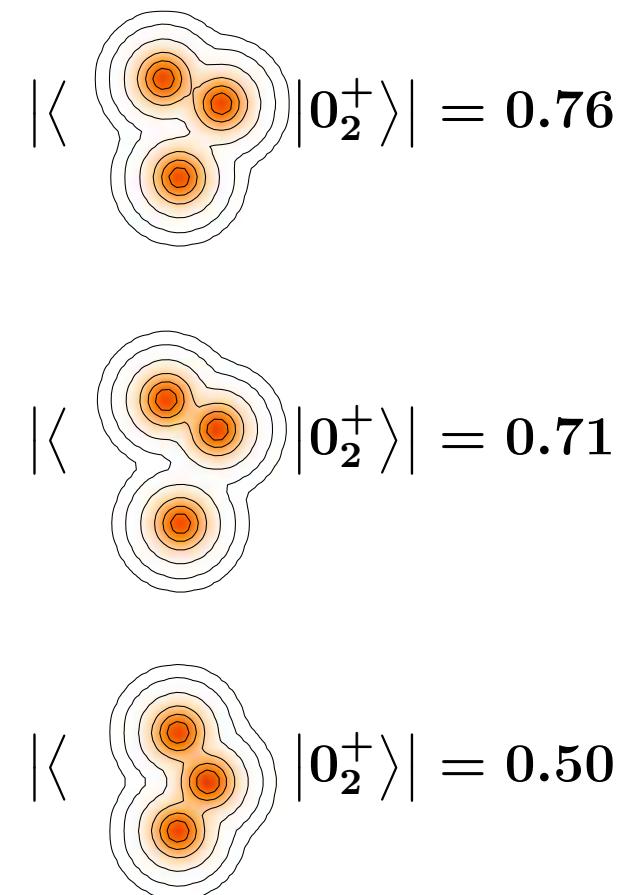
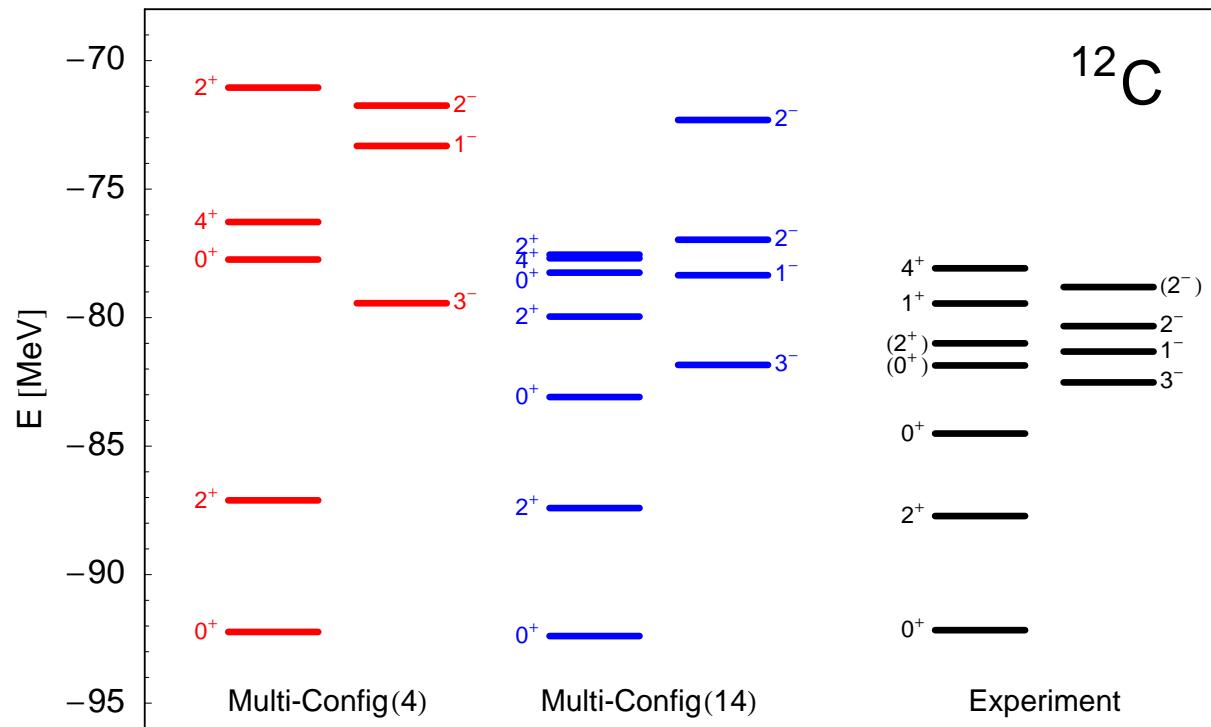


	E [MeV]	R_{ch} [fm]	$B(E2)$ [$e^2 \text{ fm}^4$]
V/PAV	81.4	2.36	-
VAP α -cluster	79.1	2.70	76.9
PAV $^\pi$	88.5	2.51	36.3
VAP	89.2	2.42	26.8
Multi-Config	92.2	2.52	42.8
Experiment	92.2	2.47	39.7 ± 3.3

Multi-Config



Structure of ^{12}C — Hoyle State



	Multi-Config	Experiment
E [MeV]	92.4	92.2
R_{ch} [fm]	2.52	2.47
$B(E2, 0_1^+ \rightarrow 2_1^+) [e^2 \text{ fm}^4]$	42.9	39.7 ± 3.3
$M(E0, 0_1^+ \rightarrow 0_2^+) [\text{fm}^2]$	5.67	5.5 ± 0.2

More References

- H. Feldmeier, T. Neff, R. Roth, and J. Schnack, Nucl. Phys. **A632**, 61 (1998)
- T. Neff, and H. Feldmeier, Nucl. Phys. **A713**, 311 (2003)
- R. Roth, T. Neff, H. Hergert, and H. Feldmeier, Nucl. Phys. **A745**, 3 (2004)
- R. Roth, H. Hergert, P. Papakonstantinou, T. Neff, and H. Feldmeier, Phys. Rev. **C72**, 034002 (2005)
- R. Roth, P. Papakonstantinou, N.Paar, H. Hergert, T. Neff, and H. Feldmeier, nucl-th/0510036, accepted for publication in Phys. Rev. C
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