



New Insights into Nuclear Clustering

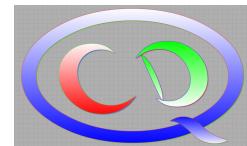
Ulf-G. Meißner, Univ. Bonn & FZ Jülich

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by DFG, SFB/TR-110

by CAS, PIFI

by Volkswagen Stiftung



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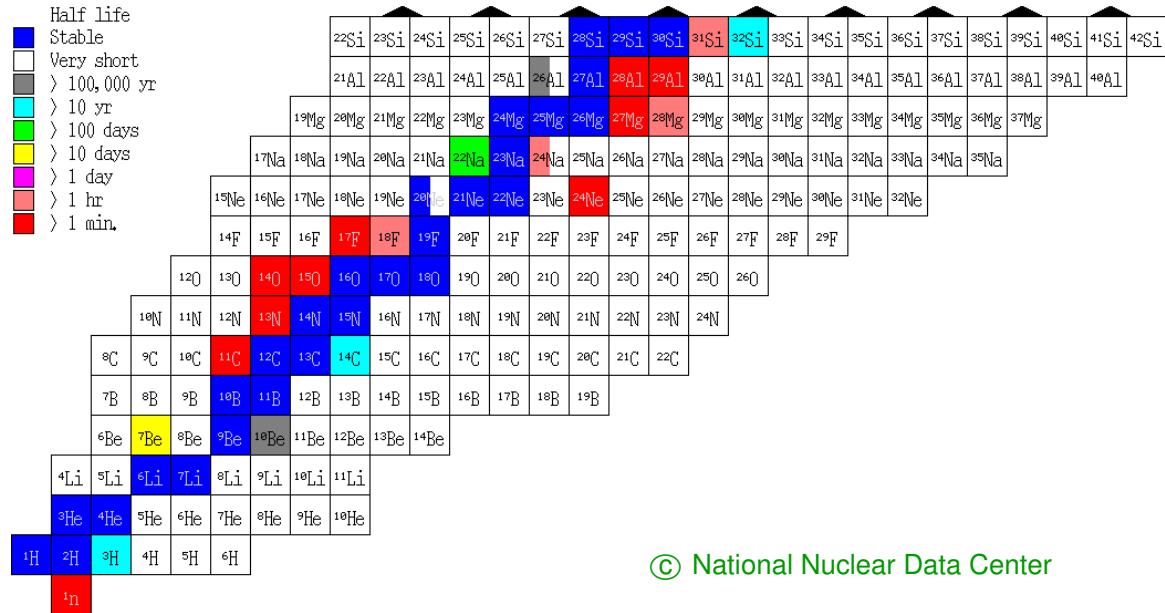
- Very brief introduction
- Basics of nuclear lattice simulations
- Results from nuclear lattice simulations
- New insights into nuclear clustering
- Summary & outlook

Very brief introduction

AB INITIO NUCLEAR STRUCTURE and SCATTERING

- Nuclear structure:

- ★ limits of stability
- ★ 3-nucleon forces
- ★ alpha-clustering
- ⋮
- this talk



- Nuclear scattering: processes relevant for nuclear astrophysics

- ★ alpha-particle scattering: $^4\text{He} + ^4\text{He} \rightarrow ^4\text{He} + ^4\text{He}$ → Dean Lee's talk
- ★ triple-alpha reaction: $^4\text{He} + ^4\text{He} + ^4\text{He} \rightarrow ^{12}\text{C} + \gamma$
- ★ alpha-capture on carbon: $^4\text{He} + ^{12}\text{C} \rightarrow ^{16}\text{O} + \gamma$
- ⋮

Basics of nuclear lattice simulations

for an easy intro, see: UGM, Nucl. Phys. News **24** (2014) 11

for an early review, see: D. Lee, Prog. Part. Nucl. Phys. **63** (2009) 117

NUCLEAR LATTICE EFFECTIVE FIELD THEORY

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Frank, Brockmann (1992), Koonin, Müller, Seki, van Kolck (2000) , Lee, Schäfer (2004), . . .
Borasoy, Krebs, Lee, UGM, Nucl. Phys. **A768** (2006) 179; Borasoy, Epelbaum, Krebs, Lee, UGM, Eur. Phys. J. **A31** (2007) 105

- new method to tackle the nuclear many-body problem

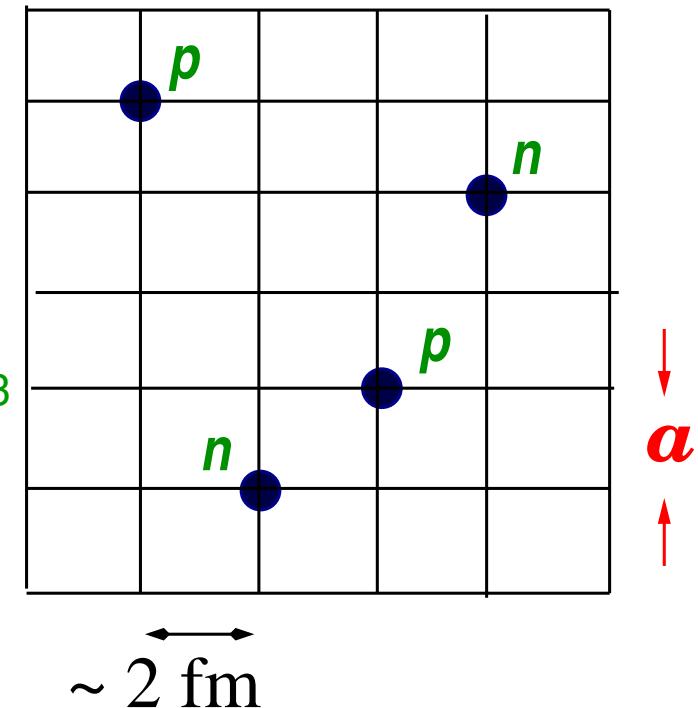
- discretize space-time $V = L_s \times L_s \times L_s \times L_t$:
nucleons are point-like particles on the sites

- discretized chiral potential w/ pion exchanges
and contact interactions + Coulomb

→ see Epelbaum, Hammer, UGM, Rev. Mod. Phys. **81** (2009) 1773

- typical lattice parameters

$$p_{\max} = \frac{\pi}{a} \simeq 314 \text{ MeV [UV cutoff]}$$



- strong suppression of sign oscillations due to approximate Wigner SU(4) symmetry

E. Wigner, Phys. Rev. **51** (1937) 106; T. Mehen et al., Phys. Rev. Lett. **83** (1999) 931; J. W. Chen et al., Phys. Rev. Lett. **93** (2004) 242302

- physics independent of the lattice spacing for $a = 1 \dots 2 \text{ fm}$

J. Alarcon et al., EPJA **53** (2017) 83

TRANSFER MATRIX METHOD

- Correlation–function for A nucleons: $Z_A(\tau) = \langle \Psi_A | \exp(-\tau H) | \Psi_A \rangle$

with Ψ_A a Slater determinant for A free nucleons
[or a more sophisticated (correlated) initial/final state]

- Transient energy

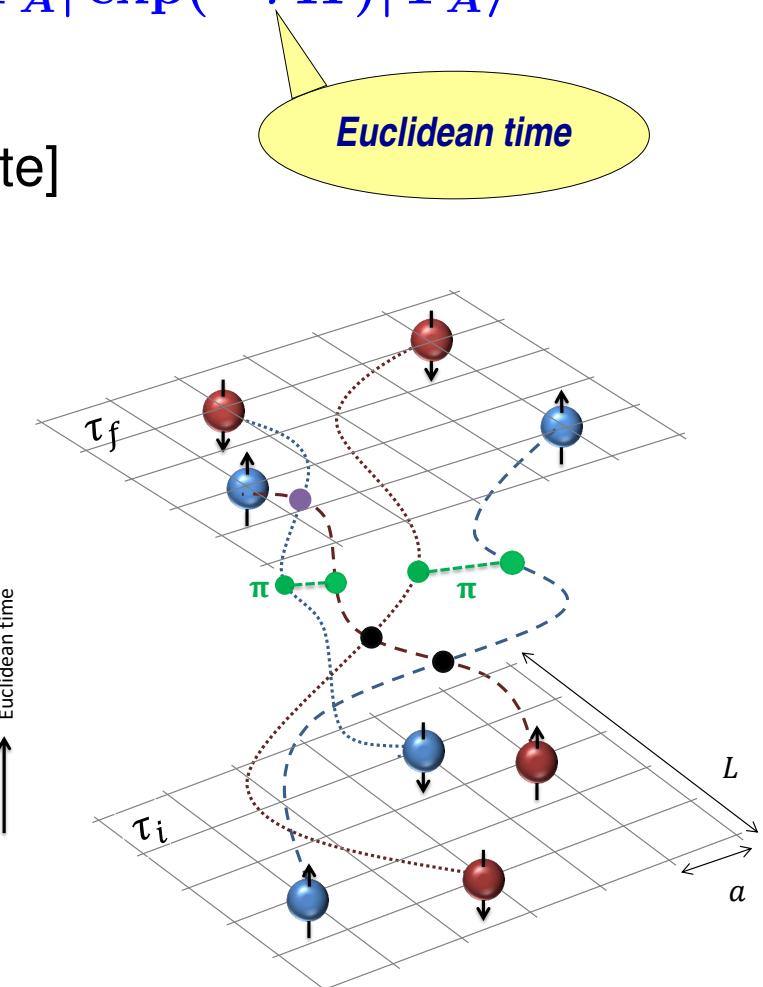
$$E_A(\tau) = -\frac{d}{d\tau} \ln Z_A(\tau)$$

→ ground state: $E_A^0 = \lim_{\tau \rightarrow \infty} E_A(\tau)$

- Exp. value of any normal–ordered operator \mathcal{O}

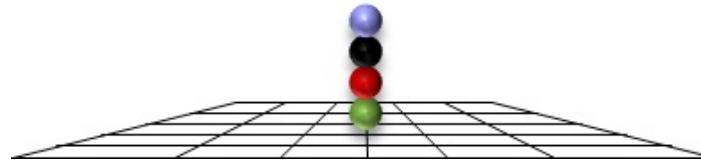
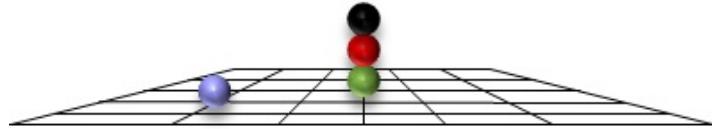
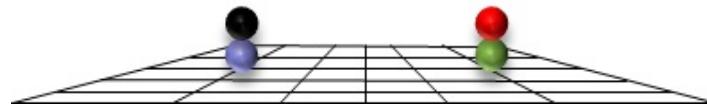
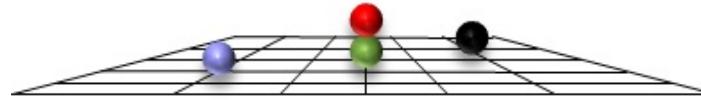
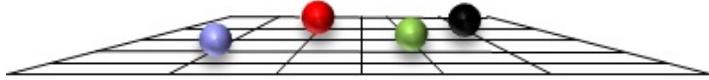
$$Z_A^\mathcal{O} = \langle \Psi_A | \exp(-\tau H/2) \mathcal{O} \exp(-\tau H/2) | \Psi_A \rangle$$

$$\lim_{\tau \rightarrow \infty} \frac{Z_A^\mathcal{O}(\tau)}{Z_A(\tau)} = \langle \Psi_A | \mathcal{O} | \Psi_A \rangle$$



CONFIGURATIONS

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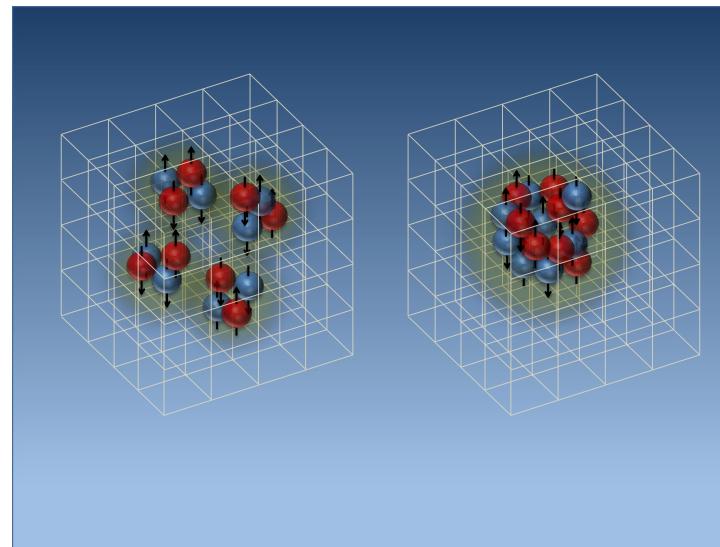
- ⇒ all *possible* configurations are sampled
- ⇒ preparation of *all possible* initial/final states
- ⇒ *clustering* emerges *naturally*

COMPUTATIONAL EQUIPMENT

- Present = JUQUEEN (BlueGene/Q)



Lattice: some results



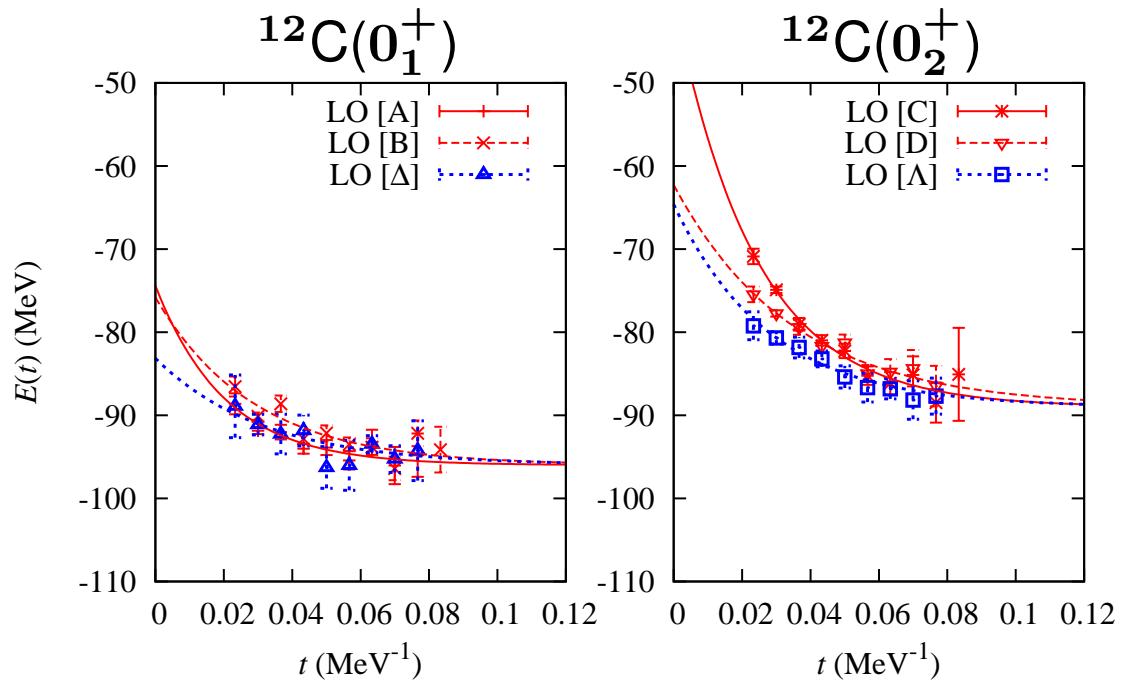
Epelbaum, Krebs, Lähde, Lee, Luu, UGM, Rupak + post-docs + students

FIXING PARAMETERS and FIRST RESULTS

Epelbaum, Krebs, Lee, UGM, Phys. Rev. Lett. **104** (2010) 142501; Eur. Phys. J. A **45** (2010) 335; ...

- some groundstate energies and differences [NNLO, 11+2 LECs]

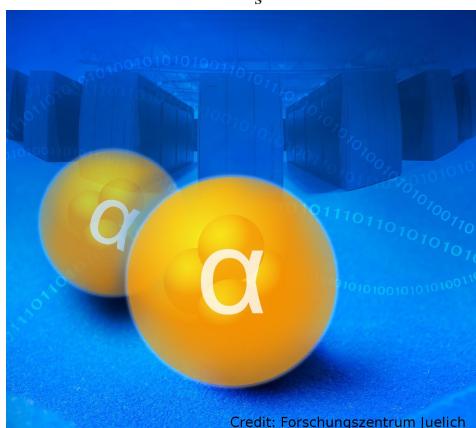
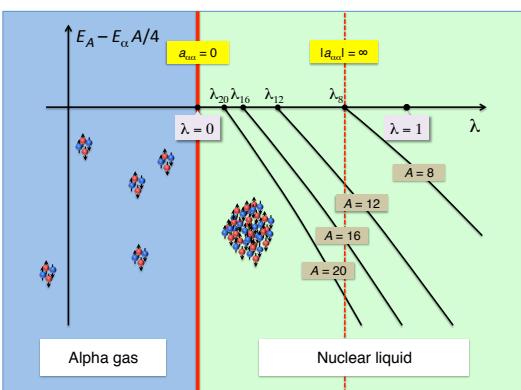
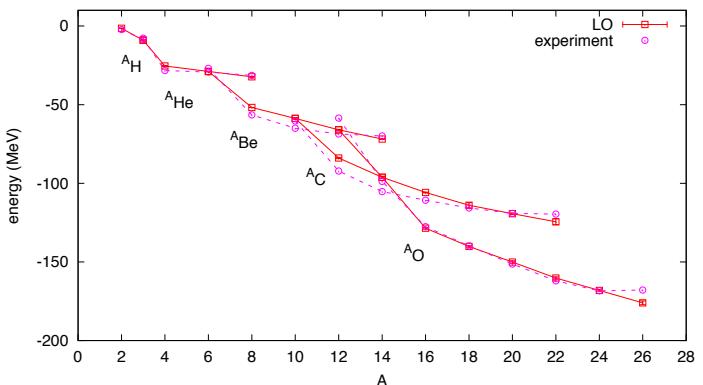
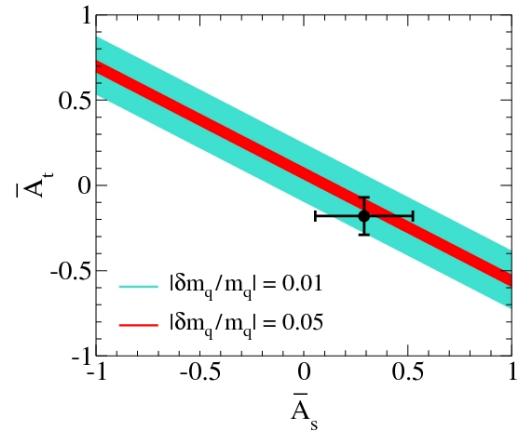
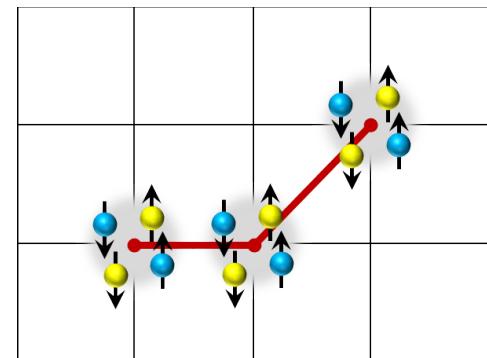
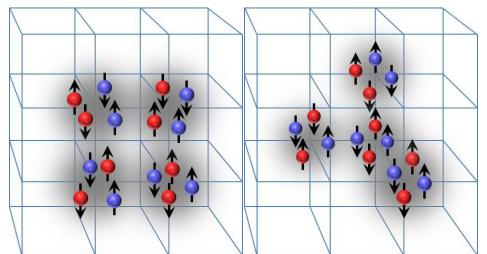
	E [MeV]	NLEFT	Exp.
old algorithm	^3He - ^3H	0.78(5)	0.76
	^4He	-28.3(6)	-28.3
	^8Be	-55(2)	-56.5
	^{12}C	-92(3)	-92.2
new algorithm	^{16}O	-131(1)	-127.6
	^{20}Ne	-166(1)	-160.6
	^{24}Mg	-198(2)	-198.3
	^{28}Si	-234(3)	-236.5



- promising results \Rightarrow uncertainties down to the 1% level
- excited states more difficult \Rightarrow projection MC method + triangulation

RESULTS from LATTICE NUCLEAR EFT

- Lattice EFT calculations for $A=3,4,6,12$ nuclei, PRL 104 (2010) 142501
- *Ab initio* calculation of the Hoyle state, PRL 106 (2011) 192501
- Structure and rotations of the Hoyle state, PRL 109 (2012) 142501
- Validity of Carbon-Based Life as a Function of the Light Quark Mass
PRL 110 (2013) 142501
- *Ab initio* calculation of the Spectrum and Structure of ^{16}O ,
PRL 112 (2014) 142501
- *Ab initio* alpha-alpha scattering, Nature 528 (2015) 111 → Dean Lee's talk
- Nuclear Binding Near a Quantum Phase Transition, PRL 117 (2016) 132501
- *Ab initio* calculations of the isotopic dependence of nuclear clustering,
arXiv:1702.05177 → this talk



Ab initio calculations of the isotopic dependence of nuclear clustering

Elhatisari, Epelbaum, Krebs, Lähde, Lee, Li, Lu, UGM, Rupak
[arXiv:1702.05117]

EARLIER RESULTS on NUCLEAR CLUSTERING

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- Already a number of intriguing results on clustering:

Ab initio calculation of the spectrum and structure of ^{12}C (esp. the Hoyle state)

Ab initio calculation of the spectrum and structure of ^{16}O

Ground state energies of α -type nuclei up to ^{28}Si within 1%

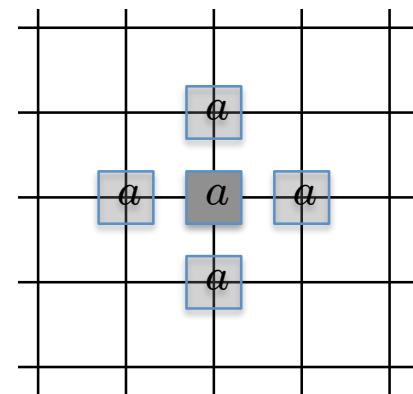
Ab initio calculation of α - α scattering

Quantum phase transition from Bose gas of α 's to nuclear liquid for α -type nuclei

- However: when adding extra neutrons/protons, the precision quickly deteriorates due to sign oscillations
- New LO action with smeared SU(4) local+non-local symmetric contact interactions & smeared one-pion exchange

$$a_{\text{NL}}(\mathbf{n}) = a(\mathbf{n}) + s_{\text{NL}} \sum_{\langle \mathbf{n}' | \mathbf{n} \rangle} a(\mathbf{n}')$$

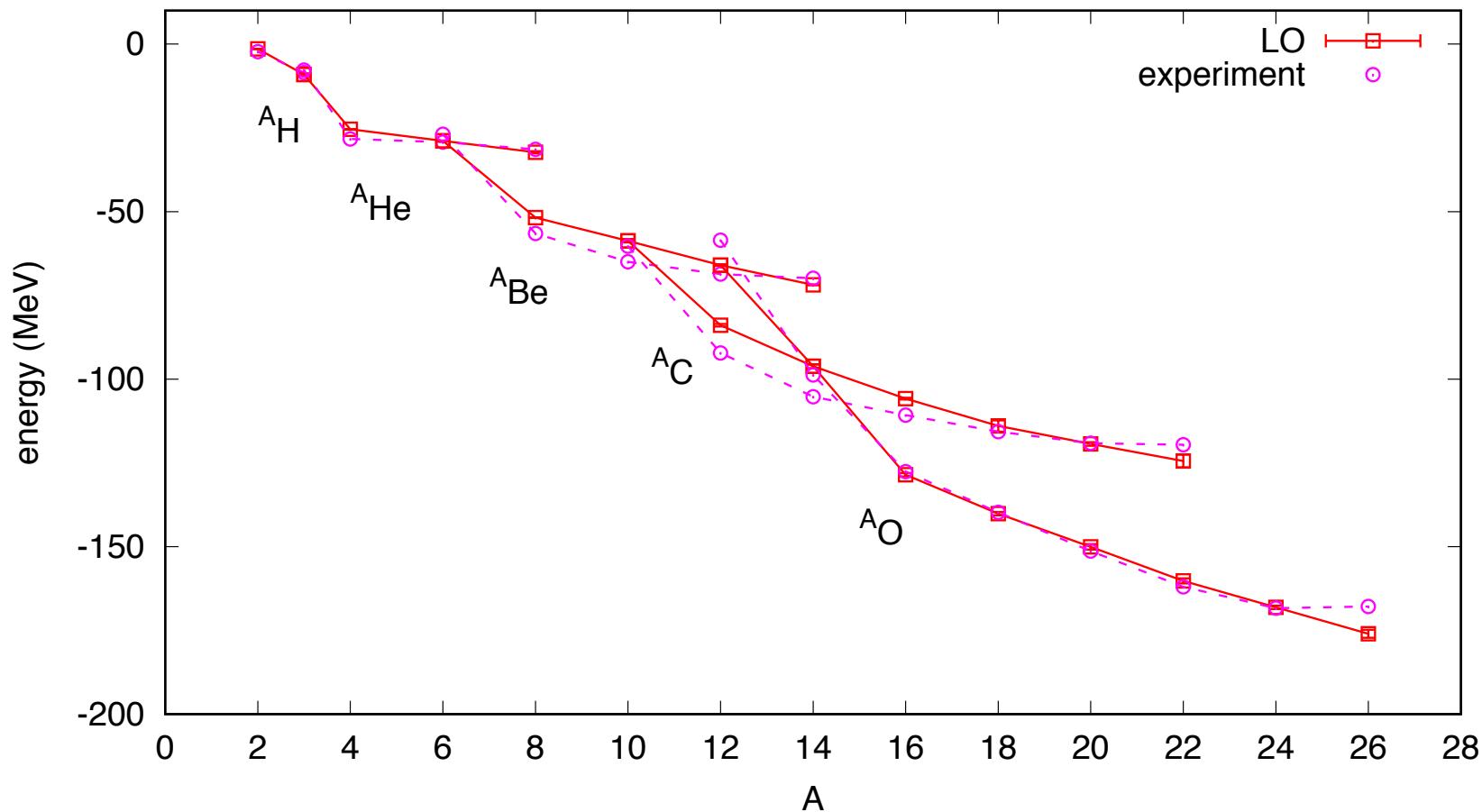
$$a_{\text{NL}}^\dagger(\mathbf{n}) = a^\dagger(\mathbf{n}) + s_{\text{NL}} \sum_{\langle \mathbf{n}' | \mathbf{n} \rangle} a^\dagger(\mathbf{n}')$$



GROUND STATE ENERGIES

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- Fit parameters to average NN S-wave scattering length and effective range and α - α S-wave scattering length
→ predict g.s. energies of H, He, Be, C and O isotopes → quite accurate (LO)



PROBING NUCLEAR CLUSTERING

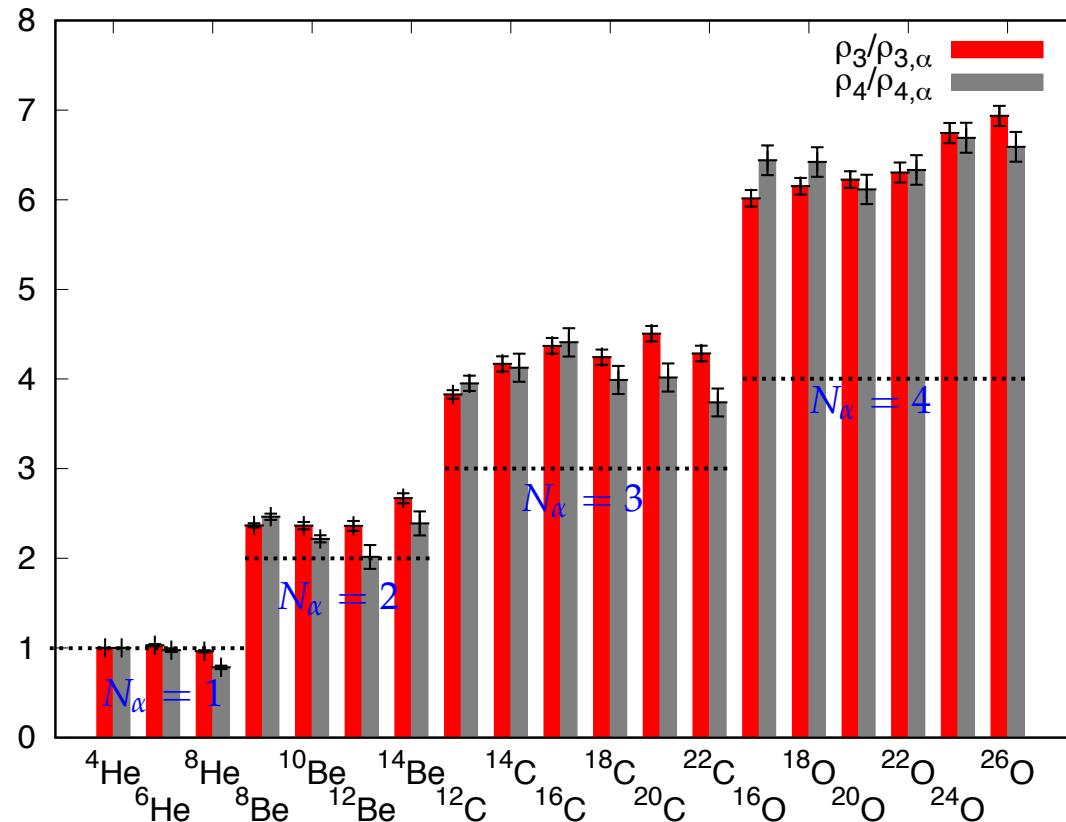
- Local densities on the lattice: $\rho(\mathbf{n})$, $\rho_p(\mathbf{n})$, $\rho_n(\mathbf{n})$
- Probe of alpha clusters: $\rho_4 = \sum_{\mathbf{n}} : \rho^4(\mathbf{n}) / 4! :$
- Another probe for $Z = N = \text{even}$ nuclei: $\rho_3 = \sum_{\mathbf{n}} : \rho^3(\mathbf{n}) / 3! :$
- ρ_4 couples to the center of the α -cluster while ρ_3 gets contributions from a wider portion of the alpha-particle wave function
- Both ρ_3 and ρ_4 depend on the regulator, a , but not on the nucleus
- The ratios $\rho_3/\rho_{3,\alpha}$ and $\rho_4/\rho_{4,\alpha}$ free of short-distance ambiguities and model-independent
- $\rho_3/\rho_{3,\alpha}$ measures the effective number of alpha-cluster N_α
 \Rightarrow Any deviation from $N_\alpha = \text{integer}$ measures the entanglement of the α -clusters in a given nucleus

PROBING NUCLEAR CLUSTERING

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- ρ_3 -entanglement of the α -clusters:

$$\frac{\Delta \rho_3}{N_\alpha} = \frac{\rho_3 / \rho_{3,\alpha}}{N_\alpha} - 1$$

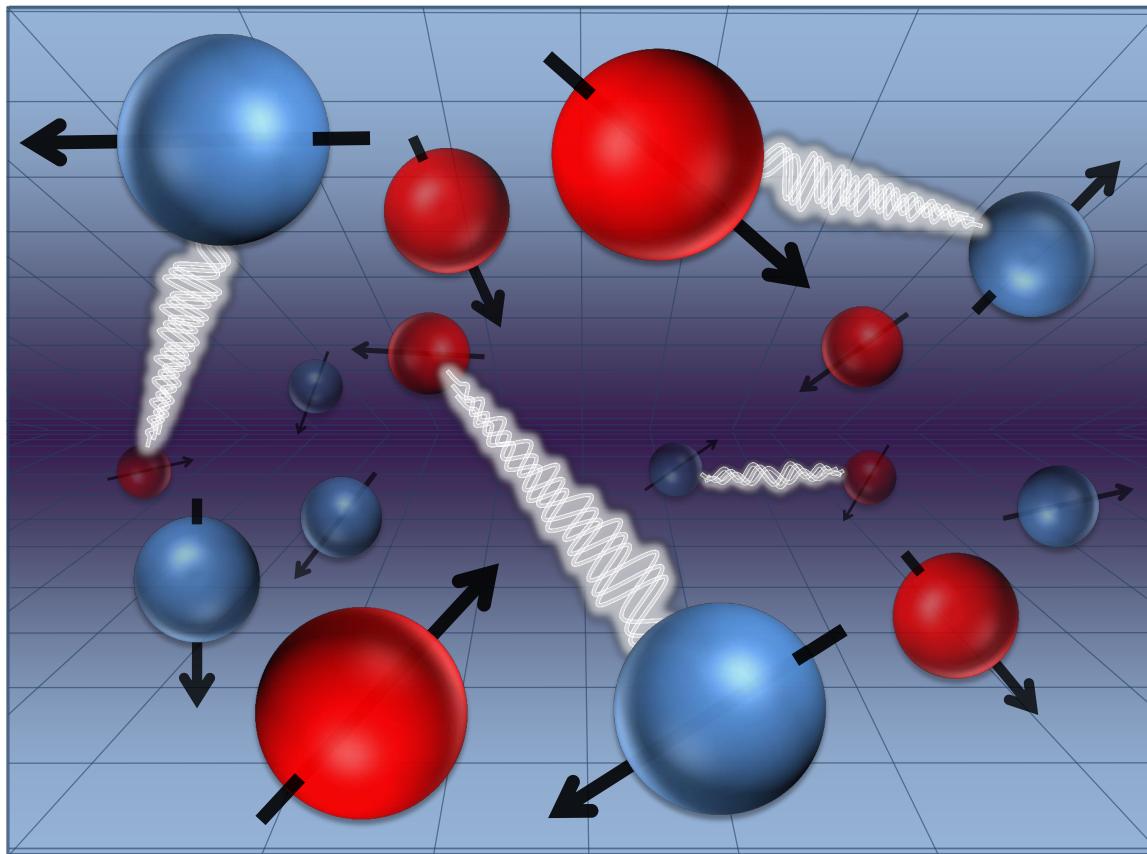


Nucleus	${}^4, {}^6, {}^8\text{He}$	${}^8, {}^{10}, {}^{12}, {}^{14}\text{Be}$	${}^{12}, {}^{14}, {}^{16}, {}^{18}, {}^{20}, {}^{22}\text{C}$	${}^{16}, {}^{18}, {}^{20}, {}^{22}, {}^{24}, {}^{26}\text{O}$
$\Delta \rho_3 / N_\alpha$	0.00 - 0.03	0.20 - 0.35	0.25 - 0.50	0.50 - 0.75

PROBING NUCLEAR CLUSTERING

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- The transition from cluster-like states in light systems to nuclear liquid-like states in heavier systems should not be viewed as a simple suppression of multi-nucleon short-distance correlations, but rather as an increasing *entanglement* of the nucleons involved in the multi-nucleon correlations.



PINHOLE ALGORITHM

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- AFQMC calculations involve states that are superpositions of many different center-of-mass positions
→ density distributions of nucleons can not be computed directly

- Insert a screen with pinholes with spin & isospin labels that allows nucleons with corresponding spin & isospin to pass = insertion of the A-body density op.:

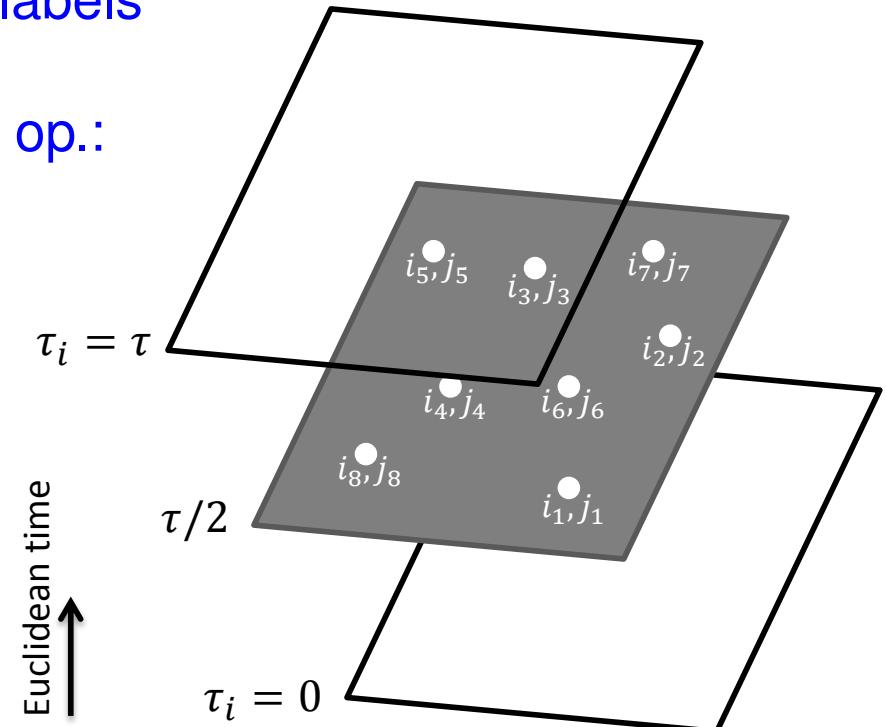
$$\rho_{i_1, j_1, \dots, i_A, j_A}(\mathbf{n}_1, \dots, \mathbf{n}_A) \\ = : \rho_{i_1, j_1}(\mathbf{n}_1) \cdots \rho_{i_A, j_A}(\mathbf{n}_A) :$$

- MC sampling of the amplitude:

$$A_{i_1, j_1, \dots, i_A, j_A}(\mathbf{n}_1, \dots, \mathbf{n}_A, L_t) \\ = \langle \psi(\tau/2) | \rho_{i_1, j_1, \dots, i_A, j_A}(\mathbf{n}_1, \dots, \mathbf{n}_A) | \psi(\tau/2) \rangle$$

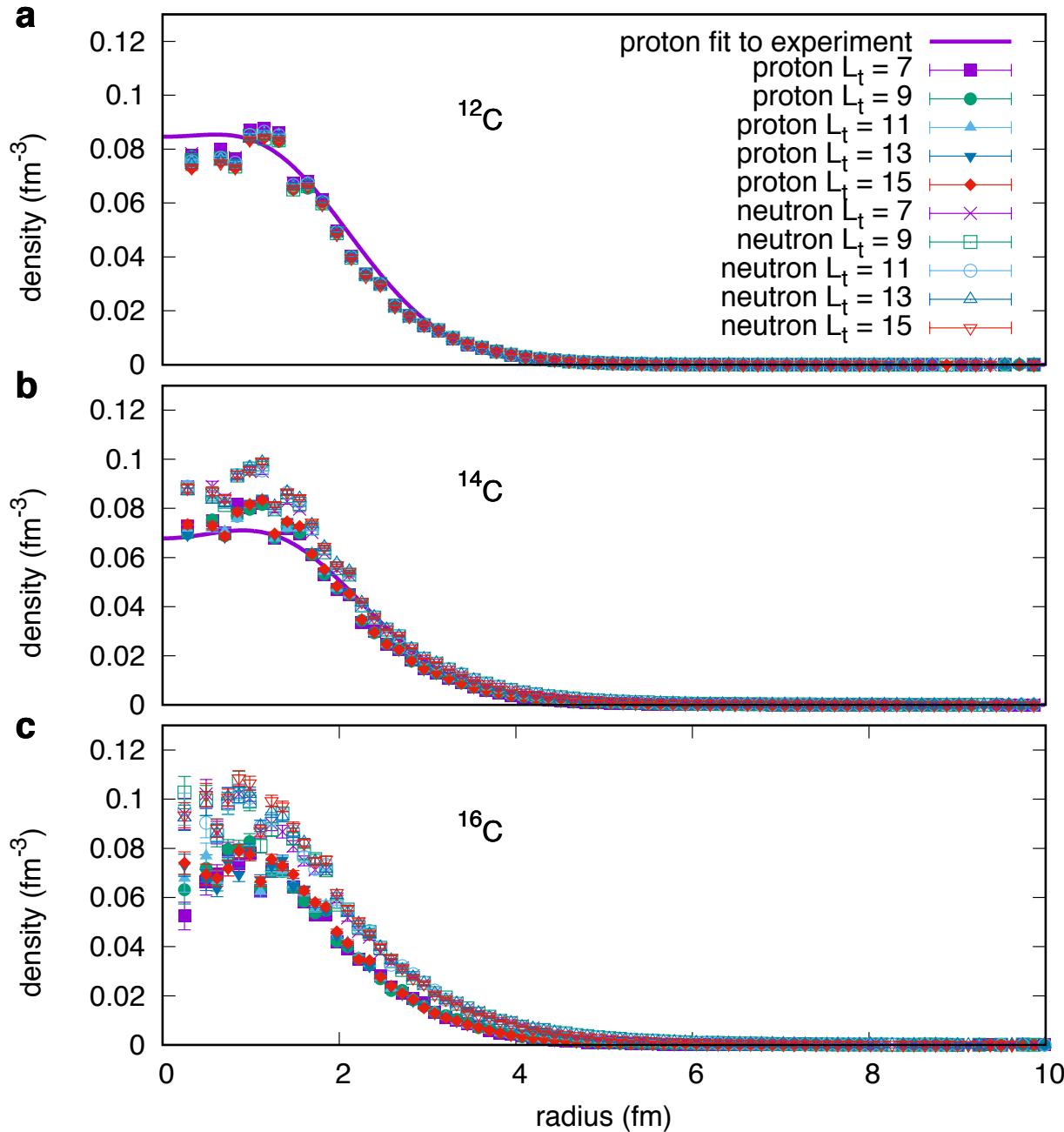
- Allows to measure proton and neutron distributions

- Resolution scale $\sim a/A$ as cm position \mathbf{r}_{cm} is an integer n_{cm} times a/A



PROTON and NEUTRON DENSITIES in CARBON

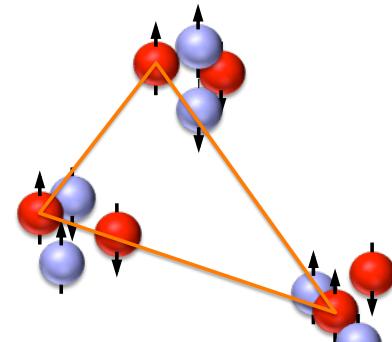
20



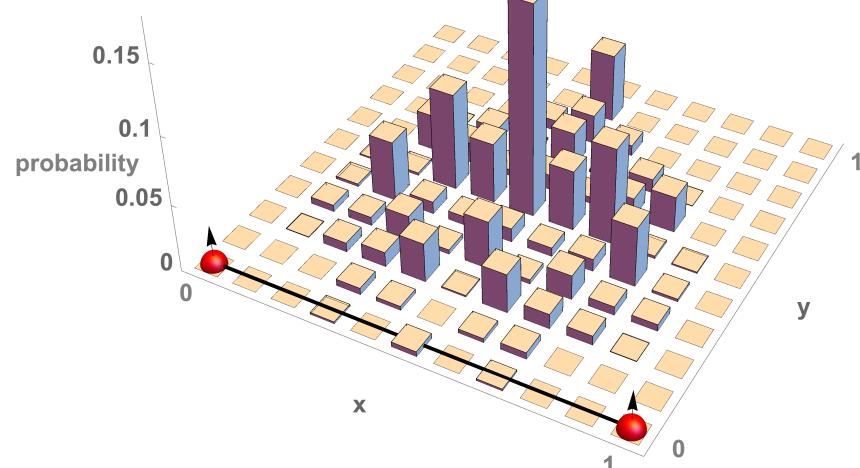
- open symbols: neutron
- closed symbols: proton
- proton size accounted for
- asymptotic properties of the distributions from the volume dependence of N-body bound states
König, Lee, [arXiv:1701.00279]
- consistent with data
- fit to data from
Kline et al., NPA209 (1973) 381

ALPHA CLUSTER GEOMETRY

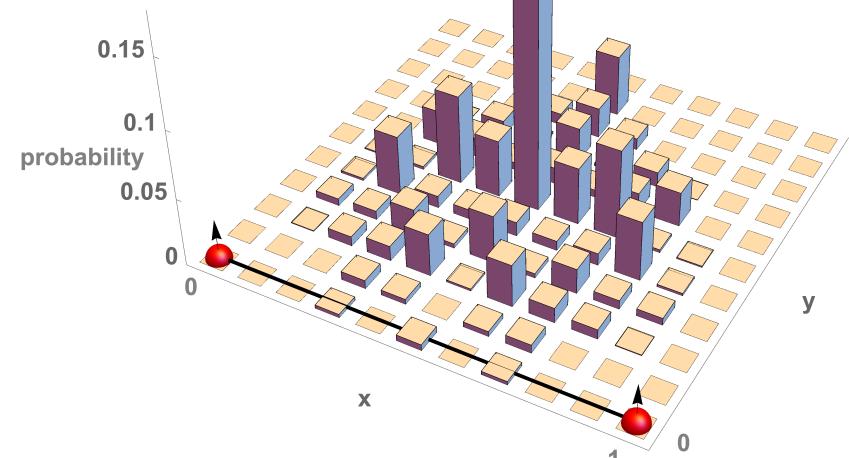
- Measuring the three spin-up protons by considering triangular shapes



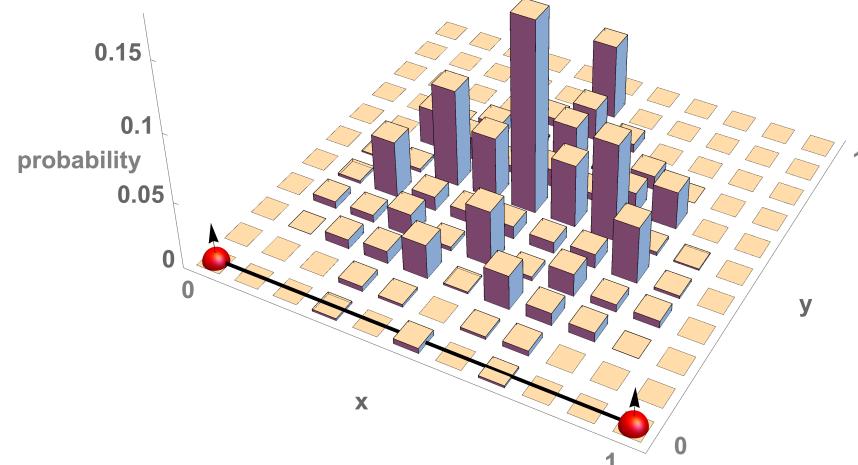
^{14}C



^{12}C



^{16}C



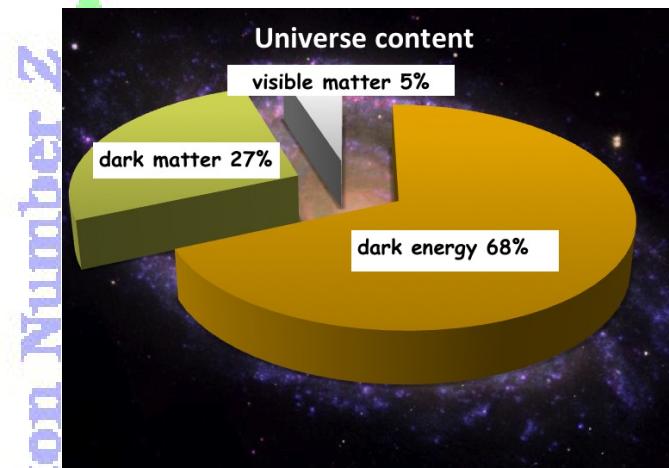
SUMMARY & OUTLOOK

- Nuclear lattice simulations: a new quantum many-body approach
 - based on the successful continuum nuclear chiral EFT
 - a number of highly visible results already obtained
- New developments
 - new highly smeared LO action
 - ↪ good description of isotopic chains of H, He, Be, C and O
 - new probes of nuclear clustering: ρ_3 and ρ_4
 - ↪ increasing entanglement between α -clusters with increasing A
 - new pinhole algorithm
 - ↪ calculation of the proton and neutron distributions in $^{12,14,16}\text{C}$
 - ↪ new method of measuring cluster geometries
- Recent review: Freer, Horiuchi, Kanada-En'yo, Lee, UGM,
 "Microscopic Clustering in Nuclei", arXiv:1705.06192

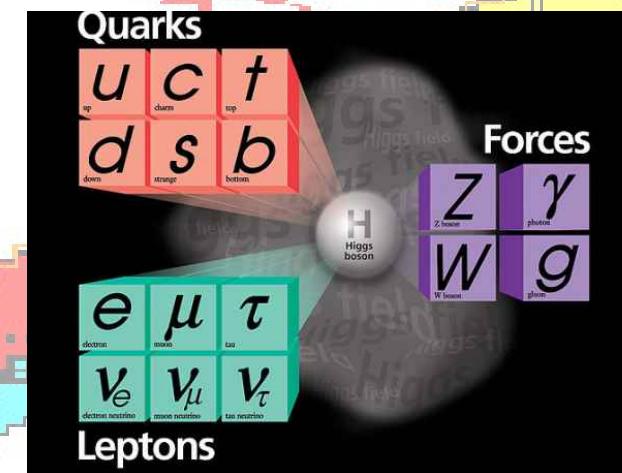
The BIG Picture

WHY NUCLEAR PHYSICS?

- The matter we are made off



- The last frontier of the SM



- Access to the Multiverse

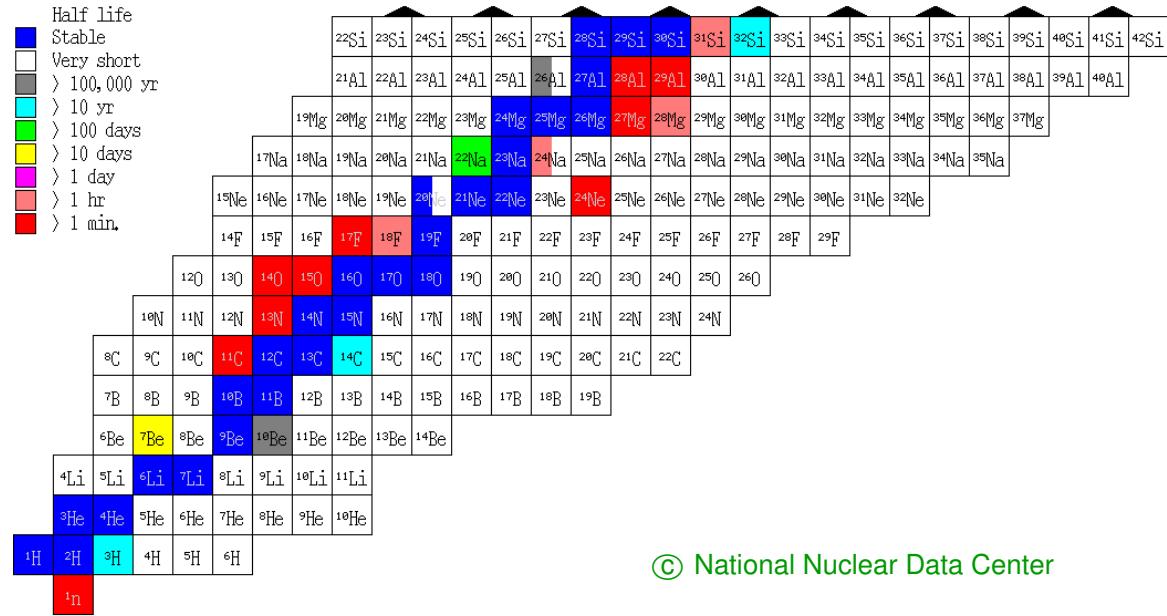


Neutron Number N

AB INITIO NUCLEAR STRUCTURE and SCATTERING

- Nuclear structure:

- ★ 3-nucleon forces
- ★ limits of stability
- ★ alpha-clustering
- ⋮



- Nuclear scattering: processes relevant for nuclear astrophysics

- ★ alpha-particle scattering: $^4\text{He} + ^4\text{He} \rightarrow ^4\text{He} + ^4\text{He}$
- ★ triple-alpha reaction: $^4\text{He} + ^4\text{He} + ^4\text{He} \rightarrow ^{12}\text{C} + \gamma$
- ★ alpha-capture on carbon: $^4\text{He} + ^{12}\text{C} \rightarrow ^{16}\text{O} + \gamma$
- ⋮

SPARES

THE NUCLEAR LANDSCAPE: AIMS & METHODS

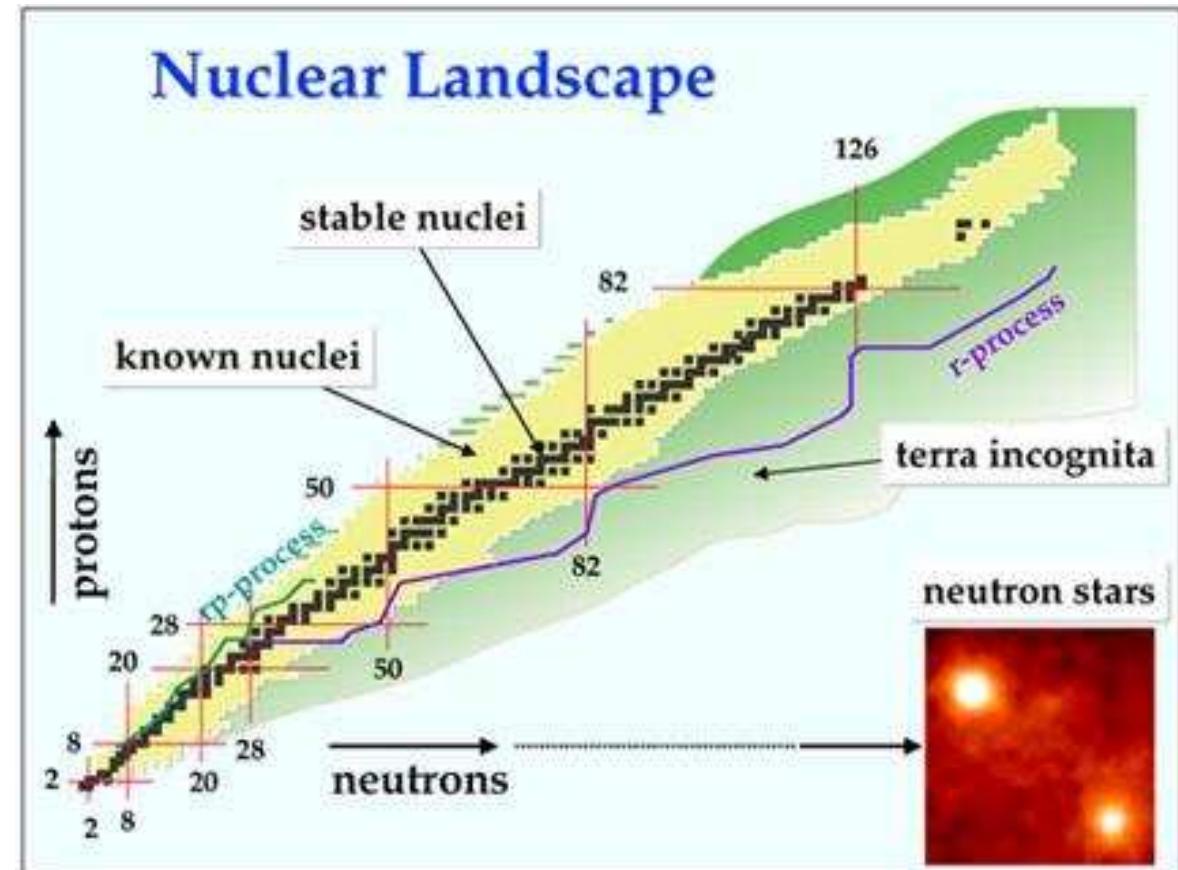
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- Theoretical methods:

- Lattice QCD: $A = 0, 1, 2, \dots$
- NCSM, Faddeev-Yakubowsky, GFMC, ... :
 $A = 3 - 16$
- coupled cluster, ... : $A = 16 - 100$
- density functional theory, ... : $A \geq 100$

- Chiral EFT:

- provides **accurate 2N, 3N and 4N forces**
- successfully applied in light nuclei with $A = 2, 3, 4$
- combine with simulations to get to larger A



⇒ Chiral Nuclear Lattice Effective Field Theory

MANY–BODY APPROACHES

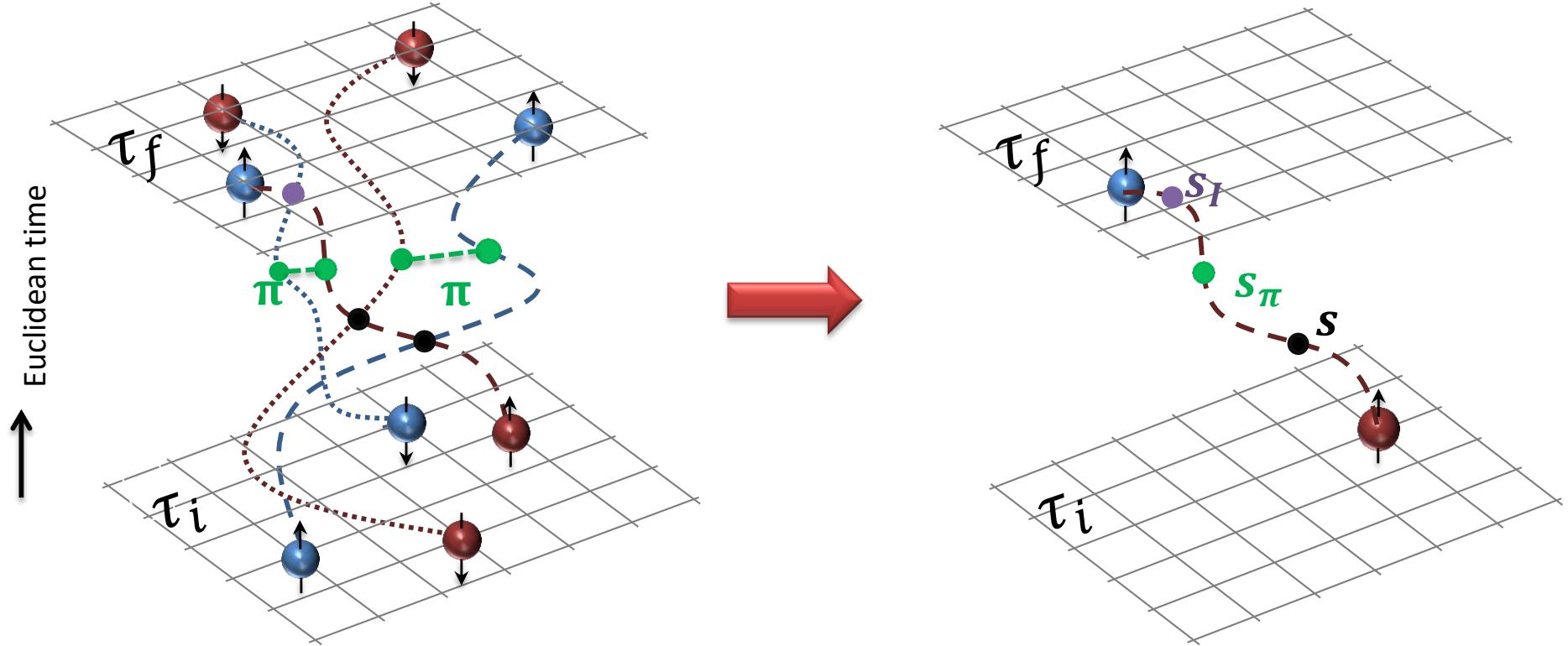
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- nuclear physics = notoriously difficult problem: strongly interacting fermions
- define *ab initio*: combine the precise and well-founded forces from chiral EFT with a many-body approach
- two different approaches followed in the literature:
 - ★ combine chiral NN(N) forces with standard many-body techniques
Dean, Hagen, Navratil, Nogga, Papenbrock, Schwenk, ...
→ successful, but problems with cluster states (SM, NCSM, CC,...)
 - ★ combine chiral forces and lattice simulations methods
→ this new method is called *nuclear lattice simulations* (NLEFT)
Borasoy, Epelbaum, Krebs, Lee, Lähde, UGM, Rupak, ...
→ rest of the talk

AUXILIARY FIELD METHOD

- Represent interactions by auxiliary fields:

$$\exp \left[-\frac{C}{2} (N^\dagger N)^2 \right] = \sqrt{\frac{1}{2\pi}} \int ds \exp \left[-\frac{s^2}{2} + \sqrt{C} s (N^\dagger N) \right]$$



EXTRACTING PHASE SHIFTS on the LATTICE

30

- Lüscher's method:

Two-body energy levels below the inelastic threshold on a periodic lattice are related to the phase shifts in the continuum

Lüscher, Comm. Math. Phys. **105** (1986) 153

Lüscher, Nucl. Phys. B **354** (1991) 531

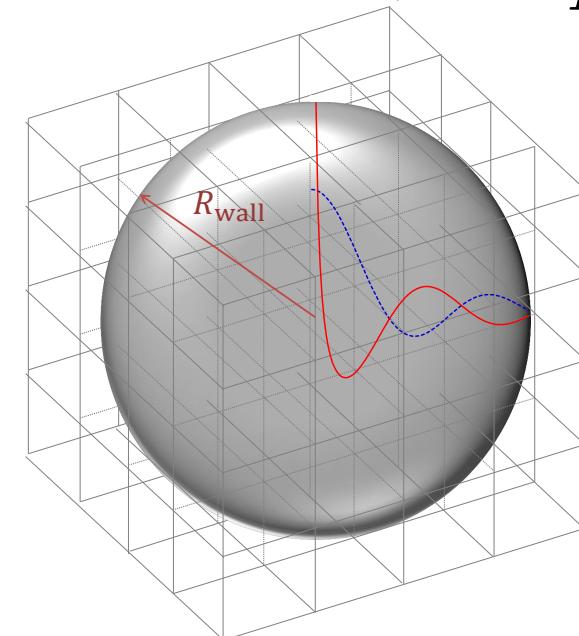
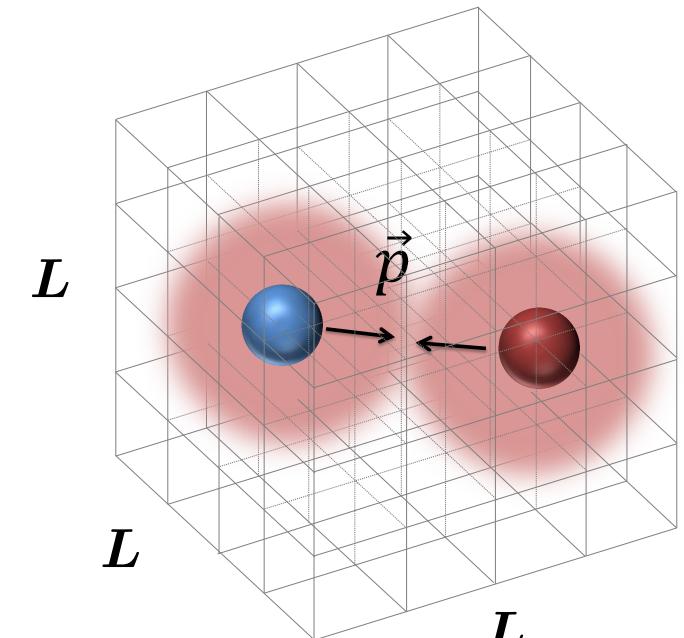
- Spherical wall method:

Impose a hard wall on the lattice and use the fact that the wave function vanishes for $r = R_{\text{wall}}$:

$$\psi_\ell(r) \sim [\cos \delta_\ell(p) F_\ell(pr) + \sin \delta_\ell(p) G_\ell(pr)]$$

Borasoy, Epelbaum, Krebs, Lee, UGM,
EPJA **34** (2007) 185

Carlson, Pandharipande, Wiringa,
NPA **424** (1984) 47



NUCLEAR FORCES: OPEN ENDS

- Why is there this hierarchy $V_{2N} \gg V_{3N} \gg V_{4N}$?
- Gauge and chiral symmetries difficult to include (meson-exchange currents)

Brown, Riska, Gari, . . .

- Connection to QCD ?

most models have one-pion-exchange, but not necessarily respect chiral symmetry

some models have two-pion exchange reconstructed via dispersion relations from $\pi N \rightarrow \pi N$

⇒ We want an approach that

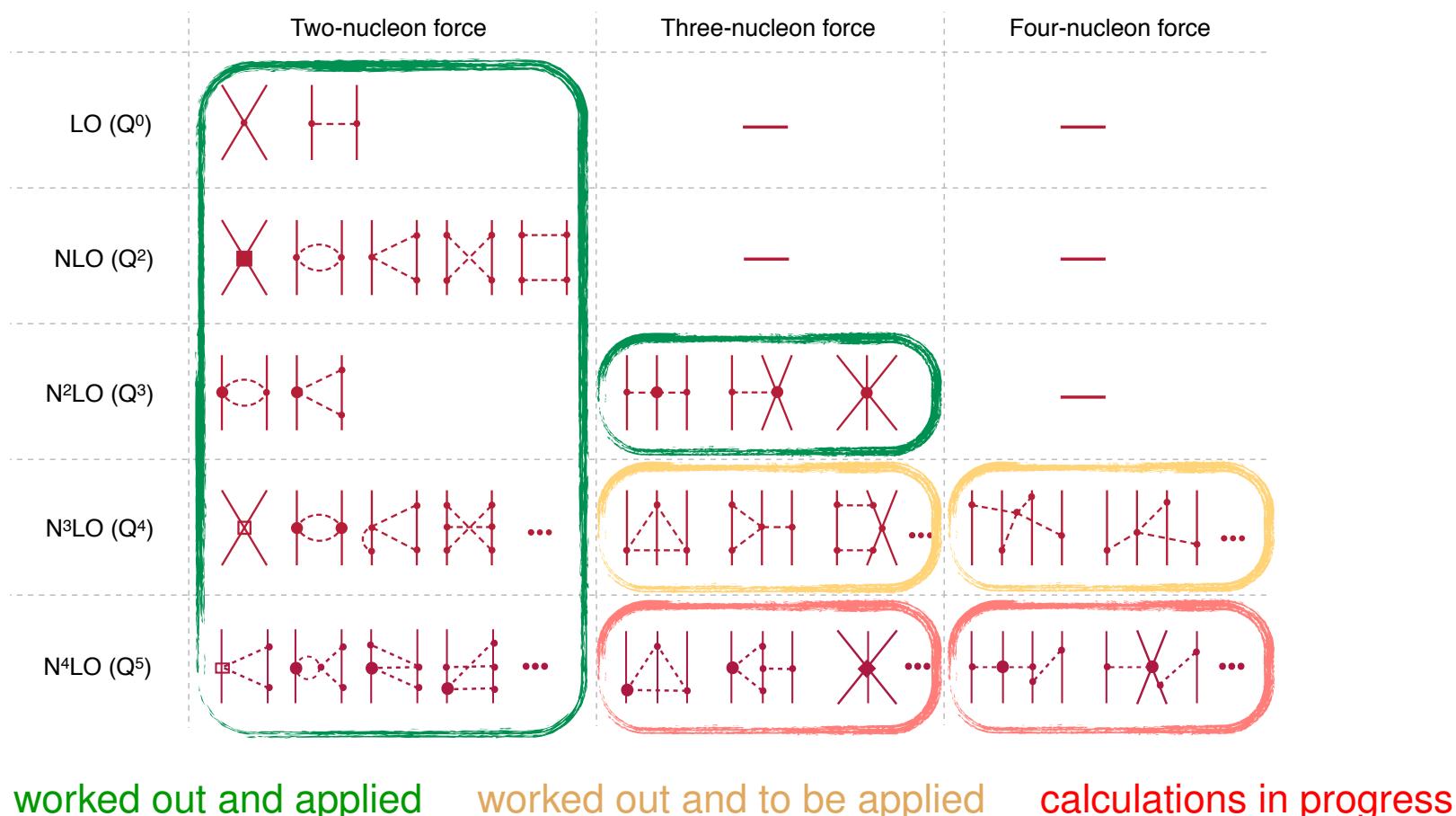
- is linked to QCD via its symmetries
- allows for systematic calc's with a controlled theoretical error
- explains the observed hierarchy of the nuclear forces
- matches nucleon structure to nuclear dynamics
- allows for a lattice formulation / chiral extrapolations
- puts nuclear physics on a sound basis

NUCLEAR FORCES in CHIRAL NUCLEAR EFT

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- expansion of the potential in powers of Q [small parameter]: $\{p/\Lambda_b, M_\pi/\Lambda_b\}$
- explains observed hierarchy of the nuclear forces
- extremely successful in few-nucleon systems

Epelbaum, Hammer, UGM, Rev. Mod. Phys. **81** (2009) 1773

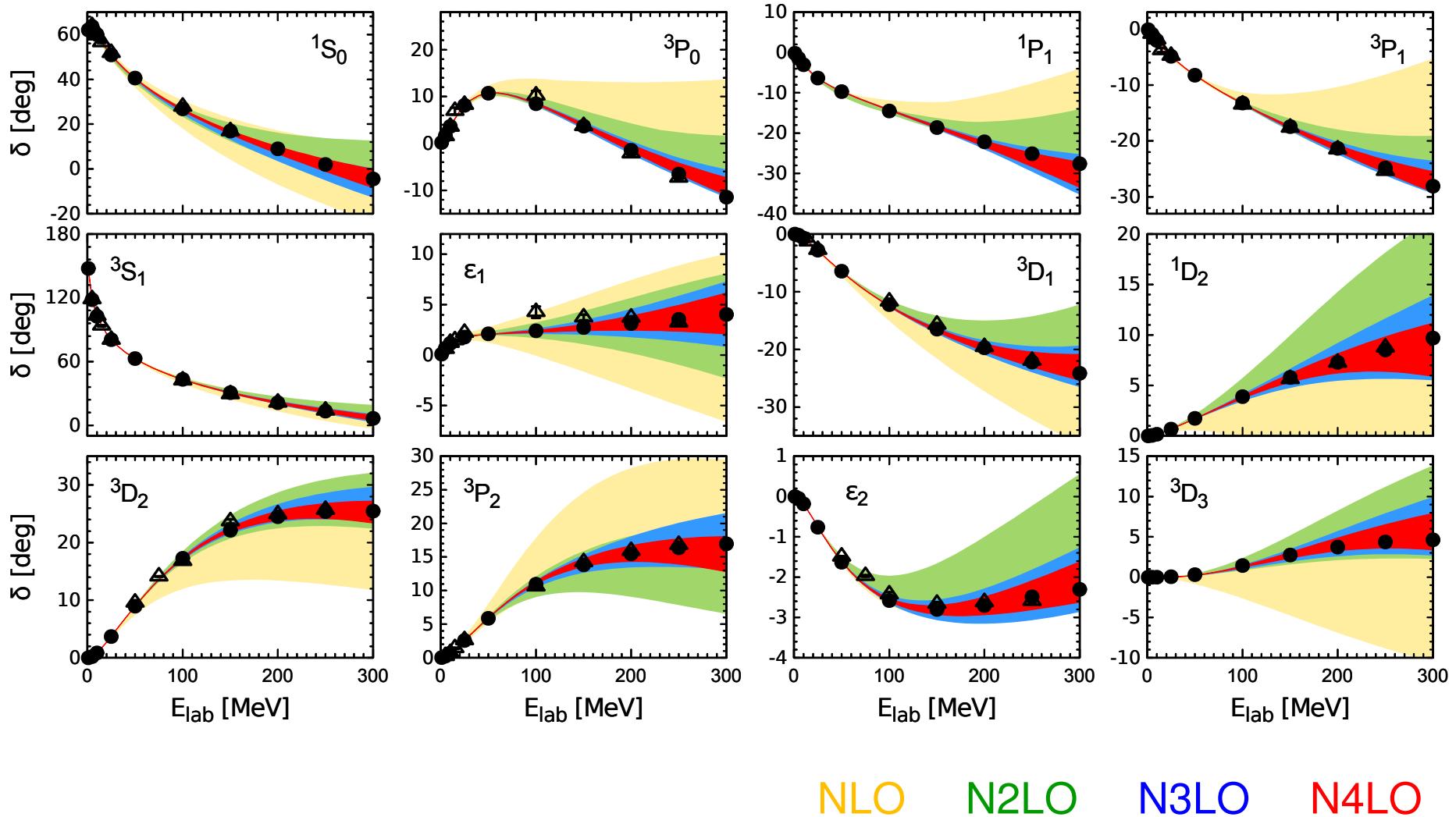


PHASE SHIFTS at N4LO

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⇒ Precision phase shifts with small uncertainties up to $E_{\text{lab}} = 300 \text{ MeV}$

Epelbaum, Krebs, UGM, Phys. Rev. Lett. **115** (2015) 122301



NLO N2LO N3LO N4LO

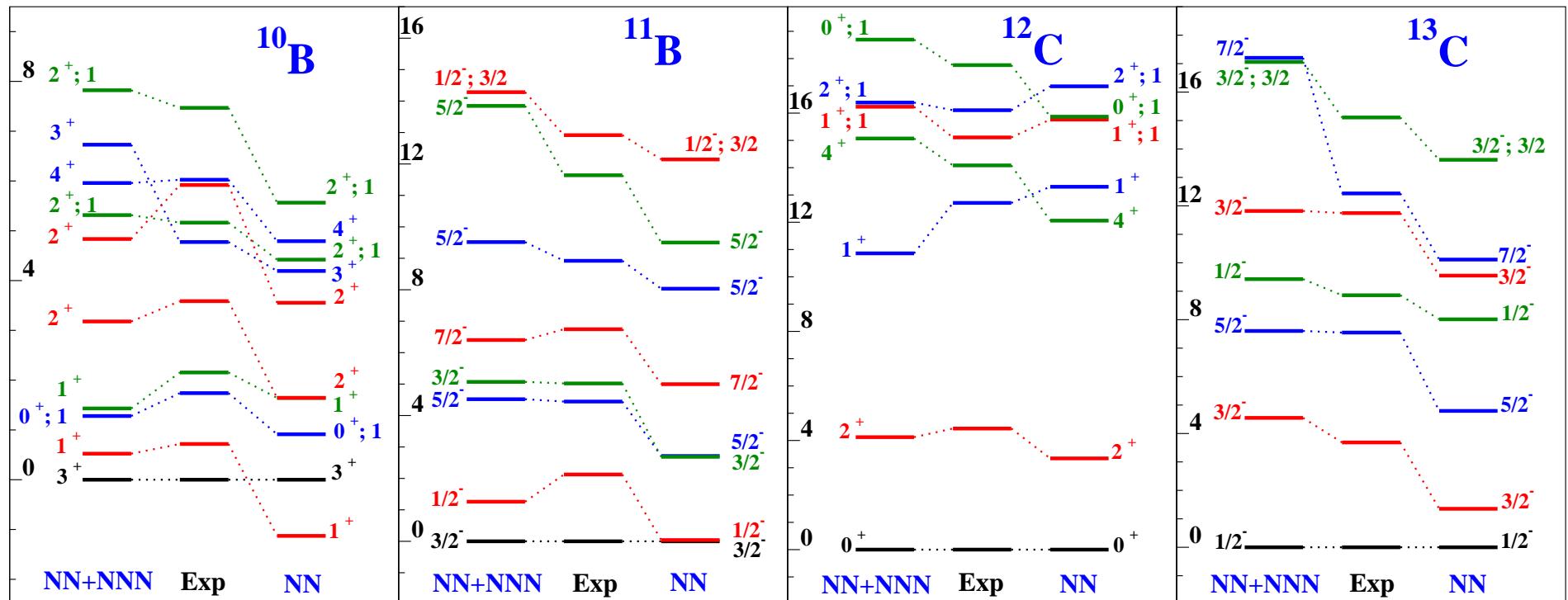
NO-CORE-SHELL MODEL: p-SHELL NUCLEI

- No-core-shell-model calculation

Navratil *et al.*, Phys. Rev. Lett. **99**, 042501 (2007)

- NN interaction at N³LO and NNN interaction at N²LO

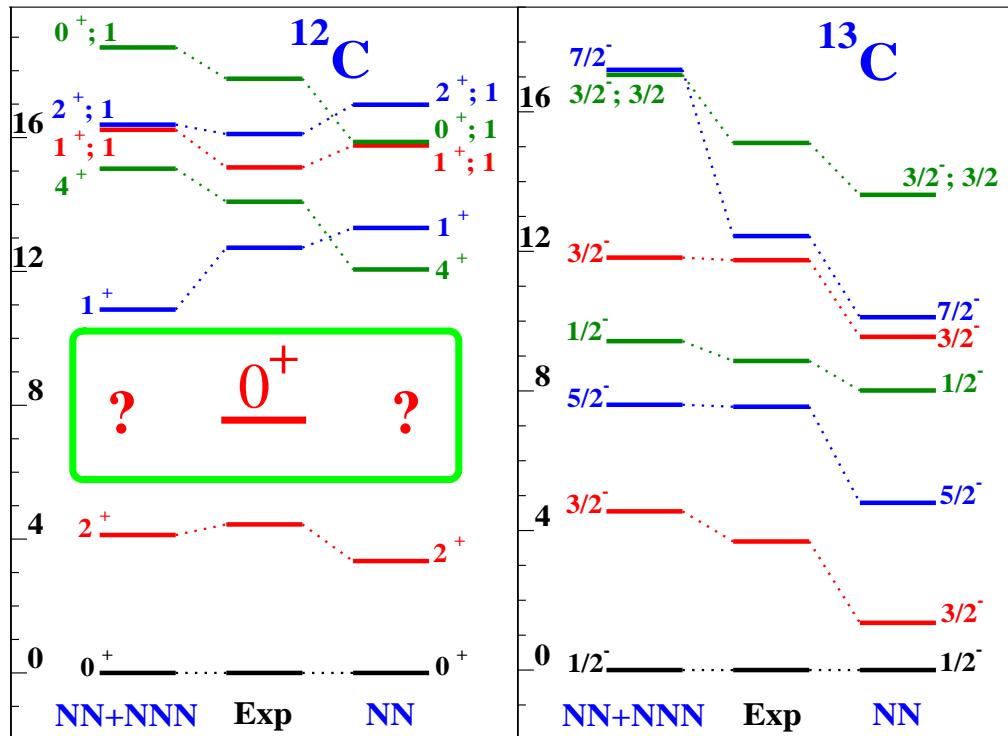
- Fix *D&E* from BE of ³H and level structure of ⁴He, ⁶Li, ^{10,11}B and ^{12,13}C



MODERN MANY-BODY THEORY and the HOYLE STATE³⁵

- one of the most sophisticated many-body theories (No-Core-Shell-Model)
- excellent description of p-shell nuclei from ^6Li to ^{13}C

P. Navratil et al., Phys. Rev. Lett. **99** (2007) 042501 + updates

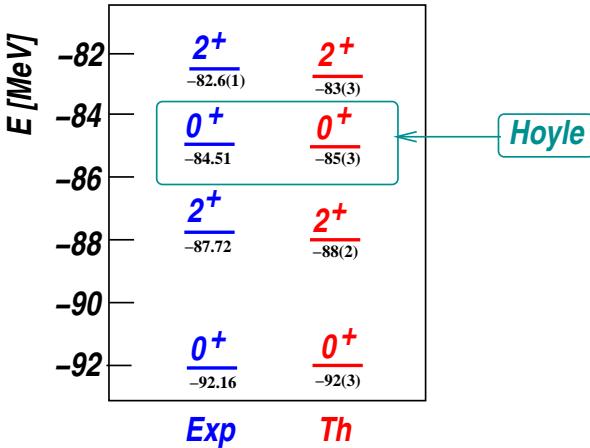


⇒ NO signal of the Hoyle state (i.g. α -cluster states)
⇒ must develop a better method

RESULTS from LATTICE NUCLEAR EFT

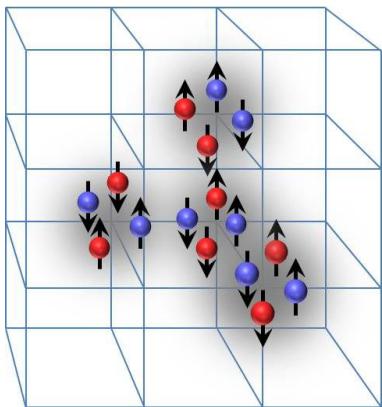
- Hoyle state in ^{12}C

PRL 106 (2011)



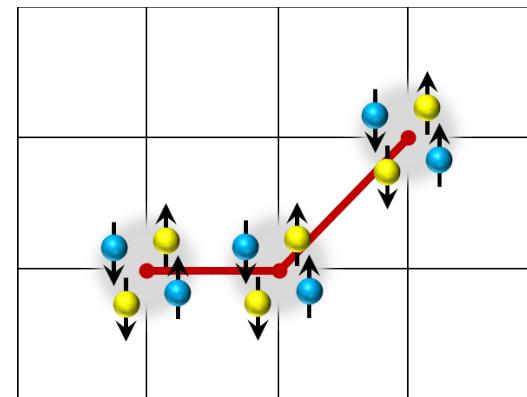
- Spectrum of ^{16}O

PRL 112 (2014)



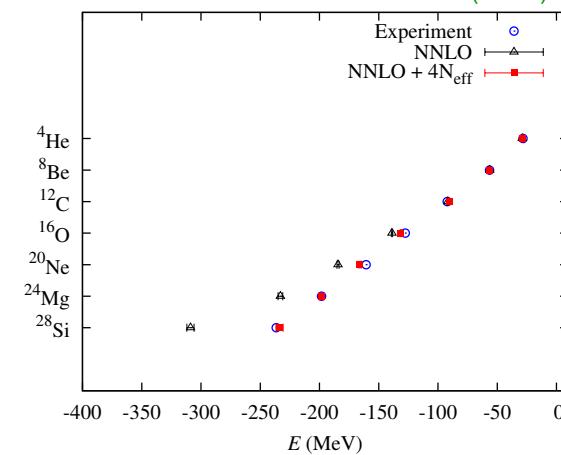
- Structure of the Hoyle state

PRL 109 (2012)



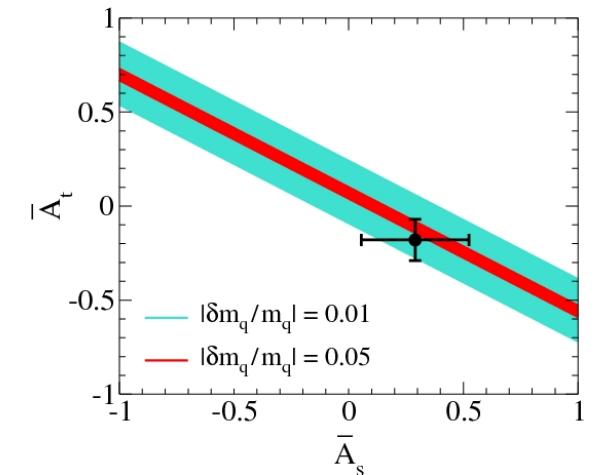
- Going up the α -chain

PLB 732 (2014)



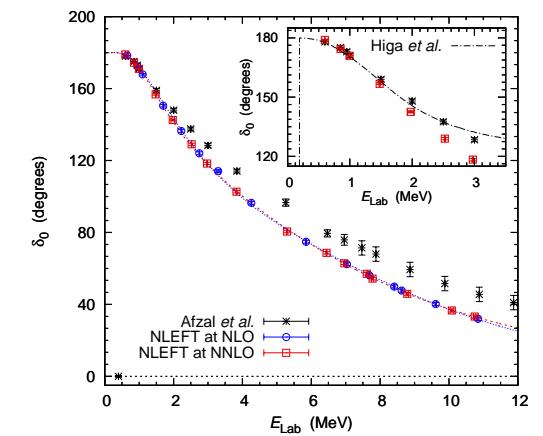
- Fate of carbon-based life

PRL 110 (2013), EPJA 49 (2013)



- Ab initio α - α scattering

Nature 528 (2015)

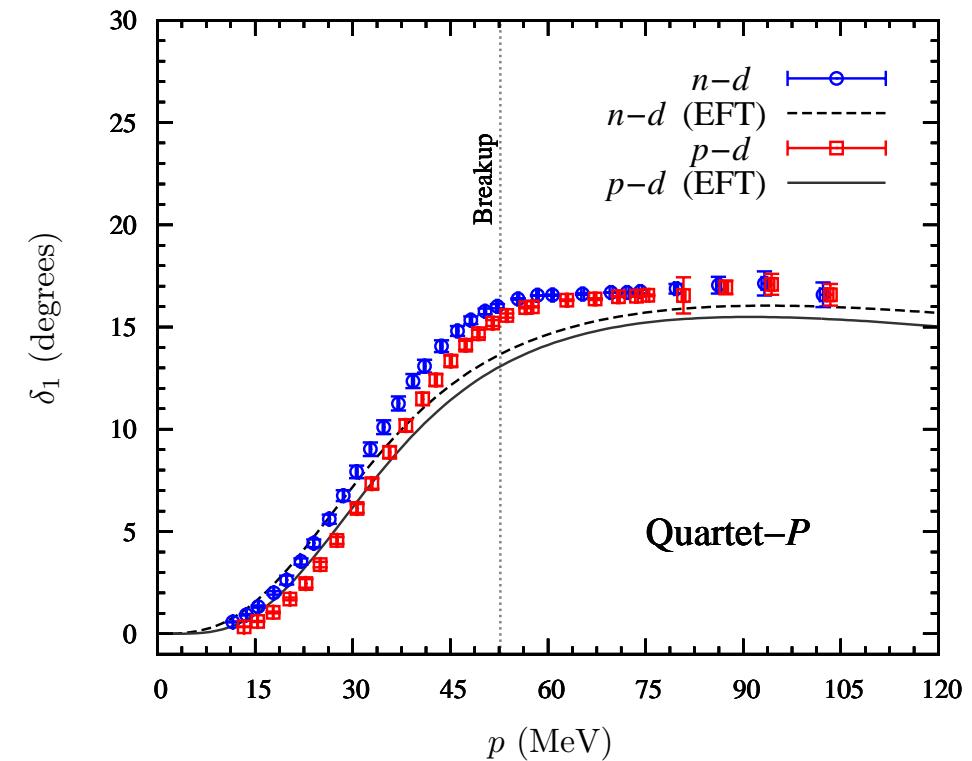
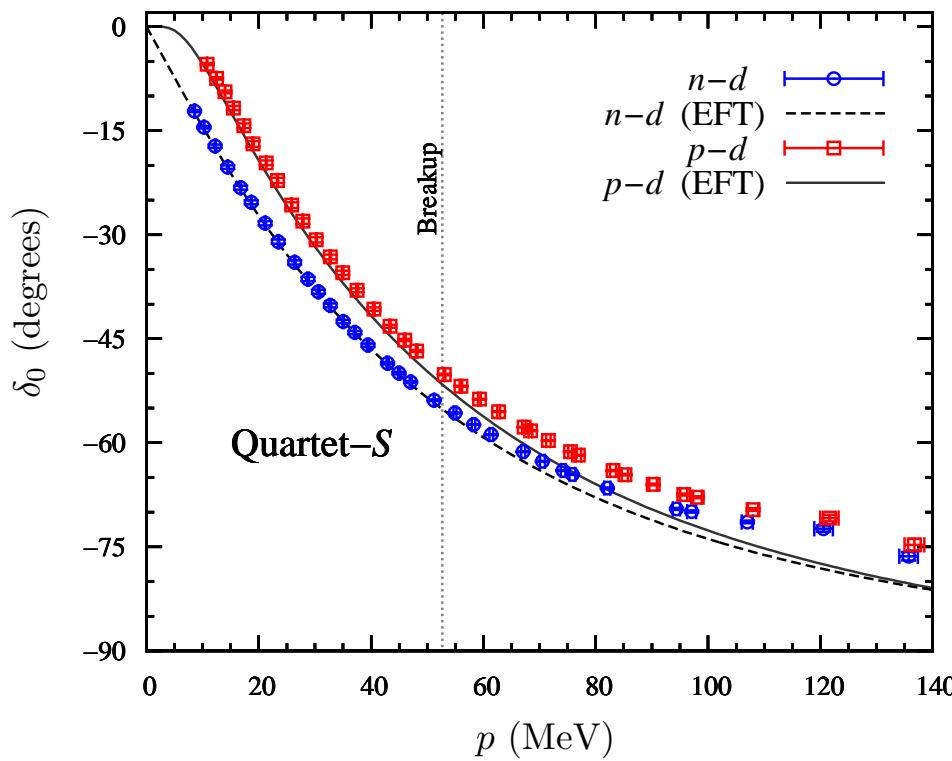


ANOTHER TEST: NUCLEON–DEUTERON SCATTERING³⁷

Elhatisari, Lee, UGM, Rupak, Eur. Phys. J. A 52 (2016) 174

- Use improved methods (cluster states projected on sph. harmonics, etc.) & algorithmic improvements
- Precision calculation of proton-deuteron and neutron-deuteron scattering

Pionless EFT: König, Hammer, Gabbiani, Bedaque, Rupak, Griesshammer, van Kolck, 1998-2011



Nuclear binding near a quantum phase transition

Elhatisari, Li, Rokash, Alarcon, Du, Klein, Lu, UGM, Epelbaum,
Krebs, Lähde, Lee, Rupak,
Phys. Rev. Lett. **117** (2016) 132501 [arXiv:1602.04539]

Editors' suggestion, featured in Physics viewpoint: D.J. Dean, Physics 9 (2016) 106

GENERAL CONSIDERATIONS

- *Ab initio* chiral EFT is an excellent theoretical framework
- not guaranteed to work well with increasing A
 - possible sources of problems:
higher-body forces, higher orders, cutoff dependence, . . .
- very many ways of formulating chiral EFT at any given order (smearing etc.)
 - use not only NN scattering and light nuclei BEs
but also light nucleus-nucleus scattering data
to pin down the pertinent interactions
 - troublesome corrections might be small
 - investigate these issues using two seemingly equivalent interactions
[not a precision study!]

LOCAL and NON-LOCAL INTERACTIONS

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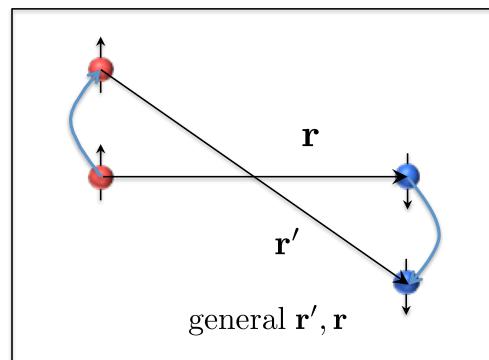
- General potential: $V(\vec{r}, \vec{r}')$

- Two types of interactions:

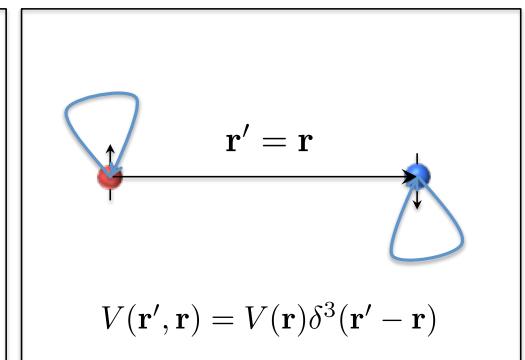
local: $\vec{r} = \vec{r}'$

non-local: $\vec{r} \neq \vec{r}'$

Nonlocal interaction



Local interaction



- Taylor two very different interactions:

Interaction A at LO (+ Coulomb)

Non-local short-range interactions
+ One-pion exchange interaction
(+ Coulomb interaction)

→ tuned to NN phase shifts

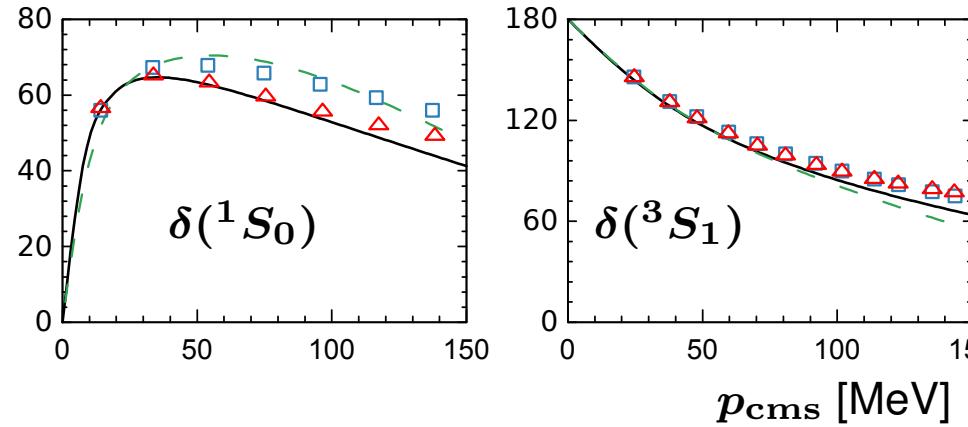
Interaction B at LO (+ Coulomb)

Non-local short-range interactions
+ Local short-range interactions
+ One-pion exchange interaction
(+ Coulomb interaction)

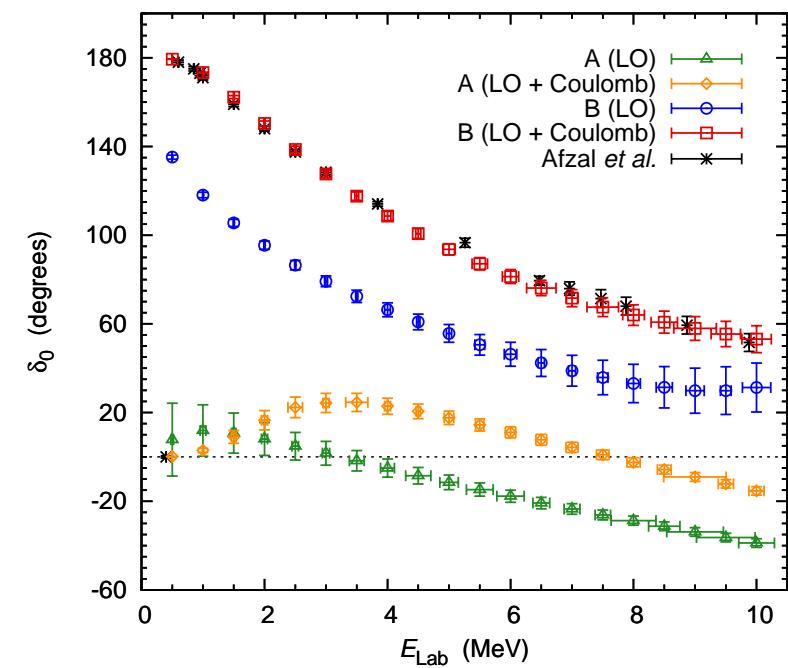
→ tuned to NN + α - α phase shifts

NN and ALPHA-ALPHA PHASE SHIFTS

- Both interactions very similar for NN but **not** for α - α phase shifts:



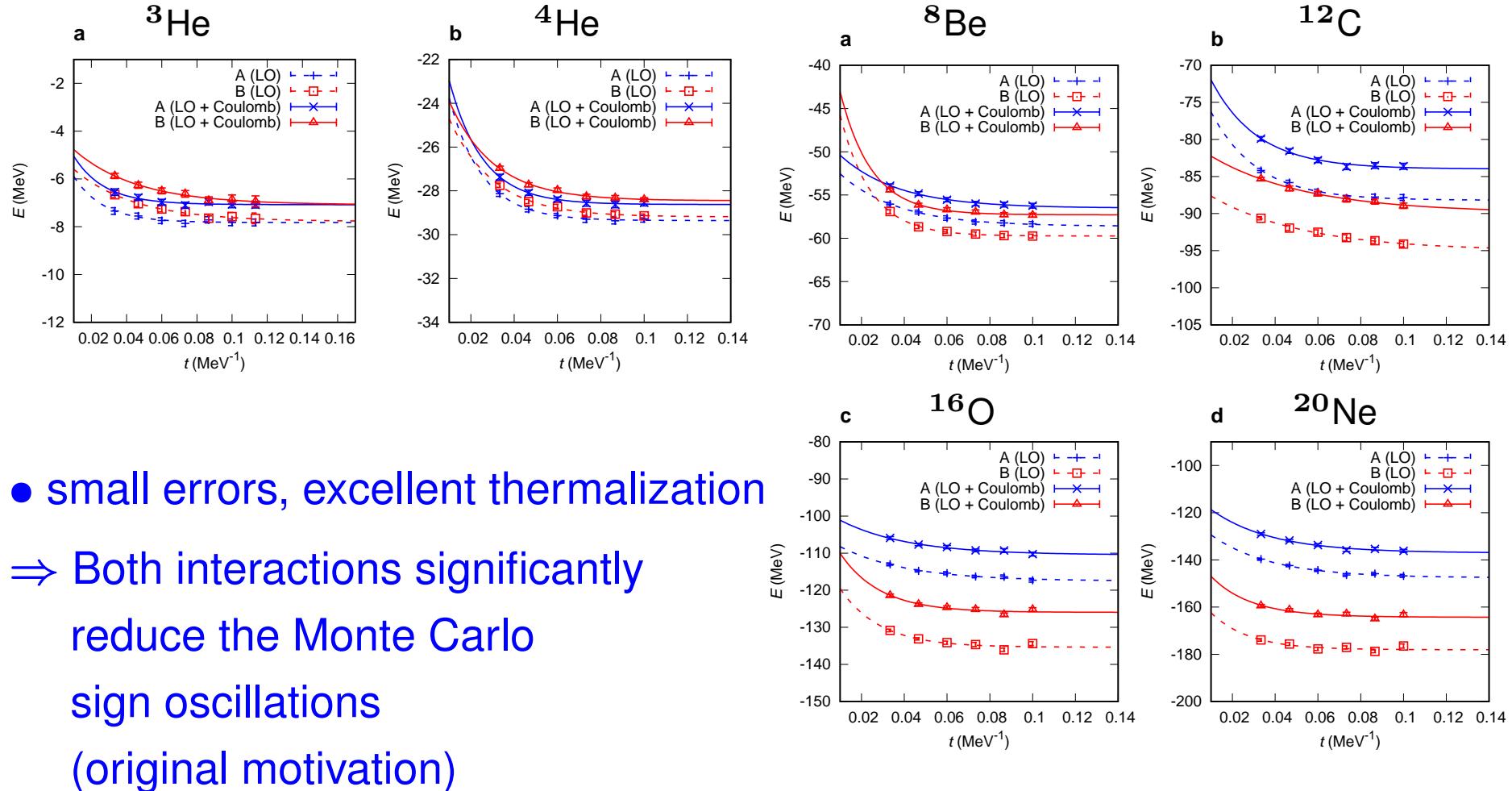
Nijmegen PWA —
 Continuum LO - -
 Lattice LO-A □
 Lattice LO-B △



- Interaction A fails, interaction B fitted
- ↪ consequences for nuclei?

GROUND STATE ENERGIES I

- Ground state energies for alpha-type nuclei plus ^3He :



- small errors, excellent thermalization
 ⇒ Both interactions significantly
 reduce the Monte Carlo
 sign oscillations
 (original motivation)

GROUND STATE ENERGIES I

- Ground state energies for alpha-type nuclei (in MeV):

	A (LO)	A (LO+C.)	B (LO)	B (LO+C.)	Exp.
⁴ He	-29.4(4)	-28.6(4)	-29.2(1)	-28.5(1)	-28.3
⁸ Be	-58.6(1)	-56.5(1)	-59.7(6)	-57.3(7)	-56.6
¹² C	-88.2(3)	-84.0(3)	-95.0(5)	-89.9(5)	-92.2
¹⁶ O	-117.5(6)	-110.5(6)	-135.4(7)	-126.0(7)	-127.6
²⁰ Ne	-148(1)	-137(1)	-178(1)	-164(1)	-160.6

- B (LO+Coulomb) quite close to experiment (within 2% or better)
- A (LO) describes a Bose condensate of particles:

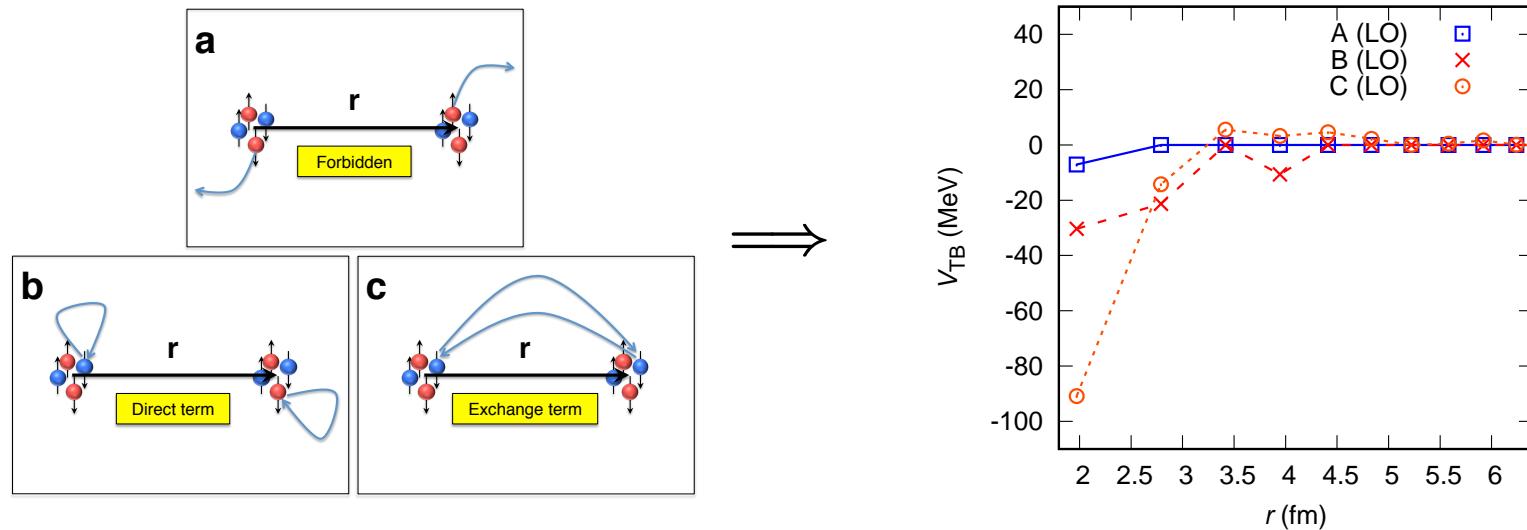
$$E(^8\text{Be})/E(^4\text{He}) = 1.997(6) \quad E(^{12}\text{C})/E(^4\text{He}) = 3.00(1)$$

$$E(^{16}\text{O})/E(^4\text{He}) = 4.00(2) \quad E(^{20}\text{Ne})/E(^4\text{He}) = 5.03(3)$$

FIRST INSIGHT

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- Interaction B was tuned to the nucleon-nucleon phase shifts, the deuteron binding energy, and the S-wave α - α phase shift
- Interaction A starts from interaction B, but *all* local short-distance interactions are switched off, then the LECs of the non-local terms are refitted to describe the nucleon-nucleon phase shifts and the deuteron binding energy
 - The alpha-alpha interaction is sensitive to the degree of locality of the NN int.
 - Qualitative understanding: tight-binding approximation (eff. α - α int.)



CONSEQUENCES for NUCLEI and NUCLEAR MATTER

- Define a one-parameter family of interactions that interpolates between the interactions A and B:

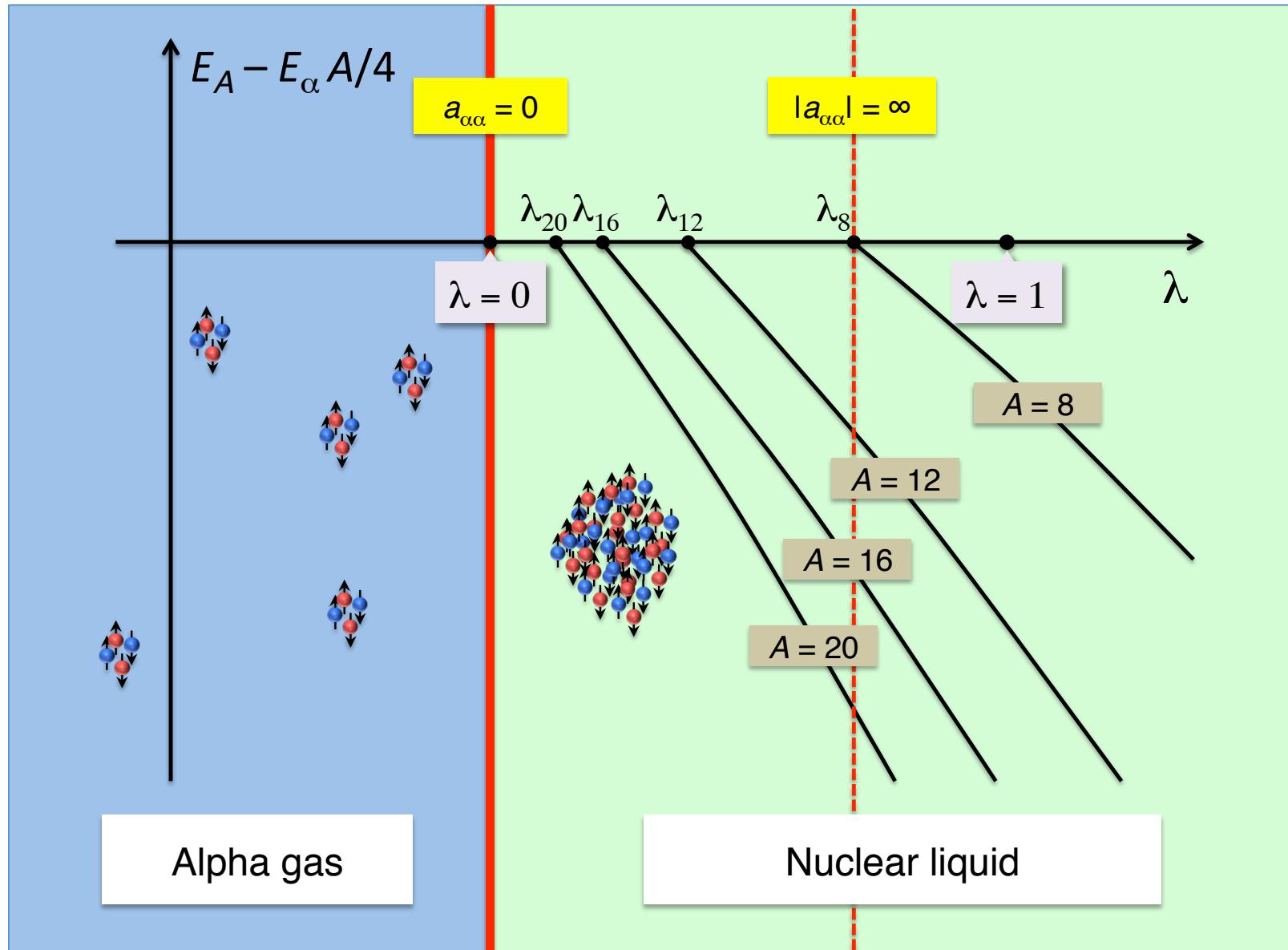
$$V_\lambda = (1 - \lambda) V_A + \lambda V_B$$

- To discuss the many-body limit, we turn off the Coulomb interaction and explore the zero-temperature phase diagram
- As a function of λ , there is a quantum phase transition at the point where the alpha-alpha scattering length vanishes

Stoff, Phys. Rev. A 49 (1994) 3824

- The transition is a first-order transition from a Bose-condensed gas of alpha particles to a nuclear liquid

ZERO-TEMPERATURE PHASE DIAGRAM

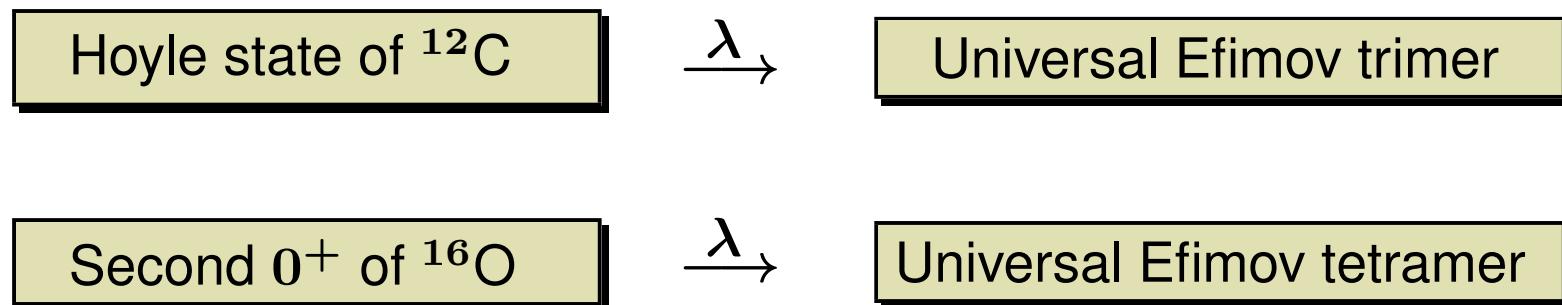


$$\begin{aligned}
 \lambda_8 &= 0.7(1) \\
 \lambda_{12} &= 0.3(1) \\
 \lambda_{16} &= 0.2(1) \\
 \lambda_{20} &= 0.2(1) \\
 \lambda_\infty &= 0.0(1)
 \end{aligned}$$

FURTHER CONSEQUENCES

- By adjusting the parameter λ in *ab initio* calculations, one can move the of any α -cluster state up and down to alpha separation thresholds.
→ This can be used as a new window to view the structure of these exotic nuclear states
- In particular, one can tune the α - α scattering length to infinity!
→ In the absence of Coulomb interactions, one can thus make contact to **universal Efimov physics**:

for a review, see Braaten, Hammer, Phys. Rept. 428 (2006) 259



LOCAL/NON-LOCAL INTERACTIONS on the LATTICE

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- Local operators/densities:

$$a(\mathbf{n}), a^\dagger(\mathbf{n}) \quad [\mathbf{n} \text{ denotes a lattice point}]$$

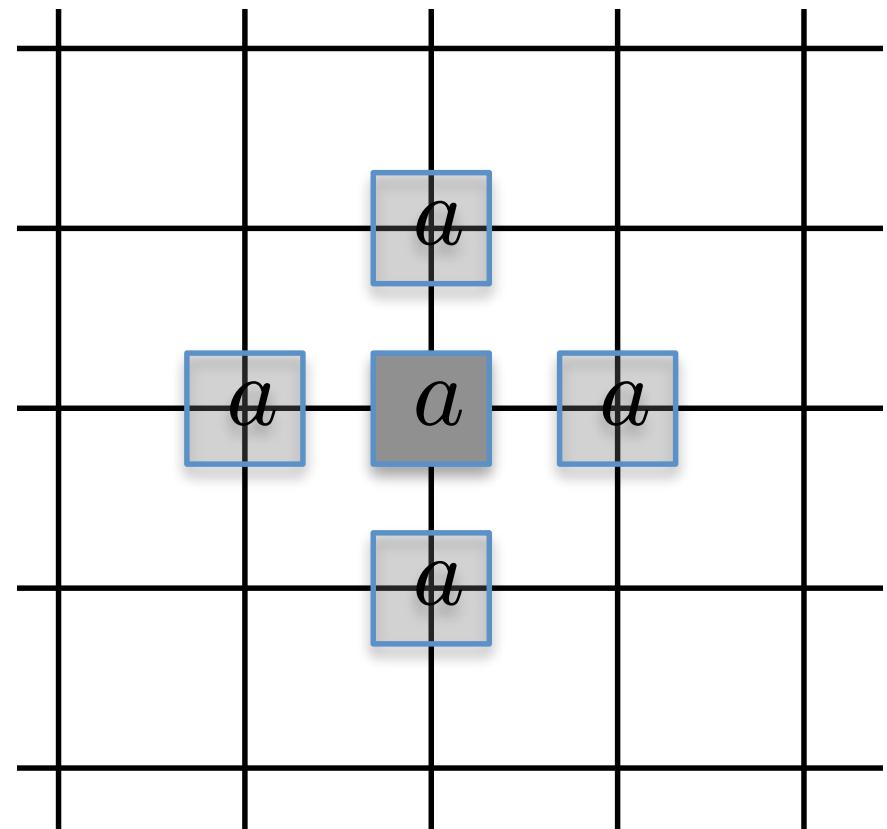
$$\rho_L(\mathbf{n}) = a^\dagger(\mathbf{n})a(\mathbf{n})$$

- Non-local operators/densities:

$$a_{NL}(\mathbf{n}) = a(\mathbf{n}) + s_{NL} \sum_{\langle \mathbf{n}' \mathbf{n} \rangle} a(\mathbf{n}')$$

$$a_{NL}^\dagger(\mathbf{n}) = a^\dagger(\mathbf{n}) + s_{NL} \sum_{\langle \mathbf{n}' \mathbf{n} \rangle} a^\dagger(\mathbf{n}')$$

$$\rho_{NL}(\mathbf{n}) = a_{NL}^\dagger(\mathbf{n})a_{NL}(\mathbf{n})$$

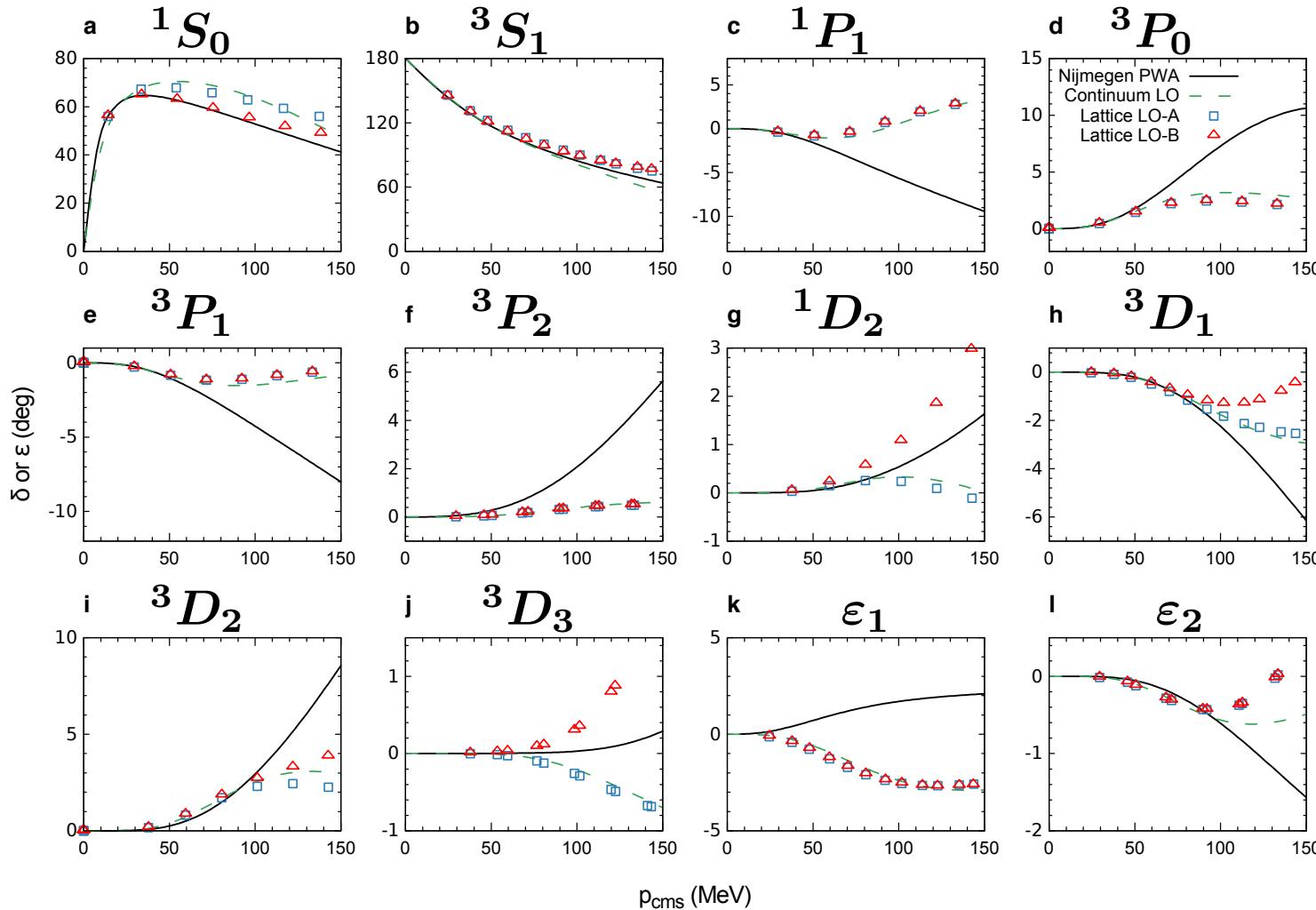


→ where $\sum_{\langle \mathbf{n}' \mathbf{n} \rangle}$ denotes the sum over nearest-neighbor lattice sites of \mathbf{n}

→ the smearing parameter s_{NL} is determined when fitting to the phase shifts

NUCLEON–NUCLEON PHASE SHIFTS

- Show results for NN [and α - α] phase shifts for both interactions:



→ both interactions very similar

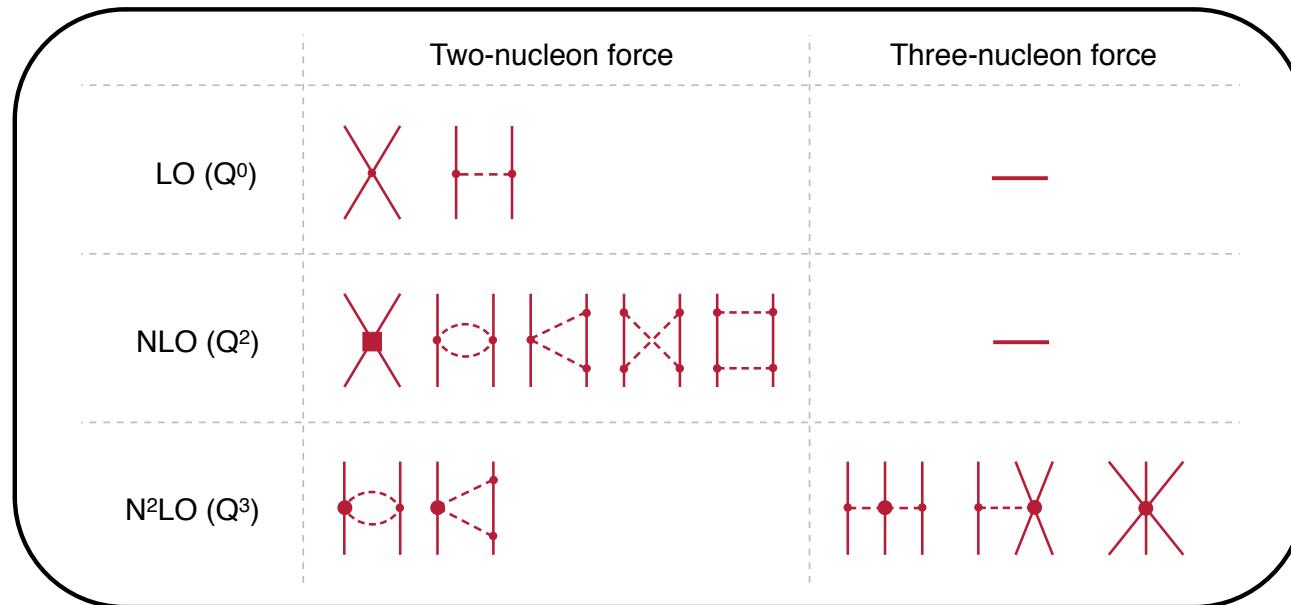
Neutron-proton scattering at NNLO for varying lattice spacings

Alarcón, Du, Klein, Lähde, Lee, Li, Luu, UGM
Eur. Phys. J. A (2017) in print [arXiv:1702.05319]

NUCLEAR FORCES at NNLO

for details, see: Epelbaum, Hammer, UGM, Rev. Mod. Phys. **81** (2009) 1773

- Potential at next-to-next-to-leading order [$Q = \{p/\Lambda, M_\pi/\Lambda\}$]:



- NN potential to NNLO [all πN and $\pi\pi N$ LECs fixed from πN scattering]:

$$\begin{aligned}
 V_{NN} &= V_{\text{LO}}^{(0)} + V_{\text{NLO}}^{(2)} + V_{\text{NNLO}}^{(3)} \\
 &= V_{\text{LO}}^{\text{cont}} + V_{\text{LO}}^{\text{OPE}} + V_{\text{NLO}}^{\text{cont}} + V_{\text{NLO}}^{\text{TPE}} + V_{\text{NNLO}}^{\text{TPE}}
 \end{aligned}$$

NUCLEAR FORCES at NNLO continued

- Analytic expressions [2+7 LECs]:

$$V_{\text{LO}}^{\text{cont}} = \mathbf{C}_S + \mathbf{C}_T (\vec{\sigma}_1 \cdot \vec{\sigma}_2)$$

$$V_{\text{LO}}^{\text{OPE}} = -\frac{g_A^2}{4F_\pi^2} \tau_1 \cdot \tau_2 \frac{(\vec{\sigma}_1 \cdot \vec{q})(\vec{\sigma}_2 \cdot \vec{q})}{q^2 + M_\pi^2}$$

\vec{q} = t-channel mom. transfer

$$\begin{aligned} V_{\text{NLO}}^{\text{cont}} = & \mathbf{C}_1 q^2 + \mathbf{C}_2 k^2 + (\mathbf{C}_3 q^2 + \mathbf{C}_4 k^2) (\vec{\sigma}_1 \cdot \vec{\sigma}_2) + i \mathbf{C}_5 \frac{1}{2} (\vec{\sigma}_1 + \vec{\sigma}_2) \cdot (\vec{q} \times \vec{k}) \\ & + \mathbf{C}_6 (\vec{\sigma}_1 \cdot \vec{q})(\vec{\sigma}_2 \cdot \vec{q}) + \mathbf{C}_7 (\vec{\sigma}_1 \cdot \vec{k})(\vec{\sigma}_2 \cdot \vec{k}) \end{aligned}$$

\vec{k} = u-channel mom. transfer

$$\begin{aligned} V_{\text{NLO}}^{\text{TPE}} = & -\frac{\tau_1 \cdot \tau_2}{384\pi^2 F_\pi^4} L(q) [4M_\pi^2 (5g_A^4 - 4g_A^2 - 1) + q^2 (23g_A^4 - 10g_A^2 - 1) \\ & + \frac{48g_A^4 M_\pi^4}{4M_\pi^2 + q^2}] - \frac{3g_A^4}{64\pi^2 F_\pi^4} L(q) [(q \cdot \vec{\sigma}_1)(q \cdot \vec{\sigma}_2) - q^2 (\vec{\sigma}_1 \cdot \vec{\sigma}_2)] \end{aligned}$$

- Loop function: $L(q) = \frac{1}{2q} \sqrt{4M_\pi^2 + q^2} \ln \frac{\sqrt{4M_\pi^2 + q^2} + q}{\sqrt{4M_\pi^2 + q^2} - q}$

$$\rightarrow 1 + \frac{1}{3} \frac{q^2}{4M_\pi^2} + \dots \text{ for } q \ll \Lambda$$

- for coarse lattices $a \simeq 2$ fm, the TPE at N(N)LO can be absorbed in the LECs C_i
- no longer true as a decreases, need to account for the TPE explicitly

A FEW DETAILS ON THE FITS

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- Fits in large & fixed volumes, vary a from 1 to 2 fm:

a^{-1} [MeV]	a [fm]	L	La [fm]
100	1.97	32	63.14
120	1.64	38	62.48
150	1.32	48	63.14
200	0.98	64	63.14

- OPE and TPE LECs completely fixed ($g_A \sim g_{\pi NN}$ and $c_{1,2,3,4}$ from RS analysis)

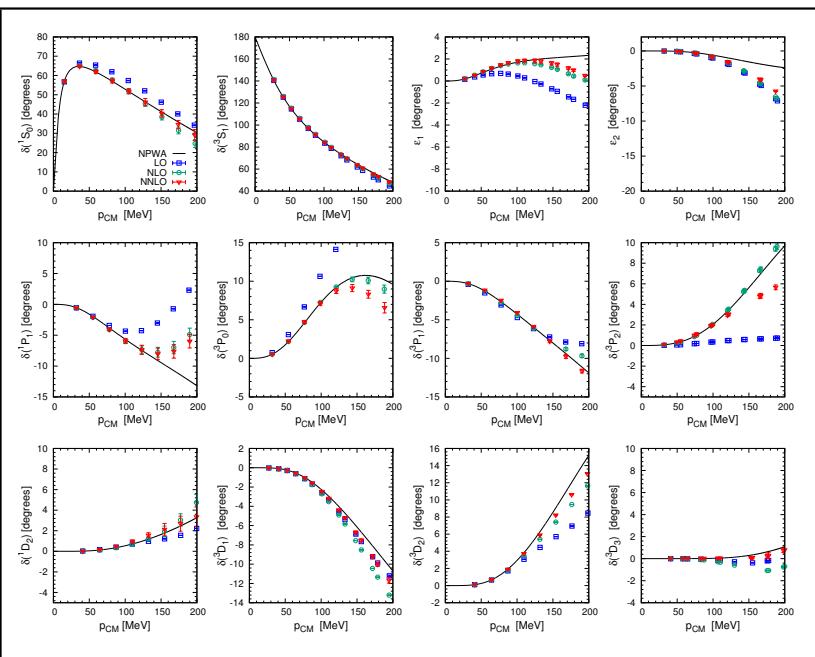
Hoferichter, Ruiz de Elvira, Kubis, UGM, Phys. Rev. Lett. **115** (2015) 092301

- Smeared LO S-wave contact interactions: $f(\vec{q}) \equiv f_0^{-1} \exp\left(-b_s \frac{\vec{q}^4}{4}\right)$
- Partial-wave projection of the contact interactions
 - fit b_s and two S-wave LECs C_i at LO up to $p_{\text{cm}} = 100$ MeV
 - w/ b_s fixed, fit two/seven S/P-wave LECs C_i at NLO/NNLO up to $p_{\text{cm}} = 150$ MeV
 - treat NLO and NNLO corrections perturbatively and non-perturbatively

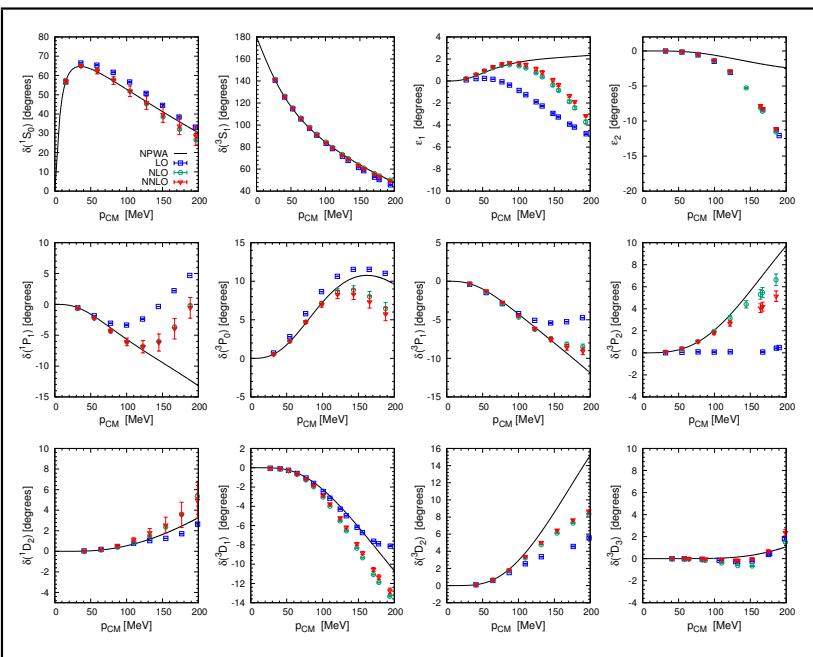
RESULTS for VARIOUS LATTICE SPACINGS - nonpert.

54

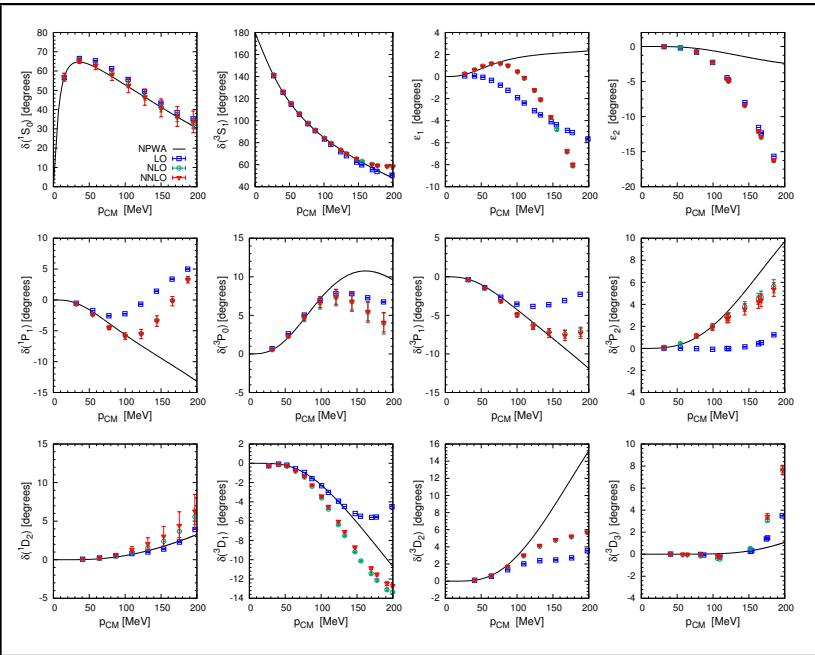
$a = 0.98 \text{ fm}$



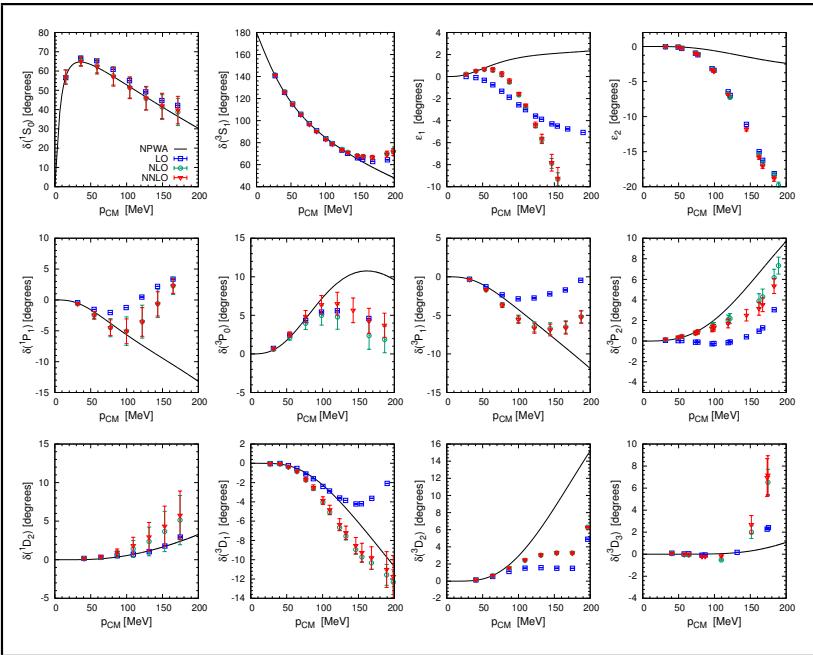
$a = 1.32 \text{ fm}$



$a = 1.64 \text{ fm}$



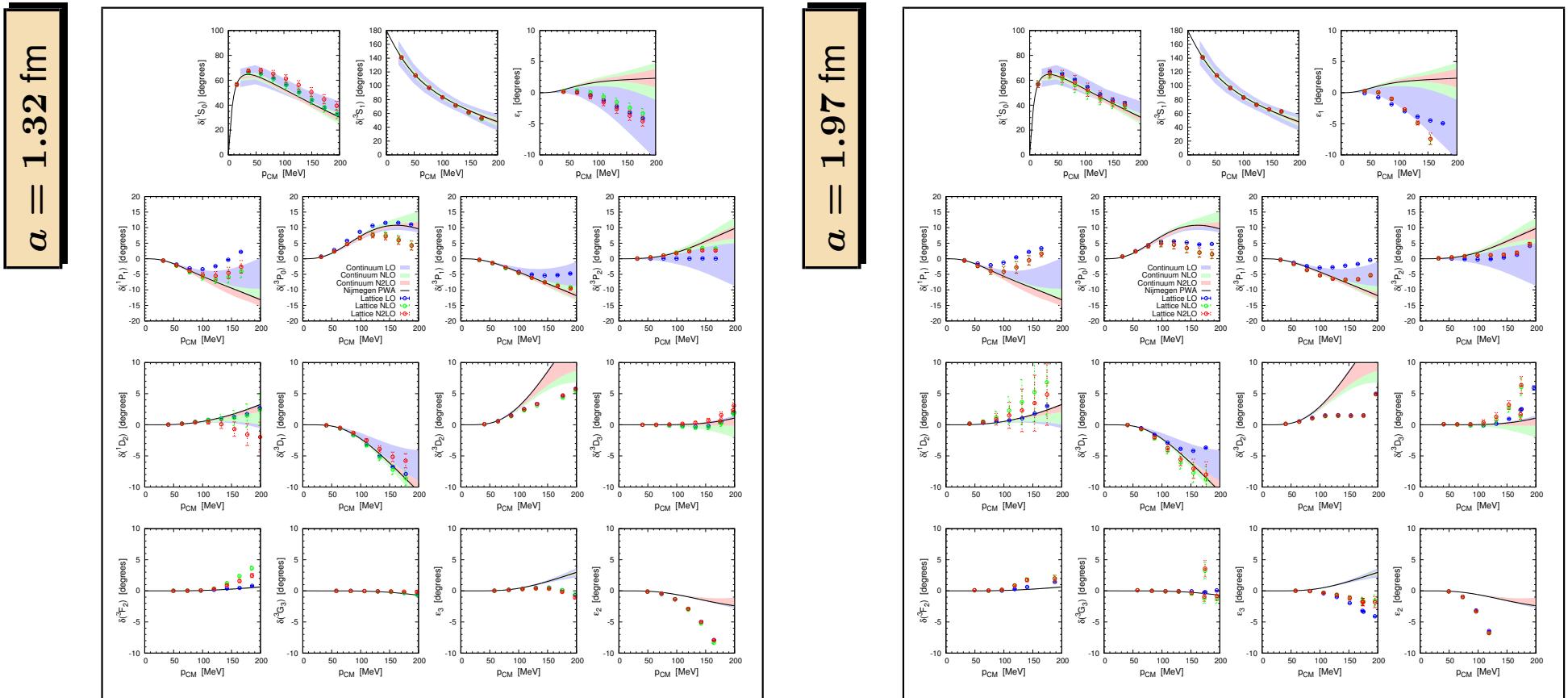
$a = 1.97 \text{ fm}$



RESULTS for VARIOUS LATTICE SPACINGS - pert.

55

- perturbative treatment of NLO and NNLO corrections



- up to $p_{\text{cm}} \simeq 150 \text{ MeV}$, physics is independent of a ✓
- description consistent with the continuum within error bands ✓
- explore this for nuclei — work in progress / stay tuned

