GROUND-STATE PROPERTIES AND MICROSCOPIC DESCRIPTION OF CLUSTERING IN ⁶HE

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Outline



Ab Initio Towards The Driplines

From quarks to nuclei

Nuclear structure



First step: Realistic interactions

Nuclear structure

Realistic nuclear interactions



Chiral EFT-based nuclear Hamiltonians

- EFT for relevant degrees of freedom (π,N) based on symmetries of QCD
- Hierarchy of consistent NN, 3N, 4N,... interactions
- See work by: Weinberg, van Kolck, Epelbaum, Meissner, Krebs, Entem, Machleidt, ...
 - In the following we use:
 - NN: N³LO of Entern & Machleidt with $\Lambda = 500$ MeV
 - ► 3NF: N²LO



Chiral EFT

- E. Epelbaum, H. Hammer, U. Meissner Rev. Mod. Phys. 81 (2009) 1773
- R. Machleidt, D. Entem, Phys. Rep. 503 (2011) 1

Advanced many-body methods

Nuclear structure

Many-body methods

Realistic nuclear interactions



A = 3,4: Several exact methods

- Faddev-Yakubovsky, SVM, GFMC, HH variational, EIHH, NCSM
- Benchmark paper: Kamada et al, Phys. Rev. C64(2001)044001
- Very important observables for testing realistic nuclear Hamiltonians

A > 4: Very few (ab initio) methods available

Intermediate Step: Unitary Transformation

Nuclear structure

Many-body methods

Unitary transformations

Realistic nuclear interactions



- adapt realistic potential to the available model space
- conserve experimentally constrained properties

Intermediate Step: Unitary Transformation

$$V(r) \propto g_1 \frac{e^{-\kappa_1 r}}{r} - g_2 \frac{e^{-\kappa_2 r}}{r}$$
$$\mathcal{F}_r(k - k')$$

Unitary transformation as a function of a flow parameter:

 $\tilde{H}_s = U_s^\dagger H U_s$

SRG evolution equation:

$$\frac{d}{ds}\tilde{H}_s = \left[\eta_s, \tilde{H}_s\right], \quad \eta_s \propto \left[T_{\rm int}, \tilde{H}_s\right]$$

SRG
S. Bogner, R. Furnstahl, A. Schwenk, Prog. Part.
Nucl. Phys. 65 (2010) 94



 $\Lambda_{HO=20MeV; n=17} \sim 3 \text{ fm}^{-1}$

Effective-interaction approaches

Bogner, Furnstahl, Schwenk, Prog. Part. Nucl. Phys. 65 (2010) 94



Many-body nuclear physics studied at low resolution scales comes at a price: the appearance of **many-body forces**.

Extremes of the (light) nuclear landscape

¹⁷Ne

109.2 ms

15O

2.03 m

 ^{14}N

stable

13C

stable

 ^{12}B

29.8 ms

¹¹Be

\$358ms

14O

70.59 s

13N

9.96 m

 ^{12}C

stable

 ^{11}B

stable

 10 Be

1.6 10⁶ a

12N

11 ms

 ^{11}C

20.38 m

 ^{10}B

stable

⁹Be

stable

17F

 $64.8 \, s$

16O

stable

15N

stable

 ^{14}C

5730 a

 ^{13}B

17.33 ms

 ^{12}Be

23.6 ms

 ^{18}F

109.7 m

17O

stable

16N

7.13 s

 ^{15}C

2.45 s

 ^{14}B

13.8 ms

18O

stable

17N

4.17 s

 ^{16}C

 $0.747 \ s$

15B

10.4 ms

 ^{14}Be

4.35 ms

19O

27.1 s

18N

0.63 s

17C

193 ms

19N

329 ms

18C

92 ms

17B

5.1 ms

Realited

19C

49 ms

Observables: anomalous trends

⁹C

126.5 ma

 ^{8}B

770 ms

⁷Be

53.29 d.

 ^{10}C

19.3 s



³He

stable

 ^{2}H

stable

n

10.25 m

 ^{1}H

stable

⁴He

stable

 ^{3}H

12.3 g



Exotic decay modes



Halo states, Borromean systems



Ground-State Properties scales in the NCSM

⁶He Ground-State Properties

 Very accurate charge-radius measurements using laser spectroscopy.

6-He references

- P. Mueller et al., Phys. Rev. Lett. **99** (2007) 252501.
- M. Brodeur et al. Phys. Rev. Lett. **108** (2012) 052504
- S. Bacca. et al. Phys. Rev. C86, (2012) 034321
- Very accurate mass measurement with a Penning trap mass spectrometer.
- Several ab initio calculations
- Most recently by S. Bacca et al
 - ▶ using EIHH and V_{lowk} NN potential based on I-N³LO.
 - Study of Vlowk cutoff-dependence and observable correlations.

Ground-state energy of ⁶He



Typical variational pattern: large $\hbar\Omega$ cuts off wf and small $\hbar\Omega$ cuts off potential. \Rightarrow trade-off between IR and UV regulators

⁶He:Two-neutron separation energy



Note that no induced three-body terms were included \Rightarrow dependence on the SRG flow-parameter in these calculations.

Changes S_{2n} by approximately +/-200 keV in [1.8, 2.2] fm⁻¹ range

C. Forssén, INT, Oct. 31, 2012

Ground-state energy of ⁶He



Ground-state energy of ⁶He



| $E(N_{\max}) = E_{\infty} + c_0 \exp\left(-c_1 N_{\max}\right)$ | | | | |
|---|-------------|---|--|--|
| E(⁶ He) | -29.221 MeV | Exp. fit , Single $\hbar\Omega$ | | |
| E(⁶ He) | -28.967 MeV | Single (N _{max} , $\hbar \Omega$) | | |

Converging sequences

Use results obtained at different HO frequencies to perform a **constrained fit**



CF, J. Vary, E. Caurier, P. Navrátil, Phys. Rev. C 77 (2008) 024301

Powerful property

Any sequence that returns the initial Hamiltonian as the model space is increased will converge to the exact result.

Ground-state energy of ⁶He



| Single (N_{max} , $\hbar \Omega$) | -28.967 MeV | E(⁶ He) |
|---------------------------------------|-----------------|---------------------|
| Exp. fit , Single $\hbar\Omega$ | -29.221 MeV | E(⁶ He) |
| Constrained fit | -29.240(93) MeV | E(⁶ He) |

Ground-state energy of ⁶He



| E(⁶ He) | -28.967 MeV | Single (N_{max} , $\hbar \Omega$) |
|---------------------|-----------------|---------------------------------------|
| E(⁶ He) | -29.221 MeV | Exp. fit , Single $\hbar\Omega$ |
| E(⁶ He) | -29.240(93) MeV | Constrained fit |

⁶He from EIHH



Use hypersperical coordinates, expand in HH and solve hyperradial equations

$$|\Psi\rangle = \sum_{[K]}^{K_{\max}} \sum_{\nu}^{\nu_{\max}} c_{[K],\nu} \mathcal{Y}_{[K]}(\Omega) e^{-\rho/2b} L_{\nu}(\rho)$$

V_{low-k} from I-N3LO with various cutoffs Λ

► Lee-Suzuki effective interaction

Study correlations of observables

6-He measurement and EIHH calculations

M. Brodeur et al. Phys. Rev. Lett.
108 (2012) 052504

• S. Bacca. et al. Phys. Rev. C**86**, (2012) 034321

⁶He: Point-Proton Radius



⁶He: Point-Proton Radius



Switching to IR and UV cutoffs as variables

* Plot $\Delta E/E_{conv}$ as a function of $\lambda_{sc} \equiv \sqrt{m\hbar\Omega/(N+3/2)}$ and $\Lambda \equiv \sqrt{m\hbar\Omega(N+3/2)}$

- Universal dependence on λ_{sc} over wide range of Δ E/E_{conv}.
- Fit shows exponential in 1 / λ_{sc}
- Plateaus to the left from UV corrections

S. Coon et al., arXiv:1205.3230.



Switching to IR and UV cutoffs as variables

* Plot $\Delta E/E_{conv}$ as a function of $\lambda_{sc} \equiv \sqrt{m\hbar\Omega/(N+3/2)}$ and $\Lambda \equiv \sqrt{m\hbar\Omega(N+3/2)}$

- Universal dependence on λ_{sc} over wide range of Δ E/E_{conv}.
- Fit shows exponential in 1 / $\lambda_{\rm sc}$
- Plateaus to the left from UV corrections



Our calculations

- not fixed Λ but color mapped

Nature and form of IR and UV corrections

Truncated basis cuts off s.p. wave functions

• First estimate of cutoffs: $\frac{1}{2}m\Omega^2 r_{\text{max}}^2 = \frac{1}{2m}p_{\text{max}}^2 = (N+3/2)\hbar\Omega$

 $\implies \Lambda_{UV} = \sqrt{2(N+3/2)}\hbar/b$ and $L_0 = \sqrt{2(N+3/2)}b$

with $b=\sqrt{\hbar/m\Omega}$ (note $\sqrt{2}$'s)

• Improved estimate for *L* from intercept of tangent at $r = L_0$:

 $L_{NLO} \approx L_0 + 0.54437 \, b \, (L_0/b)^{-1/3}$



- Square-well wave functions with mass m = 1, radius R = 1, and depth V₀ = 4
- Exact (red) is compared to HO with $\hbar\Omega = 10$ and N = 8 (blue) and to boundary condition at r = L (green) and to n = 4 wf squared (purple)

 $E_{\text{exact}} = -1.51, E_{\text{HO}} = -1.23, E_L = -1.14$ [L improved]

R.J. Furnstahl, talk at this program, 2012-10-10 R.J. Furnstahl et al., Phys. Rev. C 86(2012)031301R.

Correction for energy

Linear energy method to estimate corrections [Djajaputra]

• Let $u_E(r)$ be the radial solution regular at r = 0 with energy E, then

$$u_{L}(r) \equiv u_{E_{L}}(r) \approx u_{\infty}(r) + \Delta E_{L} \left. \frac{du_{E}(r)}{dE} \right|_{E_{\infty}} \quad \text{where} \quad E_{L} = E_{\infty} + \Delta E_{L}$$

So $u_{L}(L) = 0 \implies \Delta E_{L} \approx -u_{\infty}(L) \left(\left. \frac{du_{E}(L)}{dE} \right|_{E_{\infty}} \right)^{-1}$

- Now $u_E(r) \xrightarrow{r \gg R} A_E(e^{-k_E r} + \alpha_E e^{+k_E r})$ with $u_{\infty}(r) \xrightarrow{r \gg R} A_{\infty} e^{-k_{\infty} r}$ and k_{∞} from nucleon separation energy $S = \frac{\hbar^2 k_{\infty}^2}{2m}$
- Take the derivative and evaluate at $E = E_{\infty}$:

$$\frac{du_{E}(r)}{dE}\Big|_{E_{\infty}} = +A_{\infty} \left.\frac{d\alpha_{E}}{dE}\right|_{E_{\infty}} e^{+k_{\infty}r} + \mathcal{O}\left(e^{-k_{\infty}r}\right)$$

Substituting at r = L, we obtain our correction formula to fit:

$$\Delta E_L \approx -\left[\left.\frac{d\alpha_E}{dE}\right|_{E_{\infty}}\right]^{-1} e^{-2k_{\infty}L} + \mathcal{O}(e^{-4k_{\infty}L}) \implies E_L = E_{\infty} + a_0 e^{-2k_{\infty}L}$$

R.J. Furnstahl, talk at this program, 2012-10-10 R.J. Furnstahl et al., Phys. Rev. C 86(2012)031301R.



HO basis cutoff scales

 $\Lambda_{\rm UV} = \sqrt{2(N+3/2)}\hbar/b$

$$L_{\rm IR} = \sqrt{2(N+3/2)}b$$

 $L \approx L_{\rm IR} + 0.544b (L_{\rm IR}/b)^{-1/3}$

Extrapolations from finite HO basis • R.J. Furnstahl et al., Phys. Rev. C 86(2012)031301R

Correction to the energy due to finite HO space

 $E_L = E_{\infty} + \Delta E_L$ with $\Delta E_L = a_0 \exp(-2k_{\infty}L)$



Binding energies



S_{2n}

C. Forssén, INT, Oct. 31, 2012

1.01(1) MeV





Correction for radius

Correction for radius (or other long-distance operators)

• Use $u_L(r) \approx u_{\infty}(r) + \Delta E_L \left. \frac{du_E(r)}{dE} \right|_{E_{\infty}}$ to evaluate

$$\Delta \langle r^2 \rangle_L = \langle r^2 \rangle_L - \langle r^2 \rangle_\infty = \frac{\int_0^L |u_L(r)|^2 r^2 dr}{\int_0^L |u_L(r)|^2 dr} - \frac{\int_0^\infty |u_\infty(r)|^2 r^2 dr}{\int_0^\infty |u_\infty(r)|^2 dr}$$

• For leading *L* dependence, use $u_{\infty}(r) \longrightarrow A_{\infty}e^{-k_{\infty}r}$ and

$$\frac{du_E(r)}{dE}\Big|_{E_{\infty}} \approx -\frac{A_{\infty}}{\Delta E_L} e^{-2k_{\infty}L} e^{+k_{\infty}r} \implies \Delta \langle r^2 \rangle_L \propto \langle r^2 \rangle_{\infty} (2k_{\infty}L)^3 e^{-2k_{\infty}L}$$

• The NLO correction scales as $(2k_{\infty}L) \exp(-2k_{\infty}L)$, so

$$\langle r^2 \rangle_L \approx \langle r^2 \rangle_\infty [1 - (c_0 \beta^3 + c_1 \beta) e^{-\beta}] \text{ with } \beta \equiv 2k_\infty L$$

- $\langle r^2 \rangle_{\infty}$, c_0 , and c_1 are fit parameters while k_{∞} from energy fit
- Valid in the asymptotic regime where $\beta = 2k_{\infty}L \gtrsim 3$
- Both *E* and *r* corrections apply to *A*-body system in lab coordinates

R.J. Furnstahl, talk at this program, 2012-10-10 R.J. Furnstahl et al., Phys. Rev. C 86(2012)031301R.

⁶He: Point-Proton Radius

⁶He point-proton radius



From: R.J. Furnstahl et al., Phys. Rev. C 86(2012)031301R.





⁶He: Point-Proton Radius

| Ref. | r _{pt-p} (⁶ He) [fm] | S _{2n} (⁶ He) [fm] |
|--|---|---|
| Exp | 1.938(23) | 0.97 |
| NCSM: I-N3LO (Λ = 1.8 fm ⁻¹) | 1.81(1) | 1.17(4) |
| NCSM: I-N3LO (Λ = 2.0 fm ⁻¹) | 1.83(1) | 1.01(1) |
| NCSM: I-N3LO (Λ = 2.2 fm ⁻¹) | 1.85(3) | 0.82(2) |



Unsettled questions for the IR extrapolation

Selected from list by R.J. Furnstahl, talk at this program, 2012-10-10

- What is the optimal definition of L?
- How to make credible error estimates?
- Does the interaction matter?
 - The IR corrections are independent of the potential
 - Softer interactions mean more complete UV convergence, so larger region with IR corrections only.
- How well does extrapolation work for other operators?

 $Q(^{6}Li, |^{+})$



$B(GT): {}^{6}He \rightarrow {}^{6}Li$



Conclusions - Cutoff scales

- Introduction of UV and IR scales Combination of results from different N_{max}, ħΩ
- Optimization of run sequence BUT still need several large N_{max} computations
- Prefer UV converged results and perform IR extrapolation
- Transformation of operators

A Microscopic Description Of Clustering

⁶He as a three-body system

- Borromean nucleus
- HH and CSF three-body models with inert cluster.
 V_{nn and} V_nα
 - Core polarization needed
 r_{nα}=1.03r_{nα(free)}
 (cf. three-body force)
 - Repulsive s-wave potential ("Pauli core")
 - HH expansion
 K=2 (90%) with L=S=0 (80%) and L=S=1(10%)
- Pauli focusing.



M.V. Zhukov et al.— Phys. Rep. 231, 151 (1993),

Cluster structures in light nuclei

- Investigate clustering in NCSM wave functions
- Preserve translational invariance
- Harmonic oscillator SD many-body basis
- Transformation
 between single-particle
 and Jacobi coordinates



Three-body cluster overlap functions

$$\begin{split} u_{A-a\alpha l_{1}T_{1},a_{2}\beta l_{2}T_{2},a_{3}\gamma l_{3}T_{3};LS}(\eta_{A-a},\nu_{A-a+1}) \\ &= \left\langle A\lambda JT \right| \left. \begin{array}{l} \mathcal{A}_{A-a,a_{2},a_{3}} \Phi_{\alpha l_{1}T_{1},\beta l_{2}T_{2},\gamma l_{3}T_{3};LS}^{(A-a,a,a_{2},a_{3})JT} : \delta_{\eta_{A-a}}, \delta_{\nu_{A-a+1}} \right\rangle \\ &= \sum_{n_{1(23)},n_{23}} R_{n_{1(23)}l_{1(23}}(\eta_{A-a})R_{n_{23}l_{23}}(\nu_{A-a+1}) \\ &\times \sqrt{\frac{A!}{(A-a)!a_{2}!a_{3}!}} \left\langle A\lambda JT \right| \left. \begin{array}{l} \Phi_{\alpha l_{1}T_{1},\beta l_{2}T_{2},\gamma l_{3}T_{3};LS}^{(A-a,a,a_{2},a_{3})JT} : n_{1(23)}l_{1(23)}, n_{23}l_{23} \right\rangle \end{split}$$



Overlap function for core+N+N

Start with the **core+N+N** case:

- Do a couple of coordinate transformations
 (between relative and s.p.)
- Do a number of spin re-couplings
- Integrate over coordinates



Overlap function for core+N+N

$$\begin{split} u_{A-2\alpha l_{1}T_{1},\frac{1}{2}\frac{1}{2},\frac{1}{2}\frac{1}{2};LS}^{A\lambda JT} &= \sum \frac{R_{n_{1}(23)}l_{1}(23)}{\langle n_{1}(23)l_{1}(23)} \frac{R_{n_{23}l_{23}}(\nu_{A-1})}{\langle n_{1}(23)l_{1}(23)l_{1}(23)} \frac{R_{n_{23}l_{23}}(\nu_{A-1})}{\langle n_{1}(23)l_{1}(23)l_{1}(23)} \frac{R_{n_{23}l_{23}}(\nu_{A-1})}{\langle n_{1}(23)l_{1}(23)l_{1}(23)l_{1}(23)} \frac{R_{n_{23}l_{23}}(\nu_{A-1})}{\langle n_{1}(23)l_{1}(23)l_{1}(23)l_{1}(23)} \frac{R_{n_{23}l_{23}}(\nu_{A-1})}{\langle n_{1}(23)l_{1}(23)l_{1}(23)l_{1}(23)} \frac{R_{n_{23}l_{23}}(\nu_{A-1})}{\langle n_{1}(23)l_{1}(23)l_{1}(23)l_{1}(23)} \frac{R_{n_{23}l_{23}}(\nu_{A-1})}{\langle n_{1}(23)l_{1}(23)l_{1}(23)l_{1}(23)} \frac{R_{n_{23}l_{23}}(\nu_{A-1})}{\langle n_{1}(23)l_{1}(23)l_{1}(23)l_{1}(23)} \frac{R_{n_{23}l_{23}}(\nu_{A-1})}{\langle n_{1}(23)l_{1}(23)l_{1}(23)l_{1}(23)l_{1}(23)} \frac{R_{n_{23}l_{23}}(\nu_{A-1})}{\langle n_{1}(23)l_{1}(23)l_{1}(23)l_{1}(23)l_{1}(23)} \frac{R_{n_{23}l_{23}}(\nu_{A-1})}{\langle n_{1}(23)l_{1}(23)l_{1}(23)l_{1}(23)l_{1}(23)l_{1}(23)} \frac{R_{n_{23}l_{23}}(\nu_{A-1})}{\langle n_{1}(23)l_{1}(23)l_{1}(23)l_{1}(23)l_{1}(23)l_{1}(23)} \frac{R_{n_{23}l_{23}}(\nu_{A-1})}{\langle n_{1}(23)l_{1}(23)l_{1}(23)l_{1}(23)l_{1}(23)l_{1}(23)l_{1}(23)} \frac{R_{n_{23}l_{23}}(\nu_{A-1})}{\langle n_{1}(23)l_{1}(23)l_{1}(23)l_{1}(23)l_{1}(23)l_{1}(23)l_{1}(23)} \frac{R_{n_{23}l_{23}}(\nu_{A-1})}{\langle n_{1}(23)l$$

D. Sääf and CF, - in preparation

Ab initio $< {}^{6}\text{He} | {}^{4}\text{He}+n+n > \text{overlap}$



Ab initio $< {}^{6}$ He | 4 He+n+n > overlap

N³LO, SRG NN only, $\Lambda = 2.0$ fm⁻¹, N_{max}=14, HO=20 MeV

<⁶He (0⁺) | ⁴He (0⁺)+n+n>



L=S=0

Ab initio $< {}^{6}$ He | 4 He+n+n > overlap

N³LO, NN only SRG, $\Lambda = 2.0 \text{ fm}^{-1}$, N_{max}=8, HO=16 MeV N³LO, NN+3NF SRG, Λ = 2.0 fm⁻¹, N_{max}=8, HO=16 MeV



with 3NF

Pauli focusing

<⁶He (0⁺) | ⁴He (0⁺)+n+n>

- Dominance of the $I_{1(23)} = I_{23} = 0$ component.
- RR coefficients determine HH under coordinate-system transformation.

E.g. with
$$I_{1(23)} = I_{23} = 0$$
 we get:

- I₃₍₁₂₎= I₁₂=0 for K=0 (almost Pauli forbidden)
- Dominating $I_{3(12)} = I_{12} = I$ for K=2
- Hyperangular function for this component gives the two-peak correlation density.



$< ^{6}$ He | 4 He+n+n > overlap: N_{max} dependence

<⁶He (0⁺) | ⁴He (0⁺)+n+n>



N³LO, SRG NN only, $\Lambda = 2.0$ fm⁻¹, N_{max}=14, HO=20 MeV



Ab initio $< {}^{6}Li | {}^{4}He+n+p > overlap$

<⁶Li (|⁺) | ⁴He (0⁺)+n+n>



N³LO, SRG NN only, $\Lambda = 2.0$ fm⁻¹, N_{max}=14, HO=20 MeV

Note: different scales

Conclusion and Outlook

- Introduction of UV and IR scales Combination of results from different N_{max}, ħΩ
- Optimization of run sequence BUT still need several large N_{max} computations
- Prefer UV converged results and perform IR extrapolation
- Study of tw-body operators, and the transformation of operators
- Microscopic description of clustering
 - Calculate core swelling: rpp
 - Study projection on HH basis

Continuum effects in ab initio calculations

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THE





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Stepping Into The Continuum

Gamow states and the Berggren basis



Gamow Shell Model

• N. Michel et al Phys. Rev. Lett. **89**(2002) 042502;

- J. Rotureau et al. (2006)
- G. Hagen et al. (2010)

$$\sum_{n=b,r} |u_n\rangle \langle \bar{u}_n| + \frac{1}{\pi} \int_{L_+} |u(k)\rangle \langle u^*(k)|dk = 1$$

Gamow states and completeness

- T. Berggren, Nucl. Phys A **109**(1968)265; NPA**389**(1982)261
- T. Lind, Phys. Rev. C **47**(1993)1903

Gamow shell Model

discretization of continuum (|)contour $\sum |u_n\rangle\langle u_n| + \sum |u_{k,i}\rangle\langle u_{k,i}| \approx 1$ n=b,r(ii) construction of many-body basis $|\mathrm{SD}_i\rangle = |u_{i1}, \dots, u_{iA}\rangle$ (iii) construction of Hamiltonian matrix (complex symmetric matrix) $\langle \mathrm{SD}_i | H | \mathrm{SD}_j \rangle$

(iv) many-body spectrum contains: bound, resonant and "spurious" continuum states

Gamow Shell Model

N. Michel et al, PRL 89 (2002)
042502; PRC67 (2003) 054311; PRC70 (2004) 064313; JPG (2009) 013101
G. Hagen et al, PRC71 (2005) 044314

- J. Rotureau et al, PRL 97 (2006) 110603
- G.Papadimitriou et al, PRC(R) 84 (2011) 051304

Pole approximation is 0th order approximation:

 $H^{\text{p.a.}}|\Psi^{\text{p.a.}}\rangle = E^{\text{p.a.}}|\Psi^{\text{p.a.}}\rangle$

 Many-body resonance (bound) state have large overlap:

$$|\Psi^{\text{p.a.}}|\Psi\rangle|$$

ab initio Calculations in the Berggren Basis

NN potential

- Realistic 2b interactions V_{ij}: Argonne VI8 or chiral I-N³LO
- softened by V_{low-k}
- Single-particle states
 - s- and p-shells from HF potential
 - ▶ l > l shells from HO potential
 - Diagonalization
 - many-body Schrödinger equation

- R. Wiringa et al., Phys. Rev. C, 51 (1995)38
 D. Entern et al., Phys. Rev. C, 51 (1995)38
- D. Entem et al., Phys. Rev. C, 68(2003)041001(R)
 C. Bagrage et al., Phys. Rev. C, 68(2003)141001(R)
- S. Bogner et al., Phys. Rep., **386**(2003) |
- J. Rotureau et al., Phys. Rev. Lett., **97**(2006) | 10603



$$H_A = \frac{1}{A} \sum_{i < j}^{A} \frac{\left(\vec{p}_i - \vec{p}_j\right)^2}{2m} + \sum_{i < j}^{A} V_{NN} \left(\vec{r}_i - \vec{r}_j\right) + \dots$$

using J-coupled DMRG (basis truncation)

Ab initio Benchmark calculations in a Gamow Basis



Ab initio Benchmark calculations in a Gamow Basis

⁵He



G. Papadimitrio, J. Rotureau, B. Barrett, M. Ploszajczak, W. Nazarewicz, N. Michel
See FUSTIPEN workshop, March 2012