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It is a major challenge of nuclear theory to develop theories and algorithms that would allows us to understand the properties of these exotic systems.



<sup>6</sup>He

Closed Quantum System (nuclei near the valley of stability)



infinite well



nice mathematical properties:

Exact treatment of the c.m, analytical solution...



Theories that incorporate the continuum

### Continuum Shell Model (CSM)

- H.W.Bartz et al, NP A275 (1977) 111
- A.Volya and V.Zelevinsky PRC 74, 064314 (2006)

# Shell Model Embedded in Continuum (SMEC)

- J. Okolowicz., *et al*, PR 374, 271 (2003)
- J. Rotureau *et al*, PRL 95 042503 (2005)

## Gamow Shell Model (GSM)

- N. Michel *et al*, PRL 89 042502
- N. Michel et al., Phys. Rev. C67, 054311 (2003)
- N. Michel et al., Phys. Rev. C70, 064311 (2004
- G. Hagen et al, Phys. Rev. C71, 044314 (2005)
- N.Michel et al, J.Phys. G: Nucl.Part.Phys 36, 013101 (2009)





GSM application for He chain

PRC 70, 064313 (2004)



- $\checkmark$  Optimal basis for each nucleus via the GHF method
- $\checkmark$  Borromean nature of  $^{6,8}\text{He}$  is manifested
- $\checkmark$  Helium anomaly is well reproduced

### **GSM HAMILTONIAN**

#### We want a Hamiltonian free from spurious CM motion

Lawson method?

Jacobi coordinates?



PRC 38,1 (1988)

coordinates. No spurious states

 $\geq$ pipj matrix elements  $\langle ab | \mathbf{p}_i \mathbf{p}_j | cd \rangle = C \langle a | | \mathbf{p}_i | | c \rangle \langle b | | \mathbf{p}_j | | d \rangle$ 

x complex scaling does not apply to this particular integral...

Recoil term treatment

✓ Two methods which are equivalent from a numerical point of view

i) Transformation in momentum space

$$\frac{\hbar^2}{2\mu}k^2\psi_{nl}(k) + \int_{L_+} V_l(k,q)\psi_{nl}(q)q^2\,dq = E_{nl}\psi_{nl}(k)$$

✓ disregard numerical derivatives

 $p_i \rightarrow k_i$ 

Fourier transformation to return back to r-space

$$\phi_{nl}(r) = \sqrt{\frac{2}{\pi}} \sum_{i=1}^{N} \sqrt{\omega_i} k_i j_l(kr) \psi_{nl}(i)$$

PRC 73 (2006) 064307

ii) Expand  $p_i$  in HO basis

$$\mathbf{p}_i = \sum_{\alpha < \gamma} |\alpha\rangle \langle \alpha | \mathbf{p}_i | \gamma \rangle \langle \gamma |$$

a,γ are oscillator shells a,c are Gamow states

$$\langle a | \mathbf{p}_i | c \rangle = \sum_{\alpha < \gamma} \langle a | \alpha \rangle \langle \alpha | \mathbf{p}_i | \gamma \rangle \langle \gamma | c \rangle$$

- ✓No complex scaling is involved
- ✓ Gaussian fall-off of HO states provides convergence
- $\checkmark$  Convergence is achieved with a truncation of about  $N_{max} \sim 10$  HO quanta

### EXPERIMENTAL RADII OF 6He, 8He, 11Li

Point proton charge radii



L.B.Wang *et al*, PRL **93**, 142501 (2004) P.Mueller *et al*, PRL **99**, 252501 (2007) R.Sanchez *et al* PRL **96**, 033002 (2006) W.Nortershauser *et al* nucl-ex/0809.2607v1 (2008)

 $\rho_{pp}(^{4}He) \neq \rho_{pp}(^{6}He) \neq \rho_{pp}(^{8}He)$ 

Annu.Rev.Nucl.Part.Sci. 51, 53 (2001)

### Comparison of <sup>6</sup>He radius data with nuclear theory models



Charge radii provide a benchmark test for nuclear structure theory!

### GSM calculations for <sup>6</sup>He nucleus



#### Schematic two-body interactions employed

- 1. Separable Gaussian Interaction (GI) (PRC 71 044314)
- 2. Surface Delta Interaction (SDI) (PR 145, 830)
- 3. Surface Gaussian Interaction (SGI) (PRC 70, 064313)
  - The parameter(s) of each force is(are) fitted on the g.s energy of <sup>6</sup>He

GSM calculations for <sup>6</sup>He nucleus



Expression of charge radius in these coordinates  $\langle r_p^2(Z, A) \rangle = \underbrace{\langle r_p^2(Z, A-2) \rangle}_{core} + \underbrace{\left(\frac{2}{A}\right)^2 \frac{1}{4} \langle \vec{r_1}^2 + \vec{r_2}^2 + 2\vec{r_1} \cdot \vec{r_2} \rangle}_{center of mass correction}$ 

 $r_{c-2n} = (r_1 + r_2)/2$ 

Generalization to n-valence particles is straightforward



Renormalization of the integral based on physical arguments (density)
In our calculations we carried out the radial integration until 25fm



With an adequate number of points along the contour the fluctuations become minimal

 We "cut" when for a given number of discretization points the fluctuations are smeared out

### Results and discussion

Interaction	GI	SGI	SDI
$\left\langle r_{pp}^2 \right\rangle^{1/2}$ [fm]	1.912	1.924	1.920
n-n angle $\theta_{nn}$ [degrees]	$90.3^{o}$	$90.4^{o}$	$89.77^{o}$

charge radii and angles for a p-sd model space employed Angles estimated from the available B(E1) data and the average distances between neutrons.

PRC 76, 051602

$$\left\langle \theta_{nn} \right\rangle = 83^{\circ}{}^{+20}_{-10}$$
$$\left\langle \theta_{nn} \right\rangle = 78^{\circ}{}^{+13}_{-18}$$

#### Decomposition of the wavefunction

$(C_k)^2$	GI	SGI	SDI	
$(p_{3/2})_{res}^2$	(0.888, -0.789)	(0.886, -0.776)	(0.867, -0.747)	
$(p_{3/2})^2_{cont}$	(-0.091, 0.027)	(-0.096, 0.013)	(-0.091, 0.007)	
$(p_{3/2})^1_{res}(p_{3/2})^1_{cont}$	(0.125, 0.761)	(0.130, 0.763)	(0.115, 0.740)	J
$(s_{1/2})_{cont}^2$	0.005	0.007	0.004	
$(p_{1/2})^2_{cont}$	0.025	0.049	0.056	
$(d_{5/2})^2_{cont}$	0.0067	0.0156	0.033	
$(d_{3/2})_{cont}^2$	0.0010	0.006	0.0142	

>The p3/2 occupancy is a crucial quantity for the correct determination of the charge radius in <sup>6</sup>He

Results and discussion

- ✓ Different interactions lead to different configuration mixing.
- ✓ <sup>6</sup>He charge radius (R<sub>ch</sub>) is primarily related to the p3/2 occupation of the 2-body wavefunction.
- ✓ The recent measurements put a constraint in our GSM Hamiltonian which is related to the p3/2 occupation.



 ✓ We observe an overall weak sensitivity for both radii and the correlation angle.

#### Comparison with other structure Models



Conclusion and Future Plans

- The very precise measurements on <sup>6</sup>He, <sup>8</sup>He, <sup>11</sup>Li and <sup>11</sup>Be Halos charge radii give us the opportunity to constrain our GSM Hamiltonian.
- >The GSM description is appropriate for modelling weakly bound nuclei with large radial extension.
- The next step: charge radii <sup>8</sup>He, <sup>11</sup>Li, <sup>11</sup>Be assuming an <sup>4</sup>He core. The rapid increase in the dimensionality of the space will be handled by the GSM+DMRG method.

(J.Rotureau *et al* PRL 97 110603 (2006) and nucl-th/0810.0781.v1)

The 2<sup>+</sup> state of <sup>6</sup>He will be used to adjust the quadrupole strength V(J=2,T=1) of the interaction in <sup>8</sup>He and <sup>11</sup>Li. For <sup>11</sup>Li the T=0 channel of the interaction will be fitted to the <sup>6</sup>Li nucleus.

Develop effective interaction for GSM applications in the p and p-sd shells that will open a window for a detailed description of weakly bound systems. The effective GSM interaction depends on the valence space, but also in the position of the <u>thresholds</u> and the position of the <u>S-matrix poles</u>