Ab initio calculations for non-strange and strange few-baryon systems

Applied ab initio methods

Non-strange sector: continuum observables

Strange sector: hypernuclear bound states

Ab initio calculations for non-strange and strange few-baryon systems

- Applied ab initio methods
- Non-strange sector: continuum observables
 Resonances
 S-Factor in presence of Coulomb potential

Strange sector: hypernuclear bound states Benchmark calculation with other ab initio methods

Applied ab initio methods

- A-body system: A position vectors \mathbf{r}_{i} , removal of center of mass coordinate leads to (A-1) Jacobi vectors $\mathbf{\eta}_{i}$ (η_{i} , θ_{i} , ϕ_{i})
- Expansion of ground-state wave function or LIT state on a complete set:
 Hyperspherical Harmonics (HH)
- **3**(A-1) coordinates of HH basis: hyperradius ρ , 3(A-1) angular coordinates Θ_i and Φ_i , (A-2) hyperspherical angles: 1 hyperradius + (3A 4) angles
- HH basis states: eigen states of grand-angular momentum operator depending on the (3A - 4) angles times a hyperradial basis state
- Different HH versions: normally symmetrized basis states, but also a nonsymmetrized HH (NSHH) basis is possible
- Acceleration of convergence: effective interaction (EIHH)

Short-range two-body correlations (CHH)

Applied ab initio methods

Solve Schrödinger or LIT equation with N basis states and increase N up to the point that a sufficient convergence is obtained

LIT method

The LIT of a function R(E) is defined as follows

where the kernel \mathcal{L} is a Lorentzian,

$$\mathcal{L}(E,\sigma) = \frac{1}{(E-\sigma_R)^2 + \sigma_I^2}$$

For inclusive reactions the LIT $L(\sigma)$ is calculated by solving an equation of the form

$$(H - \sigma) \, \tilde{\Psi} = S \,,$$

where H is the Hamiltonian of the system under consideration and S is an asymptotically vanishing source term related to the operator inducing the specific reaction.

The solution Ψ is localized and the LIT is given by

$$L(\sigma) = \langle \tilde{\Psi} | \tilde{\Psi} \rangle$$
.

Alternative way:

$$L(\sigma) = -\frac{1}{\sigma_I} Im(\langle S | \frac{1}{\sigma_R + i\sigma_I - H} | S \rangle).$$

$$|S\rangle = \theta |0\rangle$$
,

where the operator θ induces a specific electroweak reaction.

The corresponding response function is given by

$$\Rightarrow R(E_f) = \int dE_f |\langle f|\theta|0\rangle|^2 \delta(E_f - E_0 - \omega)$$

Ingredients of the solution of the LIT equation via an expansion on a basis of dimension N:

$$N$$
 eigenstates with eigenenergies ϕ_n E_n

and strength

$$S_n = |\langle \phi_n | \theta | 0 \rangle|^2$$

$$L(\sigma) = \sum_{i=1}^{N} \frac{S_n}{(\sigma_R - E_n)^2 + \sigma_I^2}$$

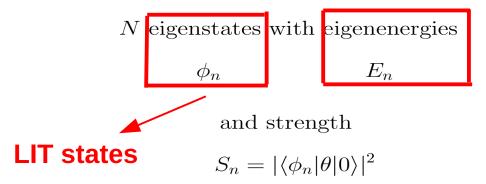
$$|S\rangle = \theta |0\rangle$$
,

where the operator θ induces a specific electroweak reaction.

The corresponding response function is given by

$$\Rightarrow R(E_f) = \int dE_f |\langle f|\theta|0\rangle|^2 \delta(E_f - E_0 - \omega)$$

Ingredients of the solution of the LIT equation via an expansion on a basis of dimension N:



$$L(\sigma) = \sum_{i=1}^{N} \frac{S_n}{(\sigma_R - E_n)^2 + \sigma_I^2}$$

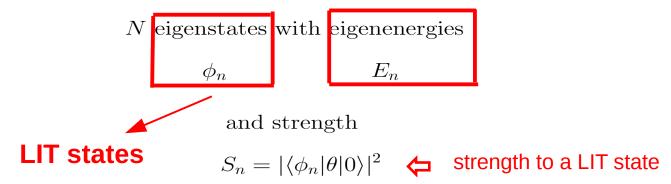
$$|S\rangle = \theta |0\rangle$$
,

where the operator θ induces a specific electroweak reaction.

The corresponding response function is given by

$$\Rightarrow R(E_f) = \int dE_f |\langle f|\theta|0\rangle|^2 \delta(E_f - E_0 - \omega)$$

Ingredients of the solution of the LIT equation via an expansion on a basis of dimension N:



$$L(\sigma) = \sum_{i=1}^{N} \frac{S_n}{(\sigma_R - E_n)^2 + \sigma_I^2}$$

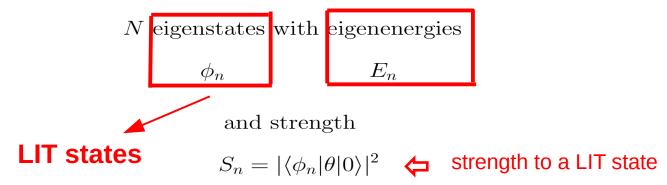
$$|S\rangle = \theta |0\rangle$$
,

where the operator θ induces a specific electroweak reaction.

The corresponding response function is given by

$$\Rightarrow R(E_f) = \int dE_f |\langle f|\theta|0\rangle|^2 \delta(E_f - E_0 - \omega)$$

Ingredients of the solution of the LIT equation via an expansion on a basis of dimension N:



Inversion of the LIT

- \bullet LIT is calculated for a fixed $\sigma_{_{\! I}}$ in many $\sigma_{_{\! R}}$ points
- Express the searched response function formally on a basis set with N basis basis functions $f_n(E)$ and open coefficients c_n with correct threshold behaviour for the $f_n(E)$ (e.g., $f_n = f_{thr}(E) \exp(-\alpha E/n)$)
- Make a LIT transform of the basis functions and determine coefficents c_n by a fit to the calculated LIT
- Increase N up to the point that a sufficient convergence is obtained (uncontrolled oscillations should not be present)

Resonances

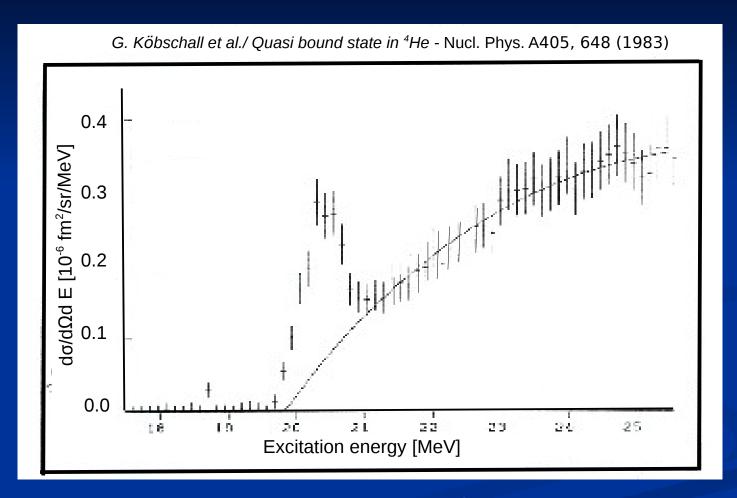
⁴He isoscalar monopole resonance

Isoscalar monopole response function $M(q, E_f = E_0 + \omega)$

with transition operator
$$\theta(q) = \frac{G_E^s(q^2)}{2} \sum_{i=1}^A j_0(qr_i)$$

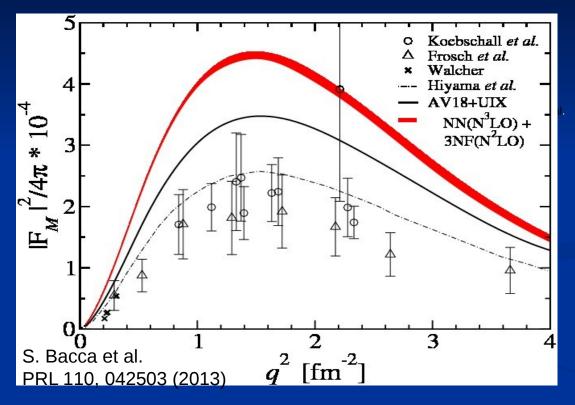
 $G_E^s(q^2)$: nucleon isoscalar electric form factor j_0 : spherical Bessel function of 0^{th} order

0+ Resonance in the ⁴He compound system



Resonance at $E_R = -8.2$ MeV, i.e. above the ³H-p threshold. Strong evidence in electron scattering off ⁴He, $\Gamma = 270\pm50$ keV

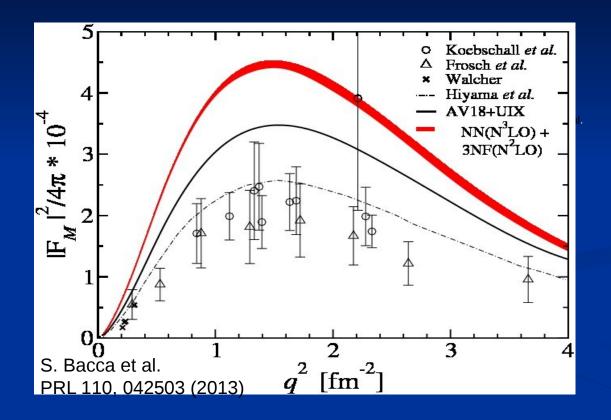
Comparison to experimental results



LIT/EIHH Calculation for AV18+UIX and Idaho-N3LO+N2LO

dash-dotted: AV8' + central 3NF (Hiyama et al.)

Comparison to experimental results



Observable is strongly dependent on potential model

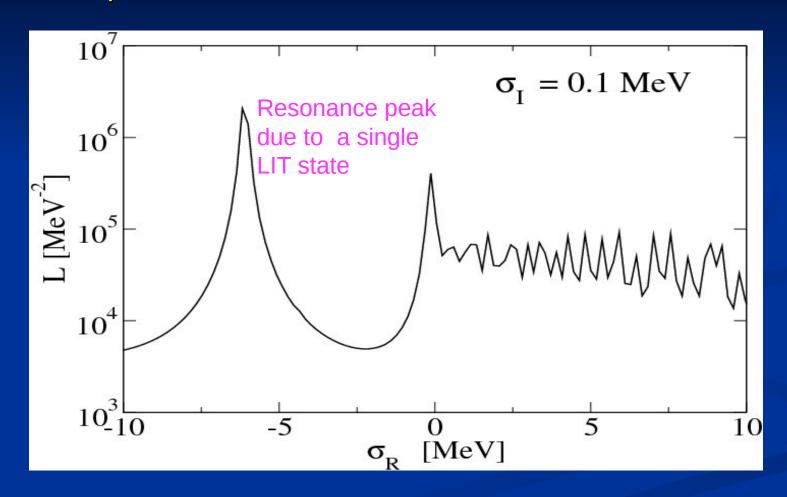
Why were we unable to determine the width of the 4He isoscalar monopole resonance?

To answer this let us check our very first LIT calculation from 1997:

⁴He(e,e') inelastic longitudinal response function

with a central NN potential

Unpublished result from a CHH calculation with the TN potential (V. Efros, WL, G. Orlandini, PRL 78,432 (1997))



To study the problem better let us consider first instead of a four-body reaction a simpler three-body reaction:

3
He + γ \longrightarrow d + p at low energies

LIT calculation with central MTI/III NN potential in unretarded dipole approximation

Aim: Increase low-energy density of LIT states

To study the problem better let us consider first instead of a four-body reaction a simpler three-body reaction:

3
He + γ \longrightarrow d + p at low energies

LIT calculation with central MTI/III NN potential in unretarded dipole approximation

Aim: Increase low-energy density of LIT states

Answer seems to be simple increase number of basis states both for hyperspherical and hyperradial states

To study the problem better let us consider first instead of a four-body reaction a simpler three-body reaction:

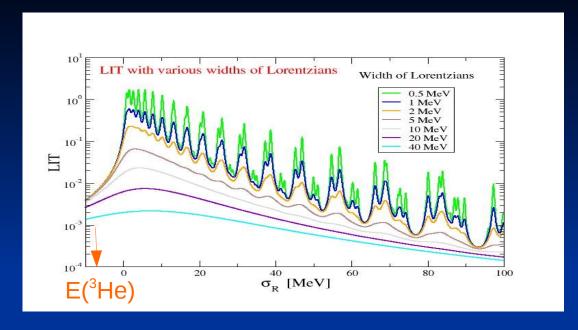
3
He + γ \longrightarrow d + p at low energies

LIT calculation with central MTI/III NN potential in unretarded dipole approximation

Aim: Increase low-energy density of LIT states

Answer seems to be simple increase number of basis states both for hyperspherical and hyperradial states

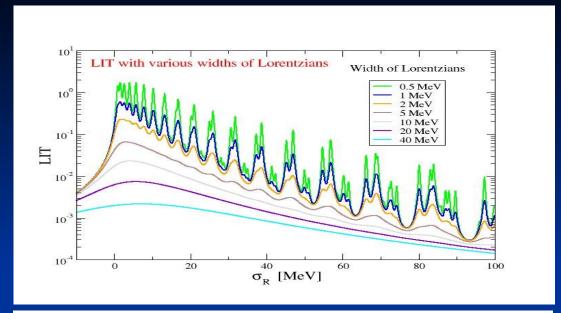
Also note: hyperradial basis states consist in an expansion on Laguerre polynomials times a spatial cutoff $\exp(-\rho/b)$ Increase of b shifts spectrum to lower energies

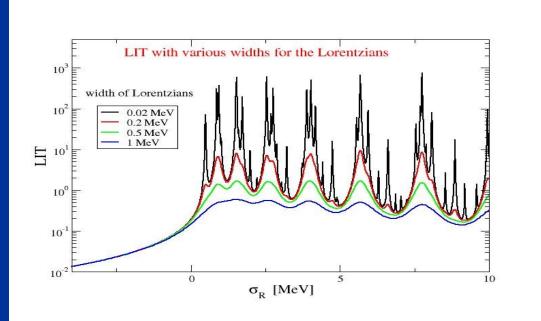


30 hyperspherical31 hyperradial

⇒ 930 basis states

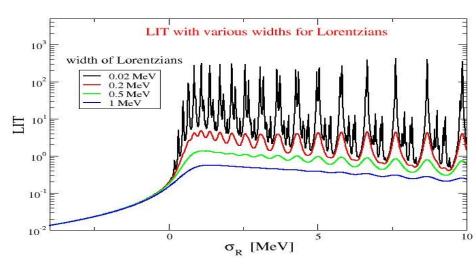
 $b = 0.6 \, fm$

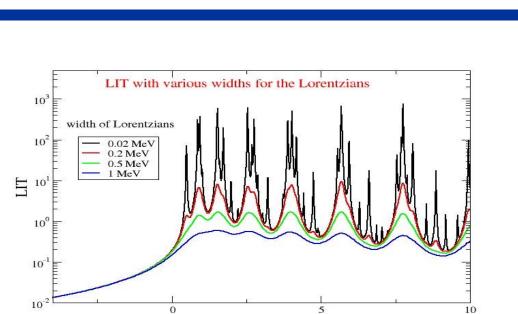




30 hyperspherical 31 hyperradial ⇒ 930 basis states b = 0.6 fm

30 hyperspherical 31 hyperradial ⇒ 930 basis states b = 0.6 fm

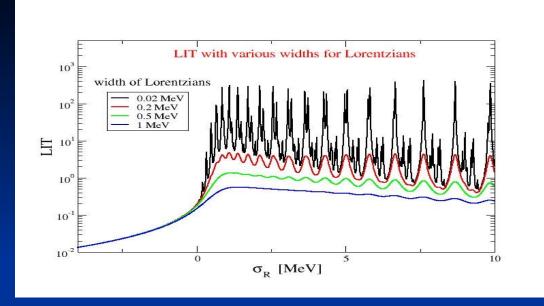


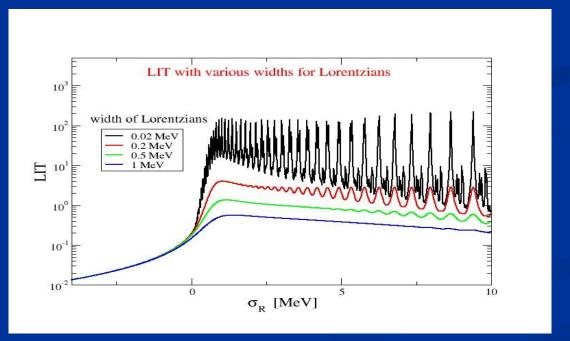


 σ_R [MeV]

40 hyperspherical 51 hyperradial ⇒ 2040 basis states b = 1 fm

30 hyperspherical 31 hyperradial ⇒ 930 basis states b = 0.6 fm





40 hyperspherical 51 hyperradial ⇒ 2040 basis states b = 1 fm

40 hyperspherical 76 hyperradial ⇒ 3040 basis states b = 2 fm

Observation

The LIT is a method with a controlled resolution

Observation

The LIT is a method with a controlled resolution

But in present LIT calculation below three-body breakup threshold not a single LIT state! Similar problem as in the previous four-body case

Observation

The LIT is a method with a controlled resolution

But in present LIT calculation below three-body breakup threshold not a single LIT state! Similar problem as in the previous four-body case

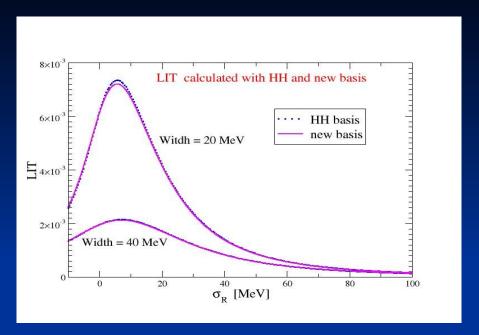
Solution: use instead of the HH basis a somewhat modified basis

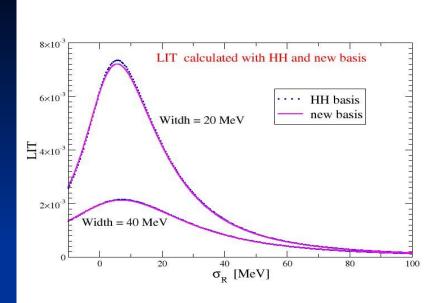
New A-body basis

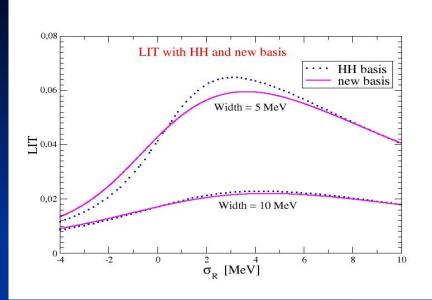
Note one of the (A-1) Jacobi vectors can be written in the following form:

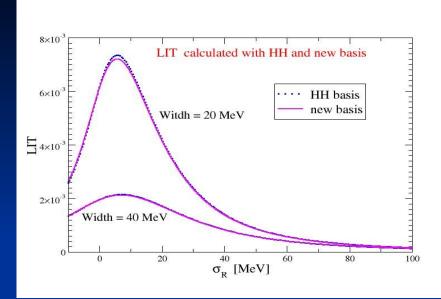
$$\eta = \mathbf{r}_{A} - \mathbf{R}_{cm}(1,2,...,A-1)$$

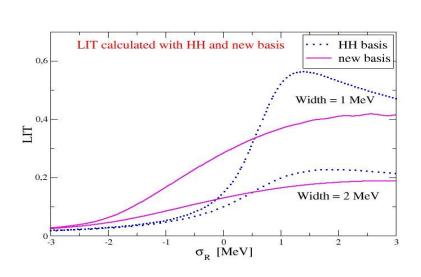
This is the coordinate one would use for the scattering of a nucleon with a (A-1)-nucleon system. In other words the relevant coordinate for a two-body breakup. Therefore

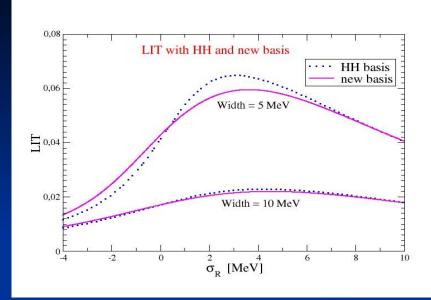


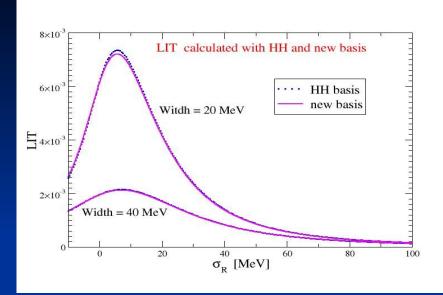


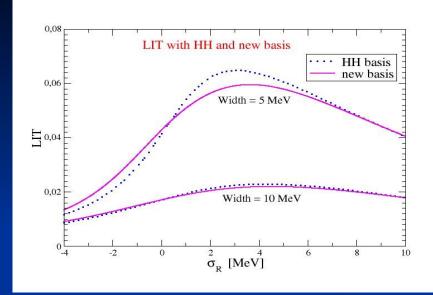


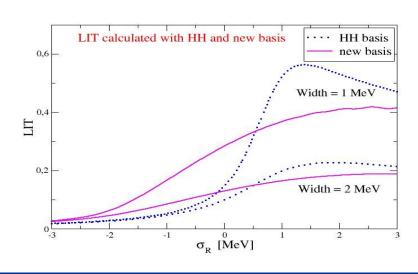


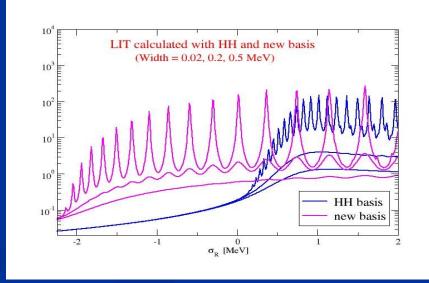








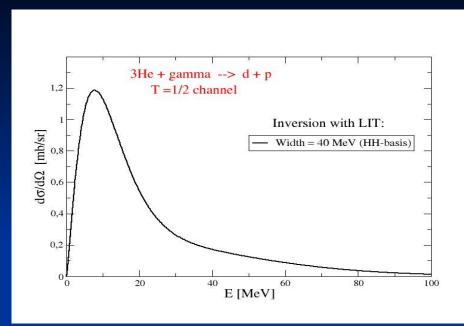


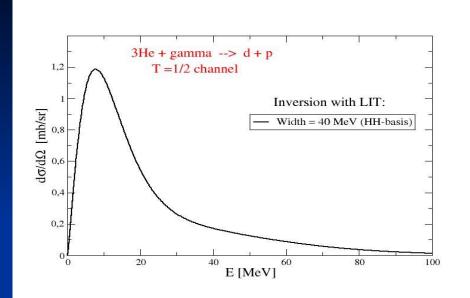


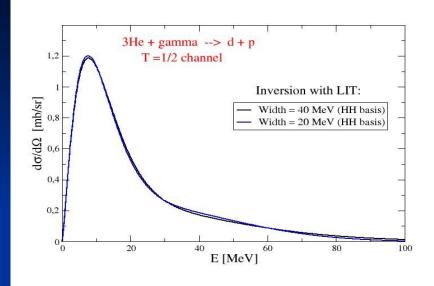
Inversions

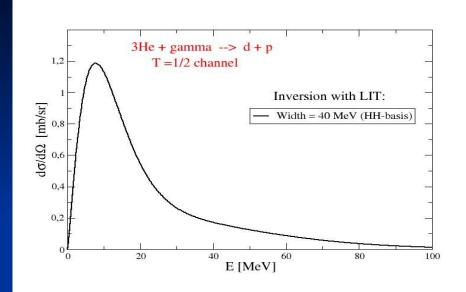
Implement correct threshold behaviour for 3 He + γ \longrightarrow d + p

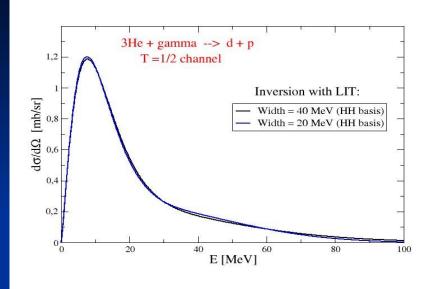
Due to Coulomb potential: usual Gamow factor

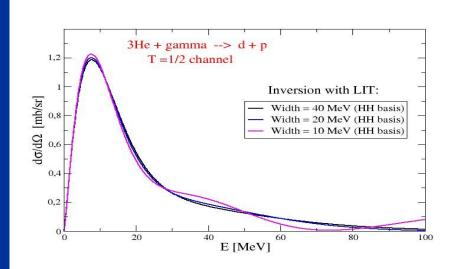


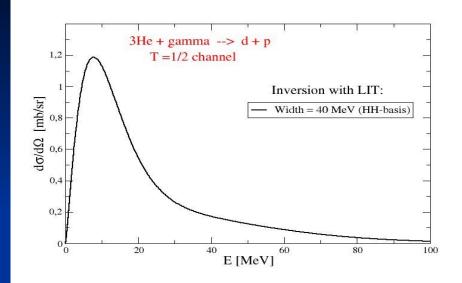


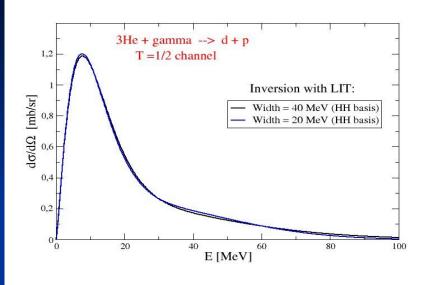


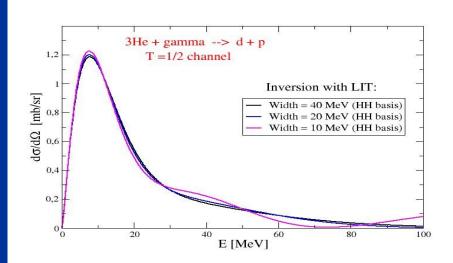


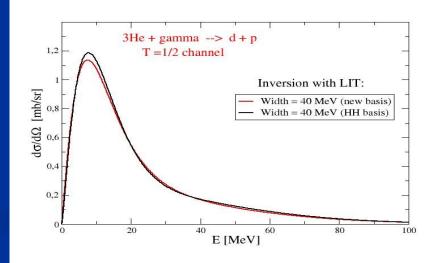


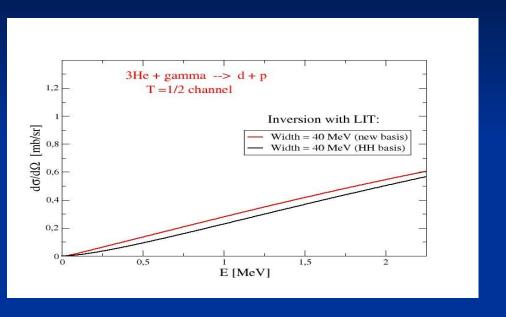


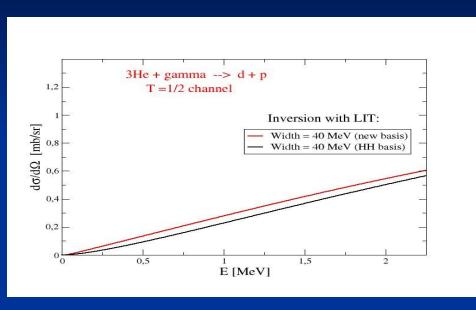


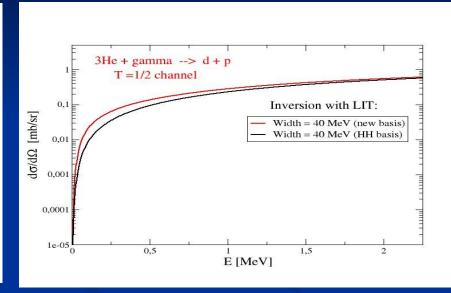


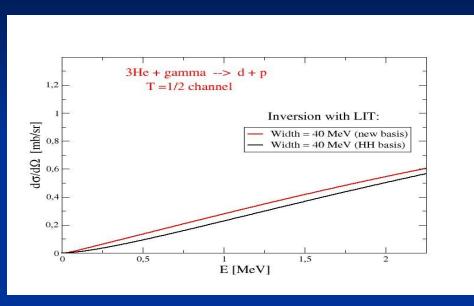


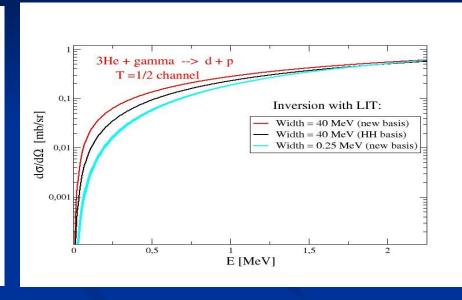


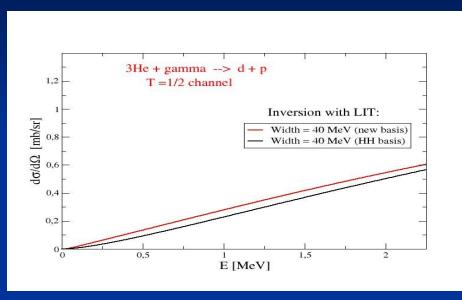


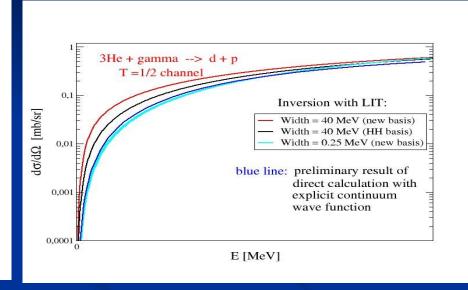




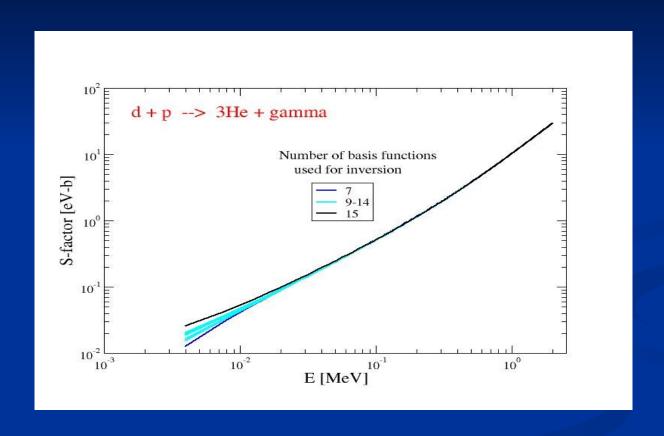






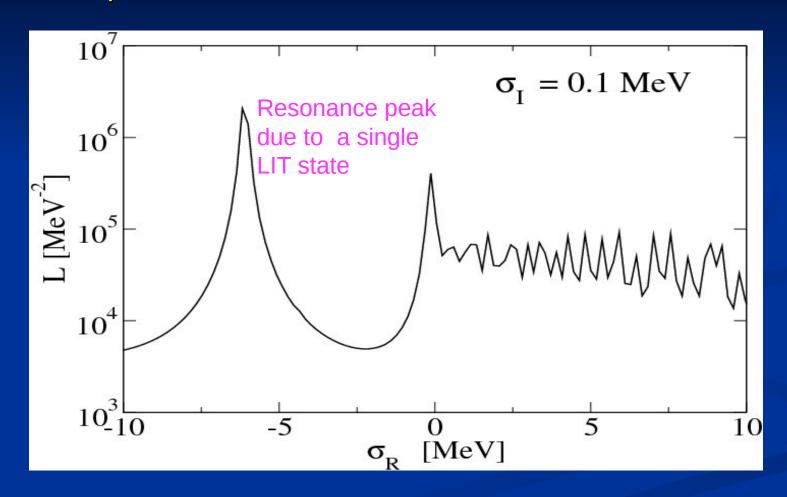


S-factor of reaction $d + p \rightarrow {}^{3}He + \gamma$

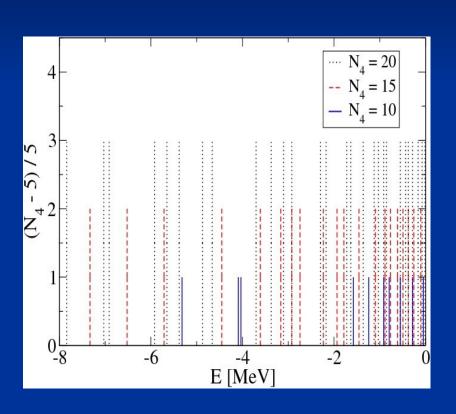


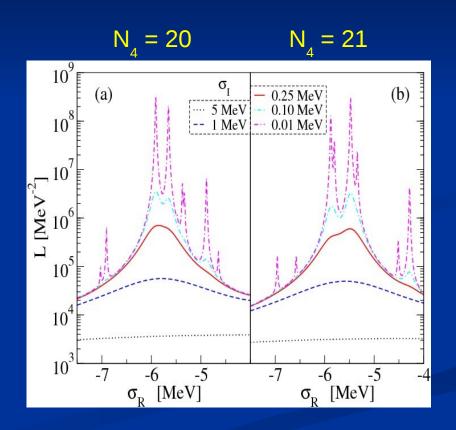
Back to the ⁴He resonance

Unpublished result from a CHH calculation with the TN potential (V. Efros, WL, G. Orlandini, PRL 78,432 (1997))



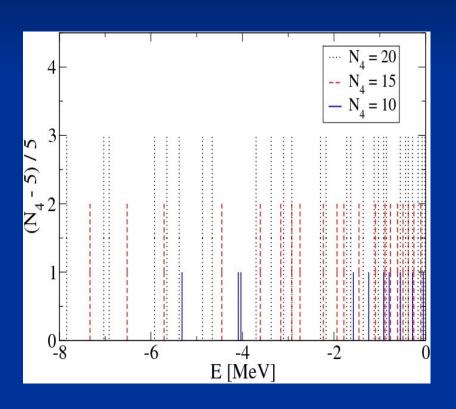
Results with new basis

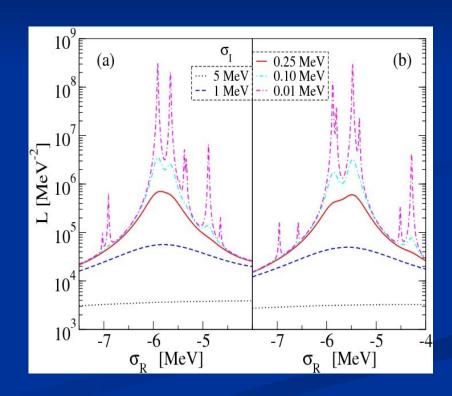






Results with new basis





Inversion: $\Gamma = 180(70)$ keV

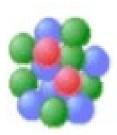
WL, PRC 91, 054001 (2015)

Benchmark calculations for hypernuclei

- Quick introduction to hypernuclei
- Short outline of our NSHH method
- Preliminary benchmark results: comparison with AFDMC (D. Lonardoni, F. Pederiva) Faddeev (A. Nogga) (GEM: E. Hiyama)

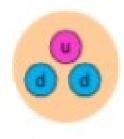
Nuclei with Strangeness

$$m_{\Lambda} = 1116 \text{ MeV}$$
 $m_{\Sigma^+} = 1189$
 $m_{\Sigma^0} = 1193$
 $m_{\Omega^-} = 1673$



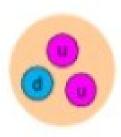
$$au_{\Lambda}=263~ps$$
 $au_{\Sigma^+}=80~ps$
 $au_{\Sigma^0}=7.4\cdot 10^{-20}s$
 $au_{\Omega^-}=82~ps$





No charge

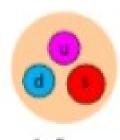
proton: 3 quarks



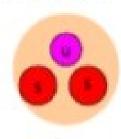
+charge

Mass: 938 MeV

hyperon: including strangeness quark



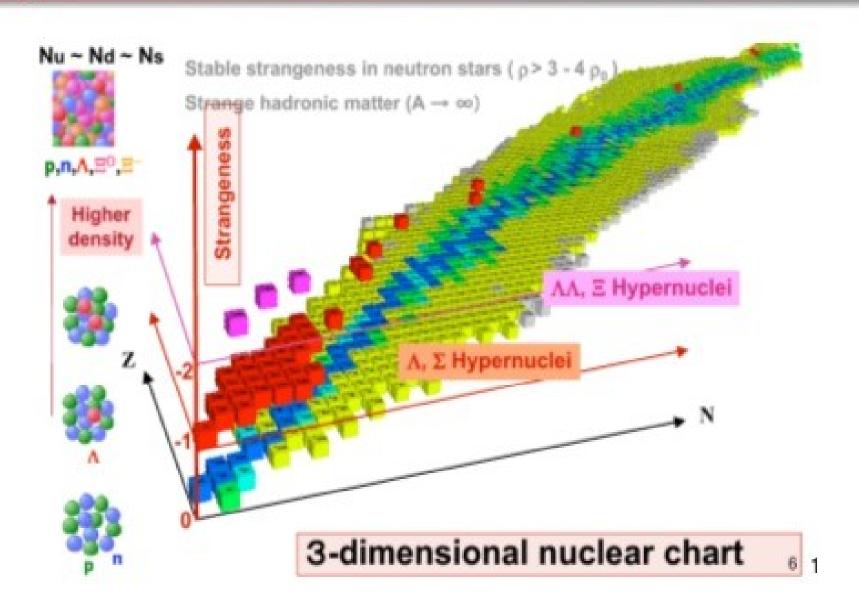






3.4

Hypernuclear Chart



Production of Hypernuclei

Strangeness exchange reaction

Associated production reaction

reaction
$$n = \begin{cases} u \\ \bar{d} \\ \bar{s} \end{cases} K^{-1}$$

$$n = \begin{cases} d \\ d \\ u \end{cases} \dots \dots \begin{cases} d \\ d \\ u \end{cases} X^{-1}$$

$$e = \begin{cases} u \\ \bar{s} \\ \bar{s} \end{cases} X^{-1}$$

Experimental Present and Future Perspectives

- Despite exensive investigations, single Λ hypernuclei knowledge is far from that of ordinary nuclei;
- Only one bound Σ-hypernucleus detected!
- No ≡ hypernuclei detected (some indications of weak attraction);
- No experimental information about Ω hypernuclei;
- Four ΛΛ-hypernuclei energies measured (⁶_{ΛΛ}He, ¹⁰_{ΛΛ}Be, ¹²_{ΛΛ}Be, ¹³_{ΛΛ}B);

Non-Symmetrized HH method

Problem: selection of **antisymmetric** states (we deal with fermions):

 \Rightarrow We add to \hat{H} the **Casimir operator** of the **permutation** group S_N , which selects "by himself" the interesting states:

$$\hat{H}' = \hat{H} + \gamma \hat{C}(A)$$
 ; $\hat{C}(A) = \sum_{i>j} \hat{P}_{ij}$

Its action on the vectors:

$$\begin{split} \hat{C}(A)\Psi_{\mathcal{S}} &= \frac{A(A-1)}{2}\Psi_{\mathcal{S}} = \lambda_{\mathcal{S}}\Psi_{\mathcal{S}}\;;\\ \hat{C}(A)\Psi_{m} &= \lambda_{m}\Psi_{m}\;;\\ \hat{C}(A)\Psi_{a} &= -\frac{A(A-1)}{2}\Psi_{a} = \lambda_{a}\Psi_{a}\;, \end{split}$$

 \Rightarrow with a proper choice of γ the g.s. energy $\mathbf{E}_{\mathcal{A}}^{0}$ becomes the **lowest eigenvalue of H**' (similar procedure for exited states).

HYP-NSHH: different particles

Hypernuclei are systems made of **two different species** of particles.

- Different masses:
 - mass weighted coordinates ⇒ dependence inside transpositions:

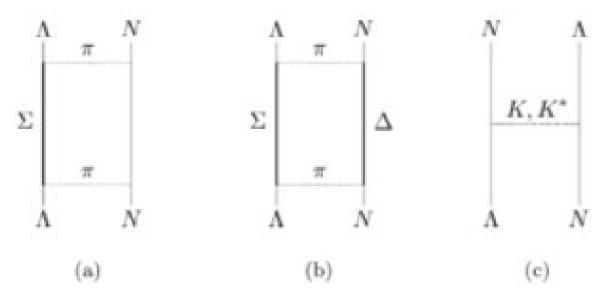
$$\mathcal{B}_{ij} \rightarrow \mathcal{B}_{ij}(m_1, m_i, m_j)$$
;

- relative coordinate rescaling mass dependence inside potential.
- Assuming 2-body potential: 3 types of different interactions:
 - NN nuclear core;
 - YN nucleon-hyperon couples;
 - YY hyperonic part.

2-body Bodmer Usmani interaction

The Λ particle has T=0 so there is no OPE term:

$$v_{\Lambda N}(r) = v_0(r) + \frac{v_{\sigma}}{4} T_{\pi}^2(r) \sigma_{\Lambda} \cdot \sigma_{N}$$



$$v_0(r) = \frac{W_c}{1 + e^{\frac{r - \bar{r}}{2}}} - \bar{v} T_{\pi}^2(r)$$

$$T_{\pi}(r) = \left[1 + \frac{3}{\mu_{\pi} r} + \frac{3}{(\mu_{\pi} r)^2}\right] \frac{e^{-\mu_{\pi} r}}{\mu_{\pi} r} (1 - e^{-cr^2})^2$$

Bodmer Usmani benchmark results

$V_{NN} + V_{YN}$	System	AFDMC	NSHH	FY
AV4'	² H	-2.245(15)	[-2.245(1)]	-2.245(1)
AV4'+U	³ H	-2.45(5)	-2.529(1)	-2.537(1)
	\hat{B}_{Λ}	0.21(5)	0.284(1)	0.292(1)
AV4'	³ H	-8.92(5)	-8.983(7)	
AV4'+U	⁴ ΛH	-11.95(5)	-12.023(3)	
	\hat{B}_{Λ}	3.03(7)	3.039(8)	
AV4'	⁴ H	-32.85(5)	-32.695(6)	
AV4'+U	5 ΛH	-39.50(5)	-39.543(10)	
	\hat{B}_{Λ}	6.65(7)	6.848(12)	

NSC97f realistic interaction

We employed the NSC97f realisitc potential³ which simulates the Nijmegen scattering phase shifts:

$$\begin{split} {}^{S}V_{NY-NY'}(r) = & \sum_{i} \left({}^{S}V_{NY-NY'}^{C} \ e^{-(r/\beta_{i})^{2}} \right. \\ & + \left. {}^{S}V_{NY-NY'}^{T} \ S_{12} \ e^{-(r/\beta_{i})^{2}} \right. \\ & + \left. {}^{S}V_{NY-NY'}^{LS} \ \mathsf{LS} \ e^{-(r/\beta_{i})^{2}} \right) \end{split}$$

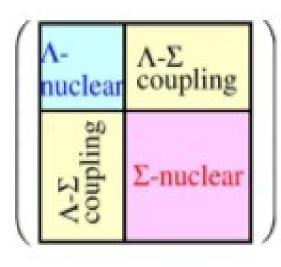
- Y→ Λ, Σ;
- C → central, T → tensor, LS → spin-orbit;
- gaussian radial functions with fitted paramenters.

Explicit use of Σ degree of freedom \Rightarrow need for extension of the HYP-NSHH method.

³E.Hiyama et al., Th.A.Rijken, Phys. Rev. C89, 061302 (2014).

Lambda-Sigma mixing

 \Rightarrow extension of the basis including Λ/Σ degree of freedom:



 definition of transformation between two Jacobi sets differing by one mass;

 extension of Lee-Suzuki procedure including Λ/Σ degree of freedom.

NSC97f interaction benchmark

$V_{NN} + V_{YN}$	System	NSHH	FY	GEM
AV8'	² H	[-2.226(1)]	-2.226(1)	
AV8'+NSC97f	³ H	-2.41(2)	-2.415(1)	
	\hat{B}_{Λ}	0.19(2)	0.189(1)	0.19(1)
AV8'	³ H	-7.76(0)		
AV8'+NSC97f	⁴ ΛH	-10.05(7)		
	\hat{B}_{Λ}	2.29(7)		2.33

Results were obtained in collaboration with

- S. Deflorian and F. Ferrari-Rufino (Trento PhD students)
- N. Barnea (Jerusalem), V. Efros (Moscow), G. Orlandini (Trento)