

Ab-Initio Study of Antiproton-Nucleus System at Low Energies

Alireza Dehghani

Supervisor: Dr. Guillaume Hupin

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IJCLab, CNRS/IN2P3 & Université Paris-Saclay, Orsay, France



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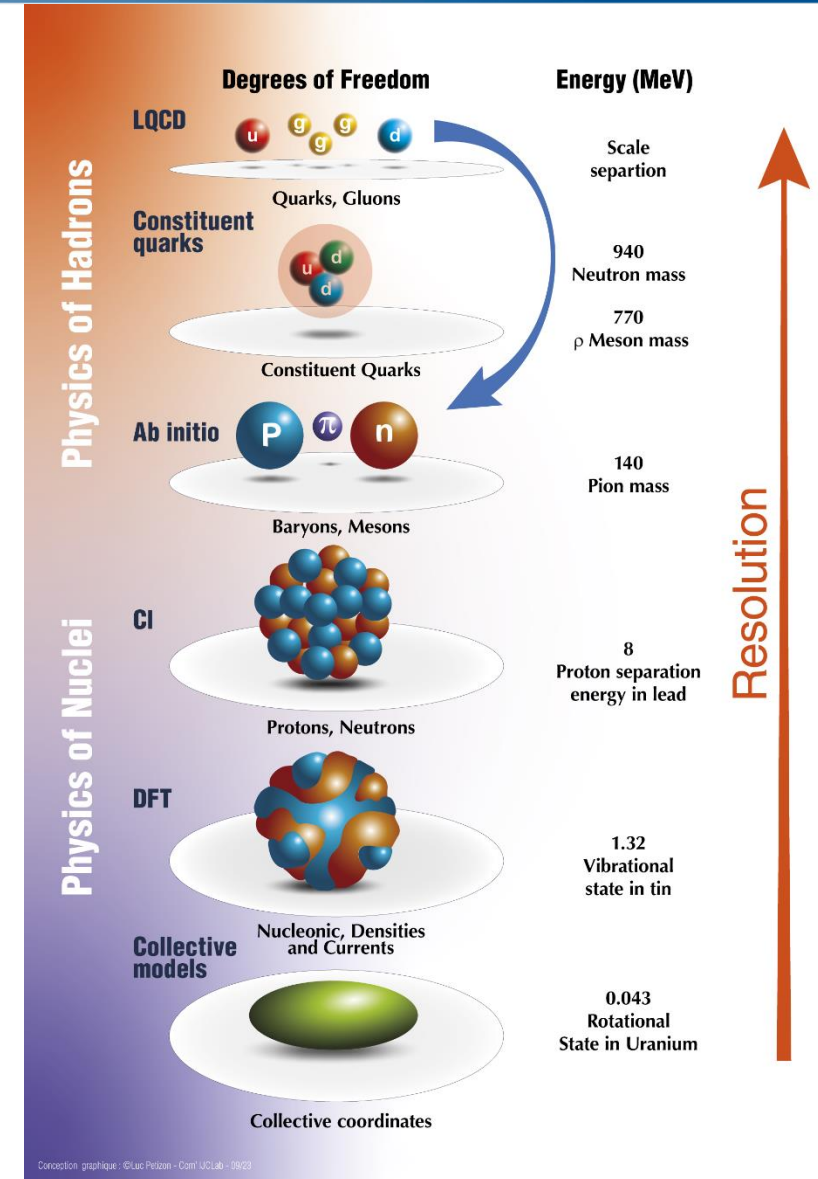
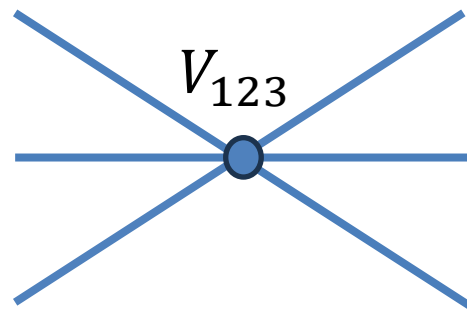
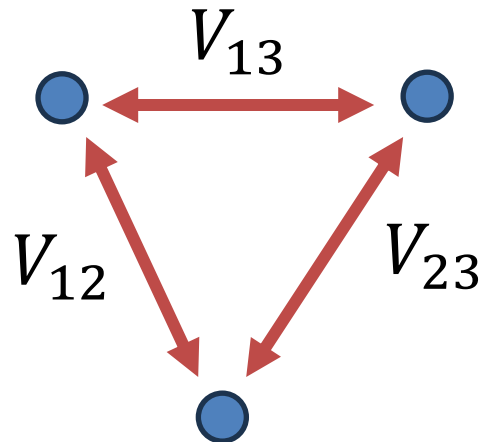




- **Goal:** Solving the Schrodinger equation (SE) for an A-body system:

$$H|\psi^{J^{\pi T}}\rangle = E|\psi^{J^{\pi T}}\rangle$$

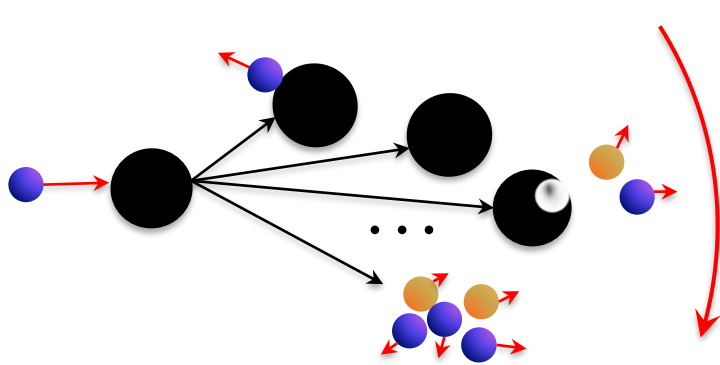
- Nucleons are considered as point-like particles.
- The SE is solved by considering two and many-body interactions between nucleons



Conception graphique: © Luc Pelton - Com UCLab - 09/23

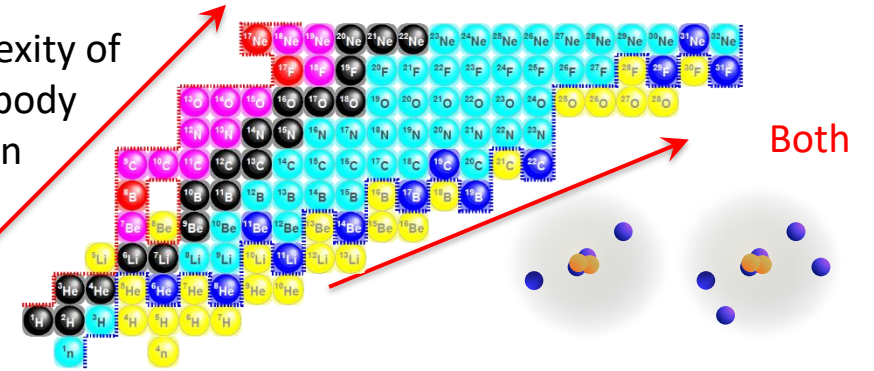


Research axis: Complex reactions with s- and p-shell nuclei



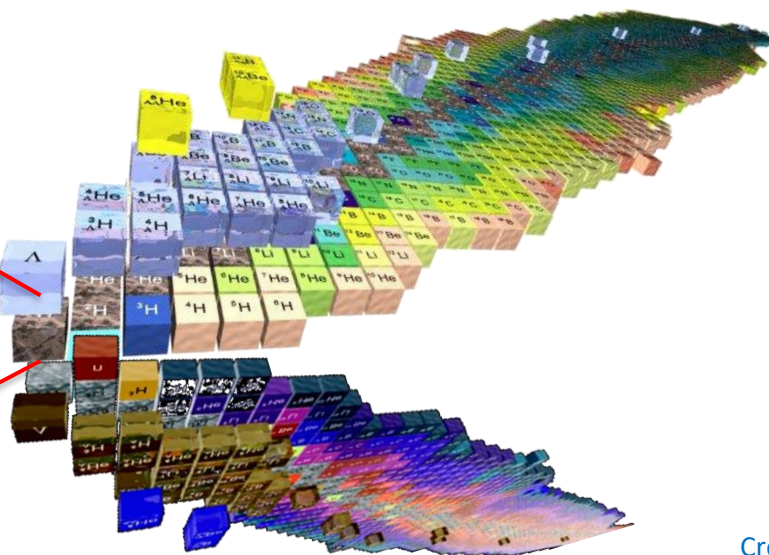
Complexity of scattering problem

Complexity of many-body solution



$\neq n, p$ particles interacting with strong force ($M_h \gg M_{n,p}$)

$$M_h \leq M_{n,p}$$



- ☹ Nuclear theory is **data driven**.
- ☹ Global optical models (NN or $N\bar{p}$ or $N\pi$ scattering) are not applicable to exotic systems.

Credits H. Lenske



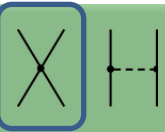
Building block - Chiral EFTs

Weinberg '90'91'92 Ordóñez and vK '92 etc...

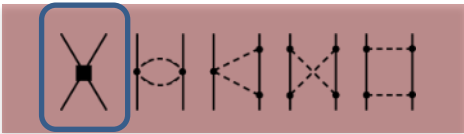
Importance



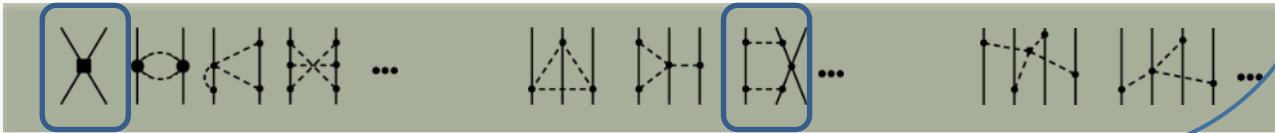
$\mathcal{O}(1)$



$\mathcal{O}(Q/M_{hi})$

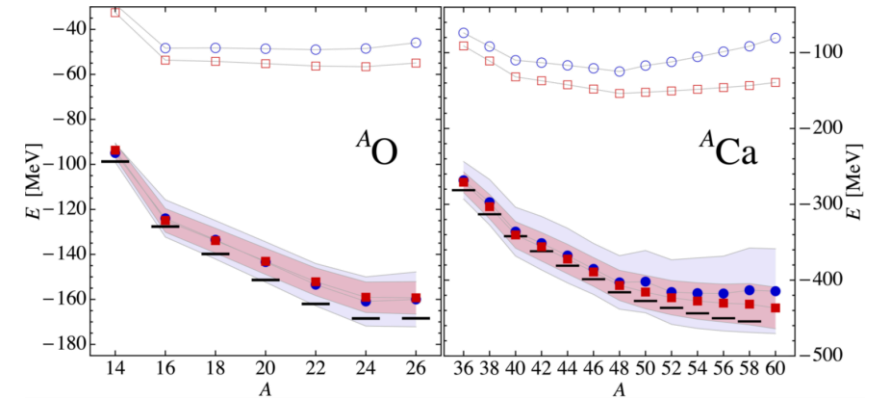


Friar '97



1. Development of improved power counting → high-precision/high-orders.
2. Stability of p-shell nuclei.

3. How does chiral expansion perform under forthcoming LQCD data?

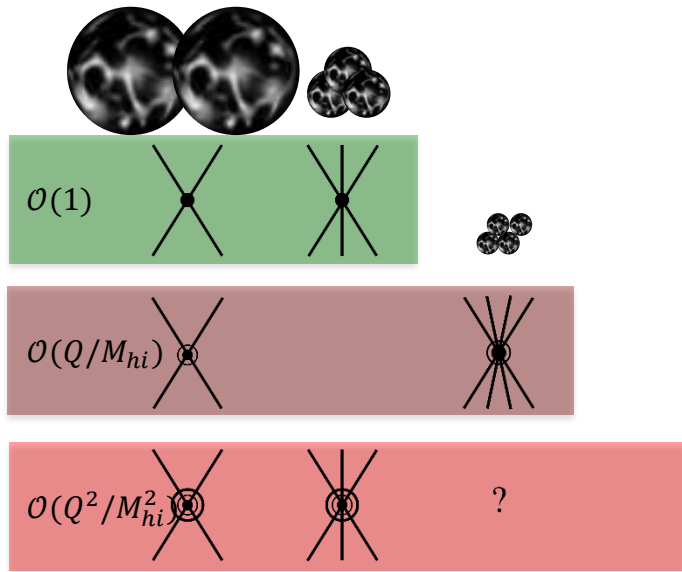


A. Tichai, et. al. Front. Phys. 8, 164 (2020)

- ☺ **High quality** nuclear interactions (at N³LO).
- ☹ **Various** fits and successes.
- ☹ Weinberg PC wrong: no **renormalizability**.
- ☺ Correct power counting: active research



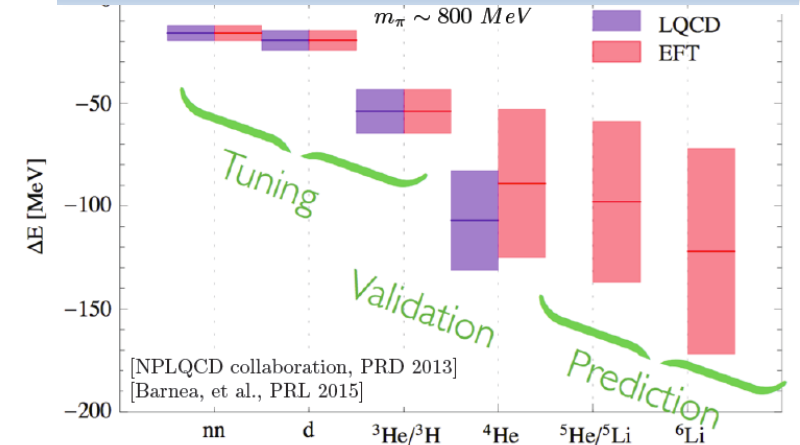
Importance



H.-W. Hammer, S. König, and U. van Kolck Rev. Mod. Phys. **92**, 025004 (2020)

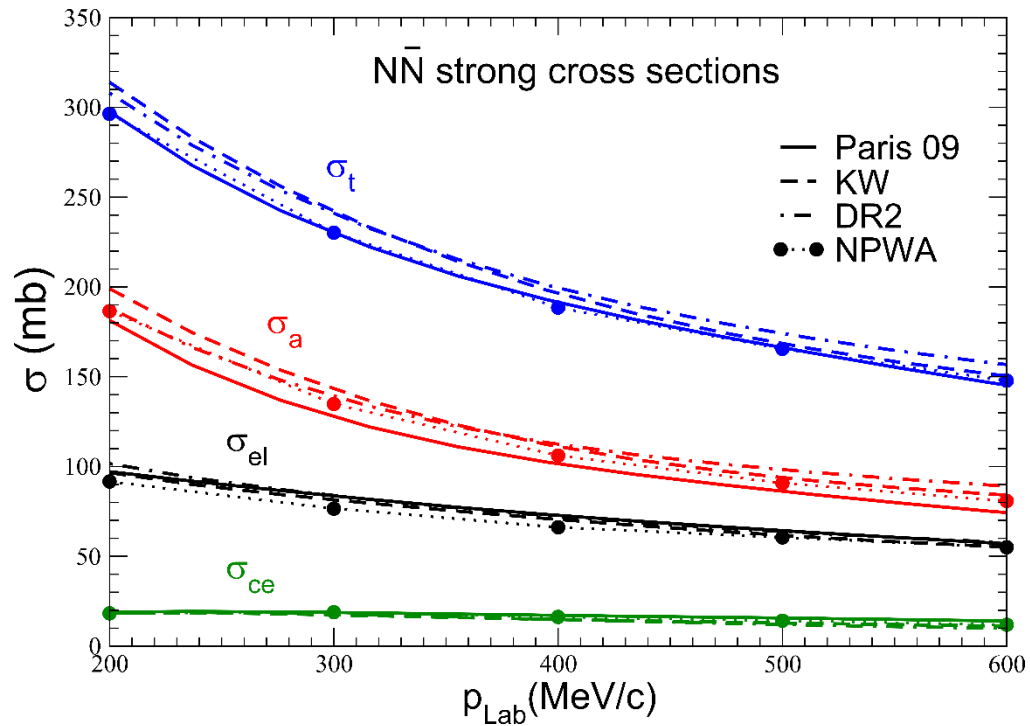
1. High-precision pionless EFT for *ab initio* nuclear studies.
2. Stability of p-shell nuclei (NLO).
3. Pionless EFT in the many-body sector: mass reach.

From LQCD to nuclear \neq EFT



☺ **Renormalizable**, connection to LQCD.

4. Pionless EFT for antinucleon systems / hypernuclei ?

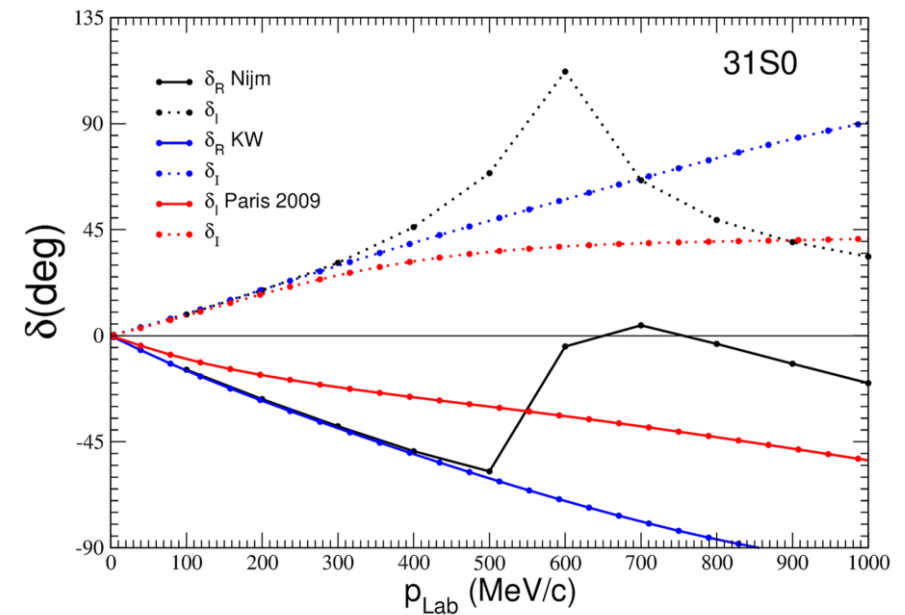
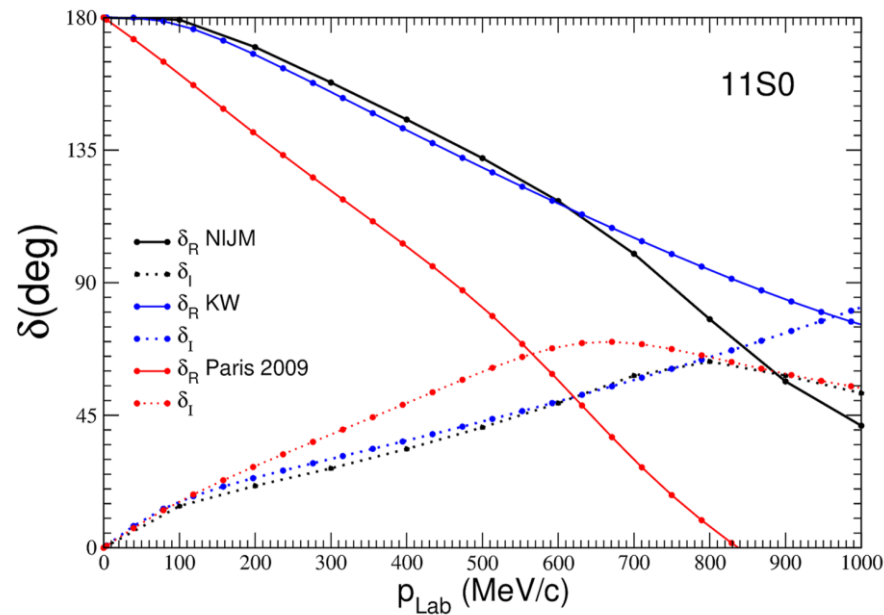


Despite significant disparities, old models, evaluation (PWA), and Chiral EFT parametrization yield **similar agreements in integrated cross-sections.**

$N\bar{N}$ strong integral cross sections for DR2 (dashed dotted line), KW (dashed line) and Paris 2009 (solid line) optical models, and the Nijmegen Partial Wave analysis.



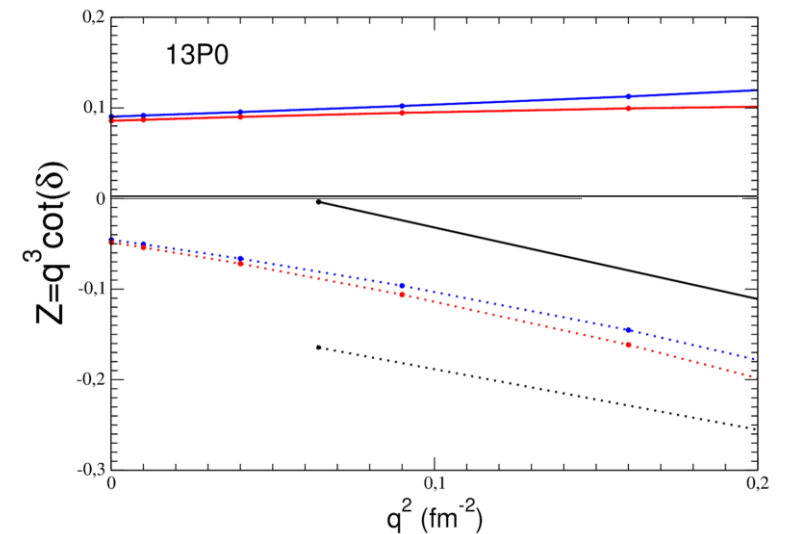
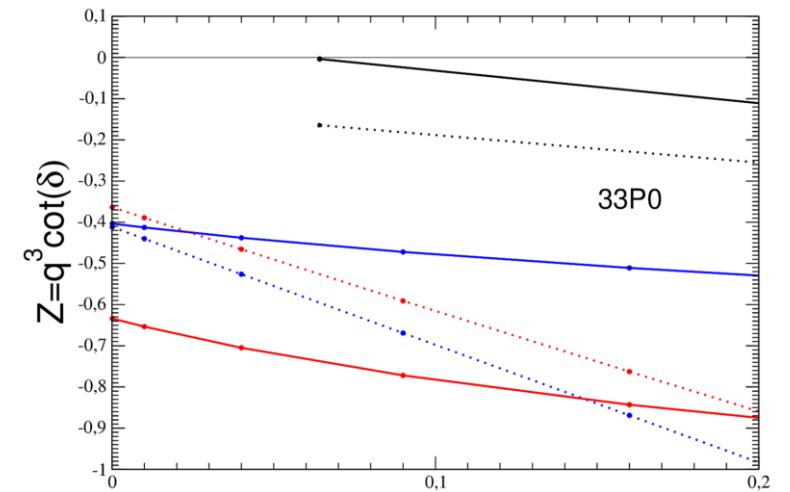
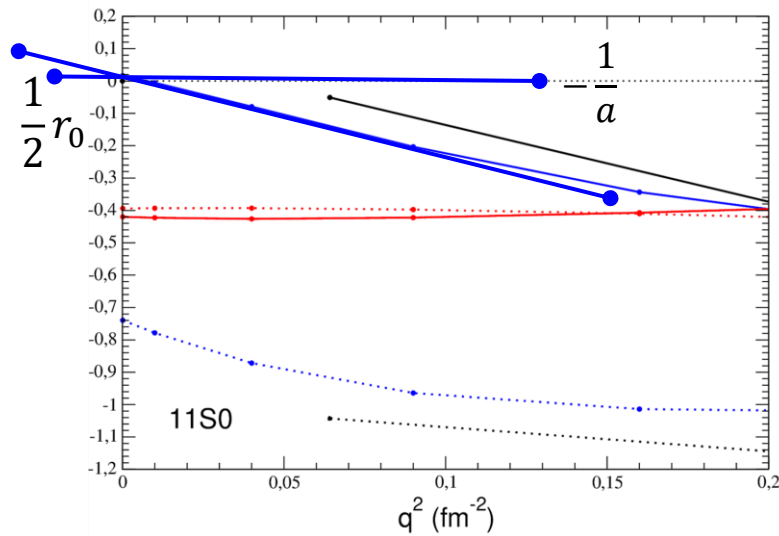
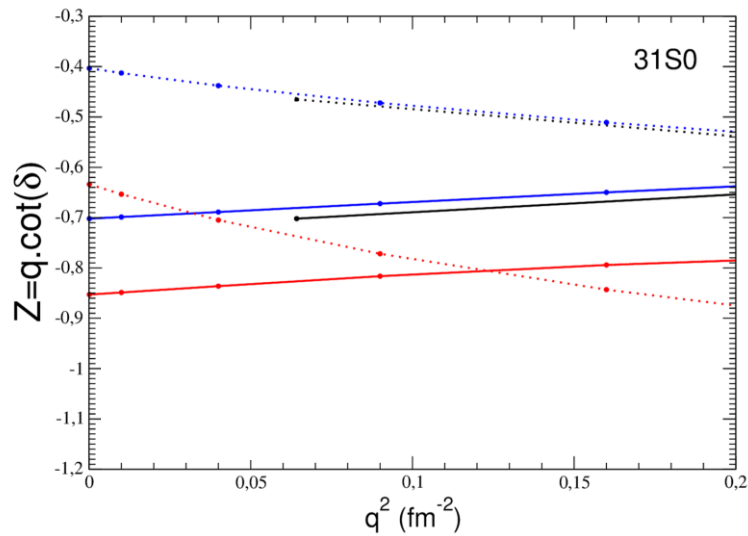
- Old models and evaluations, which were primarily fitted to intermediate energy data, may display significant discrepancies in partial wave content.
- Need for low-energy measurements ($d\sigma/d\Omega$, A_y etc ...)



$N\bar{N}$ phase-shifts comparison between two models (Paris 2009 and Kohn-Weise) and Partial Wave Analysis (PWA) of Nijmegen (data evaluation).



$N\bar{N}$: low-energy limit



- Z_r NIJM
- Z_I
- Z_r KW
- Z_I KW
- Z_r Paris 2009
- Z_I

Problems:

- $N\bar{N}$ low-energy parameters (scattering length a and effective range r_0) are weakly constrained by the present data.
- Nijmegen evaluation do not extrapolate to potential models [unlike in NN case].



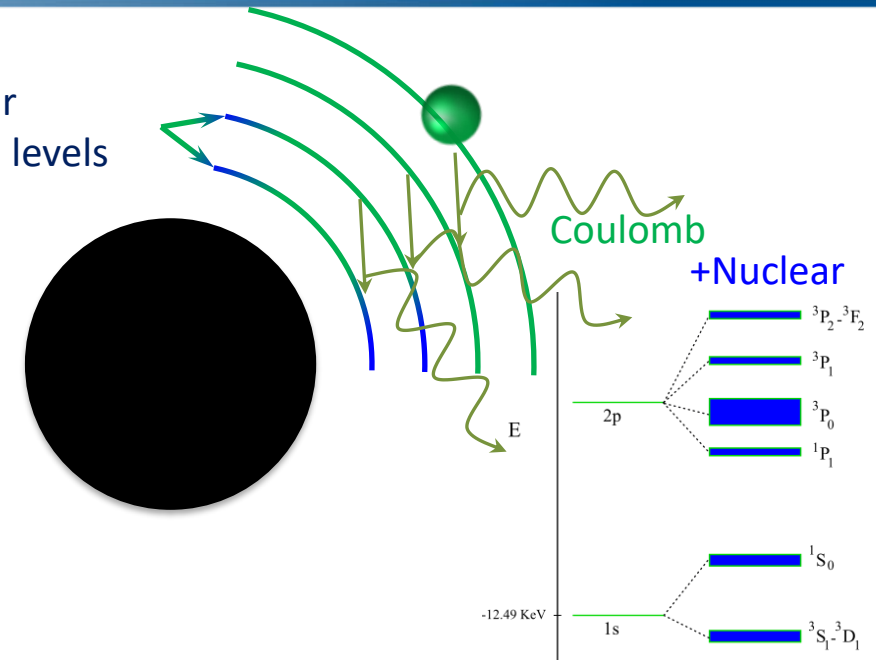
$N\bar{N}$: Effective Range Expansion parameters and protonium data

- Qualitative agreement for some waves but accuracy do not exceed 10%
- Low-energy theories need high accuracy at and below threshold

$T=0$	$^{11}P_1$		$^{13}P_0$		$^{13}P_1$		3PF_2	
Nijm*	-3.34 - 1.22i	9.3 - 1.2i	-3.06 - 7.23i	-1.7 - 1.5i	4.36 - 0.00i	-3.5 - 0.0i	--	--
Jülich	-2.87 - 0.36i	--	-2.83 - 7.82i	--	4.61 - 0.05i	--	-0.74 - 1.13i	--
Paris 09	-3.62 - 0.34i	3.8 - 0.8i	-8.78 - 4.99i	0.23 - 1.1i	5.12 - 0.02i	-3.4 - 0.02	-0.49 - 0.87i	--
KW	-3.36 - 0.62i	3.7 - 1.6i	-8.83 - 4.45i	0.25 - 0.97i	4.73 - 0.08i	-3.5 - 0.1i	-0.46 - 1.09i	--
DR2	-3.28 - 0.78i	4.2 - 2.3i	-8.53 - 3.50i	0.63 - 1.0i	5.14 - 0.09i	-3.4 - 0.1i	-0.59 - 0.85i	--

Scattering length and effective range over the years, same data.

Nuclear shifted levels



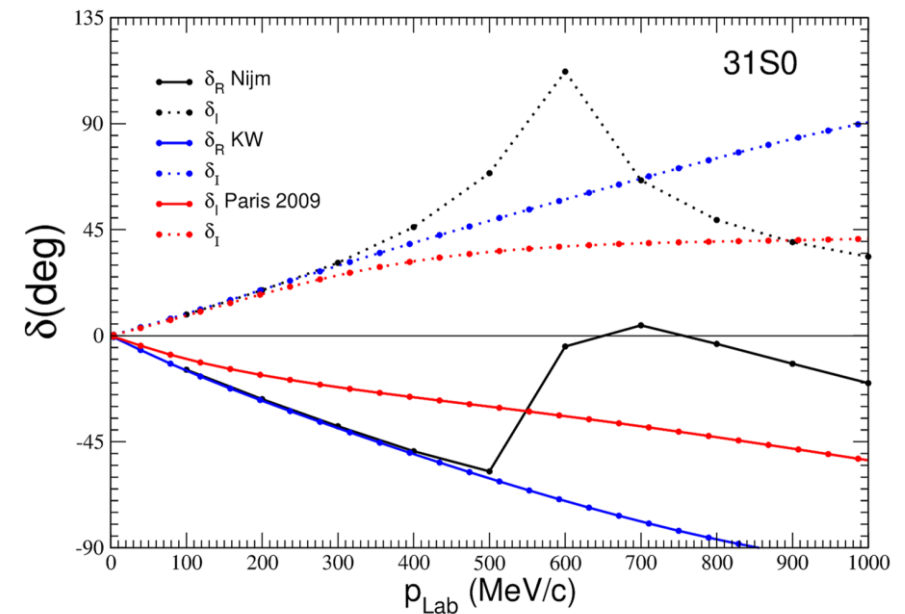
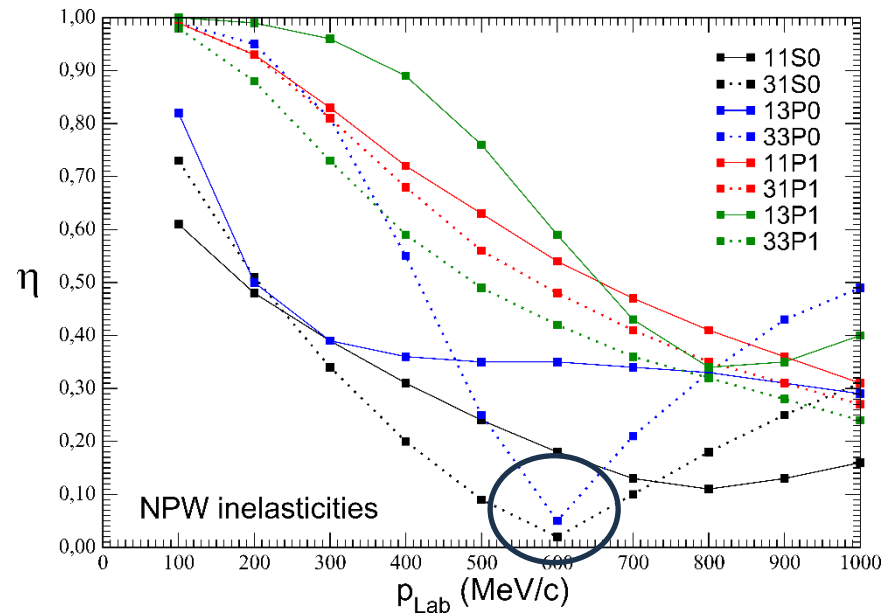
$p\bar{p}$ state	Exp	Paris 2009	Jülich	KW
$1S_0$	0.493(92) - i 0.732(146)	0.92 - i 0.67	0.50 - i 0.71	0.57 - i 0.77
3P_0	-5.68(123) - i 2.45 (49)	-2.74 - i 2.46	-0.32 - i 3.85	-2.81 - i 1.99

No agreement between average value extracted from atomic data and $N\bar{N}$



$N\bar{N}$: an example of issues to address

Fitting S-matrix components when the number data points is low, may lead to ambiguities, e.g. limit where a wave is totally absorptive and do not contribute to the elastic/c.e. cross-section. Alternative fits reproduced equally well the data (overfitting!).



$N\bar{N}$ inelasticities extracted from data (PWA).

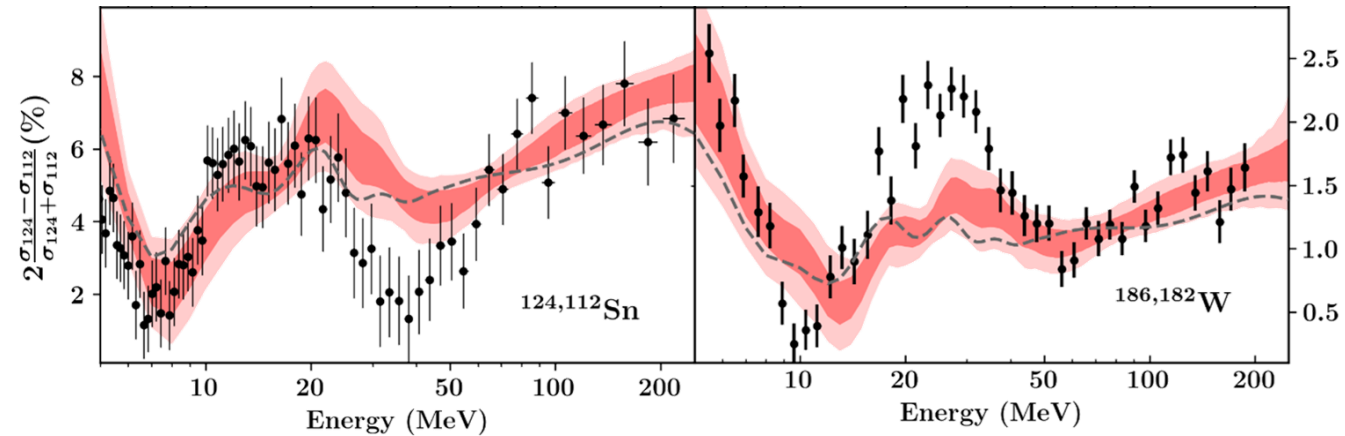
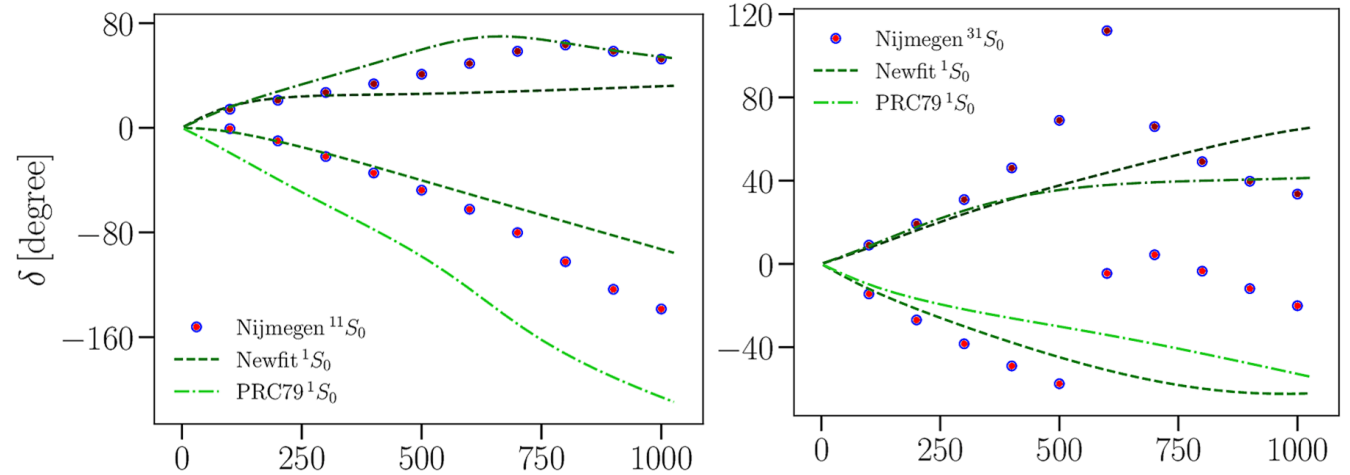


We can upgrade antediluvian $N\bar{N}$ models to match the most accurate PWA.

Yet:

1. Impossible to reproduce P-waves.
2. Unreliable extension of the partial wave analysis to low-energy.
3. Not enough information on subthreshold states.

Modern Nucleon-Nucleus optical potentials come with uncertainty quantifications, which reflect the measurements uncertainties and the limitation of the functional form of optical model



C. D. Pruitt *et al.* Phys. Rev. C **107** (2023)



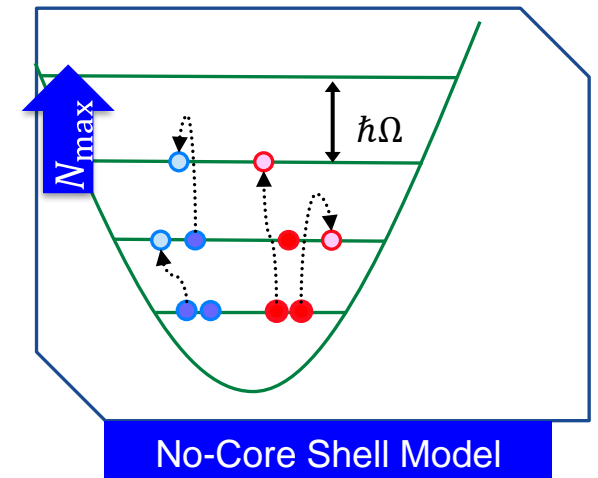
- Suitable for studying static properties of nuclei like the energy spectrum.
- In the NCSM, the wavefunction of the A-body system is expanded using A-body Harmonic oscillator (HO) basis, e. g., for A=2:

$$\Psi_{NCSM}^{(A)} = \sum_{n,l,s} c_{n,l,s} |n(l s) J^{\pi} T\rangle$$

Expansion coefficients (unknown)

Basis states

$$N_{max} = \max(2n + l)$$





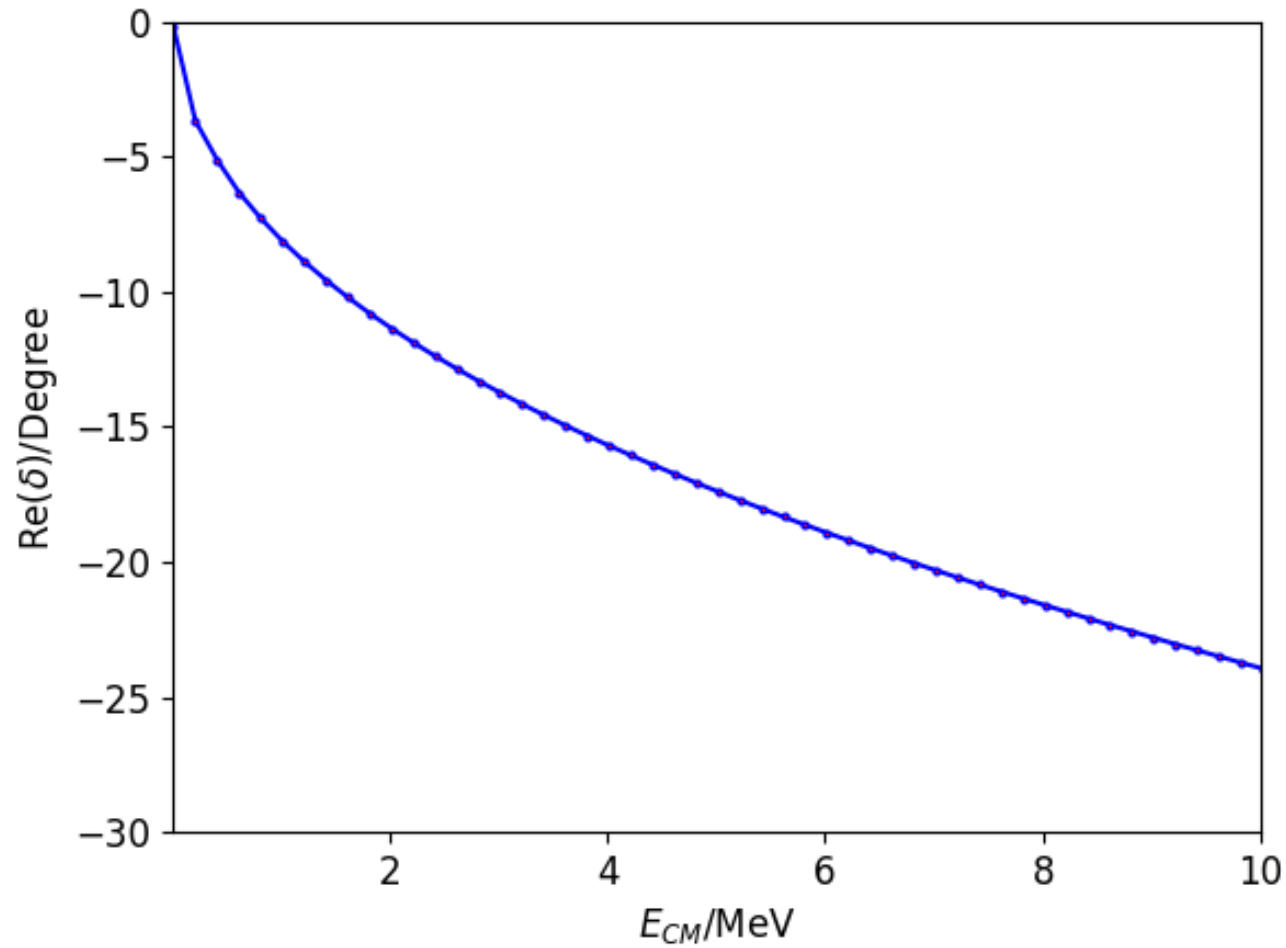
Example of No-Core Shell Model calculation: Quasi-bound states

Channel	$p\bar{n}$	$n\bar{n}$
1S_0	————	————
3SD_1	————	$-110 - 379i$
3P_0	————	————
3PF_2	————	$-242 - 345i$

Quasi-bound states (in MeV) obtained from Kohno-Weise $N\bar{N}$ potential using NCSM.
Results are obtained using $N_{\max} = 200$ and $\hbar\omega = 20$.



Example of No-Core Shell Model calculation: Scattering



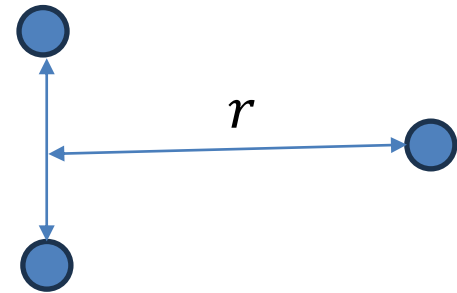
Real part of 1S_0 $\bar{p}n$ scattering phaseshift calculated using Complex Scaling method with $\theta = 20^\circ$ and $n_r = 450$.



- RGM can be used for studying nuclear reaction as well as nuclear structure.
- The A-body system is considered as two clusters, i.e., projectile and target.
- The A-body wavefunction is taken to have the following form:

$$\Psi_{RGM}^{(A)} = \sum_v \frac{g_v(r)}{r} \left| \left(\left(|\alpha \ I_1^{\pi_1} \ T_1 \rangle \right) \left(\begin{array}{cc} 1 & 1 \\ 2 & 2 \end{array} \right) s \ Y_l(\theta, \varphi) \right) J^{\pi T} \right|$$

Relative Target Projectile Relative (angular part)





The Hamiltonian (neglecting the three-body force) can be written as:

$$H = T_{rel}(r) + \bar{V}_c(r) + H_{tar} + \sum_{j=1}^{A-1} (V_{jA}^S(r_{jA}) + V_{jA}^C(r_{jA})) - \bar{V}_c(r)$$

The RGM equation then becomes:

$$\left(T_{rel}(r) + \bar{V}_c(r) - (E - E_v^{\alpha_1}) \right) \frac{g_v(r)}{r} + \sum_{v'} \int dr' W_{vv'}(r, r') \frac{g_{v'}(r')}{r'} = 0$$

$\frac{-Ze^2}{r}$ Target (binding) energy Contribution of interaction between target and projectile



For illustration, we consider a two body spinless system with a local potential:

$$(T_l + V_c + V_s - E) |u_l(E)\rangle = 0$$

We then add the Bloch operator $\mathcal{L}(B)$ to both sides:

$$(T_l + V_c + V_s - E + \mathcal{L}(B)) |u_l(E)\rangle = \mathcal{L}(B) |u_l(E)\rangle \quad , \quad \mathcal{L}(B) = \frac{\hbar^2}{2\mu} \delta(r - a) \left(\frac{d}{dr} - B \right)$$

Channel
radius

Arbitrary
parameter

For r outside the range of the potential, one can write the right hand side as:

$$\mathcal{L}(B) |u_l(E)\rangle = \mathcal{L}(B) |u_l^{ext}(E)\rangle \propto I_l(kr) + S_l(E)O_l(kr)$$



We then expand the internal wavefunction using the lagrange basis:

$$|u_l^{int}\rangle = \sum_j A_j |f_j\rangle$$

The R-Matrix (setting $B = 0$) can be defined as:

$$R_l(E) = \frac{u_l(a)}{a u_l'(a)} = \frac{u_l^{int}(a)}{a u_l^{ext'}(a)} = \frac{\hbar^2}{2\mu a} \sum_{i,j}^{n_s} f_i(a) (C^{-1})_{ij} f_j(a)$$

$$C_{ij} = \langle f_i | (T_l + V_c + V_s - E + \mathcal{L}(0)) | f_j \rangle$$

Some algebra gives:

$$S_l(E) = e^{2i\phi_l} \frac{1 - L_l^* R_l(E, 0)}{1 - L_l R_l(E, 0)}$$



$$(T_{rel}(r) + \bar{V}_c(r) - (E - E_v^{\alpha_1})) \frac{g_v(r)}{r} + \sum_{v'} \int dr' W_{vv'}(r, r') \frac{g_{v'}(r')}{r'} = 0$$

$$W_{vv'}(r, r') = \sum_{nn'}^{N_{max}} R_{n'l}(r') R_{nl}(r) \langle \phi_{v'n'}^{J^{\pi_T}} | V_{A,A-1} | \phi_{vn}^{J^{\pi_T}} \rangle \quad \left| \phi_{vn}^{J^{\pi_T}} \right\rangle \equiv \left[\left[\left(| \alpha_1 l_1^{\pi_1} T_1 \rangle \mid \frac{1}{2} \frac{1}{2} \right) s Y_l(\theta, \varphi) \right] \right]^{J^{\pi_T}} R_{nl}(r_{A,A-1})$$

Harmonic oscillator wavefunctions

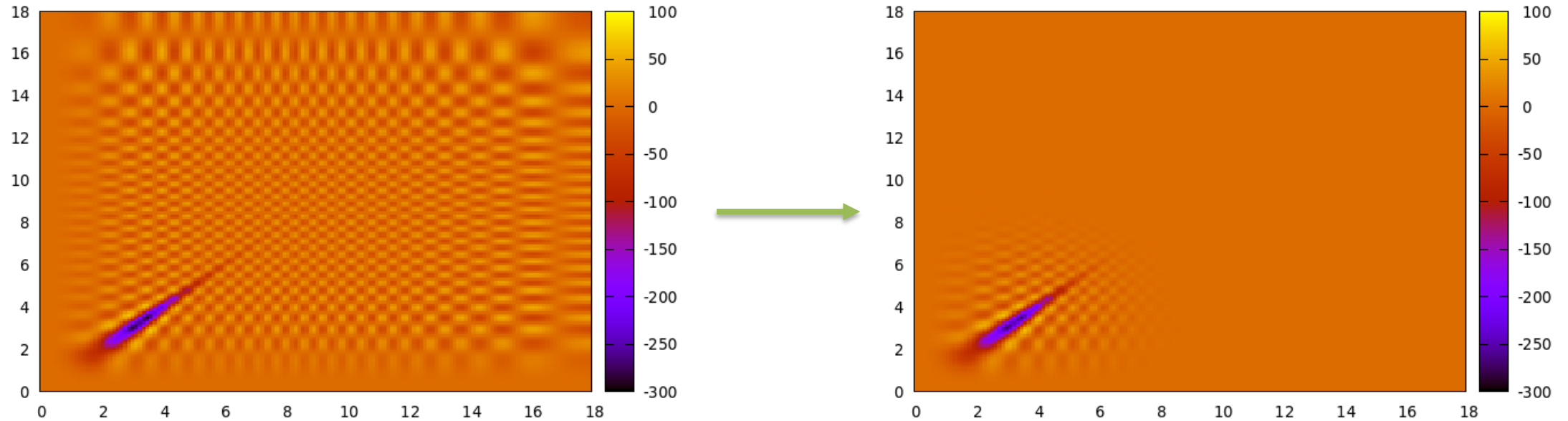
Still considerable at large n, n'

Convergence and other numerical issues



Numerical Challenges

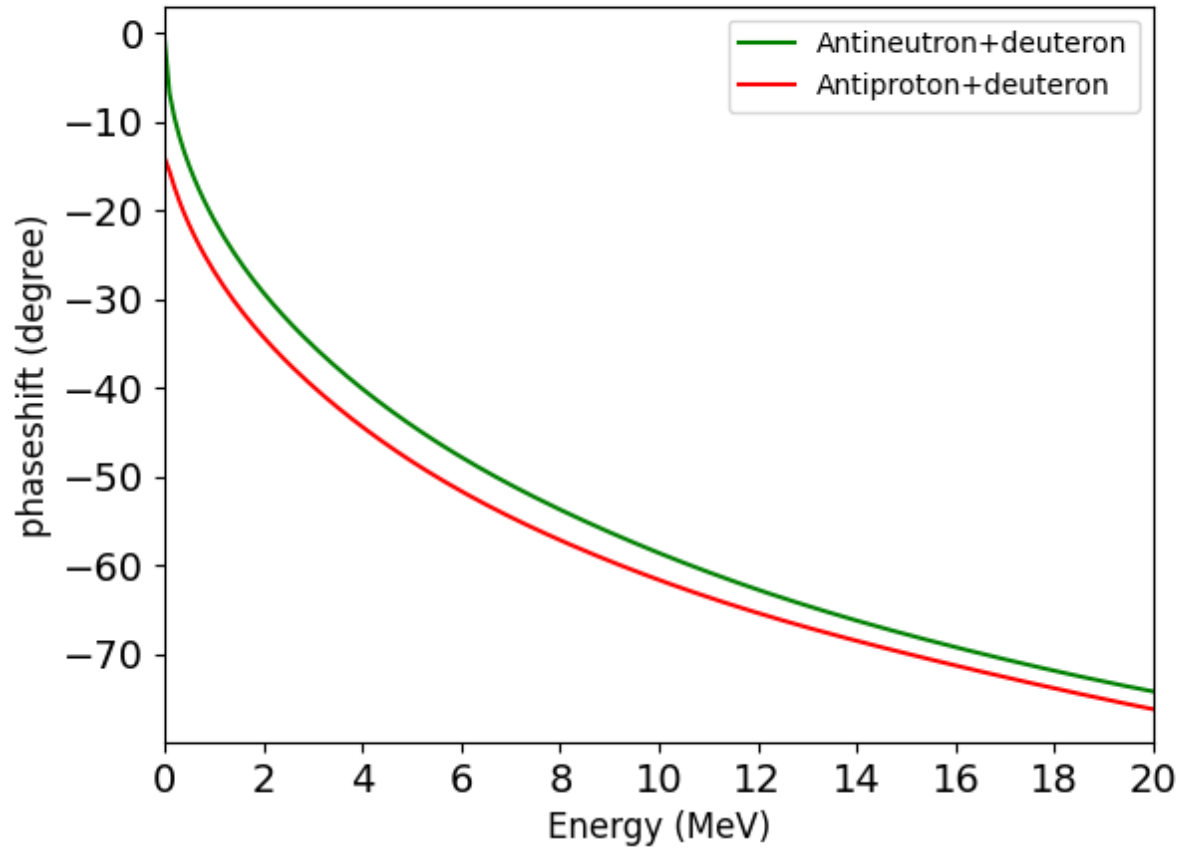
- A very high N_{max} is needed for convergence.
- The high N_{max} introduces noise in the potential:



Real part of the RGM potential for antineutron-deuteron system in ${}^2S_{1/2}$ channel before and after regularization.



Results for A=3 system (phaseshift + binding energy)

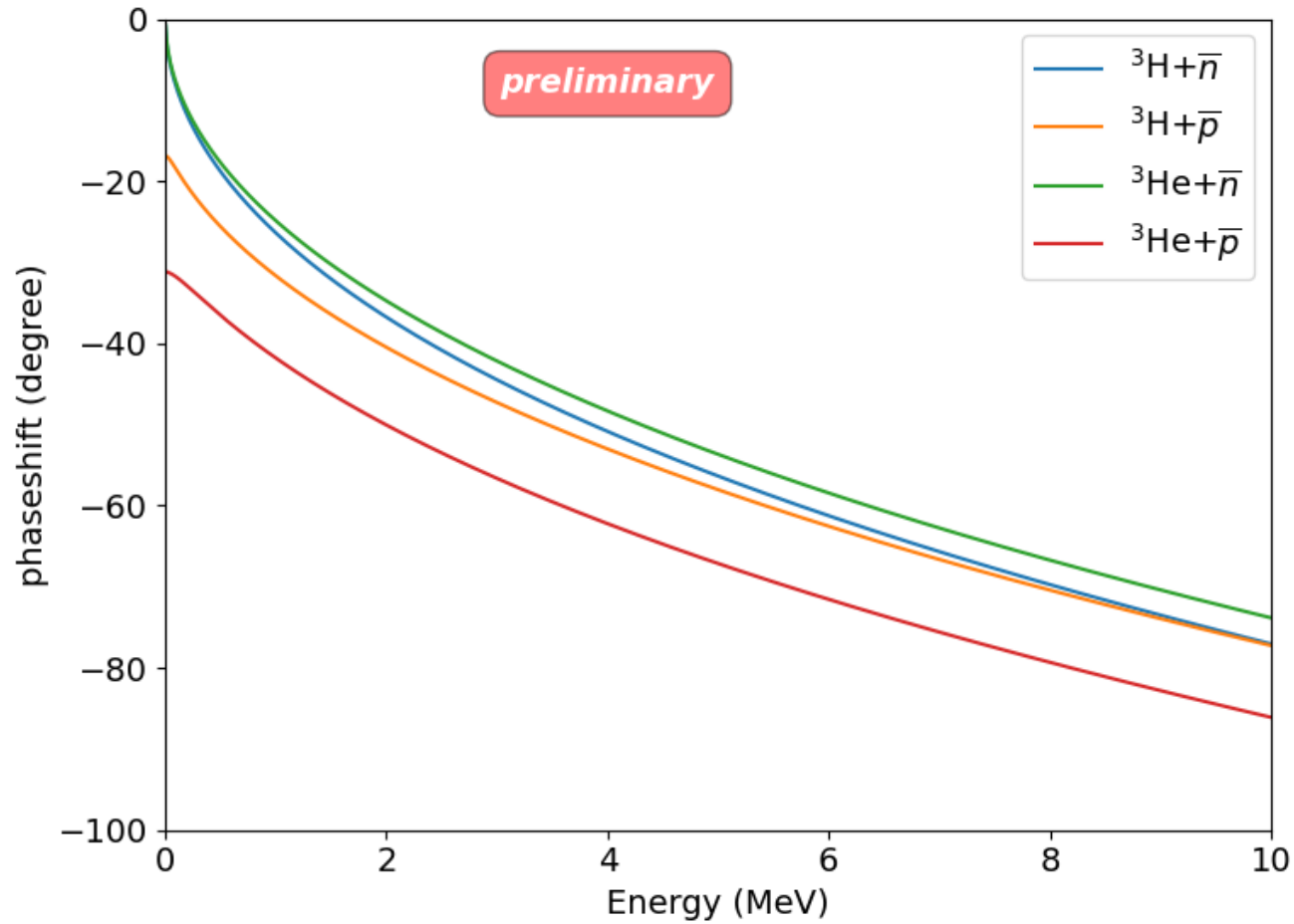


Quasi-bound state



$$E_{QB} = -21 - 153i \text{ MeV}$$

Real part of Antinucleon-deuteron phaseshift in $^2S_{1/2}$ channel with $a = 18$ fm and $N_{max} = 80$.



Real part of A=4 system phaseshift in 1S_0 channel with $a = 10$ fm and $N_{max} = 30$.



Conclusion and Outlook

- There is no consensus among current $N\bar{N}$ potentials.
- In order to develop high-quality $N\bar{N}$ potentials, one needs low-energy data from experiments.
- Currently, there are no $N\bar{N}$ observable data available at threshold.
- The advanced *ab-initio* methods designed for NN systems can be applied to $N\bar{N}$ systems.

