

TOWARDS HIGH PRECISION IN MULTI-CHANNEL REACTIONS OF COMPOSITE NUCLEI

“Always aim to convince and not primarily to impress.”

— Jürgen Habermas

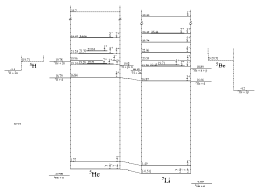
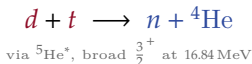
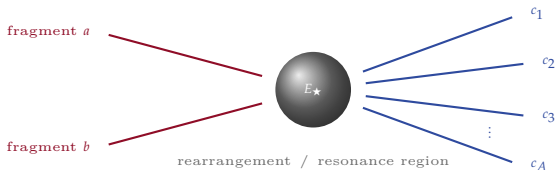
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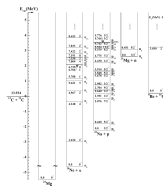
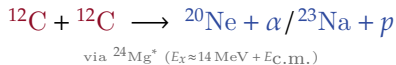
Espace de Structure et de réactions Nucléaires Théorique (ESNT)

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NCSM, RGM, GFMC, **SVM+R-matrix,&c.**

$$\Psi = \Psi_{\text{compound}} + \Psi_{\text{direct}} = \sum_{\mu} a_{\mu} \Phi_{\mu} + \mathcal{S} \left[\Phi_d \Phi_t \chi(r) \right]$$



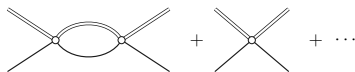
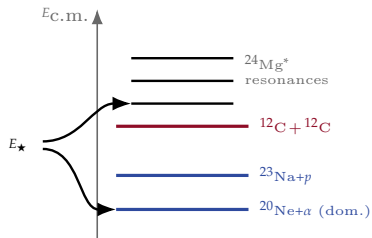
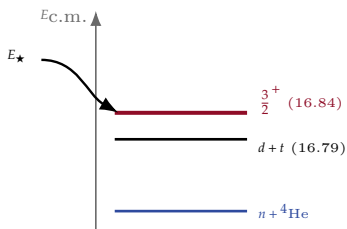
double-folding + coupled channels
(Glauber / CDCC)

$$\left[T_R + V_{cc}(R) - (E - \epsilon_c) \right] \chi_c(R) = - \sum_{c' \neq c} V_{cc'}(R) \chi_{c'}(R)$$

identical fermions: α , nucleons

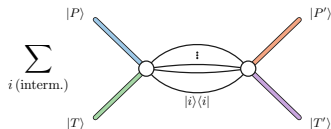
potential with long- and short-range, perturbative and non-perturbative components

defined energy interval of interest: $E_{\star}^{(\text{res/rearr})} \lesssim E_{\text{C.m.}} \ll E_{\text{exc}}^{(d,t,\alpha,N)}$

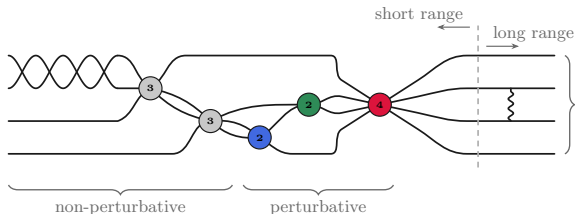


Intermediate states = the in/out fragments only
 \Rightarrow insufficient for the relevant reactions.

e.g., the $\frac{3}{2}^{+}$ resonance in ${}^5\text{He}$ is not a $d+t$ state.



In $\sum_i |i\rangle\langle i|$
 both the *interaction* **and** *solution algorithm* are active.

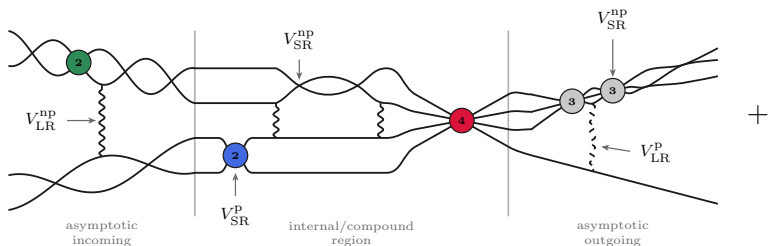


Few-body potential for non-relativistic applications

$$V = \sum_i \hat{t}_i + V_{\text{LR}}^{\text{P}} + V_{\text{LR}}^{\text{NP}} + V_{\text{SR}}^{\text{NP}} + V_{\text{SR}}^{\text{P}}$$

from a field-theoretical Hamiltonian density linking high- and low-energy scales

$$H_{\text{eff}} = \sum_i \psi_i^\dagger \left(-\frac{\nabla^2}{2m} \right) \psi_i + \tilde{C}_1 (\psi^\dagger \psi)^2 + \tilde{C}_3 (\psi^\dagger \nabla \psi) \cdot (\psi^\dagger \nabla \psi) + \dots + \tilde{D}_0 (\psi^\dagger \psi)^3 + \tilde{E}_0 (\psi^\dagger \psi)^4$$



(i) **power counting:** each vertex carries an order $v(n)$ set by its dimension & scaling

(ii) **order N :** sum *all* initial→final processes with $\sum_{\text{vertices}} v(n) \leq N$

(iii) **convergence to arbitrary accuracy:** $\Delta^{(N+1)} = \Delta^{(N)} \cdot \epsilon$, $\epsilon = \frac{\text{small scale}}{\text{large scale}} < 1$

Central challenge: Implementation

1. choose **basis-expansion** scheme \Rightarrow accurate fragment **wave functions**
(considering, at first, all iterations of all potential components)
2. **solve** the multi-channel scattering problem
(again, without regards to perturbations, but with appropriate compound structures expanded)
and map out the **functional dependence** of this solution on the strengths of all perturbative terms contributing to order N accuracy
3. adapt a **multi-dimensional power series** to this functional dependence
 \Rightarrow order- N accuracy

IMPLEMENTATION:

1. efficient representation of V

$$\begin{aligned}
 V_{\text{SR}}^{\text{n/np}}(r) &= C_1 \delta_a^{(3)}(r) + a^2 C_2 \nabla^2 \delta_a^{(3)}(r) \\
 &+ a^4 \left[C_3 \left(\frac{L(L+1)}{a^4} + \frac{2}{a^2} \left(\frac{1}{2} \nabla \right)^2 \right) \delta_a^{(3)}(r) \right. \\
 &\quad \left. + C_4 \nabla^4 \delta_a^{(3)}(r) - C_5 \nabla^2 \delta_a^{(3)}(r) \left(\frac{1}{2} \nabla \right)^2 \right]
 \end{aligned}
 \quad \leftrightarrow \quad
 \begin{aligned}
 V_{\text{SR}}^{\text{n/np}}(K, Q) &= C'_1 + C'_2 K^2 \\
 &+ C'_3 (K \times Q)^2 + C'_4 K^4 + C'_5 K^2 Q^2
 \end{aligned}$$

2. calibration of the potential

informed selection of input data to optimize convergence and minimize uncertainty already at low orders