

# Towards Faddeev-AGS equations in a Coulomb basis in momentum space

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# Focus on deuteron-induced reactions

Deuteron-induced reactions are a point of the theoretical interest  
To treat experimental results on the exotic nuclei beams with the deuterated targets in inverse kinematics.

Low energy

$E_{lab}$  from few up to 50 MeV/A.

Broad range of involved nuclei

From He isotopes up to Pb.

Main focus is on  $(d, p)$

But the tools must analyze other three-body reaction channels (e.g. deuteron breakup) on the same footing.

## $d + A$ reaction channels

### Approximated three-body:

- elastic scattering,
- direct deuteron breakup  $A(d, pn)A$  without core excitation,
- direct deuteron stripping  $(d, p)$  or  $(d, n)$  with nucleon capture without core excitation.

### Approximated three-body with excitations:

- direct deuteron breakup and stripping with core excitations,
- non-elastic direct  $(d, d')$  scattering with core excitations.

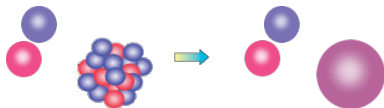
### Many-body problems:

- core breakup,
- compound-nucleus reactions...

# Deuteron-induced reactions as three-body problem

Deltuva *et al.* // Phys. Rev. C **76**, 064602 (2007)

- Neglect the internal degrees of freedom of the nucleus to get 3-body problem.
- Use Faddeev-AGS approach to treat the problem.



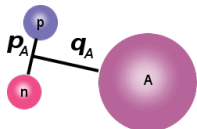
To do later:

Take into account some target degrees of freedom:

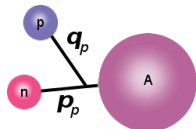
- collective core excitations,
- more complicated internal dynamics.

# Formal considerations about a 3-body system

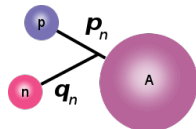
## 3-body configurations & Jacobi coordinates



$$A(pn) = d + A$$



$$p(nA)$$



$$n(Ap)$$

- Numerical indices:  $A \Leftrightarrow 1$ ,  $p \Leftrightarrow 2$ ,  $n \Leftrightarrow 3$ .
- ‘Odd man out’ configuration notation:  $a(bc)$  labeled by  $a$ .

## Denoting the two-body interactions

- $V_a \equiv V_{bc}$ , i.e.  $a$  is a spectator.

# 3-body Hamiltonian

$$H = H_0 + V_{np} + U_{nA} + v_{pA}, \quad H_0 = \mathbf{q}_a/2M_a + \mathbf{p}_a/2\mu_a.$$

Interactions between the particles

$$v_{pA} = V_{pA}^C + v_{pA}^S, \quad V_{pA}^C = \frac{Z_A \alpha^2}{r}, \quad v_{pA}^S = U_{pA} + v_{pA}^{cd},$$

- $V_{np}$ :  $NN$ -interaction potential, e.g. chiral.
- $U_{nA}, U_{pA}$ : phenomenological optical potentials, e.g. CH89.
- $v_{pA}^{cd}$ : ‘short-range Coulomb interaction’, usually, the charged sphere potential.

## Faddeev approach to a three-body problem

- $|\phi\rangle = \sum_{a=1}^3 |\phi_a\rangle [= |\phi_1\rangle]$  is initial state  $[A + (pn)]$ .

Faddeev components  $\psi_a$

$$|\Psi\rangle = |\phi\rangle + \sum_{a=1}^3 g_0^C(E) V_a^S |\Psi\rangle, \quad g_0^C(E) = (E - H_0 - V_{pA}^C + i\varepsilon)^{-1},$$

$$|\Psi\rangle = \sum_a \underbrace{\delta_{a,1} |\phi_1\rangle + g_0^C(E) V_a^S |\Psi\rangle}_{|\psi_a\rangle} = \sum_a |\psi_a\rangle.$$

Faddeev-type equations

$$\begin{cases} |\psi_1\rangle = |\phi_1\rangle + g_0^C(E) V_1^S \sum_a |\psi_a\rangle \\ |\psi_2\rangle = g_0^C(E) V_2^S \sum_a |\psi_a\rangle \\ |\psi_3\rangle = g_0^C(E) V_3^S \sum_a |\psi_a\rangle \end{cases} \Rightarrow \begin{cases} |\psi_1\rangle = |\phi_1\rangle + g_0^C(E) t_1^S \sum_{a \neq 1} |\psi_a\rangle \\ |\psi_2\rangle = g_0^C(E) t_2^S \sum_{a \neq 2} |\psi_a\rangle \\ |\psi_3\rangle = g_0^C(E) t_3^S \sum_{a \neq 3} |\psi_a\rangle \end{cases}$$

# Coulomb interaction in momentum space

Screening factor ( $g_0^C$  becomes  $g_0$ ,  $\tilde{V}^C$  goes with the other potentials)

$$\tilde{V}^C = \frac{Z_1 Z_2 \alpha^2}{r} \exp(-\mu r^n).$$

- Various short-range amplitude corrections required.



## Coulomb interaction in momentum space

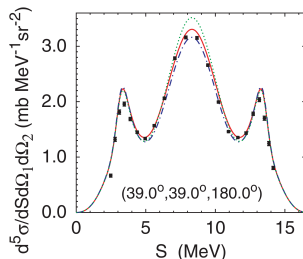
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- Various short-range amplitude corrections required.

It works for  $dp$  scattering

- Deltuva // Phys. Rev. C **80**, 064002 (2009).
- $E_{p(lab)} = 9$  MeV.
- The most advanced screening techniques to the date.
- AV18 and AN18+UIX potentials.



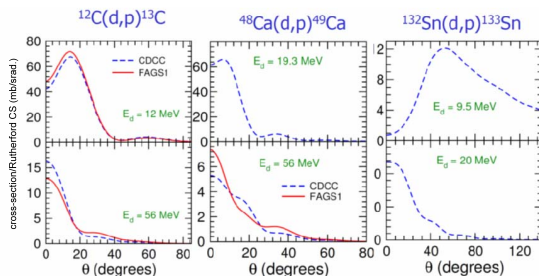
# $A(d, p)B$ reaction with screened Coulomb interaction

It works for the nuclei up to  $^{48}_{20}\text{Ca}$  (Deltuva)

\* Deltuva. // Phys. Rev. C **80**, 064002 (2009).

\* Upadhyay, Deltuva, Nunes. // Phys. Rev. C **85**, 054621 (2012).

‘The charge wall’



- Discrepancy between the results of CDCC & Faddeev calculations becomes unreasonably large as the nucleus' charge climbs up.
- Faddeev-AGS method hits the ‘charge wall’.

# Faddeev approach on the way to the highly-charged nuclei

‘Total recall’

$$\left\{ \begin{array}{l} \tilde{V}^C = \frac{Z_1 Z_2 \alpha^2}{r} \exp(-\mu r^n) \\ g_0 = [E - H_0 + i\varepsilon]^{-1} \end{array} \right. \Rightarrow \left\{ \begin{array}{l} V^C = \frac{Z_1 Z_2 \alpha^2}{r} \\ g_0^C = [E - H_0 + i\varepsilon - V_{pA}^C]^{-1} \end{array} \right.,$$

$$(H_0 = \mathbf{q}_a/2M_a + \mathbf{p}_a/2\mu_a).$$

In terms of the  $g_0$  and  $g_0^C$  proper basis states (simplified notation):

$$g_0 = \frac{|\phi\rangle\langle\phi|}{E - H_0 + i\varepsilon}, \quad \Rightarrow \quad g_0^C = \frac{|\psi_{\mathbf{k},\eta}^C\rangle\langle\psi_{\mathbf{k},\eta}^C|}{E - H_0 + i\varepsilon},$$

$$\eta = Z_A e^2 \mu_p / k.$$

## Coulomb function in momentum space

3D:

$$\langle \mathbf{p} | \psi_{\mathbf{k},\eta}^C \rangle = \psi_{\mathbf{k},\eta}^C(\mathbf{p}) = -\frac{1}{2\pi^2} \lim_{\gamma \rightarrow +0} \frac{d}{d\gamma} \left\{ \frac{[p^2 + (\gamma - ik)^2]^{i\eta}}{[\gamma^2 + (\mathbf{p} - \mathbf{k})^2]^{1+i\eta}} \right\}.$$

\* Guth and Mullin. // Phys. Rev. **83**, 667 (1951).

In partial waves:

$$\psi_{l,k,\eta}^C(p) = -\frac{2\pi e^{\pi\eta/2}}{pk} \lim_{\gamma \rightarrow +0} \frac{d}{d\gamma} \left\{ \left[ \frac{p^2 - (k + i\gamma)^2}{2pk} \right]^{i\eta} \frac{Q_l^{i\eta}(\zeta)}{(\zeta^2 - 1)^{i\eta/2}} \right\},$$

$$\eta = Z_1 Z_2 \alpha^2 \mu / k; \quad \zeta = (k^2 + p^2 + \gamma^2) / 2kp.$$

\* Dolinskii and Mukhamedzhanov. // Sov. Journ. of Nucl. Phys., vol. 3, No. 2, p. 180 (1966).

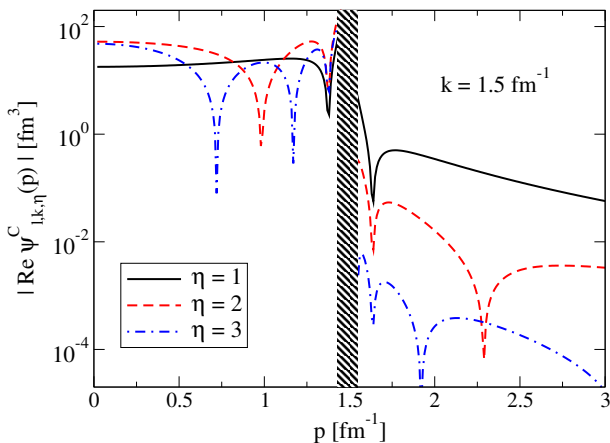
# The two representations

The ‘regular’ representation ( $p \neq k$ ):

$$\psi_{l,k,\eta}^C(p) = -\frac{4\pi\eta e^{-\pi\eta/2} q(pk)^l}{(p^2 + k^2)^{1+l+i\eta}} \lim_{\gamma \rightarrow +0} [p^2 - (k + i\gamma)^2]^{-1+i\eta} \\ \times \left[ \frac{\Gamma(1+l+i\eta)}{(1/2)_{l+1}} \right] {}_2F_1 \left( \frac{2+l+i\eta}{2}, \frac{1+l+i\eta}{2}, l+3/2; \frac{4k^2 p^2}{(p^2 + k^2)^2} \right).$$

The ‘pole-proximity’ representation ( $p \approx k$ ):

$$\psi_{l,k,\eta}^C(p) = i \frac{2\pi}{p} \exp(-\pi\eta/2 + i\sigma_l) \left[ \frac{(p+k)^2}{4pk} \right]^l \lim_{\gamma \rightarrow +0} 2 \\ \times \operatorname{Im} \left[ \frac{\Gamma(1+i\eta) e^{-i\sigma_l} (p+k)^{-1+i\eta}}{(p-k+i\gamma)^{1+i\eta}} {}_2F_1 \left( -l, -l-i\eta, 1-i\eta; \frac{(p-k)^2}{(p+k)^2} \right) \right].$$

Coulomb function in momentum space ( $l = 0$  plot)

\* Eremenko *et al.* // *Comp. Phys. Comm.* **187**, 195 (2015).

The two-body  $t$ -matrix elements in Coulomb basis

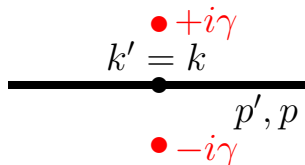
$$g_0^C(E)t_{(l)a}^S = \frac{\overbrace{|\psi_{l,k',\eta'}^C\rangle \langle \psi_{l,k',\eta'}^C| t_{(l)a}^S | \psi_{l,k,\eta}^C\rangle \langle \psi_{l,k,\eta}^C|}^{t_{a,l}^C(k',k,E)}}{E - H_0 + i\varepsilon},$$

$$t_{a,l}^C(k',k,E) = \int dp' dp \psi_{l,k',\eta'}^C(p')^\dagger t_{a,l}^S(p',p,E) \psi_{l,k,\eta}^C(p).$$

\* Some indices omitted for the simplicity.

Pinch singularity in the elastic channel:

- $E = 2\mu k^2 = 2\mu k'^2$ .
- Since  $\gamma \rightarrow +0$ , singularities of  $\psi_{l,k',\eta'}^{C\dagger}$  and  $\psi_{l,k,\eta}^C$  are pinching the integration contour.



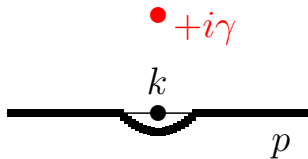
# If $t$ has a separable representation

$$t_l^S(k', k, E) = \sum_{zy} u_{l,z}(k') \lambda_{l,zy}(E) u_{l,y}(k)^\dagger,$$

$$t_l^C(k', k, E) = \sum_{zy} u_{l,z}^C(k') \lambda_{l,zy}(E) u_{l,y}^C(k)^\dagger,$$

$$u_{l,z}^C(k') = \int \frac{dp' p'^2}{2\pi^2} u_{l,z}(p') \psi_{l,k',\eta'}^C(p')^\dagger, \quad u_{l,y}^C(k)^\dagger = \int \frac{dp p^2}{2\pi^2} u_{l,y}(p)^\dagger \psi_{l,k,\eta}^C(p).$$

- Two independent integrals over  $p$  and  $p'$ .
- Cauchy's theorem.
- **No pinch singularity!**

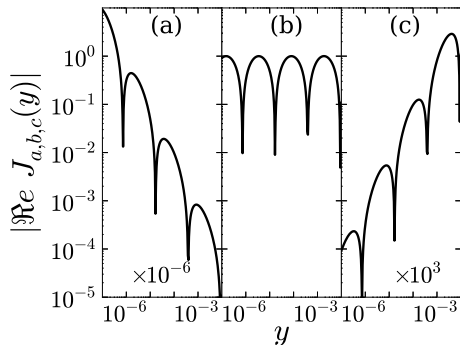




## Gel'fand-Shilov regularization for complex form-factors

$$u_l^C(k) \propto \int_{a<0}^{b>0} \frac{f(y) dy}{y^{1+i\eta}} \equiv \int_a^b dy J_a(y), \quad (\text{e.g. } f(y) = y^2 + y + 1).$$

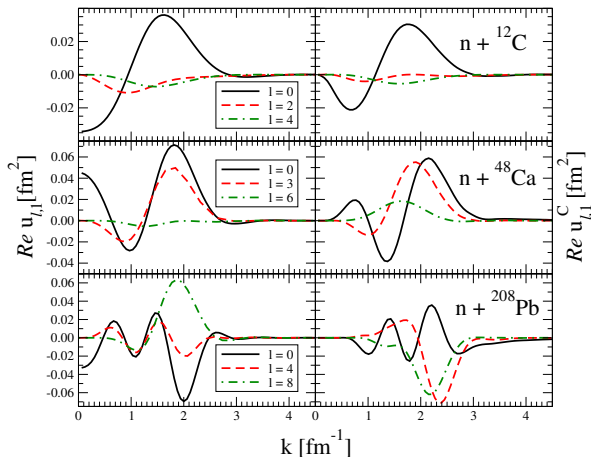
Subtract as many terms of Laurent expansion of  $f(y)$  around the integrand's special point  $y = 0$  as needed to split the integral and get the regular term, plus the analytically calculated terms.



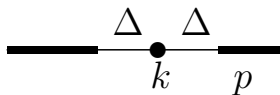
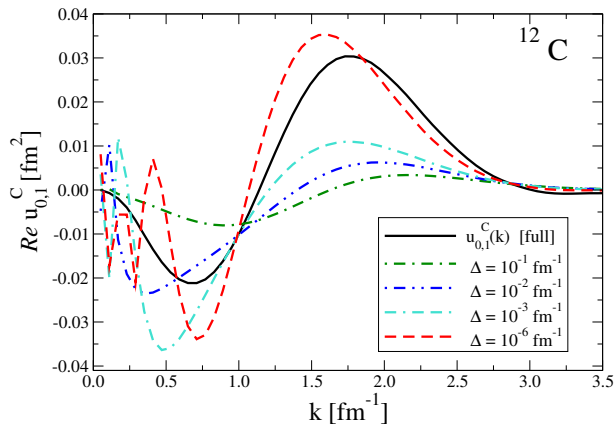
\* Upadhyay *et al.* // Phys. Rev. C **90**, 014615 (2014).

## Results: form-factors in Coulomb basis

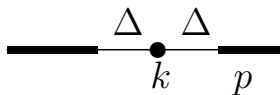
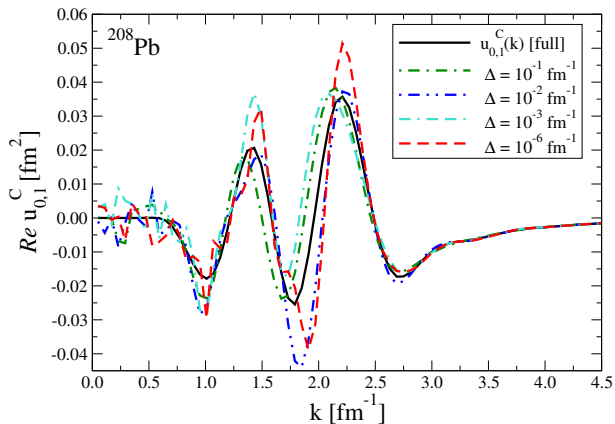
$$u_{l,1}^C(k) = \int \frac{dp p^2}{2\pi^2} u_{l,1}(p) \psi_{l,k,\eta}^C(p)^*.$$



\* Upadhyay *et al.* // Phys. Rev. C **90**, 014615 (2014).

Singularity contribution for  $n + {}^{12}\text{C}$ 

\* Upadhyay *et al.* // Phys. Rev. C **90**, 014615 (2014).

Singularity contribution for  $n + {}^{208}\text{Pb}$ 

\* Upadhyay *et al.* // Phys. Rev. C **90**, 014615 (2014).

# Faddeev-AGS equations in Coulomb basis (in progress)

## Published abstract formalism

Faddeev-AGS equations in Coulomb basis for real two-body  $t$ -matrices and spinless particles.

\* Mukhamedzhanov *et al.* // Phys. Rev. C **86**, 034001 (2012).

## To do:

- Take spin degrees of freedom into account.
- Deal with the 3-body singularity.
- Generalize the equations for the complex potentials.
- Take care about other implementation details, develop the algorithms and codes to solve the equations.

## Objective:

- Calculate scattering observables and make the code available as open source.

# Summary & Outlook

- Faddeev formalism is the theoretical tool to study deuteron-induced reactions.
- This formalism treats all possible three-body channels on the same footage.
- Momentum space is preferable due to the boundary conditions.
- Coulomb interaction can be treated properly by using the Coulomb basis in momentum space.
- Pinch singularity is avoided by choosing the two-body interactions in separable form.
- Mathematics and machinery are developed to compute Coulomb functions and matrix elements in Coulomb basis in momentum space.
- Work is in progress to cast the Faddeev-AGS equations in the Coulomb basis taking into account spin degrees of freedom and 3-body singularity in order to solve the equations numerically.