

Direct reaction studies with microscopic inputs for many-nucleon systems

... the number of the degrees of freedom excited is very small. ... even in the transition between collective states, in which infinite number of transition processes are involved, only a few collective states participate.

M. Kawai and S. Yoshida, “Nuclear Reaction Theories” (Asakura) p. 149 (Translated by KO).

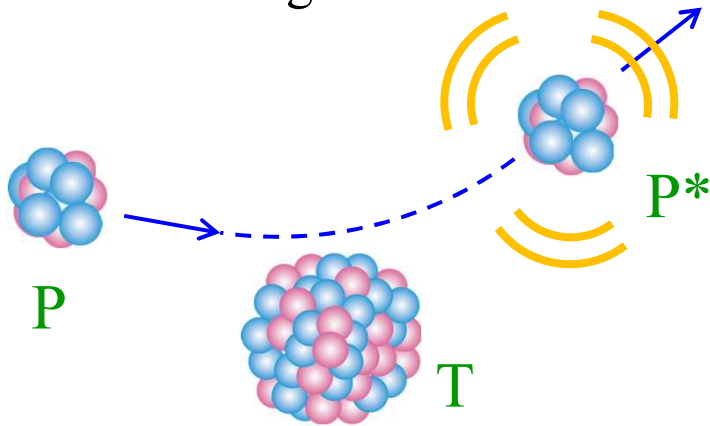
Kazuyuki Ogata

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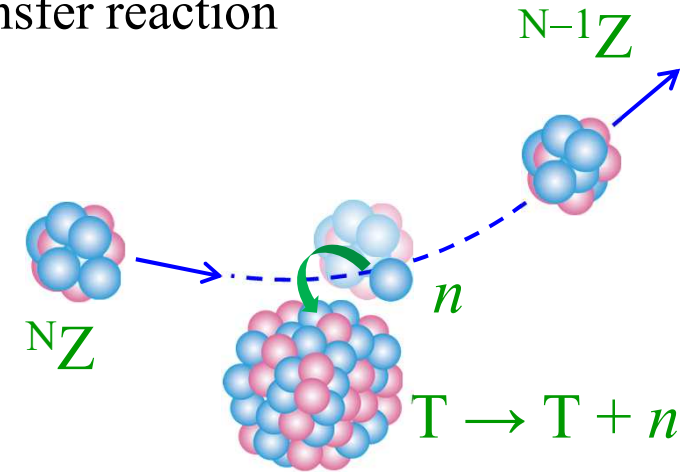
Microscopic, but not *ab initio*

Various nuclear reaction processes

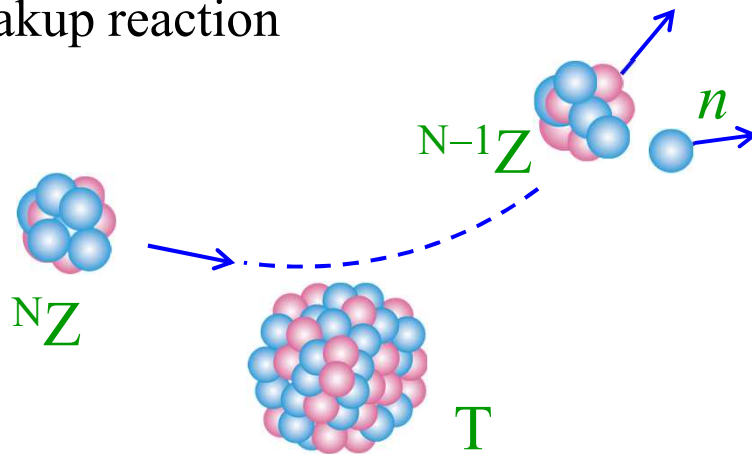
Inelastic scattering



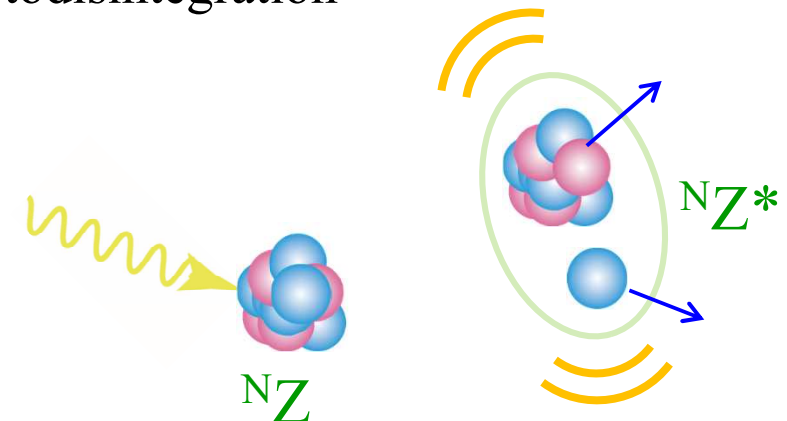
Transfer reaction



Breakup reaction



Photodisintegration



What to do?

The probability of a reaction process is given by the absolute square of the transition matrix.

$$T_{\beta\alpha} = \langle \Phi_{\beta}^{\text{free}} | V_{\beta} | \Psi_{\alpha} \rangle$$

$\Phi_{\beta}^{\text{free}}$: Internal W.Fns. of the constituents and their free relative W.Fns.

V_{β} : Interactions between the constituents

Ψ_{α} : Exact W.Fn. of the system in the initial (\neq incident) channel

Key: simplification of Ψ_{α} with keeping a microscopic picture

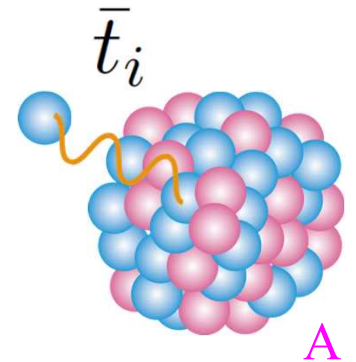
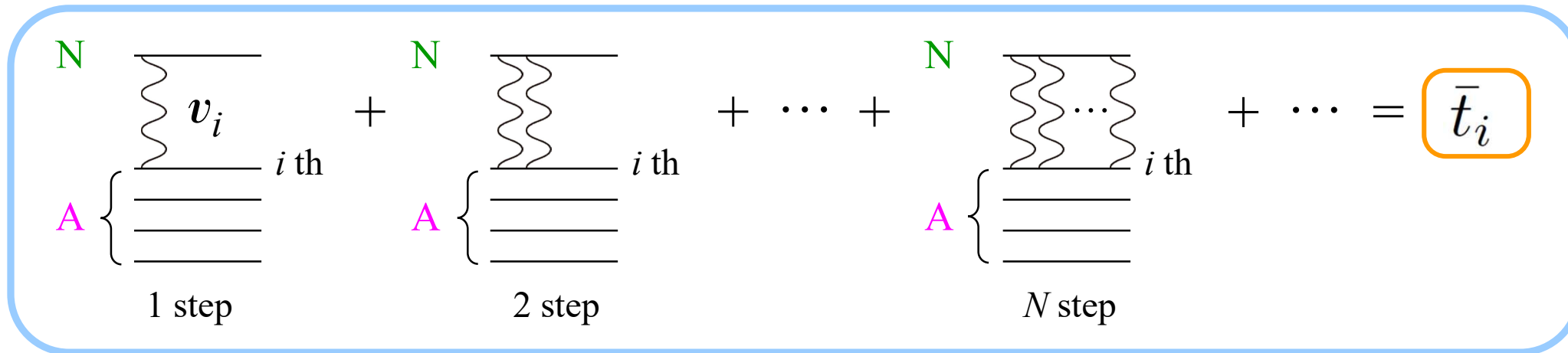
 Multiple Scattering Theory (MST)

MST

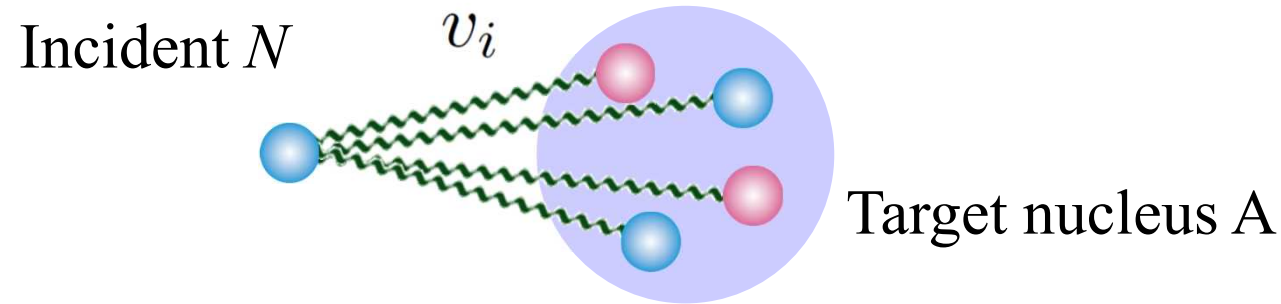
Microscopic description of optical potentials

TABLE I. Optical-Model Parameters Neutrons

NUCLIDE	ENERGY (MEV)	REAL POTENTIAL			VOL. IMAG. POTENTIAL			SURF. IMAG. POTENTIAL			SPIN-ORBIT POTENTIAL			ST	SR	FIT	NOTE	REF.
		V	R	A	W	RW	AW	WD	RD	AD	VSO	RSO	ASO					
AL	1.	40.	1.25*	0.65*				5.0G*	1.25*	0.98*	10.*	1.25*	0.65*	3520	1340	S3	15	GIL63
AL	1.5	47.4	1.25*	0.46				6.3G	1.25*	0.98*	10.*	1.25*	0.46	3204		S1	10	KOR68
AL	2.47	48.0	1.14	0.65				8.42	1.19	0.48*	8.0*	1.14	0.65	2530	1270	S2	2	HOL71
AL	3.00	47.9	1.13	0.72				7.35	1.08	0.48*	8.0*	1.13	0.72	2520	1250	S2	2	HOL71
AL	3.49	48.7	1.18	0.61				8.46	1.29	0.48*	8.0*	1.18	0.61	2360	1130	S1	2	HOL71
AL	4.00	49.1	1.20	0.62				7.99	1.26	0.48*	8.0*	1.20	0.62	2290	1090	S2	2	HOL71
AL	4.56	50.2	1.18	0.59				8.38	1.26	0.48*	8.0*	1.18	0.59	2060	1020	S1	2	HOL71
AL	6.09	47.8	1.20	0.67				8.23	1.23	0.48*	8.0*	1.20	0.67	1880	1070	S3	2	HOL71
AL	7.	45.5	1.25*	0.65*				9.5G	1.25*	0.98*	8.6	1.25*	0.65*			X3		BJO58
AL	7.05	49.1	1.20	0.68				7.90	1.20	0.48*	8.0*	1.20	0.68	1800	1040	S2	2	HOL71
AL	7.97	49.4	1.20	0.69				12.1	1.30	0.41	9.8	1.20	0.69			S1	2	BRA72



MST description of nucleon-nucleus (NA) scattering



Difficult to handle

$$T = \sum_i \boxed{v_i} + \sum_i v_i G_0^{(+)} \sum_j v_j + \sum_i v_i G_0^{(+)} \sum_j v_j G_0^{(+)} \sum_k v_k + \dots = \sum_i \Lambda_i$$

“free” propagator

Kinetic energy Op.

Hamiltoniam of A

$$G_0^{(+)} = \frac{1}{E - (T_{\text{NA}} + H_A) + i\eta}$$

Resummation (1/3)

$$\Lambda_i \equiv v_i + v_i G_0^{(+)} \sum_j v_j + v_i G_0^{(+)} \sum_j v_j G_0^{(+)} \sum_k v_k + \dots$$

$$\Lambda_i = v_i + v_i G_0^{(+)} v_i + v_i G_0^{(+)} v_i G_0^{(+)} \sum_k v_k + \dots$$

$$+ v_i G_0^{(+)} \sum_{j \neq i} v_j + v_i G_0^{(+)} \sum_{j \neq i} v_j G_0^{(+)} \sum_k v_k + \dots$$

$$= v_i + v_i G_0^{(+)} \left(v_i + v_i G_0^{(+)} \sum_k v_k + \dots \right)$$

$$+ v_i G_0^{(+)} \sum_{j \neq i} \left(v_j + v_j G_0^{(+)} \sum_k v_k + \dots \right)$$

N interacts last with i

$$\boxed{\Lambda_i} = v_i + v_i G_0^{(+)} \boxed{\Lambda_i} + v_i G_0^{(+)} \sum_{j \neq i} \boxed{\Lambda_j}$$

N interacts last with

$j \neq i$

Resummation (2/3)

$$\Lambda_i = v_i + v_i G_0^{(+)} \Lambda_i + v_i G_0^{(+)} \sum_{j \neq i} \Lambda_j$$

$$\Lambda_i = \frac{1}{1 - v_i G_0^{(+)}} v_i + \frac{1}{1 - v_i G_0^{(+)}} v_i G_0^{(+)} \sum_{j \neq i} \Lambda_j$$

$$\left(t_i = v_i + v_i G_0^{(+)} t_i \right) \quad t_i \equiv \frac{1}{1 - v_i G_0^{(+)}} v_i \quad \longrightarrow \quad \Lambda_i = t_i + t_i G_0^{(+)} \sum_{j \neq i} \Lambda_j$$

T-matrix in terms of an effective interaction

$$T = \sum_i t_i + \sum_i t_i G_0^{(+)} \sum_{j \neq i} t_j + \sum_i t_i G_0^{(+)} \sum_{j \neq i} t_j G_0^{(+)} \sum_{k \neq j} t_k + \dots$$

Resummation (3/3)

If the A -body W.Fn. is antisymmetrized, the label i has no specific meaning:

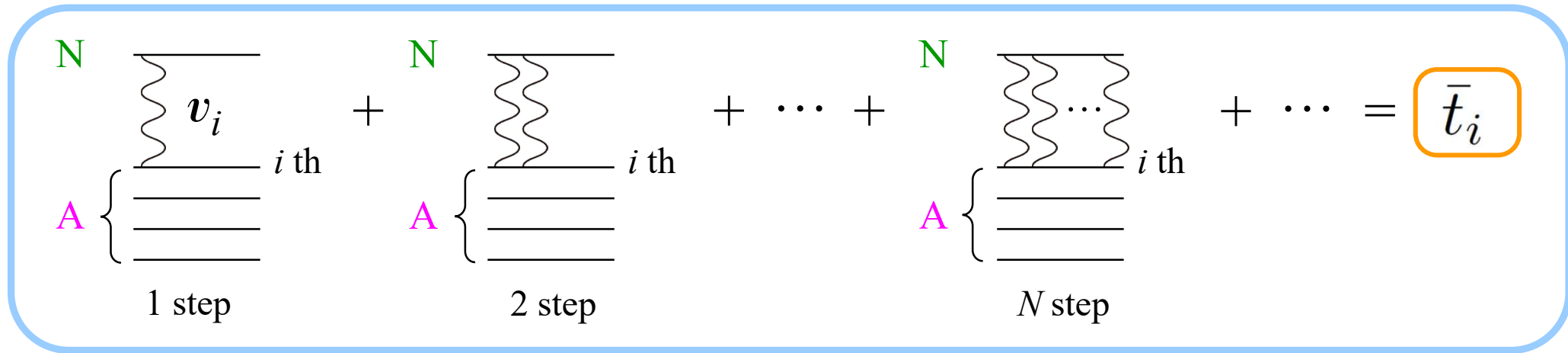
$$\sum_{j \neq i} t_j \rightarrow \frac{A-1}{A} \sum_j t_j \equiv \sum_j \bar{t}_j$$

$$T = \frac{A}{A-1} \bar{T}$$

$$\bar{T} \equiv \sum_i \bar{t}_i + \sum_i \bar{t}_i G_0^{(+)} \sum_j \bar{t}_j + \sum_i \bar{t}_i G_0^{(+)} \sum_j \bar{t}_j G_0^{(+)} \sum_k \bar{t}_k + \dots$$

One can use **an effective interaction** instead of **the bare NN interaction**.

Multiple Scattering Theory (MST)



$$\left(T_{\text{NA}} + \sum_i v_i + H_A - E\right) \Psi = 0 \xrightarrow{\text{Resummation}} \left(T_{\text{NA}} + \sum_i \bar{t}_i + H_A - E\right) \bar{\Psi} = 0$$

Note: we have restricted the b.c.

$$T = \sum_i v_i + \sum_i v_i G_0^{(+)} \sum_j v_j + \dots$$

$$\bar{t}_i = \frac{A-1}{A} t_i, \quad t_i = v_i + v_i G_0^{(+)} t_i$$

L. L. Foldy, Phys. Rev. **67**, 107 (1945); K. M. Watson, Phys. Rev. **89**, 115 (1953).
 A. K. Kerman, H. McManus, and R. M. Thaler, Ann. Phys. (NY) **8**, 551 (1959).
 M. Yahiro, K. Minomo, KO, and M. Kawai, PTP **120**, 767 (2008).

Impulse approximation to t

$$t_i = v_i + v_i \frac{1}{E - (T_{\text{NA}} + H_{\text{A}}) + i\eta} t_i$$

$$t_i = v_i + v_i \frac{1}{\boxed{E} - T_{\text{NA}} - T_{i\text{A}} \boxed{-H_{\text{B}} - V_{i\text{B}}} + i\eta} t_i$$

$\approx E'$



$$t_i^{\text{IA}} = v_i + v_i \frac{1}{E' - T_{\text{NA}} - T_{i\text{A}} + i\eta} t_i^{\text{IA}}$$

At high energies, the effect from the nucleons except the interacting one may be disregarded.

g matrix approximation to t

NN interaction is obtained by solving the Bethe-Goldstone Eq.
assuming infinite nuclear matter of density ρ :

$$g(\mathbf{q}', \mathbf{q}; \mathbf{K}) = v(\mathbf{q}', \mathbf{q}) + \int d\mathbf{k}' v(\mathbf{q}', \mathbf{k}') \frac{\overbrace{Q(\mathbf{k}', \mathbf{K}; k_F)}^{\text{Pauli's Op.}}}{E(\mathbf{k}_0, \mathbf{K}) - E(\mathbf{k}', \mathbf{K})} g(\mathbf{k}', \mathbf{q}; \mathbf{K}),$$
$$E(\mathbf{k}_0, \mathbf{K}) - E(\mathbf{k}', \mathbf{K}) = \frac{\hbar^2}{m} (k_0^2 - k'^2) + \underbrace{U(|\mathbf{k}_0 + \mathbf{K}|) + U(|\mathbf{k}_0 - \mathbf{K}|) - U(|\mathbf{k}' + \mathbf{K}|) - U(|\mathbf{k}' - \mathbf{K}|)}_{\text{Auxiliary pot.}}.$$

Elastic scattering

Microscopic optical potential with MST

$$\left[T_{\text{NA}} + U^{\text{dr}}(R) + U^{\text{ex}}(R) - E_{\text{cm}} \right] \chi(\mathbf{R}) = 0$$

Direct term:

$$U^{\text{dr}}(R) = \frac{A-1}{A} \int g^{\text{dr}}(s, k_{\text{F}}(|\mathbf{R} + \mathbf{s}/2|)) \rho(|\mathbf{R} + \mathbf{s}|) d\mathbf{s},$$

Nuclear one-body density

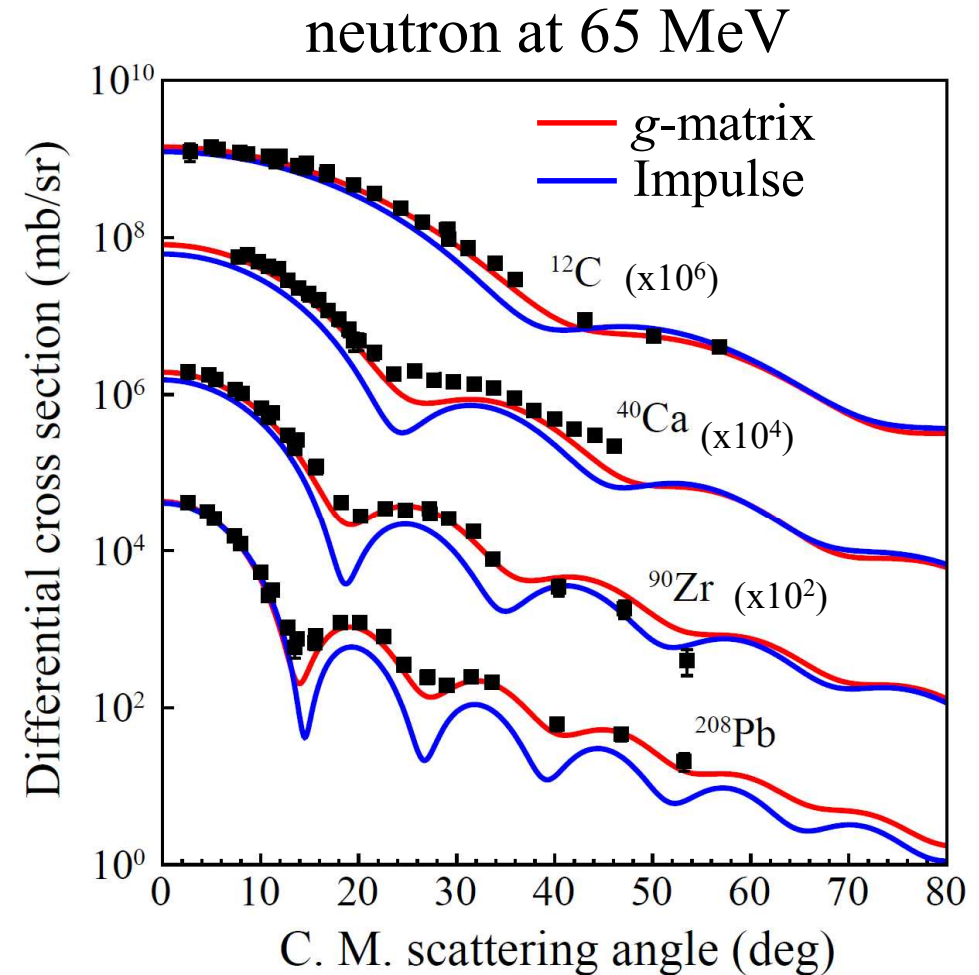
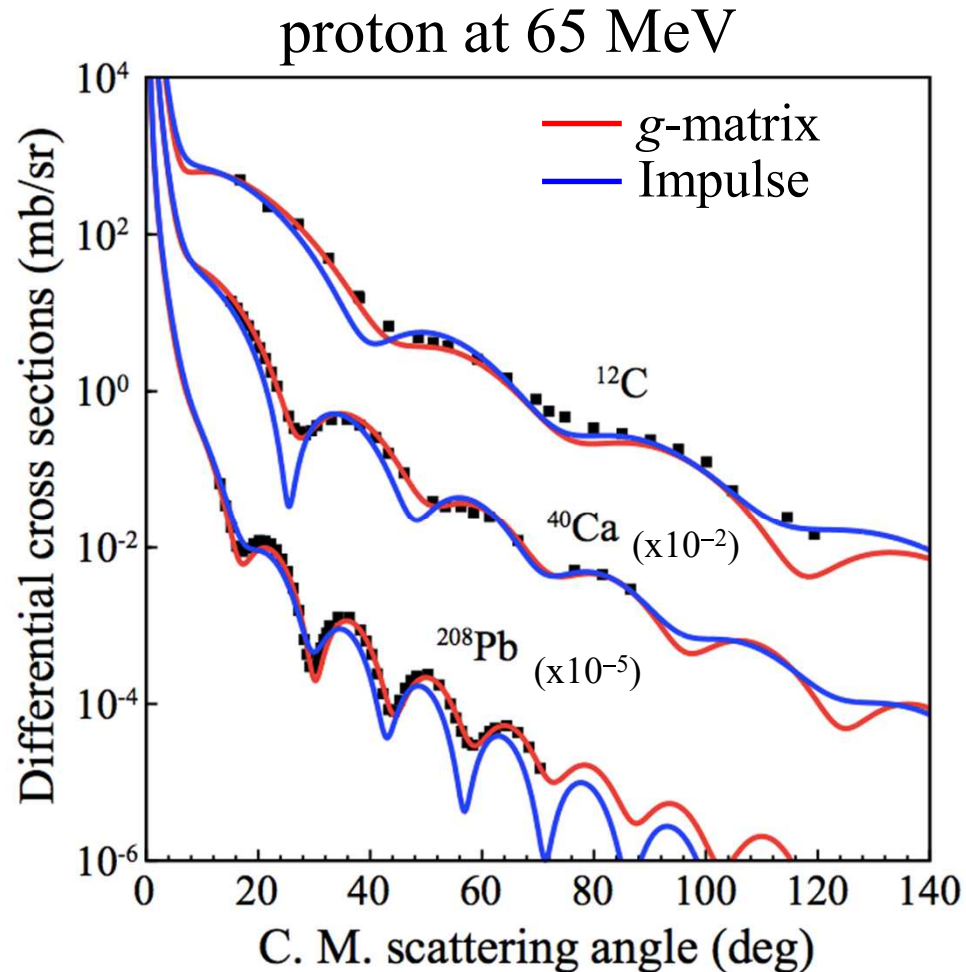
Knock-on exchange term with Brieva-Rook localization:

F. A. Brieva and J. R. Rook, Nucl. Phys. A 291, 317 (1977).

$$U^{\text{ex}}(R) = \frac{A-1}{A} \int \frac{3j_1(k_{\text{F}}(|\mathbf{R} + \mathbf{s}/2|)s)}{k_{\text{F}}(|\mathbf{R} + \mathbf{s}/2|)s} j_0(K(R)s) \\ \times g^{\text{ex}}(s, k_{\text{F}}(|\mathbf{R} + \mathbf{s}/2|)) \rho(|\mathbf{R} + \mathbf{s}/2|) d\mathbf{s}.$$

Nuclear one-body density

Application of MST to NA elastic scattering



cf. K. Amos+, *Adv. Nucl. Phys.* **25**, 275 (2000).

T. Furumoto+, *PRC* **78**, 044610 (2008).

M. Toyokawa+, *PRC* **92**, 024618 (2015).

Inelastic scattering

MST-based description of inelastic scattering

Hagino, Moro, and O, Prog. Part. Nucl. Phys. 125, 10395 (2022).

$$\left(T_{\text{NA}} + \sum_i \bar{t}_i + H_A - E\right) \bar{\Psi} = 0$$

$$\bar{\Psi} = \sum_c \chi_c \phi_c^A \quad H_A \phi_c^A = \varepsilon_c \phi_c^A \quad \langle \phi_{c'}^A | \phi_c^A \rangle = \delta_{c'c}$$

Microscopic coupled-channel equations:

$$\left(T_{\text{NA}} + \underbrace{\int \phi_c^{*A} \sum_{i \in A} \bar{t}_i \phi_c^A d\xi_A}_{\text{Diagonal pot.}} - E_c\right) \chi_c = - \underbrace{\sum_{c' \neq c} \left(\int \phi_c^{*A} \sum_{i \in A} \bar{t}_i \phi_{c'}^A d\xi_A \right) \chi_{c'}}_{\text{Coupling pot.}}$$

Diagonal pot.

Coupling pot.

(Processes through channel c')

Coupled-Channel equations/potentials

Hagino, O, Moro, PPNP 125, 103951 (2022).

□ CC Eqs.
$$\left[-\frac{\hbar^2}{2\mu} \frac{d^2}{dR^2} + \frac{\hbar^2}{2\mu} \frac{(L')^2}{R^2} - E_{n'I'} \right] \chi_{c'c_0}^{(J_T)}(k_{n'I'}, R) = - \sum_c U_{c'c}^{(J_T)}(R) \chi_{cc_0}^{(J_T)}(k_{nl}, R)$$

□ CC Pots.
$$U_{c'c, J_T}^{\text{dr}}(R) = \int d\hat{\mathbf{R}} \sum_{m'_I m'_L} \langle I' m'_I L' m'_L | J_T M_T \rangle i^{-P'} i^{-L'} Y_{L' m'_L}^*(\hat{\mathbf{R}}) \\ \times \sum_{m_I m_L} \langle I m_I L m_L | J_T M_T \rangle i^P i^L Y_{L m_L}(\hat{\mathbf{R}}) \sum_{\lambda \mu} \langle I m_I \lambda \mu | I' m'_I \rangle \mathcal{U}_{n'I'nl, \lambda}^{\text{dr}}(\mathbf{R}),$$

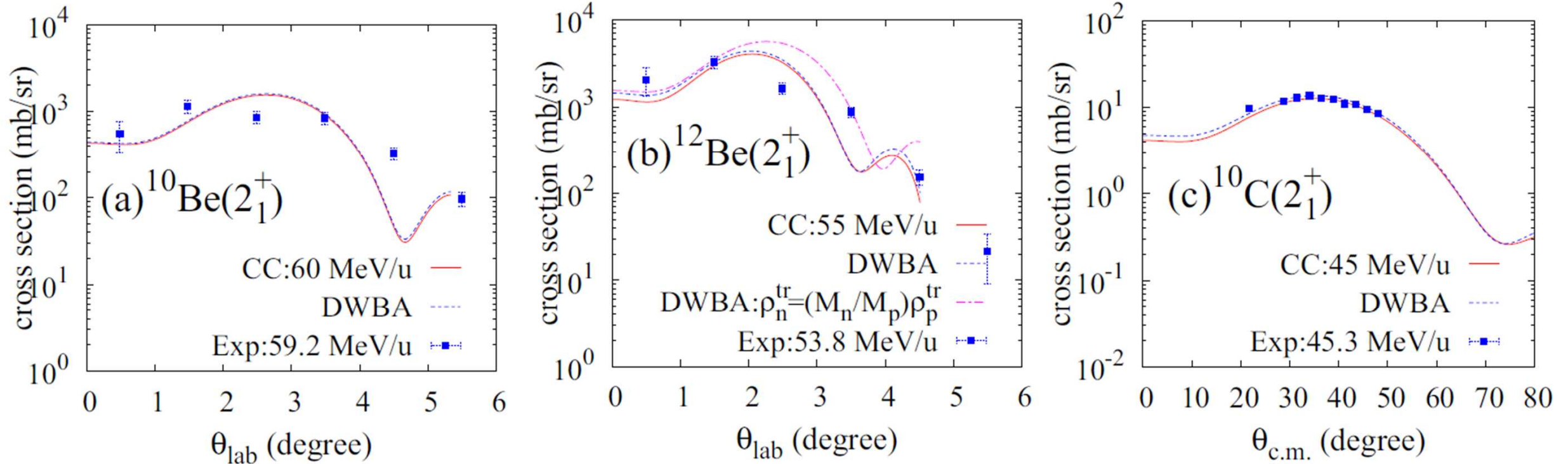
$$\mathcal{U}_{n'I'nl, \lambda}^{\text{dr}}(\mathbf{R}) \equiv \int \bar{g}^{\text{dr}}(s, k_{F; n'I'nl}(|\mathbf{R} + \mathbf{s}/2|)) \rho_{n'I'nl, \lambda}^{\text{tr}}(r_A) Y_{\lambda \mu}^*(\hat{\mathbf{r}}_A) d\mathbf{r}_A$$

One-body transition density

$$\rho_{n'I' m'_I n l m_I}^{\text{tr}}(\mathbf{r}_A) \equiv \int \Phi_{n'I' m'_I}^*(\xi) \sum_{i=1}^A \delta(\mathbf{r}_i - \mathbf{r}_A) \Phi_{n l m_I}(\xi) d\xi = \sum_{\lambda \mu} \langle I m_I \lambda \mu | I' m'_I \rangle \rho_{n'I'nl, \lambda}^{\text{tr}}(r_A) Y_{\lambda \mu}^*(\hat{\mathbf{r}}_A)$$

Application to p-inelastic scattering off ^{10}Be , ^{12}Be , ^{10}C

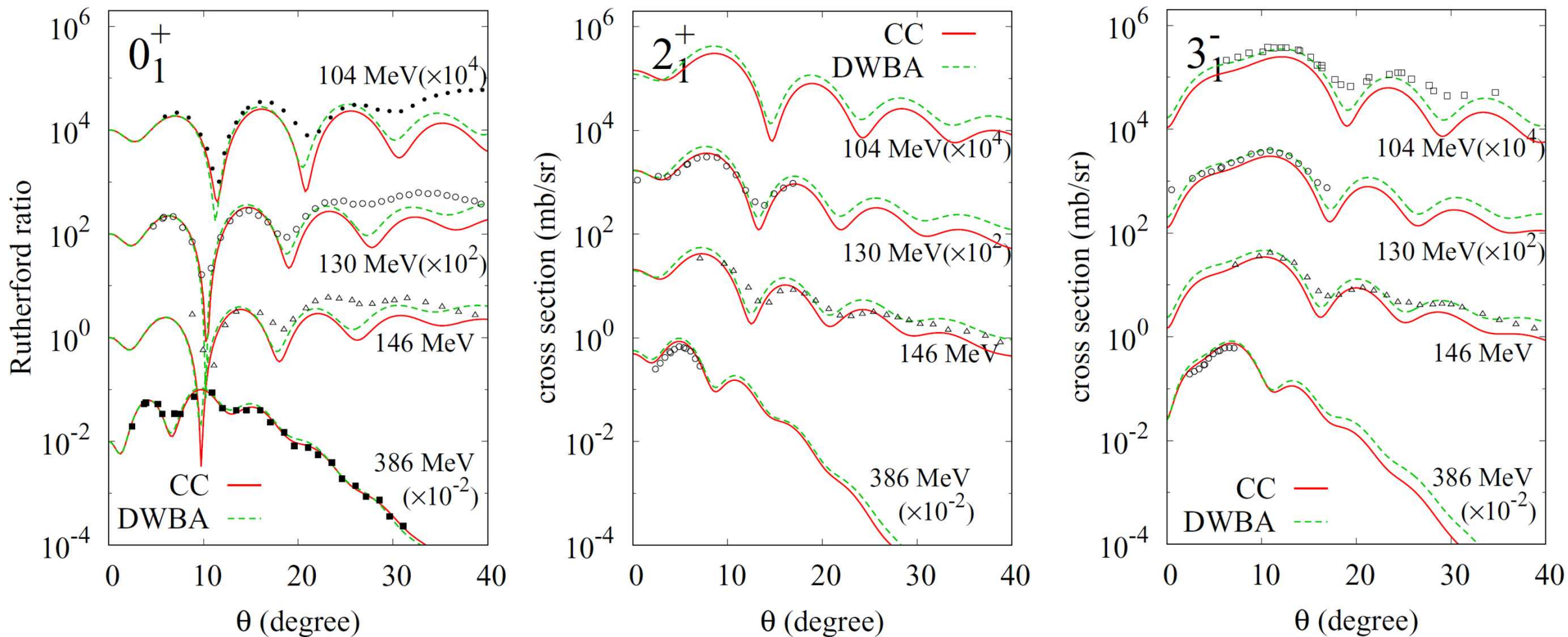
Kanada-En'yo and O, Phys. Rev. C 100, 064616 (2019).



AMD (Antisymmetrized Molecular Dynamics) densities are used.

Application to α - ^{16}O scattering

Kanada-En'yo and O, Phys. Rev. C 99, 064608 (2019).



α -A CC pots. are obtained by folding NA CC pots. with a one-body density of α .

Breakup reaction

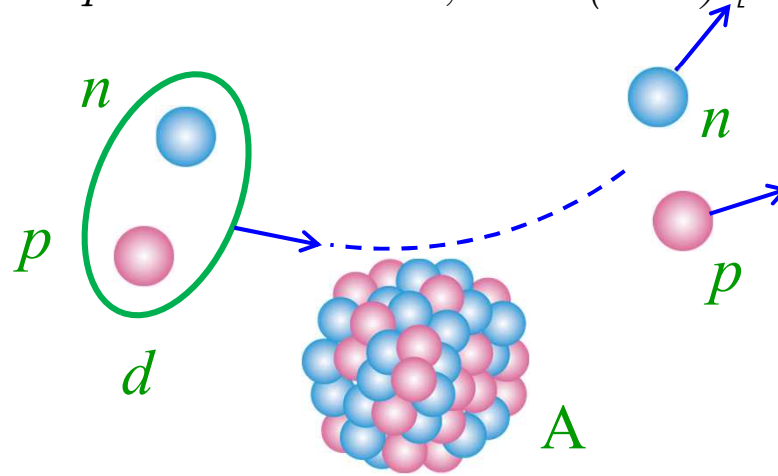
Two questions

1. *Can we use a single Lippmann-Schwinger Eq.?*

2. *How to implement the continuum st. of many-body system?*

The Faddeev theory

*L. D. Faddeev, Zh. Eksp. Theor. Fiz. **39**, 1459 (1960) [Sov. Phys. JETP **12**, 1014 (1961)].*



$$[E - K - V_{pn} - V_{pA} - V_{nA}] \Psi = 0, \quad \Psi = \Psi_d + \Psi_p + \Psi_n.$$

Faddeev Eqs.

$$[E - K - V_{pn}] \Psi_d = V_{pn} (\Psi_p + \Psi_n),$$

$$[E - K - V_{nA}] \Psi_n = V_{nA} \Psi_d + V_{nA} \Psi_p,$$

$$[E - K - V_{pA}] \Psi_p = V_{pA} \Psi_d + V_{pA} \Psi_n.$$

Three-body theory in a model space

*N. Austern, M. Yahiro, and M. Kawai, Phys. Rev. Lett. **63**, 2649 (1989);*

*N. Austern, M. Kawai, and M. Yahiro, Phys. Rev. C **53**, 314 (1996).*

$$[E - K - V_{pn} - V_{pA} - V_{nA}] \Psi = 0, \quad \Psi = \Psi_d + \Psi_p + \Psi_n.$$

Faddeev Eqs.

not pair int. but 3-body int.

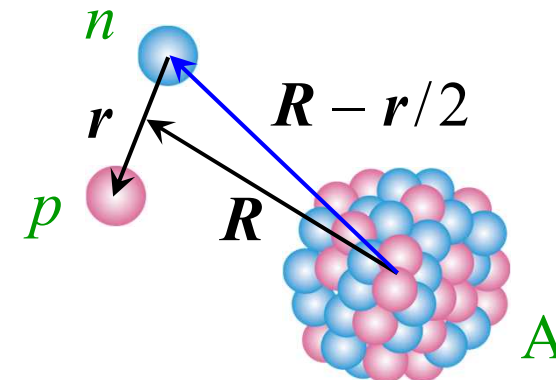
$$[E - K - V_{pn} - \mathcal{P}_{l_{\max}} (V_{nA} + V_{pA}) \mathcal{P}_{l_{\max}}] \Psi_d = V_{pn} (\Psi_p + \Psi_n),$$

$$[E - K - V_{nA}] \Psi_n = (V_{nA} - \mathcal{P}_{l_{\max}} V_{nA} \mathcal{P}_{l_{\max}}) \Psi_d + V_{nA} \Psi_p,$$

$$[E - K - V_{pA}] \Psi_p = (V_{pA} - \mathcal{P}_{l_{\max}} V_{pA} \mathcal{P}_{l_{\max}}) \Psi_d + V_{pA} \Psi_n.$$

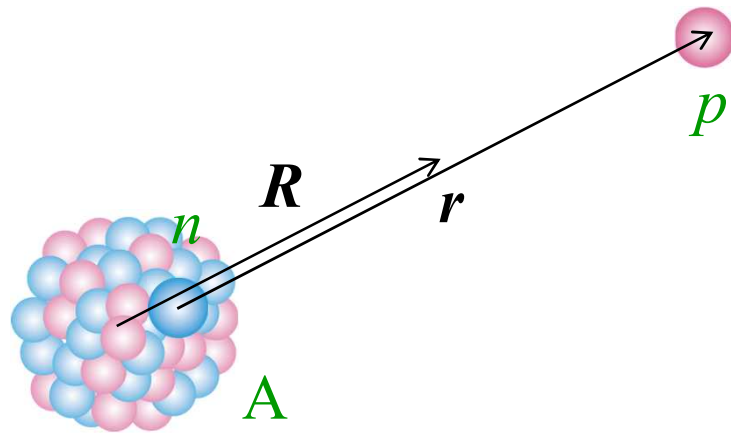
$$\mathcal{P}_{l_{\max}} = \int d\hat{r}' \sum_{l \leq l_{\max}} \sum_m Y_{lm}(\hat{r}) Y_{lm}^*(\hat{r}')$$

$$\mathcal{P}_0 e^{-\mu(\mathbf{R} - \mathbf{r}/2)^2} \rightarrow e^{-\mu R^2} e^{-\mu r^2/4}$$



l -truncation

*N. Austern, M. Yahiro, and M. Kawai, Phys. Rev. Lett. **63**, 2649 (1989);
N. Austern, M. Kawai, and M. Yahiro, Phys. Rev. C **53**, 314 (1996).*



$$\mathcal{P}_{l_{\max}} = \int d\hat{\mathbf{r}}' \sum_{l \leq l_{\max}} \sum_m Y_{lm}(\hat{\mathbf{r}}) Y_{lm}^*(\hat{\mathbf{r}}')$$

$\mathcal{P}_{l_{\max}}$ smears out $\hat{\mathbf{r}}$ with the resolution of $1/l_{\max}$.

[If $l_{\max} \rightarrow \infty$, it means $\delta(\hat{\mathbf{r}}' - \hat{\mathbf{r}})$.]

- We have no rearrangement-like channel in the asymptotic region because of $\mathcal{P}_{l_{\max}}$.
- As l_{\max} increases, the coupling between the 1st Eq. and the other two becomes weaker.

Three-body theory in a model space

*N. Austern, M. Yahiro, and M. Kawai, Phys. Rev. Lett. **63**, 2649 (1989);*

*N. Austern, M. Kawai, and M. Yahiro, Phys. Rev. C **53**, 314 (1996).*

$$[E - K - V_{pn} - V_{pA} - V_{nA}] \Psi = 0, \quad \Psi = \Psi_d + \Psi_p + \Psi_n.$$

Faddeev Eqs.

not pair int. but 3-body int.

→ 0

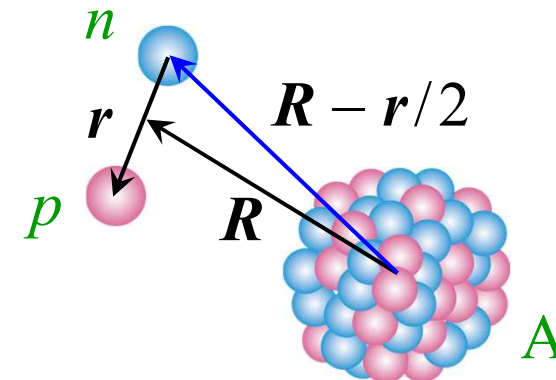
$$[E - K - V_{pn} - \mathcal{P}_{l_{\max}} (V_{nA} + V_{pA}) \mathcal{P}_{l_{\max}}] \Psi_d = V_{pn} (\Psi_p + \Psi_n)$$

$$[E - K - V_{nA}] \Psi_n = (V_{nA} - \mathcal{P}_{l_{\max}} V_{nA} \mathcal{P}_{l_{\max}}) \Psi_d + V_{nA} \Psi_p,$$

$$[E - K - V_{pA}] \Psi_p = (V_{pA} - \mathcal{P}_{l_{\max}} V_{pA} \mathcal{P}_{l_{\max}}) \Psi_d + V_{pA} \Psi_n.$$

$$\mathcal{P}_{l_{\max}} = \int d\hat{r}' \sum_{l \leq l_{\max}} \sum_m Y_{lm}(\hat{r}) Y_{lm}^*(\hat{r}')$$

$$\mathcal{P}_0 e^{-\mu(\mathbf{R} - \mathbf{r}/2)^2} \rightarrow e^{-\mu R^2} e^{-\mu r^2/4}$$



The Continuum-Discretized Coupled-Channels method

*N. Austern, M. Yahiro, and M. Kawai, Phys. Rev. Lett. **63**, 2649 (1989);*

*N. Austern, M. Kawai, and M. Yahiro, Phys. Rev. C **53**, 314 (1996).*

CDCC solves the following LS eq.:

$$\Psi^{\text{CDCC}} = e^{i\mathbf{K}\cdot\mathbf{R}}\phi_d + \frac{1}{E - H_d + i\varepsilon} \mathcal{P}_{l_{\max}} (V_{nA} + V_{pA}) \mathcal{P}_{l_{\max}} \Psi^{\text{CDCC}}.$$

CDCC gives a proper solution to a three-body scattering problem
if the solution converges with respect to l .

- Continuum-Discretization has nothing to do with the justification of CDCC.
- l -truncation allows one to truncate also r and k .
- Convergence for other quantities (r_{\max} , k_{\max} , etc.) must be confirmed.

Faddeev–Alt-Grassberger-Sandhas (FAGS) vs. CDCC

N. J. Upadhyay, A. Deluva, F. M. Nunes, PRC 85, 054621 (2012).

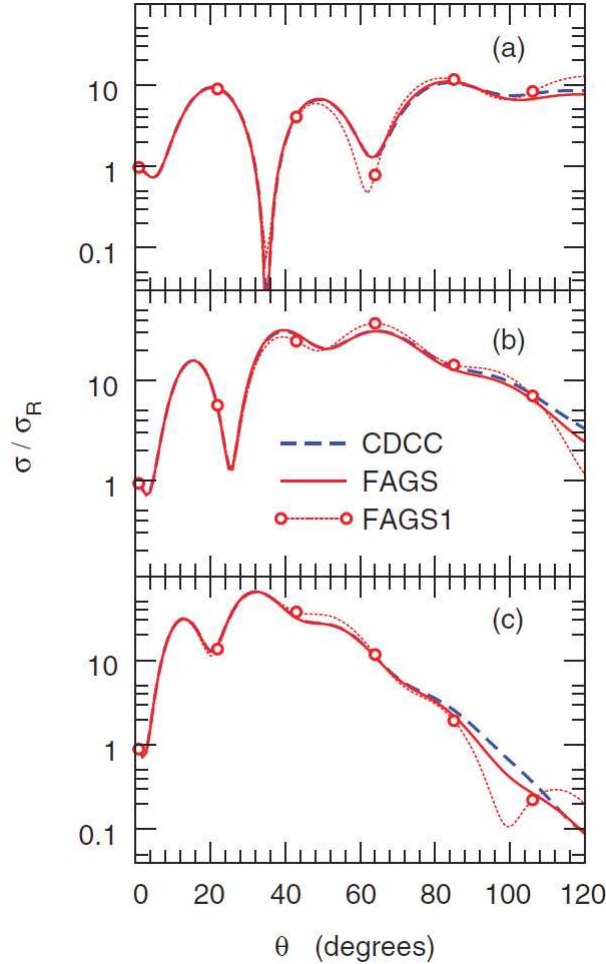


FIG. 2. (Color online) Elastic cross section for $d+^{10}\text{Be}$: (a) $E_d = 21.4$ MeV, (b) $E_d = 40.9$ MeV, and (c) $E_d = 71$ MeV.

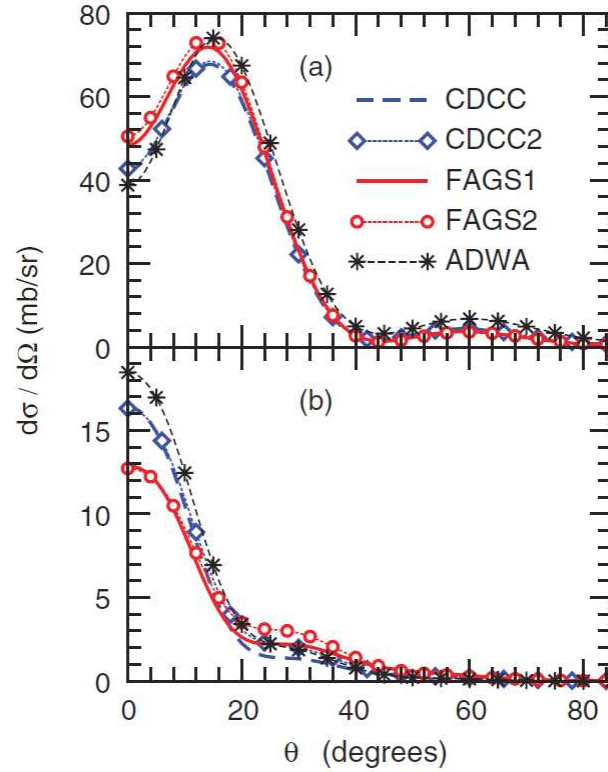


FIG. 6. (Color online) Angular distribution for $^{12}\text{C}(d, p)^{13}\text{C}$: (a) $E_d = 12$ MeV and (b) $E_d = 56$ MeV.

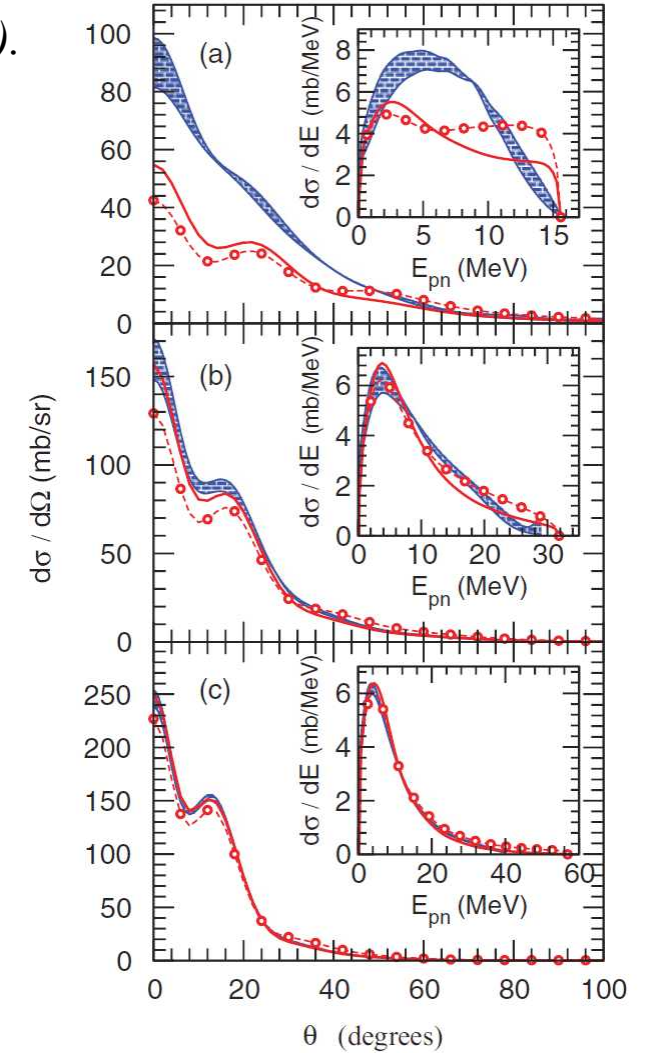
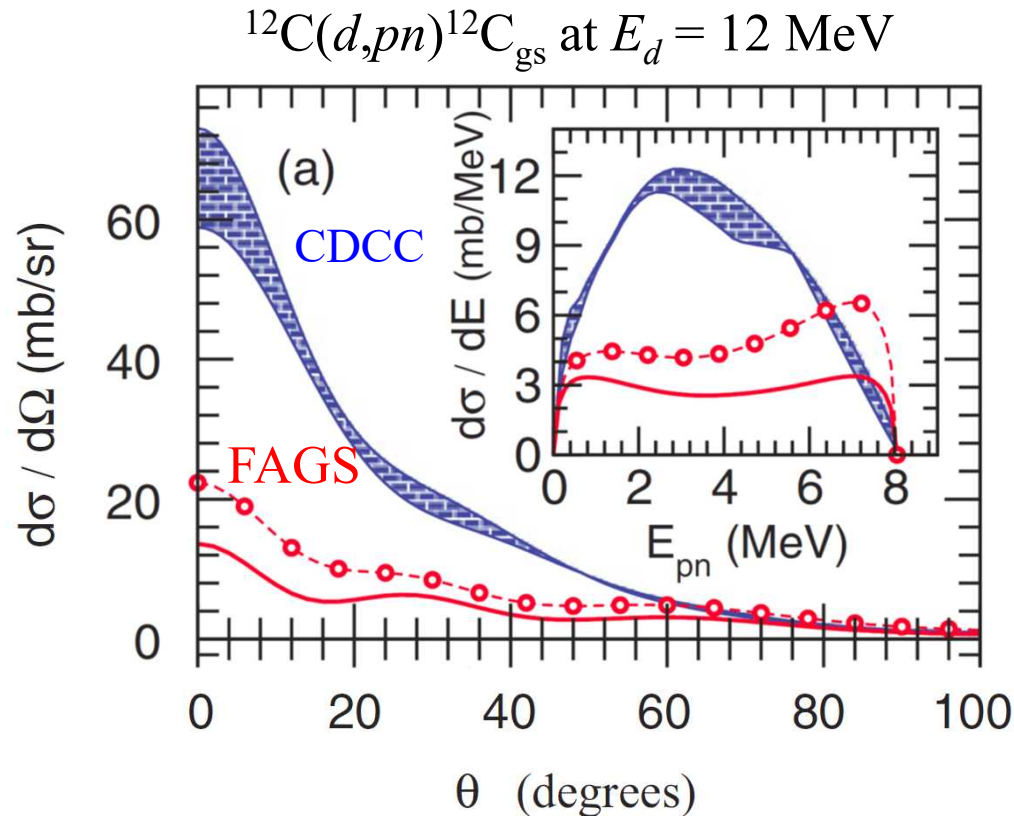


FIG. 8. (Color online) Breakup distributions for the $^{10}\text{Be}(d, pn)^{10}\text{Be}$ reaction at (a) $E_d = 21$ MeV, (b) $E_d = 40.9$ MeV, and (c) $E_d = 71$ MeV. Results for CDCC (hatched band), FAGS (solid), and FAGS1 (circles).

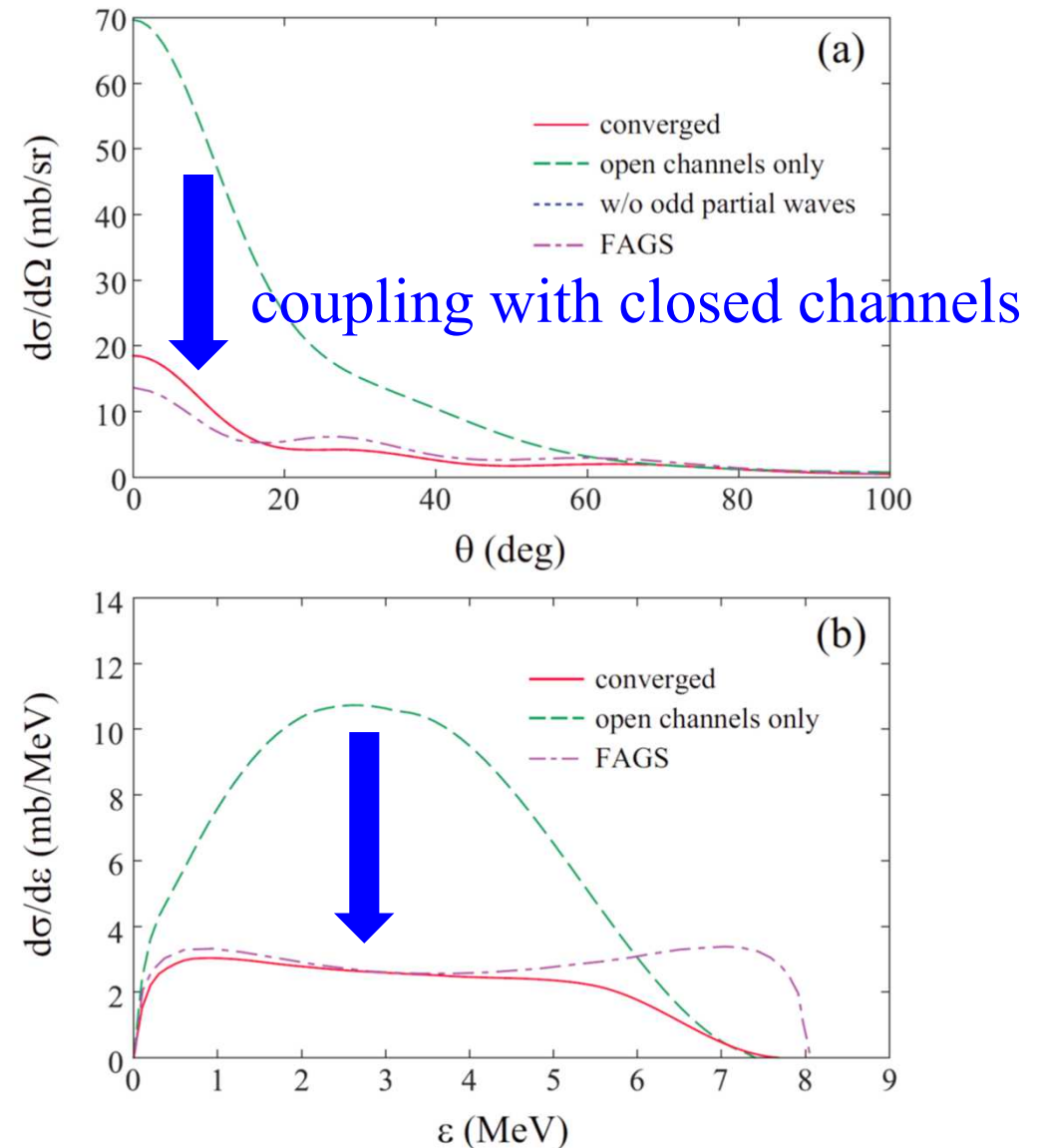
Applicability of CDCC to low energy BU process

N. J. Upadhyay et al., PRC 85, 054621 (2012).



CDCC severely overshoots the result of FAGS.

KO and K. Yoshida, PRC 94, 051603(R) (2016).



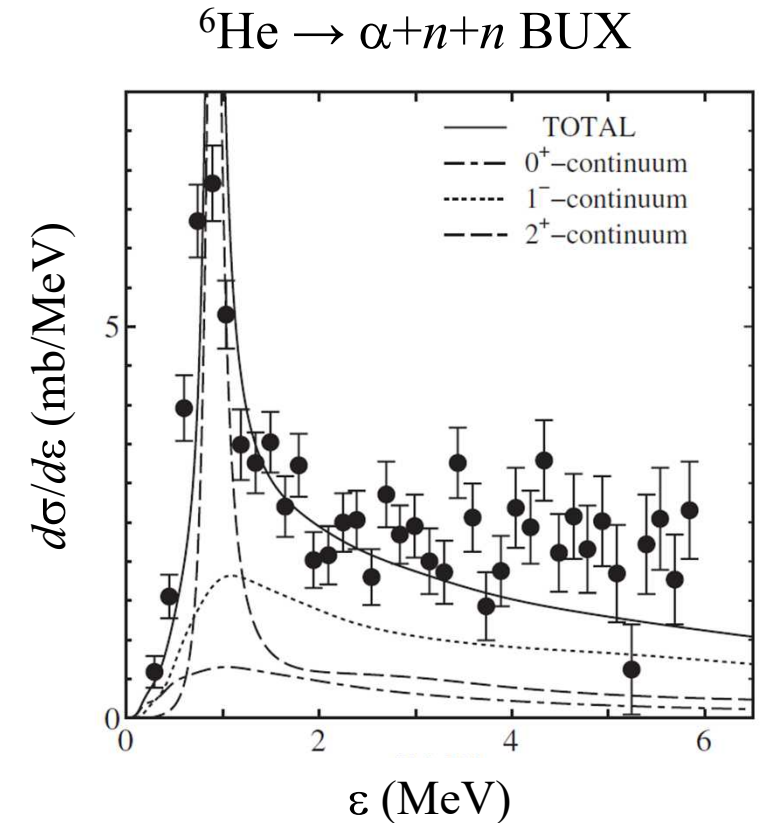
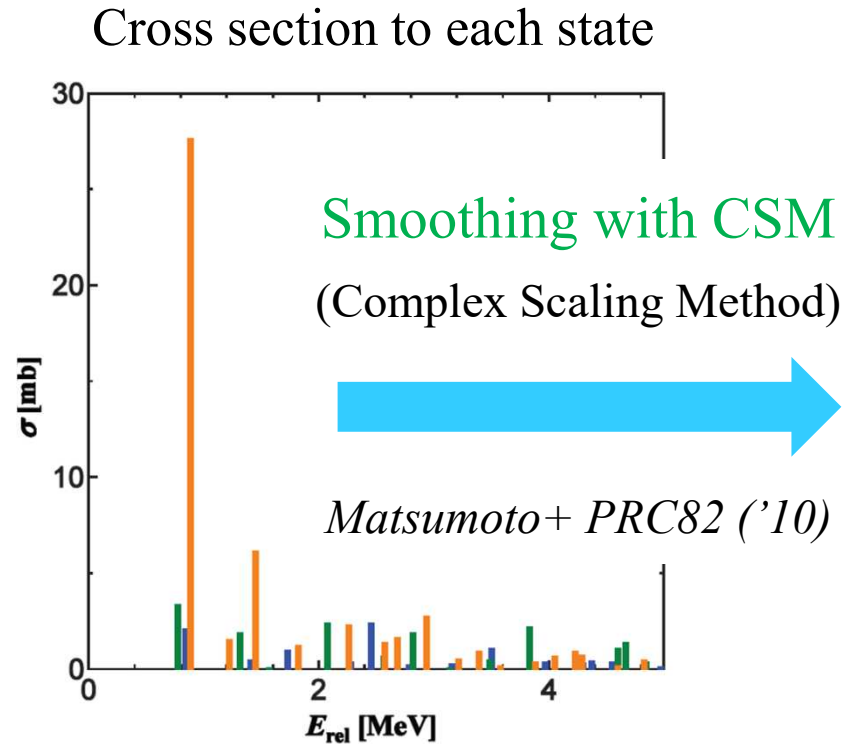
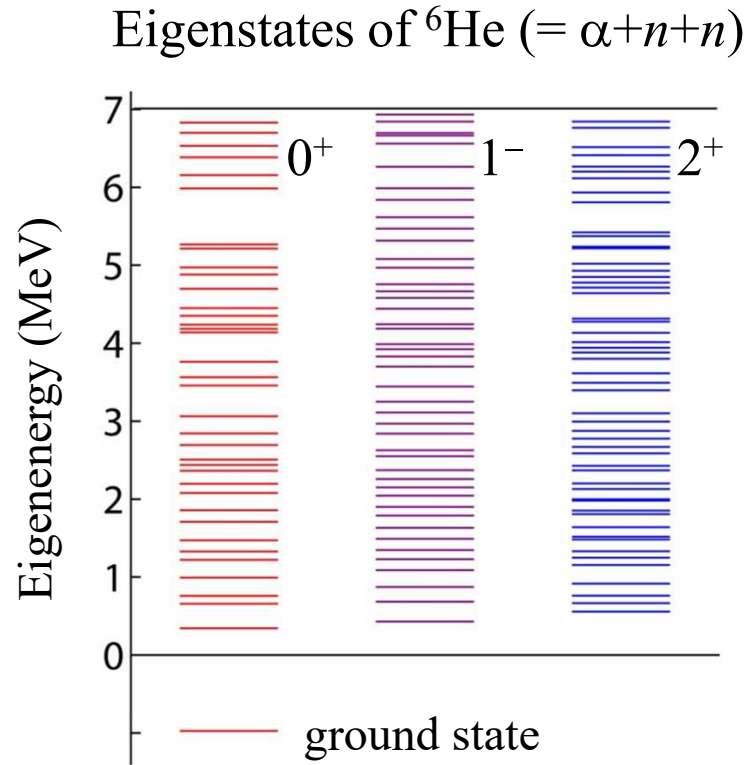
Two questions

✓. *Can we use a single Lippmann-Schwinger Eq.?*

2. *How to implement the continuum st. of many-body system?*

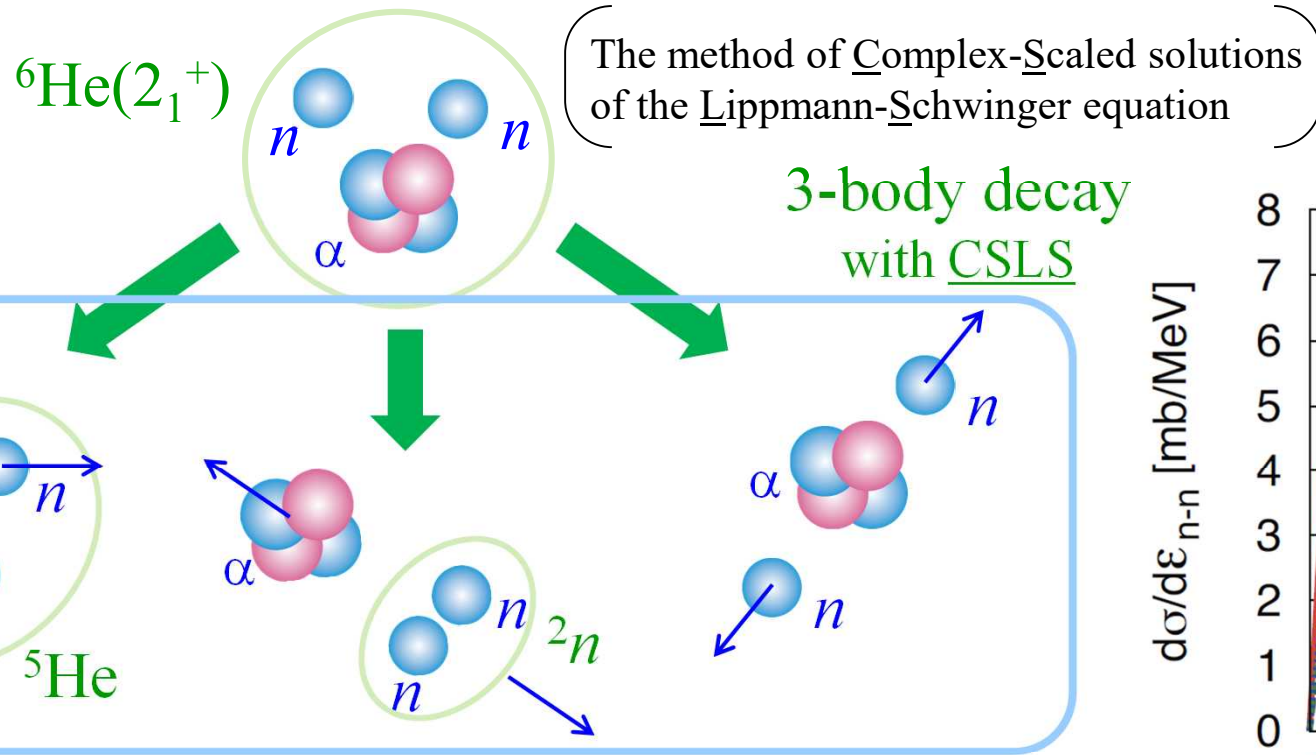
Gaussian Expansion Method (GEM) for ${}^6\text{He}$

Yahiro+, Prog. Theor. Exp. Phys. 2012, 01A206.

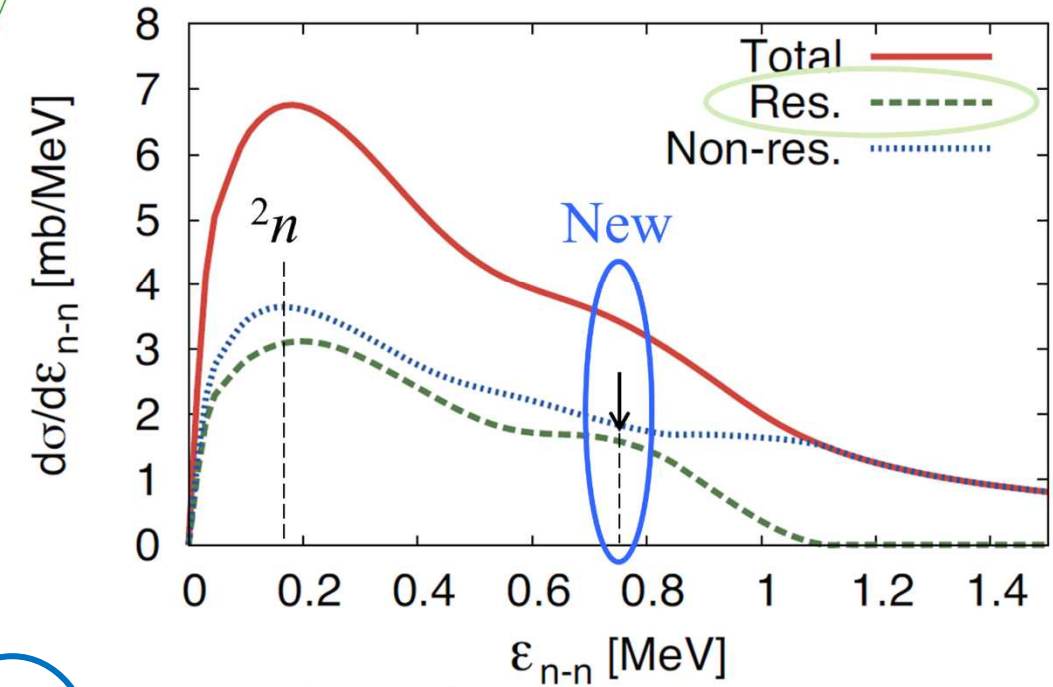


Hiyama+, Prog. Part. Nucl. Phys. 51, 223 (2003).

Specification of the final state



Kikuchi+ PTP122 ('09), PRC88 ('13)

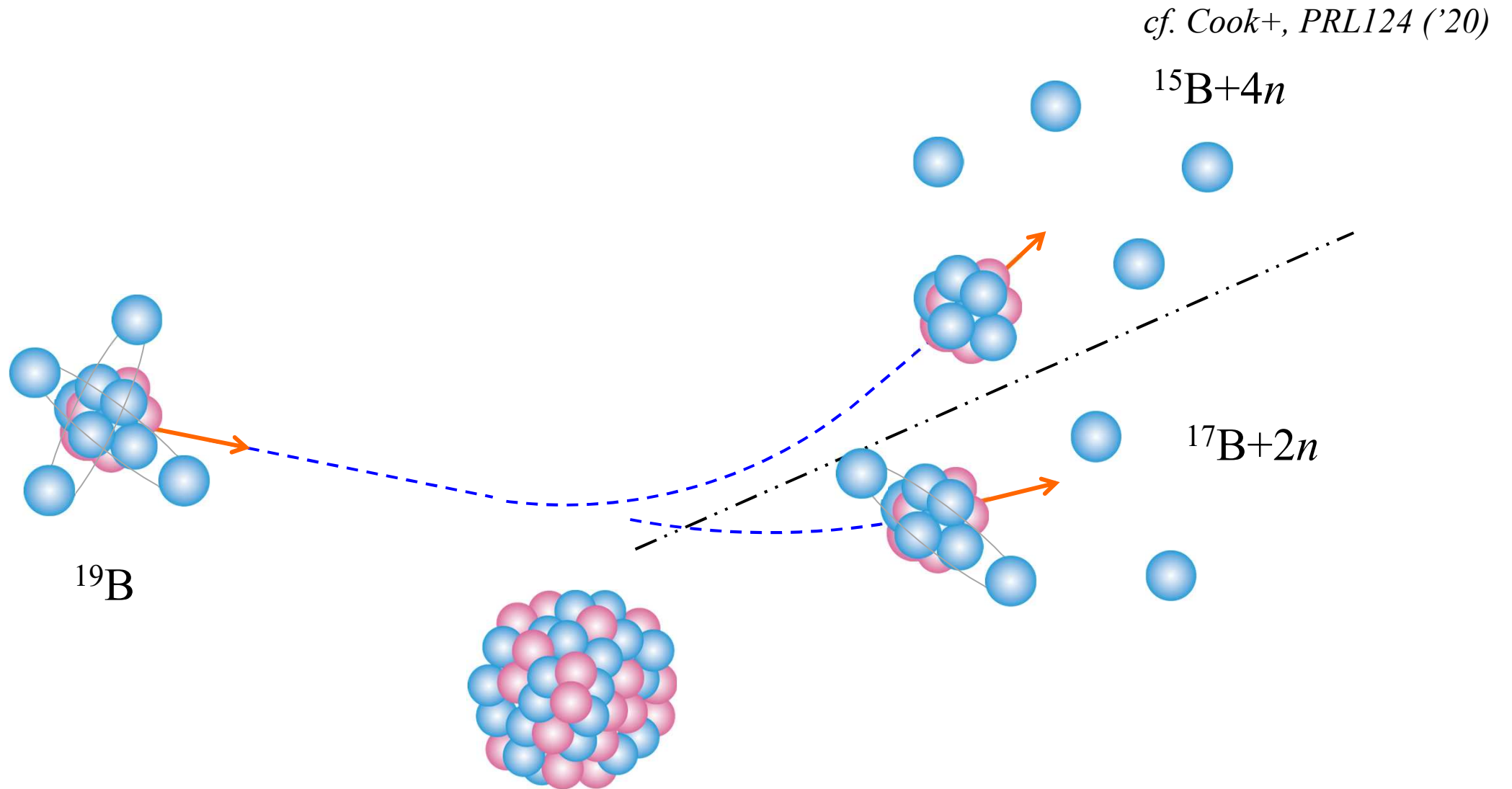


$$|\psi^{(+)}\rangle = |\phi_{\text{free}}\rangle + \sum_n U^{-1}(\theta) |\Phi_n^\theta\rangle \frac{1}{\varepsilon - \varepsilon_n^\theta} \langle \tilde{\Phi}_n^\theta | U(\theta) V | \phi_{\text{free}} \rangle$$

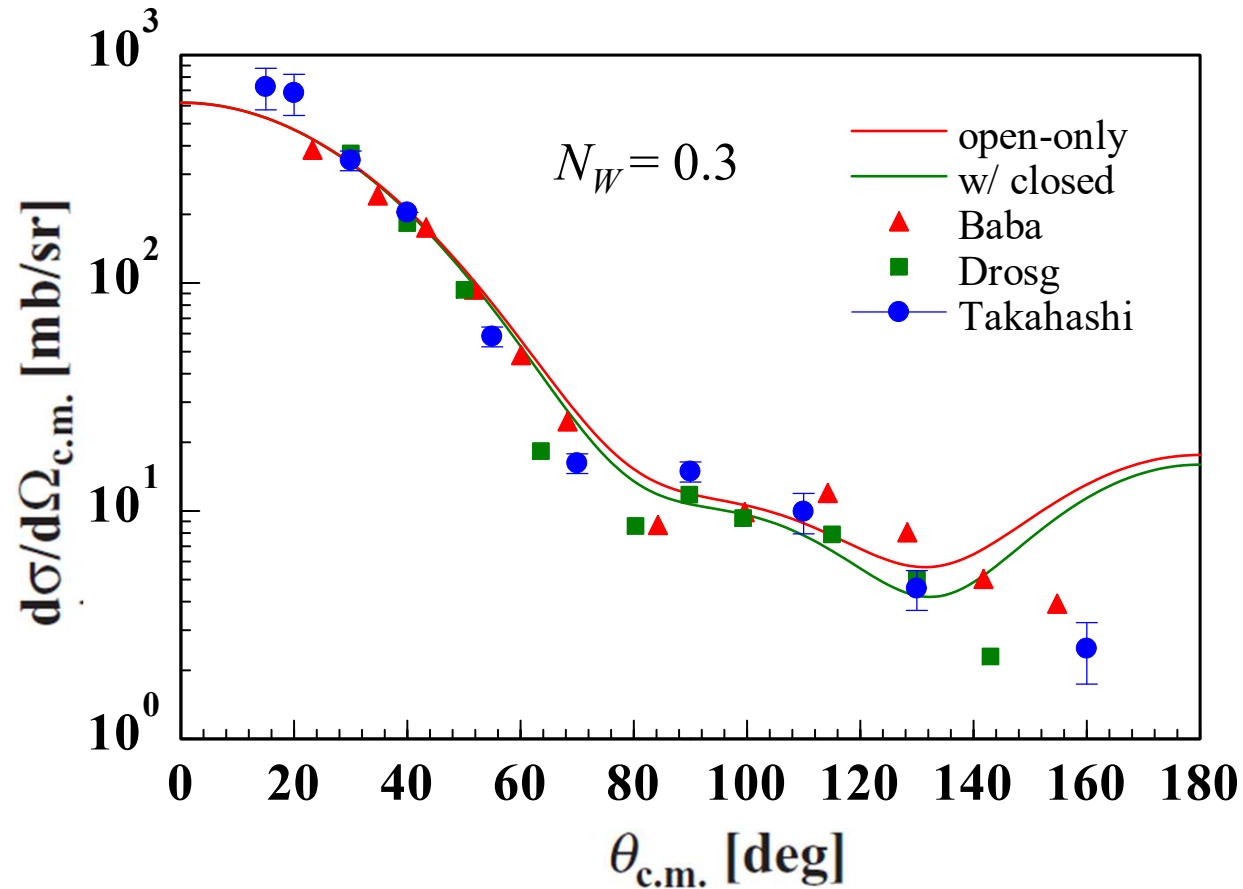
Eigenstate of a CS Hamiltonian

cf. A practical FS specification: Watanabe+ PRC103 ('21)

Possible future study



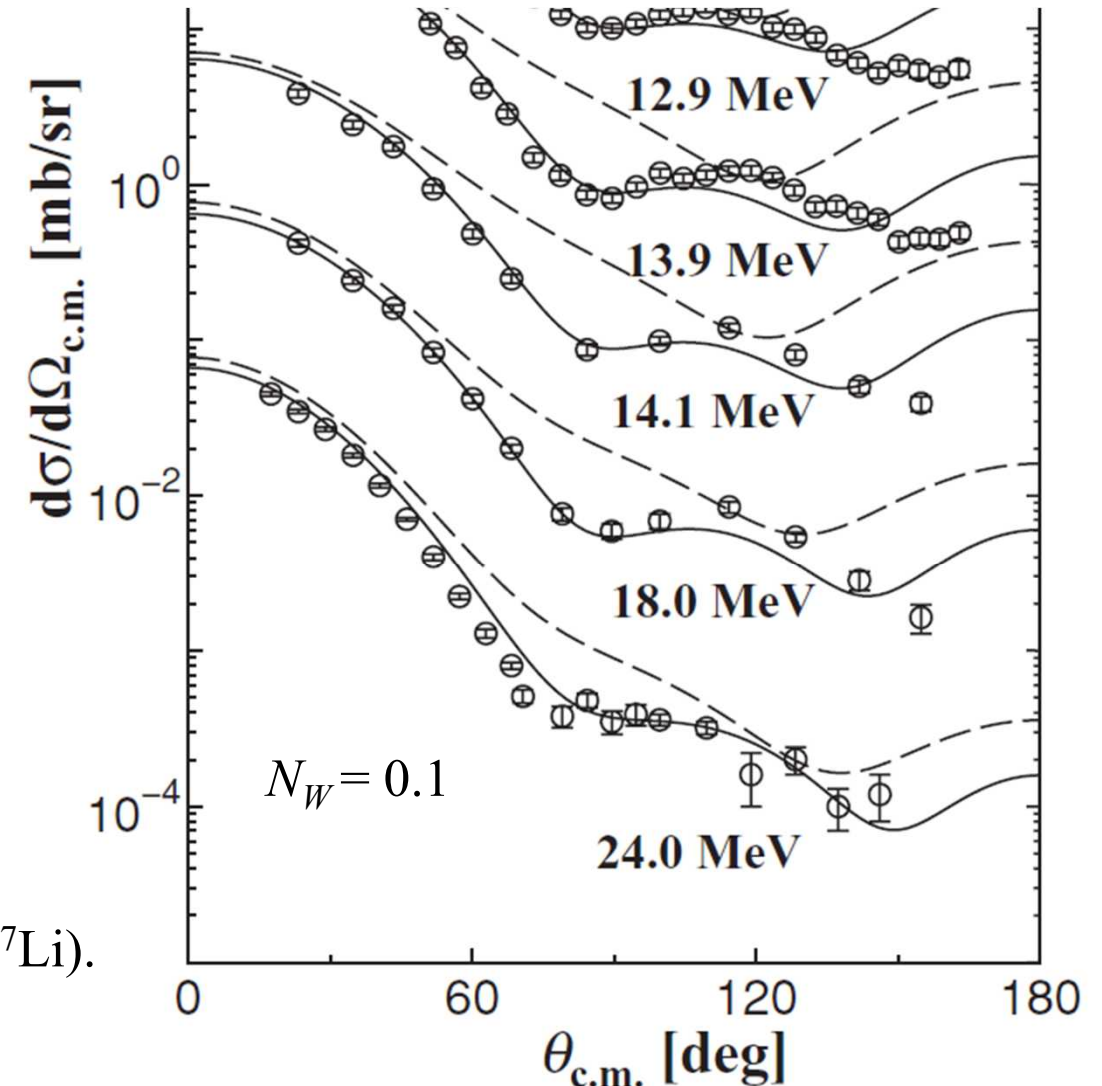
(My) 1st 4-body CDCC calc. for n-⁶Li elas. scat.



Physics case: t production around ITER

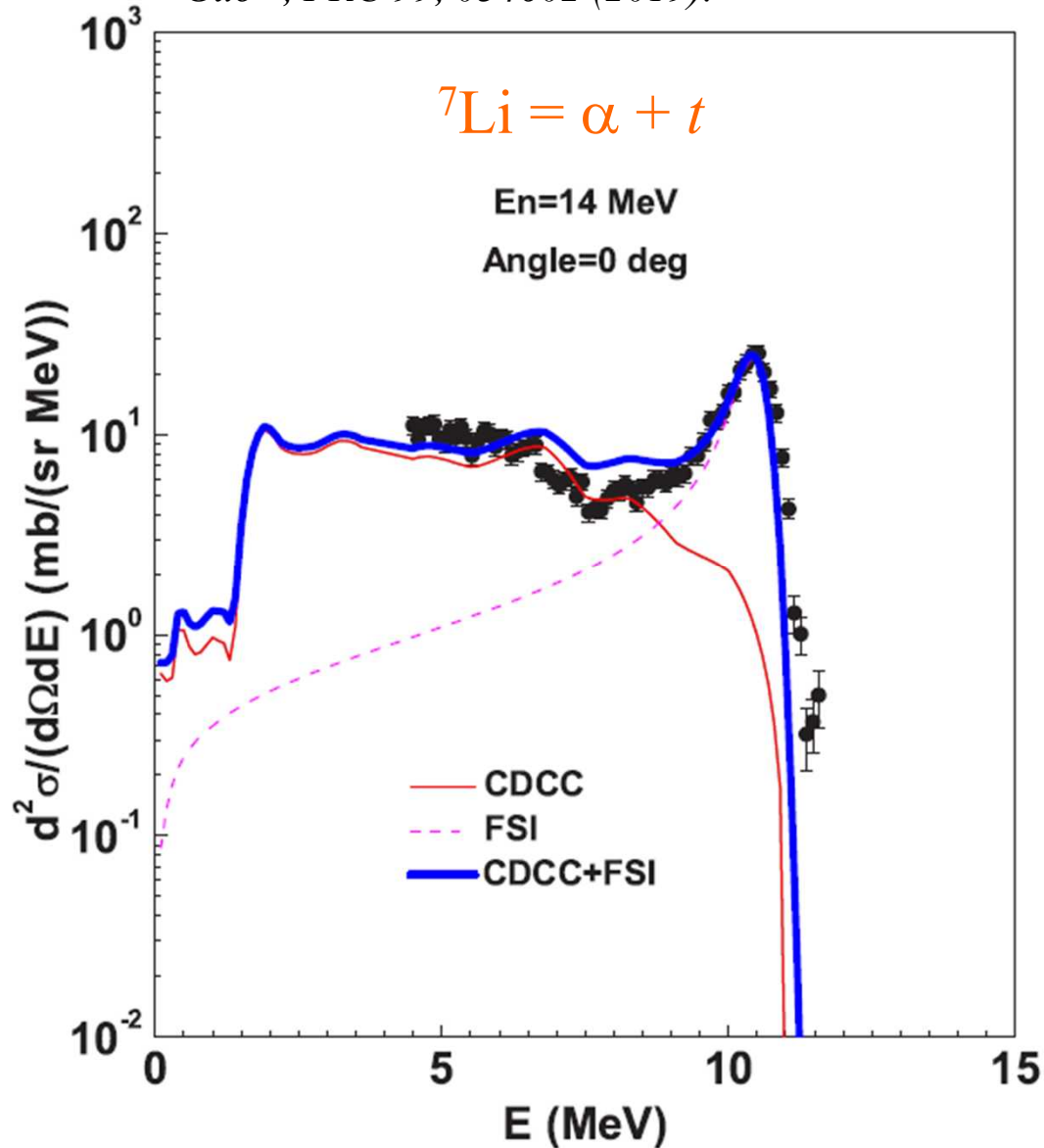
- ✓ In the preceding studies, $\alpha+d$ ($\alpha+t$) is assumed for ${}^6\text{Li}$ (${}^7\text{Li}$).
- ✓ A simplified g-matrix int. (JLM) is employed, with a normalization of its imaginary part.

Matsumoto+, PRC 83, 064611 (2011).



t production from n-⁷Li reaction

Guo+, PRC 99, 034602 (2019).



How does the result change when ⁷Li is described with $\alpha+p+n+n$?

Note:

The specification of the final channel with CSLS is necessary.

Summary

- ✓ I have introduced an MST-based microscopic approach to nuclear reactions.
 - ✓ NN eff. int. is a key ingredient; g-matrix approach works quite well.
 - ✓ One-body transition densities, ideally with 3NF effect, are expected to be provided by B1 and B2 collaborators.
- ✓ CDCC, a possible alternative to Faddeev-AGS, is the key reaction model.
 - ✓ Its theoretical foundation was reviewed.
 - ✓ Specification of the final channel with CSLS is crucial.
- ✓ Description of t-production from N-^{6,7}Li is the first physics case of B3.
 - ✓ GEM-CDCC calc., in collaboration with Prof. Hiyama, is getting ready for N-⁶Li.
 - ✓ A comparison using different W.Fns. will be performed in collaboration with Prof. W. Horiuchi (OMU).
 - ✓ After N-⁶Li studies, we will consider how to proceed with ⁷Li.