



# Systematics of reaction cross sections and geometrical parameters from double folding and single folding optical potentials

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#### Work Plan

- o Motivation
- o Basic equations of eikonal formulae
- o Numerical results
- o Conclusion

#### Motivation

o Breakup reaction is one of the main tools for the study of exotic nuclei.

o The optical potential model can be obtained microscopically through a folding model approach: single folding and double folding.

o The single-folded optical potential is more accurate than the double-folded optical potential.

#### In this work

A systematic comparison of calculated reaction cross sections on a  ${}^9Be$  target via a single folding versus a double folding optical potential.

❖ Comparison of the strong absorption radius parameter extracted from the S matrices for single folding and double folding results.

## Basic equations of eikonal formalism

The eikonal reaction cross section:

$$\sigma_R = 2\pi \int_0^\infty bdb \left( 1 - \left| S_{PT}(\mathbf{b}) \right|^2 \right) \tag{1}$$

Where

$$\left|S_{PT}(\mathbf{b})\right|^2 = e^{2\chi_I(b)} \tag{2}$$

■ The imaginary part of the eikonal phase shift:

$$\chi_I(\mathbf{b}) = \frac{1}{\hbar v} \int dz W^{PT}(\mathbf{b}, z) \tag{3}$$

## Basic equations of eikonal formalism

#### Single folding potential:

#### Double folding potential:

$$W_{s.f.}^{PT}(\mathbf{r}) = \int d\mathbf{b}_1 W^{nT} \left( \mathbf{b}_1 - \mathbf{b}, z \right) \int dz_1 \rho_P \left( \mathbf{b}_1, z_1 \right)$$
 (4)

 $W_{d.f.}^{PT}(\mathbf{r}) = -\frac{1}{2}\hbar v \sigma_{nn} \left[ d\mathbf{b}_1 \rho_T \left( \mathbf{b}_1 - \mathbf{b}, z \right) \right] dz_1 \rho_P \left( \mathbf{b}_1, z_1 \right)$  (5)

n+9Be phenomenological nucleon-target potential (AB)

• A. Bonaccorso, R.J. Charity, Phys. Rev. C 89 (2014) 024619.

Energy-dependent nucleon-nucleon (nn) cross section

**Densities: HFB code** 

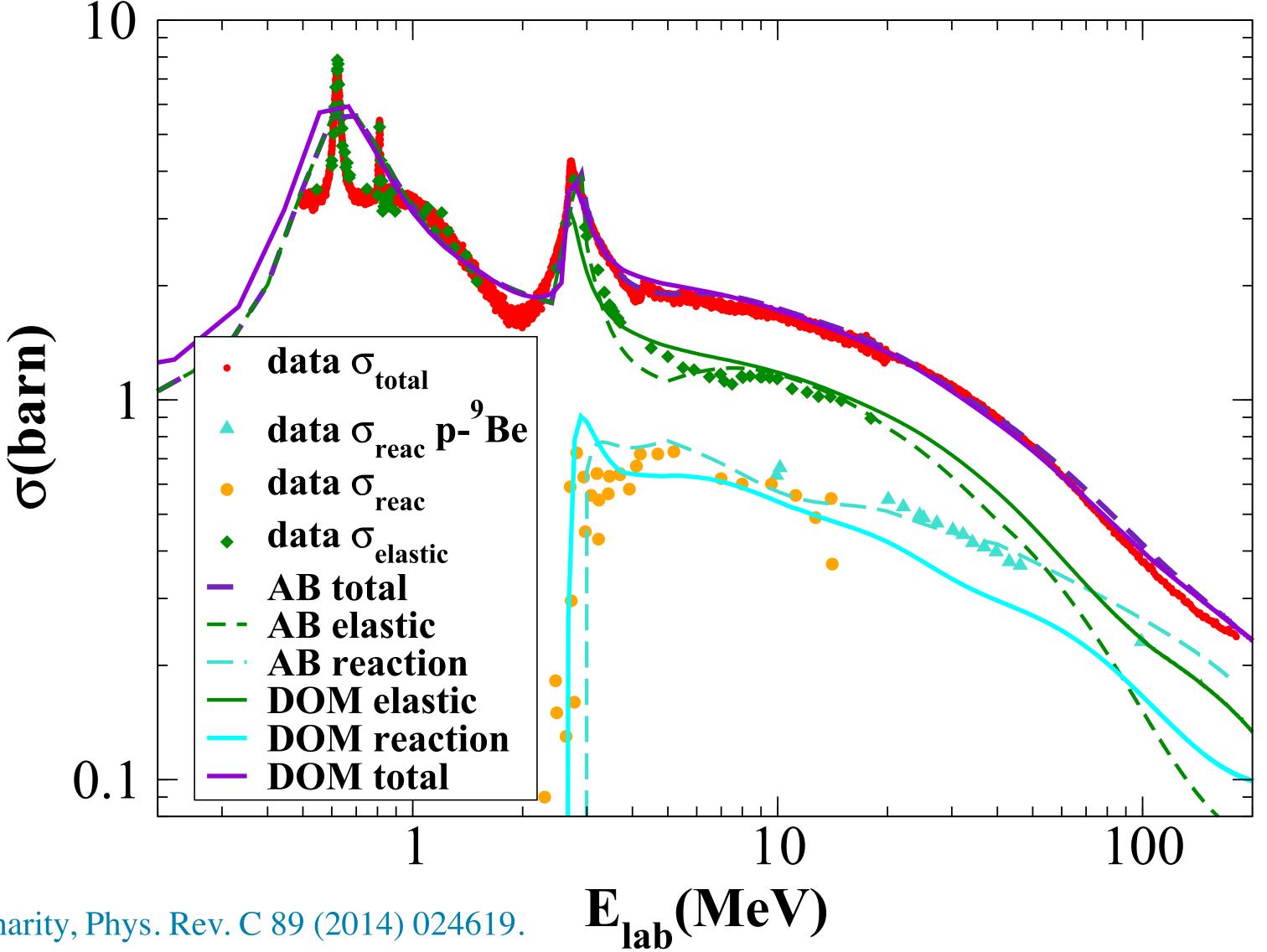
- C.A. Bertulani, C. De Conti, Phys. Rev. C 81 (2010) 064603.
- → Eq. (5) can be given the same structure as Eq. (4) by defining

$$W^{nT}(\mathbf{r}) = -\frac{1}{2}\hbar v \sigma_{nn} \rho_T(\mathbf{r})$$
 (6)

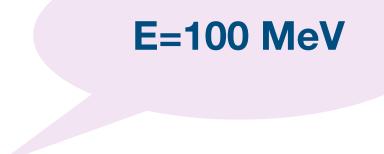
#### Justification

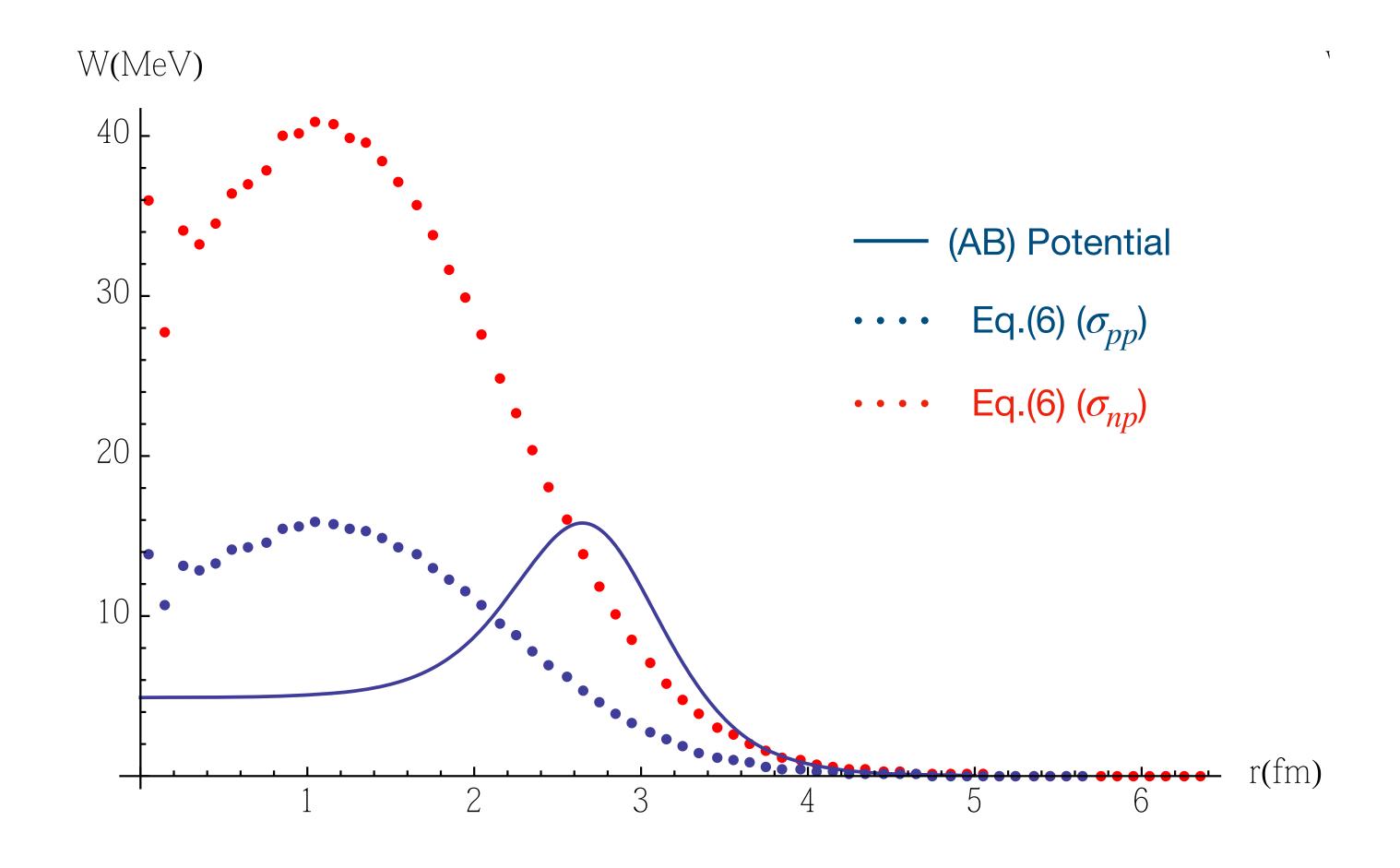
#### Reminder:

$$W_{s.f.}^{PT}(\mathbf{r}) = \int d\mathbf{b_1} W^{nT} \left( \mathbf{b_1} - \mathbf{b}, z \right) \int dz_1 \rho_P \left( \mathbf{b_1}, z_1 \right)$$
 (4)



#### Justification



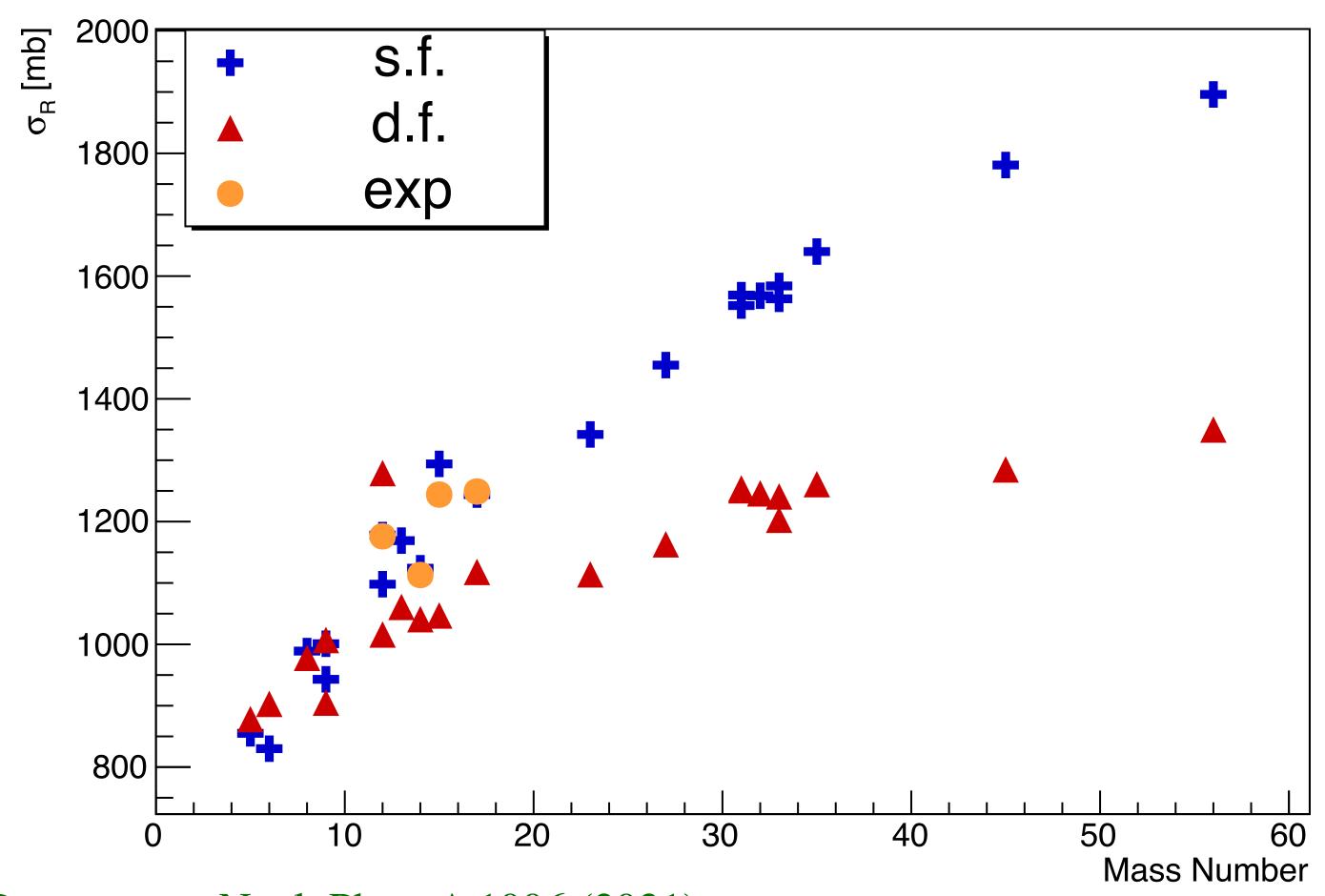


$$W^{nT}(\mathbf{r}) = -\frac{1}{2}\hbar v \sigma_{nn} \rho_T(\mathbf{r})$$
 (6)

• A. Bonaccorso, et al. Few-Body Syst (2016) 57:331-226.

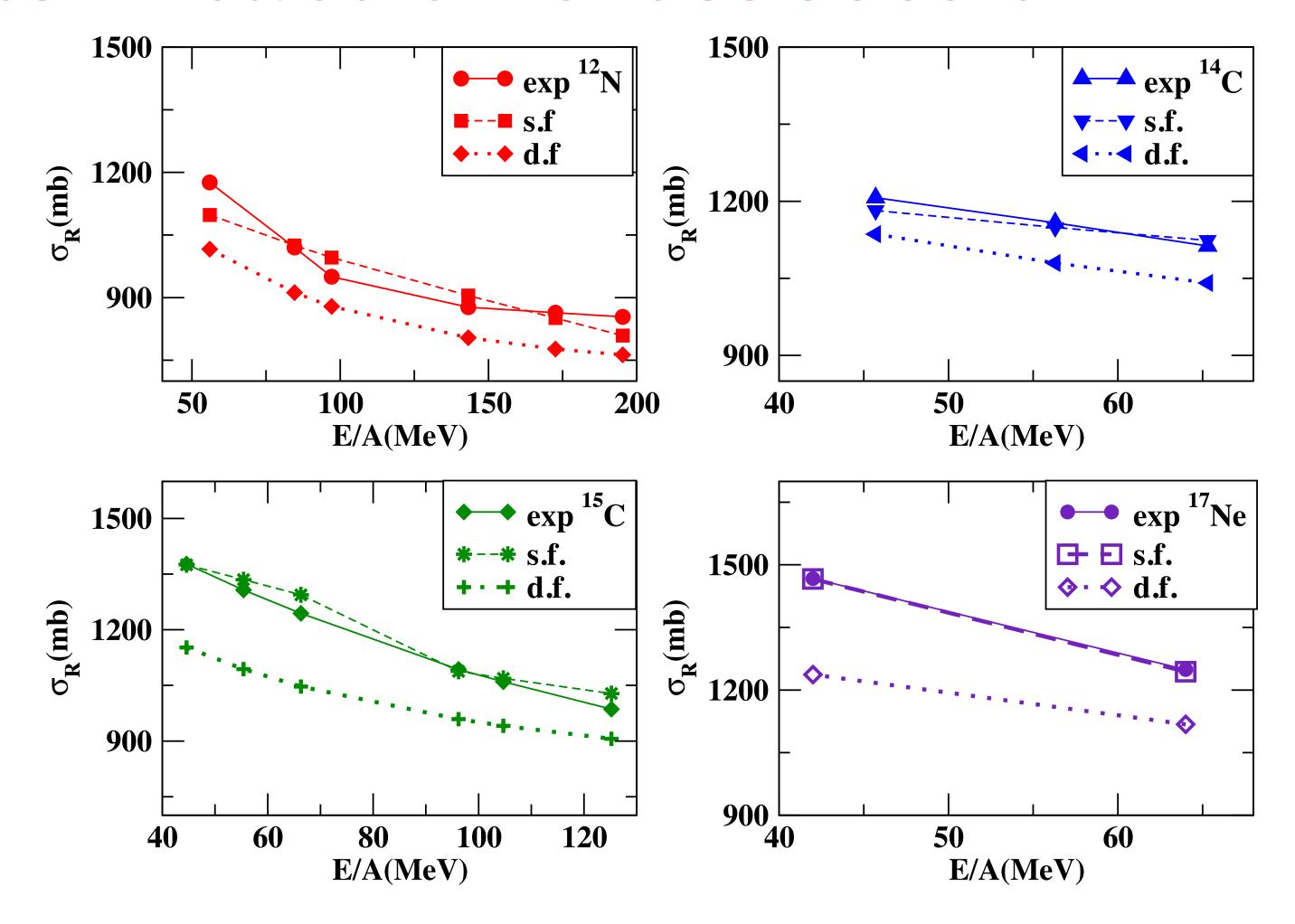
#### Results: Reaction cross section





I. Moumene and A. Bonaccorso, Nucl. Phys. A 1006 (2021)

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<sup>&</sup>lt;sup>12</sup>N: M. Fukuda, Y. Morita et al., JPS Conf. Proc. 6, 030103 (2015).

<sup>&</sup>lt;sup>15</sup>C: H. Du, M. Fukuda et al., Acta Physica Polonica B 48, 473 (2017). J. A 25 (s01) (2005) 221.

<sup>&</sup>lt;sup>17</sup>Ne: K. Tanaka, M. Fukuda, et al., Eur. Phys. J. A 25 (s01) (2005) 221.

<sup>&</sup>lt;sup>14</sup>*C*: M. Fukuda, private communication.

## Strong absorption radius

The strong absorption radius:

$$|S_{PT}(R_S)|^2 = \frac{1}{2}$$
 (7)

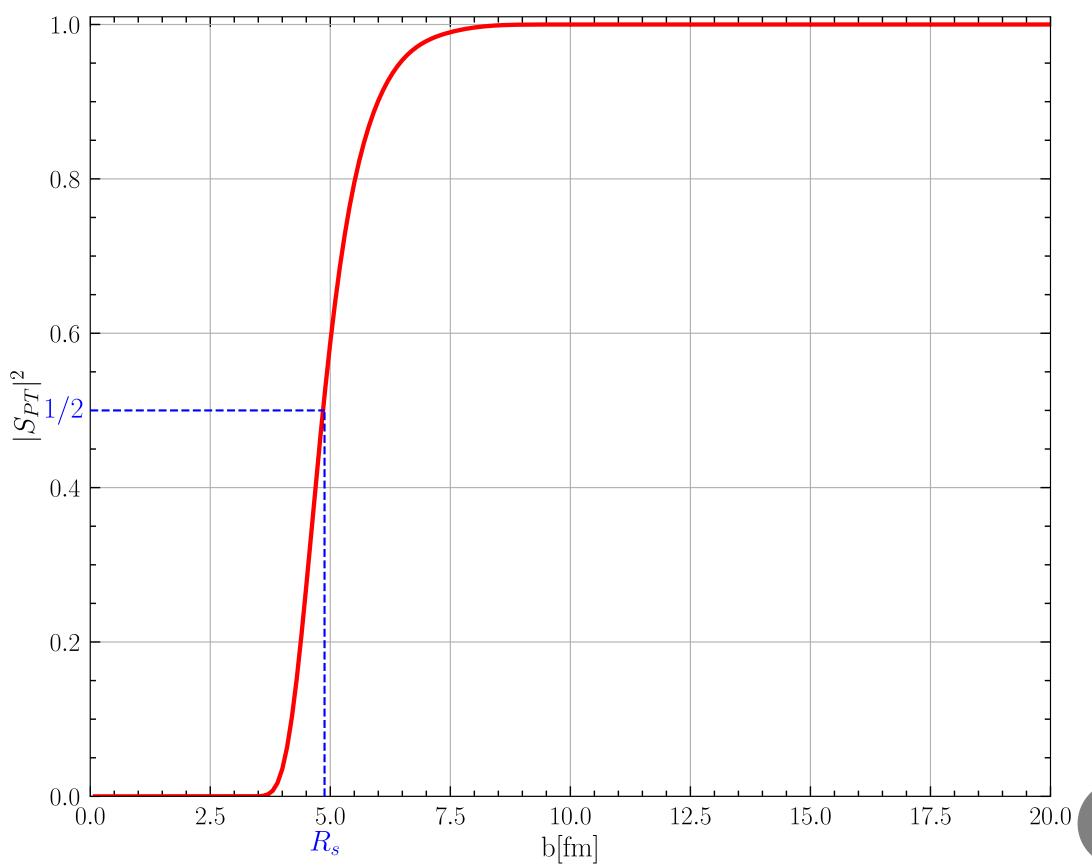
With

$$R_s = r_s \left( E_{inc} \right) \left( A_P^{1/3} + A_T^{1/3} \right)$$
 (8)

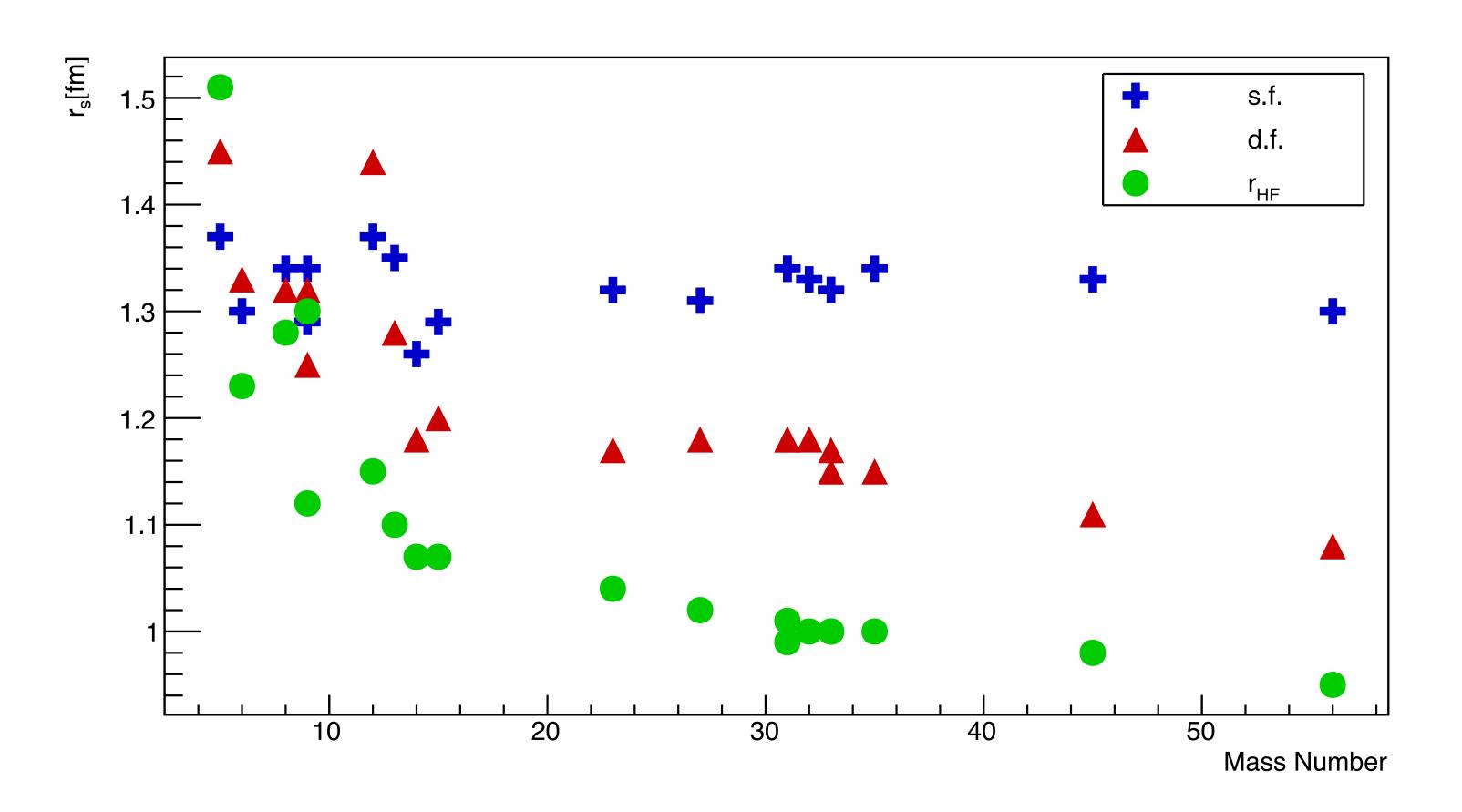
 $r_{s}$ : Determine the range of impact parameters for which surface reaction dominate the core-target interaction from regions in which the strophysical absorption regime applies.

#### **Reminder:**

$$\sigma_R = 2\pi \int_0^\infty bdb \left( 1 - \left| S_{PT}(\mathbf{b}) \right|^2 \right)$$
$$\left| S_{PT}(\mathbf{b}) \right|^2 = e^{2\chi_I(b)}$$

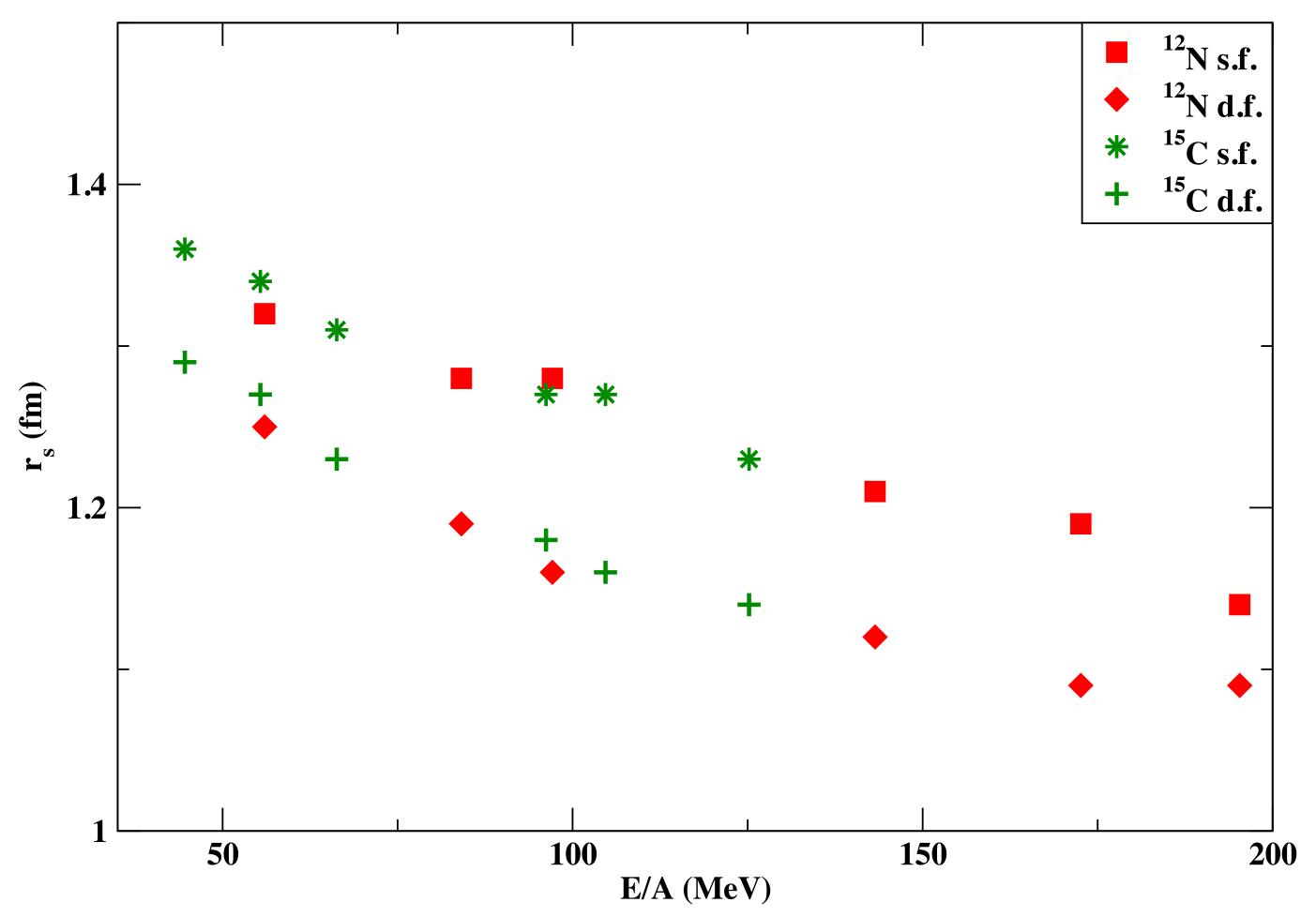


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#### Conclusion

- The single folding model is more reliable than the double folding model in describing the total reaction cross sections for several nuclei.
- ❖ The d.f. gives smaller reaction cross sections and in turn it is expected to produce larger breakup cross sections.
- The fact that the single folding model has provided very stable values of the strong absorption radius parameter,  $r_s=1.3-1.4$  fm, confirms the validity and potentiality of our s.f. approach for future studies with breakup and knockout reactions.
- Study the reaction cross section using densities from coupled cluster model (In progress)

## Thank you for your attention