# Uncertainty in nuclear-reaction calculations involving halo nuclei

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- Introduction
  - Halo nuclei
  - Reaction model
- Uncertainties in nuclear-reaction modelling
  - in halo-nucleus description
  - solving the problem with Halo-EFT
  - in optical potentials
  - looking for a solution from xEFT N-N interaction
- Summary

#### Halo nuclei

Exotic nuclear structures are found far from stability In particular halo nuclei with peculiar quantal structure:

- Light, n-rich nuclei
- Low  $S_n$  or  $S_{2n}$

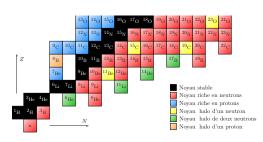
Exhibit large matter radius

due to strongly clusterised structure :

neutrons tunnel far from the core and form a halo

## One-neutron halo

<sup>11</sup>Be ≡ <sup>10</sup>Be + n <sup>15</sup>C ≡ <sup>14</sup>C + n Two-neutron halo <sup>6</sup>He ≡ <sup>4</sup>He + n + n <sup>11</sup>Li ≡ <sup>9</sup>Li + n + n



#### Reactions with halo nuclei

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Halo nuclei are fascinating objects but difficult to study [\tau_{1/2}(^{11}Be)=13 \text{ s}] \Rightarrow require indirect techniques, new probes, like reactions:
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Breakup ≡ dissociation of halo from core
by interaction with target
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Need good understanding of the reaction mechanism
(i.e. a good reaction model)
to know to what the probe is sensitive
(i.e. what nuclear-structure information it provides)
have reliable inputs for the model
(i.e. optical potentials to describe the interactions with target)
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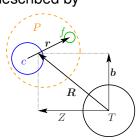
Here I look at the uncertainty related to the last two points For the former, a precise reaction models is coupled with Halo EFT For the latter, I do what I can...

#### Reaction model

Projectile (P) modelled as a two-body quantum system : core (c)+loosely bound nucleon (f) described by

$$H_0 = T_r + V_{cf}(\mathbf{r})$$

 $V_{cf}$  effective interaction describes the projectile structure with ground state  $\Phi_0$ 



Target T assumed structureless

Interaction with target simulated by optical potentials  $U_{cT}$  and  $U_{fT}$   $\Rightarrow$  breakup reduces to three-body scattering problem :

$$\left[T_R + H_0 + U_{cT} + U_{fT}\right] \Psi(\mathbf{r}, \mathbf{R}) = E_T \Psi(\mathbf{r}, \mathbf{R})$$

with initial condition  $\Psi(\mathbf{r},\mathbf{R}) \underset{Z \to -\infty}{\longrightarrow} e^{iKZ} \Phi_0(\mathbf{r})$ 

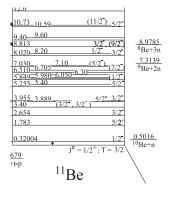
We use the Dynamical Eikonal Approximation (DEA)

[Baye, P. C., Goldstein, PRL 95, 082502 (2005)]

## Sources of uncertainty

- $\bullet$   $V_{cf}$ 
  - Describes the projectile structure
  - Fitted to external data (binding energy...)
  - Not an observable in the sense of J. Oakley but it still has meaning outside the reaction code
- $\bullet$   $U_{cT}$  and  $U_{fT}$ 
  - Optical potentials simulating the c-T and f-T interaction
  - Found in the literature
  - Not observable either
  - Very uncertain
- Numerical parameters
  - Control inputs
  - Value chosen to ensure convergence of numerical scheme

## $^{11}$ Be $\equiv ^{10}$ Be $\otimes$ n

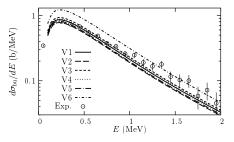


- $\frac{1}{2}^+$  ground state :  $\epsilon_{\frac{1}{2}^+} = -0.503$  MeV In our model, seen as  $1s_{\frac{1}{2}}$  neutron bound to  $^{10}$ Be( $0^+$ )
- $\frac{1}{2}$  bound excited state :  $\epsilon_{\frac{1}{2}} = -0.184$  MeV In our model, seen as  $0p_{\frac{1}{2}}$  neutron bound to  $^{10}$ Be( $0^+$ )

## Uncertainty in halo-nucleus description

 $^{11}$ Be+Pb  $\rightarrow$   $^{10}$ Be+n+Pb @ 69AMeV

We have used phenomenological  $V_{cf}$  with different geometries Fitted to reproduce  $^{11}$ Be low-energy spectrum



Exp. : [Fukuda *et al.* PRC 70, 054606 (2004)] Th. : [P.C., Nunes, PRC 73, 014615 (2006)]

 $\Rightarrow$  significant variation in  $d\sigma_{\rm bu}/dE$ 

Can we control this uncertainty?

## Halo EFT description of <sup>11</sup>Be

Use Halo EFT : clear separation of scales

⇒ provides an expansion parameter (small scale / large scale) upon which the low-energy behaviour is expanded (cf. S. Wesolowski)

[H.-W. Hammer, C. Ji, D. R. Phillips JPG 44, 103002 (2017)]

Use narrow Gaussian potentials :  $V_{cf}(r) = V_0 \ e^{-\frac{r^2}{2\sigma^2}} + V_2 \ r^2 e^{-\frac{r^2}{2\sigma^2}}$ 

Fit  $V_0$  and  $V_2$  to reproduce known experimental values (and *ab initio* predictions [A. Calci *et al.* PRL 117, 242501 (2016)])

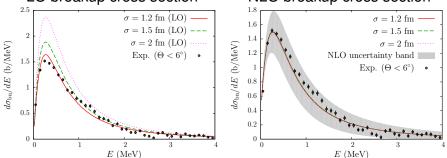
 $\sigma$  = 1.2, 1.5 or 2 fm is a parameter used to evaluate the sensitivity of the calculations to short-range physics

@LO:  $V_2 = 0 \& V_0$  adjusted to fit  $\epsilon_{1/2^+}$  in s wave  $(V_{cf} = 0 \ \forall l > 0)$ 

@NLO :  $V_2$  &  $V_0$  adjusted in s and p waves to fit  $\epsilon_{1/2^+}$  &  $\epsilon_{1/2^-}$ , ANCs and  $\delta_{p3/2}$  (ab initio predictions)  $(V_{cf}=0 \ \forall l>1)$ 

[PC, D. R. Phillips, H.-W. Hammer PRC 98, 034610 (2018)]

## LO & NLO analyses of <sup>11</sup>Be+Pb→<sup>10</sup>Be+n+Pb @ 69AMeV LO breakup cross section NLO breakup cross section



Exp. : [Fukuda et al. PRC 70, 054606 (2004)]

Th.: [PC, D. R. Phillips, H.-W. Hammer PRC 98, 034610 (2018)]

- LO: Large differences due to differences in ANC
- NLO : All calculations identical (ANCs &  $\delta_p$  fitted) Excellent agreement with data
- ⇒ Halo EFT shows which degrees of freedom matter

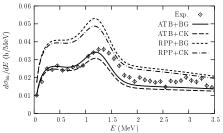
i.e. which structure observables are probed in reaction provides an estimate of the uncertainty due to truncation

## Uncertainty in optical potentials

The reaction model requires optical potentials  $U_{cT}$  and  $U_{fT}$ 

$$\left[T_R + H_0 + U_{cT} + U_{fT}\right] \Psi(\mathbf{r}, \mathbf{R}) = E_T \Psi(\mathbf{r}, \mathbf{R})$$

How can one estimate the uncertainty related to these interactions? Example :  $^{11}\text{Be+C} \rightarrow ^{10}\text{Be+n+C} \oslash 67A\text{MeV}$ 



Exp. : [Fukuda *et al.* PRC 70, 054606 (2004)] Th. : [P. C., Goldstein, Baye PRC 70, 064605 (2004)]

⇒ that uncertainty can be huge

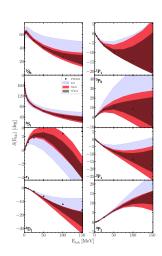
Problem: reaction calculations are expensive ⇒ use emulator?

## Idea: build our own optical potential

Idea : using a double-folding procedure with accurate NN interactions from  $\chi {\sf EFT}$ 

Gezerlis *et al.* have developed local NN interactions up to N<sup>2</sup>LO [PRL 111, 032501 (2013), PRC 90, 054323 (2014)]

Based on this formalism, build a double-folding potential  $V_F$ 

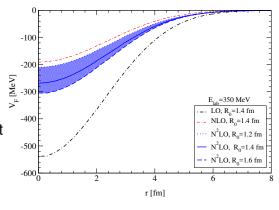


## <sup>16</sup>O-<sup>16</sup>O potential

Potential built

- at different orders
- for different cutoffs

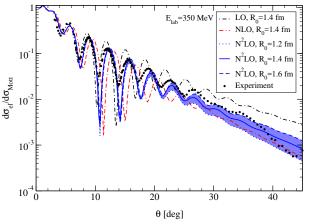
Calculations by V. Durant PLB 782, 668 (2018)



The imaginary part is assumed proportional to  $V_F$ 

$$U_F(r) = (1 + N_W i) V_F(r)$$
 with  $N_W = 0.6 - 0.8$ 

## <sup>16</sup>O-<sup>16</sup>O elastic scattering @350 MeV



- Fair agreement with experiment (no fitting parameter)
- Systematic order-by-order behaviour
- Small uncertainty related to the cutoff
- Larger uncertainty to N<sub>W</sub>

## Summary and prospect

- Halo nuclei studied mostly through reactions
- Mechanism of reactions with halo nuclei understood
- Halo EFT
  - emulates ab initio structure calculations in reaction codes
  - shows which degrees of freedom matter
  - provides an estimate of the uncertainty
- Optical potentials have huge impacts on reaction calculations
  - Could we estimate this uncertainty using an emulator?
  - Can be built by double-folding from  $\chi$ EFT NN interactions

#### Thanks...

to you for your attention and to my collaborators

**Daniel Phillips** 



Victoria Durant Lukas Huth Hans-Werner Hammer Achim Schwenk



Daniel Baye Gerald Goldstein

