

# Breakup Project

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## Reaction Project

During the exercise sessions, you'll be computing and analysing **breakup cross sections** of **halo nuclei**

- You'll be using the Fortran code **Chaconne.f** (see Indico)  
That code implements the Coulomb Corrected Eikonal (**CCE**)
  - ▶ runs fast (a few minutes at most)
  - ▶ accounts for the  $P$ - $T$  interaction at all orders
  - ▶ includes a 1st order correction of the Coulomb interaction
- You pick one (or two, or all. . .) of the reactions
  - ▶  $^{11}\text{Be} + \text{Pb} \rightarrow ^{10}\text{Be} + n + \text{Pb} @69\text{A MeV}$   
[Fukuda *et al.* PRC 70, 054606 (2004)]
  - ▶  $^{11}\text{Be} + \text{C} \rightarrow ^{10}\text{Be} + n + \text{C} @67\text{A MeV}$   
[Fukuda *et al.* PRC 70, 054606 (2004)]
  - ▶  $^{15}\text{C} + \text{Pb} \rightarrow ^{15}\text{C} + n + \text{Pb} @68\text{A MeV}$   
[Nakamura *et al.* PRC 79, 035805 (2009)]
  - ▶  $^{19}\text{C} + \text{Pb} \rightarrow ^{19}\text{C} + n + \text{Pb} @67\text{A MeV}$   
[Nakamura *et al.* PRL 83, 1112 (1999)]

## Goal Project

- Study the reaction :
  - ▶ develop a  $V_{cf}$  interaction (within Halo EFT)  
(use the code `Boscov.f` to fit the interaction, see Indico)
  - ▶ find suitable **optical potentials**  $V_{cT}$  and  $V_{fT}$
  - ▶ check the **convergence**
  - ▶ compare to existing **data** (available on Indico)  
There are energy and angular distributions  
Don't forget to account for the experimental resolution
  - ▶ analyse the agreement/differences with experiment
- Work in **groups of 4**  
(make sure that one of you has a computer to run the code)
- Friday morning, present the results of your study to the others
- This afternoon session is to decide on the system and set  $V_{cf}$

## Resources on Indico

- Codes **Boscov.f** (structure) and **Chaconne.f** (reaction) with short **user's manuals** and examples of input files (\*.dat files) and output files (\*.dep and \*.sdE files)
- Experimental data (\*.dat and \*.rtf files)
  - ▶ projectile and target are self-explanatory
  - ▶ `ere1_*.*` are energy distributions ( $d\sigma_{bu}/dE$ ) obtained after integration over angular range
  - ▶ `angle_*.*` are angular distributions ( $d\sigma_{bu}/d\Omega$ ) obtained after integration over a definite energy range

Details about the beam energy, experimental resolution etc. can be found in the original articles, which are provided in that same folder.