

## From ab initio structure predictions to reaction calculations via effective field theory

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Halo nuclei exhibit an uncommon nuclear structure with a larger matter radius compared to stable nuclei [1]. This large size is qualitatively understood as due to the loose binding of one or two valence neutrons, which have then a high probability of presence at a large distance from the other nucleons. They thus form a sort of halo around the compact core of the nucleus. The best known examples are  $^{11}\text{Be}$ , with a one-neutron halo, and  $^{11}\text{Li}$ , with a two-neutron halo. Due to their short lifetime, these nuclei are mostly studied through reactions like breakup [2]. In order to extract valuable structure information from measured cross sections, a precise model of the reaction coupled to a reliable description of the projectile is needed. Many such models have been developed (see Ref. [3] for a recent review). However, they mostly rely on a simple two- or three-body description of the nucleus. Recently, some of these exotic nuclei have become accessible to ab initio calculations [4]. Unfortunately, such  $A$ -body descriptions are too computationally demanding to be directly included within existing reaction models. In the present work, we use the outputs of an ab initio calculations of  $^{11}\text{Be}$  as inputs to the description of that nucleus within a reliable breakup model [5]. That description is inspired by an effective field theory treatment of  $^{11}\text{Be}$  [6] (see Ref. [7] for a recent review). Our calculations of the breakup of  $^{11}\text{Be}$  on Pb and C at about 70A MeV are in very good agreement with experimental measurements [2]. These excellent results prove the feasibility of incorporating results from ab initio calculations in reaction theory in this way. More importantly, they confirm the results for important aspects of  $^{11}\text{Be}$  obtained by the calculations of Calci et al. [4], hence improving our understanding of the nuclear structure far from stability.

### References

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### Summary

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