Halo effective field theory analysis of one-neutron knockout reactions of ¹¹Be and ¹⁵C

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Background: One-nucleon knockout reactions provide insightful information on the single-particle structure of nuclei. When applied to one-neutron halo nuclei, they are purely peripheral, suggesting that they could be properly modeled by describing the projectile within a halo effective field theory (halo-EFT).

Purpose: We reanalyze the one-neutron knockout measurements of ¹¹Be and ¹⁵C—both one-neutron halo nuclei—on beryllium at about 60 MeV/nucleon. We consider halo-EFT descriptions of these nuclei which already provide excellent agreement with breakup and transfer data.

Method: We include a halo-EFT description of the projectile within an eikonal-based model of the reaction and compare its outcome to existing data.

Results: Excellent agreement with experiment is found for both nuclei. The asymptotic normalization coefficients inferred from this comparison confirm predictions from *ab initio* nuclear-structure calculations and values deduced from transfer data.

Conclusions: Halo-EFT can be reliably used to analyze one-neutron knockout reactions measured for halo nuclei and test predictions from state-of-the-art nuclear structure models on these experimental data.

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I. INTRODUCTION

Halo nuclei exhibit an unusually large matter radius compared to their isobars. These very neutron-rich nuclei are found far from stability, close to or at the neutron dripline. There, the binding of one or two valence nucleons is so loose that they can tunnel far into the classically forbidden region to form an extended diffuse halo around the compact core of the nucleus [1]. Example of halo nuclei are ¹¹Be and ¹⁵C seen as a one-neutron halo bound to a ¹⁰Be or ¹⁴C core, and ⁶He and ¹¹Li with two neutrons in their halo. Because of their peculiar structure, halo nuclei have been the subject of many studies since their discovery in the mid-1980s [2,3].

Due to their short lifetime, halo nuclei are often studied through reactions such as breakup [4–7], transfer [8–11], and knockout [12–15]. In the first reaction, the halo dissociates from the core during its interaction with the target. The breakup hence reveals the internal structure of the nucleus. Thanks to the fragile nature of the projectile, the cross sections are high. In transfer reactions, the halo state is populated through a (d, p) transfer reaction often measured in inverse kinematics. Knockout, which is the focus of the present study, corresponds to the removal of the halo neutron from the nucleus. This reaction therefore contains both the aforementioned breakup—often called *diffractive breakup*—and the absorption of the valence neutron by the target, known as *stripping*. Because it does not require the detection of a neutron in coincidence with the core, the knockout measurement exhibits higher statistics than breakup. For that reason it is favored over the latter at low beam intensity [15,16].

The typical observable measured in knockout reactions is the momentum distribution of the core after the collision [12-15]. When measured at high-enough energy, the reaction is sudden and that observable retains the memory of the momentum distribution the core had within the nucleus. Because of the large spatial extension of the core-halo wave function and following Heisenberg's uncertainty principle, that distribution is narrow, hence providing a stringent probe of the halo structure [12-15].

Because of their high statistics, the use of knockout reactions has been extended to study the single-particle structure of more deeply-bound exotic nuclei [15,16]. Experimental data are usually compared to eikonal calculations of the reaction which include shell-model predictions for the spectroscopic factor of the projectile overlap wave function. In such analyses, the experimental cross section is usually smaller than what theory predicts [15–19]. The quenching factor R_S , i.e., the ratio between experiment and theory, is therefore interpreted as quantifying the deviations from shell-model predictions caused by missing correlations in the truncated model space used in structure calculations. Surprisingly, R_S decreases with the binding energy of the knocked-out nucleon: $R_S \approx 1$ for the removal of weaklybound nucleons, as in halo nuclei and $R_S \approx 0.3$ in the

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knockout of strongly-bound nucleons [17–19]. Since this feature is not observed in the analysis of transfer or quasi-free scattering reactions [16,20–22], it remains unclear if it is due to a bias in the interpretation of the knockout data or more profound differences between these reaction mechanisms and how they should be analyzed and compared to one another.

In a previous analysis, we have studied the sensitivity of the knockout cross sections of one-neutron halo nuclei to the description of their structure [23]. We have shown that these reactions are peripheral, in the sense that they probe only the tail of the core-halo wave function. Moreover, they are quite insensitive to the description of the core-neutron continuum. Accordingly, to properly reproduce experimental data, the model of the reaction should include a description of the projectile that reproduces its binding energy, the correct asymptotic normalization coefficient (ANC), and excited bound states and their ANC. For this goal, the halo effective field theory (halo-EFT [24,25], see Ref. [26] for a recent review) at next-to-leading order (NLO) includes all the relevant structure information. Interestingly, there already exist such halo-EFT descriptions for ¹¹Be [27] and ¹⁵C [28]. They are constrained with experimental one-neutron separation energies and the ANCs of the bound states predicted by ab initio calculations [29,30] or inferred from analyses of transfer reactions [28,31]. These descriptions successfully reproduce both breakup [27,28,32] and transfer [28,31] data. In this work, we use these halo-EFT descriptions to reanalyze knockout parallel-momentum distributions and integrated cross sections for ¹¹Be and ¹⁵C on a ⁹Be target at, respectively, 60 MeV/nucleon and 54 MeV/nucleon [13,14].

We briefly present in Sec. II the reaction model, including the halo-EFT descriptions of ¹¹Be and ¹⁵C and our choices of optical potentials. In Sec. III, we compare our theoretical predictions with experimental data and investigate the uncertainty associated with the choice of optical potentials. We conclude in Sec. IV.

II. FEW-BODY MODEL OF KNOCKOUT

A. Reaction model

To model the one-neutron knockout of halo nuclei, we rely on the usual few-body framework [15,33]. The projectile is described as a two-body system: a core c, assumed to be in its 0^+ ground state, to which the halo neutron n is loosely bound. The structure of the target T is neglected. The c-ninteraction is modeled by an effective single-particle potential whose parameters are fitted to reproduce the lower-energy spectrum of the projectile: the energy, spin and parity of its bound states and of some of its resonances above the c-nseparation threshold. The interaction between the projectile constituents and the target is simulated by optical potentials, which include an imaginary term that accounts for all the inelastic channels not explicitly accounted for by the model.

Within this three-body framework, we evaluate the knockout cross sections using eikonal-based models of the collision [34]. Two processes contribute to the knockout observables: the diffractive breakup, in which both the core and the halo neutron survive the collision, and the stripping, in which the neutron is absorbed by the target. The diffractive-breakup contribution is computed within the dynamical eikonal approximation (DEA) [35,36], which provides a proper dynamical description of the collision. Because the DEA has not yet been extended to stripping observables, we compute these contributions at the usual eikonal approximation [15,37,38]. Contrary to the DEA, the usual eikonal model relies on the adiabatic approximation, which sees the internal coordinates of the projectile as frozen during the collision. This approximation neglects the dynamical effects in the stripping.

All calculations are obtained with the same model spaces as the ones detailed in Refs. [23,39].

B. Halo-EFT description of ¹¹Be and ¹⁵C

We describe ¹¹Be and ¹⁵C within halo-EFT. This approach exploits the separation of scales between the size of the halo and that of the core to expand the projectile Hamiltonian upon the small parameter $R_{\rm core}/R_{\rm halo}$. In this expansion, both the core and the halo neutron are considered structureless and the breakdown scale is set by the size of the core [24–26]. The short-range physics is absorbed by a contact interaction and its derivatives, whose parameters are adjusted in each partial wave to reproduce known long-range properties of the nucleus, such as its one-neutron separation energy and the ANC of its bound states. For practical handling within reaction codes, these interactions are regulated with a Gaussian (see Eq. (13) of Ref. [27]).

Both ¹¹Be and ¹⁵C exhibit a similar structure. They both have a $1/2^+$ ground state, which is seen as a $1s_{1/2}$ neutron bound to their ¹⁰Be or ¹⁴C core in its 0^+ ground state. Their spectra also include a subthreshold excited state: a $1/2^{-}$ state for ¹¹Be and a $5/2^+$ state for ¹⁵C. These states are also described as a neutron bound to the 0^+ core, in either the $0p_{1/2}$ or the $0d_{5/2}$ orbital, respectively. Following Refs. [27,28], we describe ¹¹Be and ¹⁵C at NLO. For ¹¹Be, we set an effective *c-n* potential in the $s_{1/2}$ and $p_{1/2}$ waves, which reproduces the experimental one-neutron separation energies in the $1/2^+$ and $1/2^{-}$ states, as well as the ANCs predicted by the *ab initio* calculations of Calci et al. [29]. In the present work, we consider the potential with the Gaussian range of 1.2 fm, for which the depths can be found in Tables I and II of Ref. [27]. The uncertainty associated with the choice of this range is negligible, i.e., less than 1.5% for integrated knockout cross sections. We do not include any interaction in the $p_{3/2}$ or higher partial waves, first to reproduce the nearly nil phaseshifts predicted ab initio at low ¹⁰Be -n energy [29] and second because the details of the core-neutron interaction in the continuum do not matter in such calculations [23].

In the case of ¹⁵C, we use an effective potential in the $s_{1/2}$ partial wave to reproduce the ground state and go beyond NLO by including an interaction within the $d_{5/2}$ wave to describe the bound excited state, that can affect knockout calculations [23]. In this article, we use the parameters of these interactions for a Gaussian range of 1.2 fm, which are given in Table III of Ref. [28]. As for the ¹¹Be case, the uncertainty associated with this range are negligible. These potentials reproduce the ANCs inferred from an analysis of transfer observables [28], and agree very well with those

TABLE I. Parameters of the optical potentials (1) used to simulate the c^{-9} Be and n^{-9} Be interactions for the one-neutron knockout of ¹¹Be and ¹⁵C on ⁹Be at 60 MeV/nucleon and 54 MeV/nucleon. To estimate the uncertainty related to this choice, two potentials have been selected for each interaction.

Interacti	on	E _{beam} (MeV)	V _R (MeV)	<i>r_R</i> (fm)	a_R (fm)	W _I (MeV)	<i>r</i> _{<i>I</i>} (fm)	<i>aI</i> (fm)	W _D (MeV)	<i>r</i> _D (fm)	<i>a</i> _D (fm)	<i>r</i> _C (fm)	Ref.
<i>c</i> - ⁹ Be	V_{cT}^1		123.0	0.75	0.80	65.0	0.78	0.80	0	0	0	1.2	[40]
	V_{cT}^2		127.0	0.80	0.78	13.90	1.25	0.70	0	0	0	1.0	[42]
<i>n</i> - ⁹ Be	V_{nT}^1		33.08	1.14	0.65	4.15	1.14	0.65	9.18	2.83	0.18		[43]
	V_{nT}^2	60	22.60	1.37	0.29	5.30	1.30	0.30	15.25	1.30	0.30		[44]
		54	23.47	1.40	0.29	5.36	1.30	0.30	15.55	1.30	0.30		

predicted by an *ab initio* calculation [30]. As in Ref. [28], we assume a nil c-n interaction in the other partial waves, in particular the p ones. First, we do not have any reliable prediction for the corresponding phaseshifts and, second, we know from our previous study that the knockout cross sections are insensitive to the particulars of the c-n continuum [23].

Once coupled with accurate reaction models, these halo-EFT descriptions of ¹¹Be and ¹⁵C lead to an excellent agreement with breakup and transfer data [27,28,31,32]. The goal of the present study is to see if, following our previous study [23], they can also explain the knockout cross sections measured in Refs. [13,14].

C. Optical potentials

To analyze one-neutron knockout reactions of ¹¹Be and ¹⁵C on ⁹Be at 60 MeV/nucleon and 54 MeV/nucleon, respectively [13,14], we need optical potentials to simulate the interaction between the projectile constituents and the target. To estimate the uncertainty related to that choice, we select in the literature two different optical potentials for each interaction.

The first *c*-⁹Be potential V_{cT}^1 , developed in Ref. [40], reproduces elastic-scattering data of ¹⁰Be on ¹²C at a beam energy $E_{\text{beam}} = 59.4 \text{ MeV/nucleon}$ [41]. Although it does not correspond to the exact same target, it provides a fair estimate for the ¹⁰Be - ⁹Be interaction at the right beam energy. Not having found a more adequate potential for the ¹⁴C - ⁹Be interaction, we use this potential also in that latter case. The second *c*- ⁹Be potential V_{cT}^2 has been adjusted on elastic-scattering data for ⁹Be on ¹³C at 40 MeV [42]. Although it has been fitted at a beam energy much lower than those considered here, it is the second most realistic potential we could find in the literature.

For the first n^{-9} Be interaction V_{nT}^1 , we adopt the global optical potential developed by Weppner [43]. It has been fitted to elastic-scattering angular distributions and polarization data for a nucleon off a nucleus with mass number $A \leq 13$ at energies between 65 and 75 MeV. This potential is therefore well adapted for the collisions studied here. The second n^{-9} Be potential V_{nT}^2 , parametrized by Bonaccorso and Charity, reproduces the total, elastic, and reaction cross sections for a neutron on ⁹Be at energies between 1 and 100 MeV [44]. We compute its parameters for the two beam energies considered here.

The nuclear part of all these potentials exhibit the usual expression

$$V(R) = -V_R f_{WS}(R, R_R, a_R) - i W_I f_{WS}(R, R_I, a_I) + i 4 a_D W_D \frac{d}{dR} f_{WS}(R, R_D, a_D),$$
(1)

with $f_{WS}(R, R_X, a_X) = \frac{1}{1+e^{\frac{R-R_X}{a_X}}}$ the Woods-Saxon form. For the *c*-*T* interaction, the radii are parametrized as $R_x = r_x(A_c^{1/3} + A_T^{1/3})$, with A_c and A_T the core and target mass numbers, respectively. The radii for the *n*-*T* interactions are obtained with $R_x = r_x \times A_T^{1/3}$. The Coulomb *c*-*T* interaction is simulated by the potential generated by a uniformly charged sphere of radius $R_C = r_C \times (A_c^{1/3} + A_T^{1/3})$. The parameters of the optical potentials considered in this study are listed in Table I.

III. ANALYSIS OF THE ONE-NEUTRON KNOCKOUT FROM ¹¹Be AND ¹⁵C

To see if halo-EFT descriptions of ¹¹Be and ¹⁵C at NLO correctly reproduce one-neutron knockout, we perform reaction calculations within the model described in Sec. II and compare our results with the data of Refs. [13,14] obtained on a ⁹Be target at 60 MeV/nucleon and 54 MeV/nucleon, respectively. To estimate the uncertainty related to the simulation of the interaction between the projectile constituents and the target, we consider the four possible combinations of the optical potentials listed in Table I. For each of these choices, the total one-neutron knockout cross section (σ_{ko}) is provided in Table II, as well as its two contributions: diffractive breakup (σ_{bu}) and stripping (σ_{str}). The experimental values of Refs. [13,14] are also listed. The knockout cross sections expressed as a function of the momentum of the core parallel to the beam axis $p_{c\parallel}$ are displayed in Fig. 1. For the first combination of optical potentials, the diffractivebreakup and stripping contributions are shown separately. The experimental data (red points) are extracted from Figs. 11 and 12 of Ref. [14]. The error bars correspond to the 15% and 12% experimental uncertainties reported in Refs. [13,14], respectively. For a proper comparison between theory and experiment, following Ref. [14], we have adjusted the position of the center of each parallel-momentum distributions to the data. Beside that minor adjustment, no parameter has been

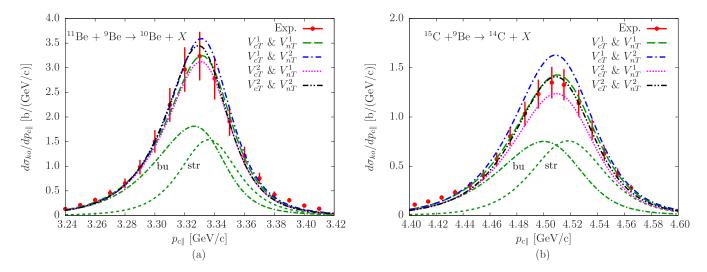


FIG. 1. Parallel-momentum distributions for the one-neutron knockout on ${}^{9}\text{Be}$ of (a) ${}^{11}\text{Be}$ at 60 MeV/nucleon, and (b) ${}^{15}\text{C}$ at 54 MeV/nucleon. Calculations obtained with the different optical potentials of Table I are compared to the experimental data of Refs. [13,14]. The experimental error bars correspond to the 15% and 12% uncertainty cited in Refs. [13,14], respectively. For the first choice of optical potentials, the diffractive-breakup (bu, double-dash-dotted green line) and stripping (str, short-dashed green line) contributions to the cross section are shown separately.

fitted to the data; in particular, the magnitude of the cross sections is the direct output of the calculations.

We observe an excellent agreement between theory and experiment for both knockout observables. This is first seen in the magnitude of the cross section. In both Table II and Fig. 1, but for two exceptions, all calculations fall within one standard deviation of the experimental value. In addition, our calculations reproduce very well the experimental parallel-momentum distribution. As deduced from the initial studies, this confirms the clear one-neutron halo structure of ¹¹Be and ¹⁵C [13–15].

Because our results are nearly insensitive to the details of the halo-EFT description of the projectile, the dominant uncertainty in the reaction model resides in the choice of the optical potentials. This choice affects mostly the magnitude of the cross section, leading to about 11% difference in the ¹¹Be case and up to 26% in the ¹⁵C calculations, which is much larger than the 1.5% uncertainty on the *c-n* interaction we have observed. For the former nucleus, most of that uncertainty comes from V_{nT} , whereas for the latter both *c-T* and *n-T* interactions share an equal role in the changes in the cross section. Reducing that uncertainty requires the development of new, more accurate, optical potentials, e.g., derived from first principles [45–48]. A systematic study within a Bayesian approach, such as done in Refs. [49–51] for transfer reactions, would help us better understand the influence of each interaction on the reaction process and provide a more reliable uncertainty. Although in a much less significant way, the optical potentials also influence the shape of the parallel-momentum distribution. The n-⁹Be optical potential of Bonaccorso and Charity [44] produces a slightly less asymmetric peak than Weppner's [43]. Albeit not statistically significant, this hints at the influence of the n-T interaction in those reactions dynamics [52].

To study further the role of the optical potentials, we separate the knockout cross section into its diffractivebreakup (bu) and stripping (str) components in Table II. As in Ref. [14], we note that each component contributes for roughly half of the knockout cross section. The actual division into diffractive breakup and stripping depends mostly on V_{nT} : The less absorptive V_{nT}^1 of Ref. [43] leads to a larger σ_{bu} and smaller σ_{str} than the more absorptive V_{nT}^2 of Ref. [44]. The *c*-*T* interaction plays only a minor role in this division.

TABLE II. Integrated diffractive-breakup (σ_{bu}), stripping (σ_{str}), and knockout ($\sigma_{ko} = \sigma_{bu} + \sigma_{str}$) cross sections in mb for ¹¹Be and ¹⁵C on ⁹Be at 60 MeV/nucleon and 54 MeV/nucleon, respectively. Theoretical results obtained with the different potentials listed in Table I are compared to the experimental values from Refs. [13,14].

		Expt.	V_{cT}^{1} & V_{nT}^{1}	V_{cT}^1 & V_{nT}^2	V_{cT}^2 & V_{nT}^1	$V_{cT}^2 \& V_{nT}^2$
	$\sigma_{ m bu}$		113	89	110	89
¹¹ Be	$\sigma_{ m str}$		81	118	74	108
	$\sigma_{ m ko}$	203 ± 31	194	207	185	197
	$\sigma_{ m bu}$		60	48	52	43
¹⁵ C	$\sigma_{ m str}$		51	75	43	63
	$\sigma_{ m ko}$	109 ± 13	111	124	95	105

The separation of the $p_{c\parallel}$ distribution into its two contributions is illustrated in Fig. 1 for the first choice of optical potentials (the other choices lead to similar results, but for the relative magnitude of the two contributions, see Table II). The diffractive-breakup contribution is plotted in double-dash-dotted green line and the stripping one in shortdashed green line. The former exhibits a notably asymmetric shape with a tail extending to the low-momentum side. This asymmetry is due to the dynamics of the reaction [14,39], which is correctly accounted for within the DEA [35,36]. The stripping cross section, on the contrary, is purely symmetric around its maximum because it is computed within the usual eikonal approximation, i.e., including the adiabatic approximation [14]. Although their sum reproduces very well the experimental cross section, our calculations slightly underestimate the data in the low-momentum tail, especially for the knockout of ¹⁵C. Our description of the reaction therefore misses some dynamical effects. To account for the reaction dynamics in the stripping process, the authors of Ref. [14] assume the same shape of that parallelmomentum distribution as the diffractive-breakup one. Their methodology hence includes more dynamical effects than ours; accordingly the difference between their calculations and the data at low momenta is smaller than ours. A more precise treatment of the dynamics within the stripping reaction is thus needed to properly describe the full reaction process. Nevertheless, but for this minor issue, our treatment of the reaction mechanism is accurate enough to analyze these measurements.

Since knockout is purely peripheral for halo nuclei, the cross section scales very well with the square of the ANC of the ground state [23]. From our calculations, we can infer an ANC by adjusting the magnitude of our parallel-momentum distributions to the data. For ¹¹Be we obtain an ANC² = $0.62 \pm 0.06 \pm 0.09$ fm⁻¹ and for ¹⁵C, an ANC² = $1.57 \pm 0.30 \pm 0.18$ fm⁻¹. The first quoted uncertainty is associated with the choice of optical potentials, while the second is due to the experimental error. Because they are of the same order of magnitude, reducing the uncertainty on the ANC inferred from knockout data would require better constraints on the optical potentials and more precise data.

These ANCs inferred from the knockout data are in excellent agreement with the ones predicted by ab initio calculations, viz. 0.618 fm⁻¹ for ¹¹Be [29] and 1.644 fm⁻¹ for ^{15}C [30], and with halo-EFT analyses of transfer reactions, viz. $0.616 \pm 0.001 \text{ fm}^{-1}$ for ¹¹Be [31] and $1.59 \pm 0.06 \text{ fm}^{-1}$ for ¹⁵C [28]. Although in the lower end of the spectrum, the ANC for ¹⁵C agrees well with the values extracted from other analyses [9,53–58] (see list in Table II of Ref. [28]). The values listed above have been used to constrain the halo-EFT *c*-*n* interactions considered in this analysis as well as in Refs. [27,28,32]: They lead to an excellent agreement with data for breakup at both intermediate [27,28] and high [28,32] energies, transfer [28,31], and the radiative capture ${}^{14}C(n, \gamma)$ at energy of astrophysical interest [28]. This result is thus the last piece of a puzzle, which demonstrates that with one halo-EFT description of halo nuclei, we can reproduce various sets of experimental data using accurate reaction models. Accordingly, halo-EFT properly identifies the nuclear-structure observables that matter most in these reactions, viz. the binding energy of the halo neutron, its orbital angular momentum and the ANC of the projectile ground state. In particular, the norm of the overlap wave function does not influence the results of the calculations, as long as the ANC is properly reproduced [23,27].

IV. CONCLUSIONS

Knockout reactions provide a clean probe of halo nuclei because their cross section reveals clearly the spatial extension of the nuclear wave function and they exhibit high statistics [12–15]. Our previous analysis [23] has shown that the one-neutron knockout of halo nuclei is sensitive only to the asymptotics of the ground-state wave function while being insensitive to the norm of that wave function, i.e., the spectroscopic factor of the corresponding overlap wave function. Moreover, knockout cross sections do not depend on the description of the *c*-*n* continuum. Accordingly, a halo-EFT description at NLO of the projectile should include all the relevant nuclear-structure information to reproduce such data. In this article, we test this idea by reanalyzing measurements of one-neutron knockout of ¹¹Be [13] and ¹⁵C [14] using halo-EFT descriptions reproducing well data on breakup [27,28,32] and transfer [28,31].

For both nuclei, our calculated parallel-momentum distributions and integrated knockout cross sections are in excellent agreement with the data. In particular, the typical narrow width of the distributions, their slight asymmetry, and the magnitude at the peak are well reproduced. Using four different optical potentials, we have estimated the uncertainty of the reaction model and suggested that more accurate optical potentials would reduce that uncertainty [45–48].

Comparing our calculations with the data, we have inferred the ground-state ANC for both projectiles. These values agree well with *ab initio* predictions [29,30] and transfer analyses [28,31]. Our results complete a series of previous studies and demonstrate that one description of ¹¹Be and ¹⁵C based on halo-EFT reproduces independent experimental data for breakup [27,28,32], transfer [28,31], radiative-capture [28], and, from this study, knockout [13,14] measurements. Since our results are independent of the normalization of the groundstate wave function [23], they also suggest that the quenching factor $R_S \approx 1$ obtained for halo nuclei in previous analyses of knockout data [17–19] is due more to the use of realistic ANCs in reaction calculations than to accurate spectroscopic factors. This is not surprising since, for neutrons loosely bound in an s wave, the ANC of the wave function is predominantly determined by the binding energy, with little sensitivity to the *c*-*n* potential geometry [59].

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