

Electric dipole polarizability of ^{40}Ca

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The electric dipole strength distribution in ^{40}Ca between 5 and 25 MeV has been determined at RCNP, Osaka, from proton inelastic scattering experiments at very forward angles. Combined with total photoabsorption data at higher excitation energy, this enables an extraction of the electric dipole polarizability $\alpha_D(^{40}\text{Ca}) = 1.92(17) \text{ fm}^3$. Together with the measured α_D in ^{48}Ca , it provides a stringent test of modern theoretical approaches, including coupled-cluster calculations with chiral effective field theory interactions and state-of-the-art energy density functionals. The emerging picture is that for this medium-mass region dipole polarizabilities are well described theoretically, with important constraints for the neutron skin in ^{48}Ca and related equation of state quantities.

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Introduction. The nuclear equation of state (EOS) not only determines basic properties of nuclei [1] but also plays a key role for the properties of neutron stars and the dynamics of core-collapse supernovae and neutron star mergers [2]. New observations from neutron stars and mergers provides constraints for the EOS of neutron-rich matter that can be compared with those derived from nuclear physics (see, e.g., Refs. [3–5]). However, while the EOS of symmetric nuclear matter is well determined around saturation density, the properties of neutron-rich matter are less explored experimentally. The latter depends on the symmetry energy, whose properties are typically encoded in an expansion around saturation density n_0 , with the symmetry energy at saturation density $J(n_0)$ and its density dependence $L = 3n_0 \partial J(n_0) / \partial n$.

Theoretically, a model-dependent correlation between L and the neutron-skin thickness r_{skin} in nuclei with neutron excess has been established [6–9]. This correlation was also

recently confirmed in *ab initio* computations of the neutron skin in ^{208}Pb [10]. Experimental attempts to determine the neutron skin thickness have been performed with a variety of probes (see, e.g., Ref. [11] and references therein), but many of them suffer from systematic uncertainties entering in the description of the reaction processes. Parity-violating elastic electron scattering (a weak process mediated by the Z^0 boson) can be used for a nearly model-independent extraction of the neutron distribution in nuclei and, by comparison with accurately measured charge radii, the neutron skin thickness. Recently, results with this technique have been reported by the CREX and PREX collaborations for ^{48}Ca [12] and ^{208}Pb [13], respectively. The r_{skin} values inferred with selected nuclear models favor a comparatively small neutron skin in the former and a large skin in the latter case.

Alternatively, the electric dipole polarizability α_D has been established as a possible measure of the neutron skin, based on the strong correlation with r_{skin} [8,14]. Data for α_D extracted from proton inelastic scattering experiments at extreme forward angles have been presented for both ^{48}Ca [15] and ^{208}Pb [16]. In these papers, two theoretical approaches have been used to describe α_D : *ab initio* coupled-cluster (CC) calculations [17,18] starting from chiral two- and three-nucleon interactions [19,20] and energy density functional (EDF) theory [21].

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Attempts to simultaneously describe $\alpha_D(^{208}\text{Pb})$ and the parity-violating asymmetry from PREX and CREX with EDF models have shown limited success [22–25]. The values derived for r_{skin} [13] and L [26] from PREX are in tension with EDFs capable of describing [27] the presently available results on α_D in ^{48}Ca [15], ^{68}Ni [28], ^{120}Sn [29,30], and ^{208}Pb [16]. While the CREX results is in excellent agreement with *ab initio* predictions [18], the PREX result is in mild tension with the recent *ab initio* computations of ^{208}Pb [10].

Correlations between experimental observables and symmetry energy properties are well explored in EDF theory [6–8,14], but predictions for isovector observables like α_D are less well constrained. On the other hand, *ab initio* calculations provide a direct link to the EOS, as nuclear matter properties can be calculated based on the same chiral interactions [10,18,19,31,32]. Results presented here are based on the set of two- and three-nucleon interactions from Refs. [19,20] applied to study α_D in ^{48}Ca [15,18]. The calculations of the $E1$ response are based on merging the Lorentz integral transform approach with CC theory, as described in Refs. [33,34]. Recent work has extended the original two-particle–two-hole (2p-2h) CC truncation to include correlations up to three-particle–three-hole (3p-3h), so-called triples corrections, in the computation of α_D [35]. Their inclusion leads to a reduction of the predictions for $\alpha_D(^{48}\text{Ca})$ of the order of 10%, allowing an improved simultaneous description of the charge radius [35]. A similar improvement was achieved for ^{68}Ni [36].

In this Letter, we present the measurement of the dipole polarizability for ^{40}Ca and confront it with CC and EDF calculations. This tests the emerging picture that nuclear theory can describe very well the neutron skin in medium-mass nuclei and related observables.

Experiment. Cross sections for the $^{40}\text{Ca}(p, p')$ reaction have been measured at RCNP, Osaka, at an incident proton energy of 295 MeV. Data were taken with the Grand Raiden spectrometer [37] in a laboratory scattering angle range 0.4° – 14.0° and for excitation energies in the range 5–25 MeV. Dispersion matching techniques were applied to achieve an energy resolution of about 30 keV (full width at half maximum). The experimental techniques and the raw data analysis are described in Ref. [38].

In the top panel of Fig. 1 we show representative energy spectra measured at laboratory scattering angles $\Theta_{\text{lab}} = 0.4^\circ$, 1.74° , 3.18° , and 5.15° . The predominant cross sections lie in the energy region above 10 MeV. $M1$ strength in ^{40}Ca is known to be concentrated in a single prominent transition at 10.32 MeV [39]. The cross sections above 10 MeV show a broad resonance structure peaking at about 19 MeV increasing towards 0° . The angular dependence is consistent with relativistic Coulomb excitation of $E1$ transitions. We identify this resonance structure as the isovector giant dipole resonance.

The various contributions to the spectra were separated using a multipole decomposition analysis (MDA) as described in Ref. [40]. Results for the most forward angle measured are presented in the bottom part of Fig. 1 as example, where the spectrum was rebinned to 200 keV. Theoretical angular distributions for the relevant multipoles were obtained from distorted-wave Born approximation calculations with transition amplitudes from quasiparticle-phonon-model calcu-

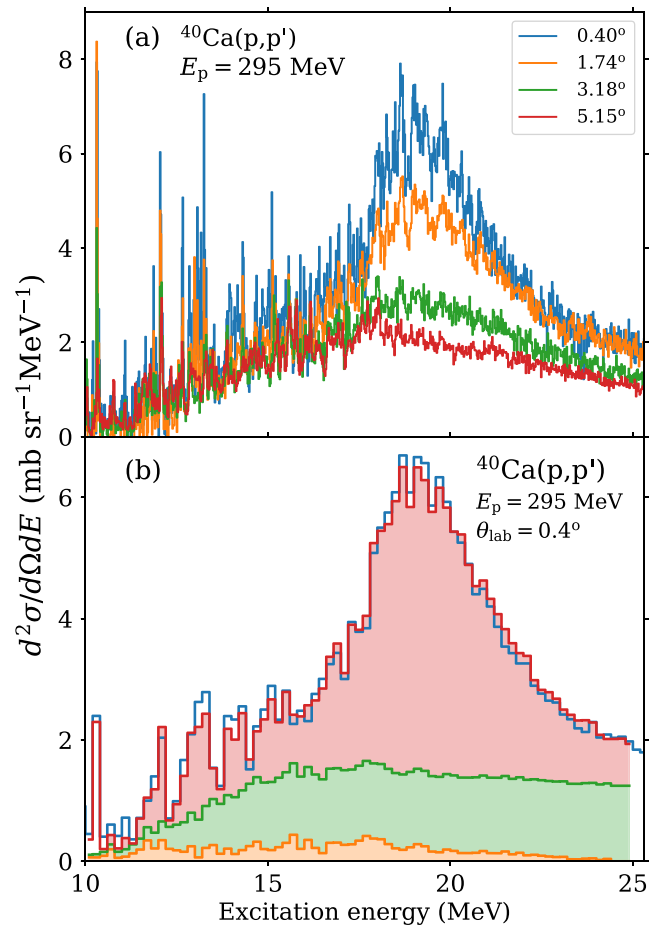


FIG. 1. (a) Spectra of the $^{40}\text{Ca}(p, p')$ reaction at $E_0 = 295$ MeV and scattering angles $\Theta_{\text{lab}} = 0.4^\circ$, 1.74° , 3.18° , and 5.15° . (b) Example of the MDA of the spectrum at $\Theta_{\text{lab}} = 0.4^\circ$ in 200 keV bins (blue) and decomposition into contributions of $\lambda \neq 1$ multipoles (orange), continuum background (green), and $E1$ (red).

lations similar to the analysis of ^{48}Ca [15]. Additionally, a background due to pre-equilibrium multistep scattering was considered. Its angular dependence was taken from experimental systematics [41,42] while the amplitude was derived by two means. Initially, an unconstrained fit was done at each energy bin of the set of spectra. The resulting cross sections could be well approximated by a simple Fermi function but showed strong fluctuations for certain excitation energy bins due to the similarity to some of the $E1$ theoretical angular distributions. Thus, in the final analysis, the continuum contribution was determined by fitting a Fermi function to the unconstrained excitation energy dependence.

Photoabsorption cross sections and dipole polarizability. The $E1$ cross sections resulting from the MDA were converted into equivalent photoabsorption cross sections using the virtual photon method [43]. The virtual photon spectrum was calculated in an eikonal approach [44] to Coulomb excitation, integrated over the distribution of scattering angles covered in the solid angle of each angular bin. The photoabsorption spectra derived from scattering data at 0.40° and 1.00° were essentially identical, and that at 1.74° deviated only slightly, consistent with an estimate of the grazing angle (1.33°) at

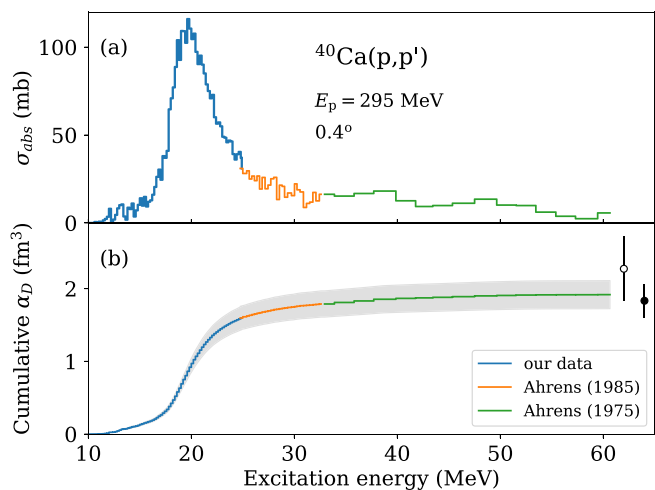


FIG. 2. (a) Photoabsorption cross section derived at a scattering angle of 0.40° using the virtual photon method. (b) Electric dipole polarizability α_D derived from the photoabsorption cross sections. The blue curve shows the present data, while the orange and green curves show the extrapolation to higher energies using the data of Refs. [45,46]. The open (full) black circles are the CC results for the NNLO_{sat} interaction including up to doubles (triples) contributions in the cluster expansion.

which Coulomb-nuclear interference becomes relevant. The resulting photoabsorption cross section is displayed as blue histogram in Fig. 2(a).

The electric dipole polarizability α_D was obtained from the photoabsorption cross section over the energy range 10–25 MeV leading to a contribution $1.60(14) \text{ fm}^3$. The integration was extended to 60 MeV, where the cumulative sum plotted in Fig. 2(b) shows saturation. The data at higher excitation energies were taken for 25–31 MeV from Ref. [46] and for 31–60 MeV from Ref. [45] to obtain the total $\alpha_D(^{40}\text{Ca}) = 1.92(17) \text{ fm}^3$. The uncertainty considers systematic errors of (i) the absolute cross sections, (ii) the MDA (determined as described, e.g., in Ref. [47]), and (iii) the parametrization of the continuum background (evaluated by varying the amplitude of the Fermi function), added in quadrature. The latter, dominating the total uncertainty budget, was estimated by the variation needed to change the χ^2 value of the MDA fit by one. Statistical errors turned out to be negligible. A detailed breakdown of the error contributions is given in Table I.

TABLE I. Budget of error contributions to $\alpha_D(^{40}\text{Ca})$.

Source	Value (%)
Trigger efficiency	0.1
Drift chamber efficiency	0.8
Charge collection	0.3
Target thickness	1.0
Determination of solid angle	3.0
MDA	1.2
Background parametrization	8.3
Total	9.0

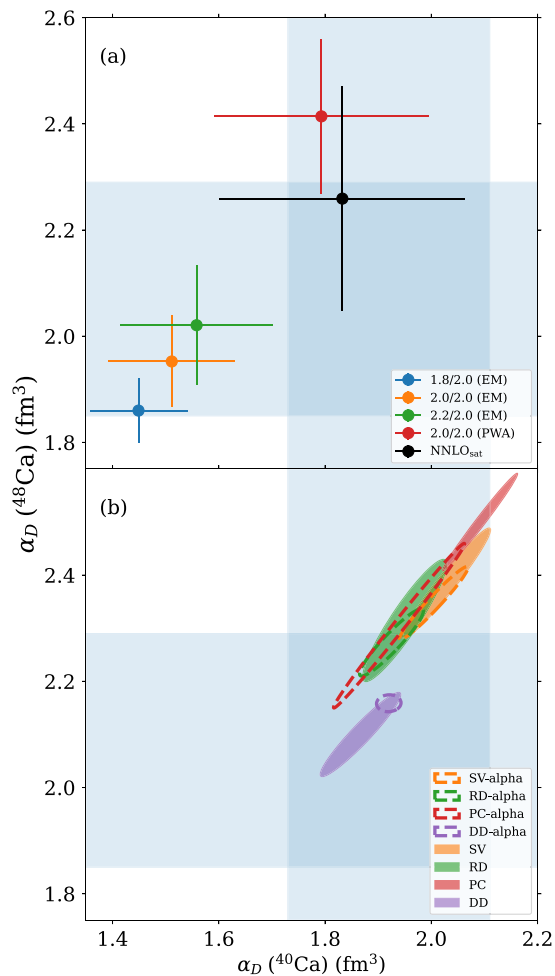


FIG. 3. Comparison of the experimental dipole polarizabilities of ^{40}Ca (present work) and ^{48}Ca [15] shown as blue bands with (a) CC calculations with different interactions, including triples contributions, and (b) EDF calculations with different energy density functionals [22]. For details see text.

Comparison with coupled-cluster calculations. The extracted value of α_D serves as a benchmark for CC theory [18,33–35]. Coupled-cluster calculations were recently performed for the dipole polarizability of ^{48}Ca [15] and ^{68}Ni [36], which led to an improved understanding of the neutron and proton distributions in nuclei, as well as their difference encoded in the neutron skin. We have performed CC computations of α_D in ^{40}Ca starting from a Hartree-Fock reference state considering a basis of 15 major harmonic oscillator shells. To gauge the convergence of our results we varied the oscillator frequency in the range $\hbar\omega = 12\text{--}16 \text{ MeV}$. Three-nucleon contributions had an additional energy cut of $E_{3\text{max}} = 16\hbar\omega$.

Figure 3 explores the correlation between α_D for ^{40}Ca and ^{48}Ca as predicted by theory. Figure 3(a) shows the CC results including triples contributions, not available for ^{40}Ca so far. The theoretical uncertainties for the different Hamiltonians stem from the truncation of the CC expansion and the residual dependence on CC convergence parameters, calculated as described in Ref. [34]. Similarly to ^{48}Ca , we find

that the inclusion of 3p-3h correlations reduces the value of $\alpha_D(^{40}\text{Ca})$ by an amount varying between 10% and 20% for different interactions. While the EM and PWA interactions [19] are not simultaneously compatible with both ^{40}Ca and ^{48}Ca experimental data, the set of employed interactions shows an approximately linear trend between the two quantities overlapping with both experimental results. A particular improvement in the reproduction of both $\alpha_D(^{48}\text{Ca})$ and $\alpha_D(^{40}\text{Ca})$ is seen for the NNLO_{sat} interaction [20], which is capable of accurately describing binding energies and radii of nuclei up to ^{40}Ca as well the saturation point of symmetric nuclear matter. The different interactions predict a range of symmetry energy parameters $J = 27\text{--}33$ MeV, $L = 41\text{--}49$ MeV [18], with the NNLO_{sat} values at the lower end ($J = 27$ MeV, $L = 41$ MeV).

Comparison with EDF approaches. Recently, it was investigated whether the dipole polarizability and the parity-violating asymmetry A_{PV} for ^{208}Pb and ^{48}Ca can be simultaneously accounted for with modern EDFs [24]. We use the four representative forms of functionals from that study: nonrelativistic Skyrme functionals SV [48] and RD [49], the latter with different forms of density dependence, and relativistic functionals DD [50] with finite-range meson-exchange coupling and PC [51] with point coupling. All four have been calibrated to the same set of ground-state data to determine the model parameters. With these sets, it was shown that PREX and CREX results for A_{PV} (and r_{skin}) cannot be consistently explained within the model uncertainties while the α_D were reproduced. Hence, the present result in ^{40}Ca provides an important test of the global predictive power of these EDFs.

Figure 3(b) displays the EDF results for α_D with 1σ error ellipses (for their definition see Refs. [22,24]). The parametrizations as given from the ground-state fits are shown by filled ellipses. The DD functional performs rather well. The other predictions tend to slightly overestimate the experimental mean values of both ^{40}Ca and ^{48}Ca , while their 1σ error ellipses do overlap with the experimental bands, except for PC. In all cases, the two α_D values are highly correlated. We note that the same behavior is found for the description of $\alpha_D(^{208}\text{Pb})$ [16] after correction for the quasideuteron contribution [27]. Thus, all the models are capable to account for the mass dependence of the polarizability.

The dashed ellipses show results from a refit where additionally the experimental α_D value of ^{208}Pb [16] corrected for the quasideuteron part [27] was included yielding the functionals SV-alpha; RD-alpha, PC-alpha, and DD-alpha [22,24]. This improves the agreement with experiment,

particularly for the PC model, and shrinks most error ellipsoids. The uncertainty reduction is especially large for the DD model because this functional has the least isovector freedom. The linear trend shown by the different theoretical approaches in Fig. 3 is similar although the CC calculations tend to underestimate the α_D in ^{40}Ca and perform nicely for ^{48}Ca . The bulk symmetry energies range from $J = 30$ MeV for DD to 35 MeV for PC and from 32 to 82 MeV for L , respectively. The fits which include also α_D in ^{208}Pb narrow the prediction to $J = 30\text{--}32$ MeV and $L = 35\text{--}52$ MeV which correlates nicely to the narrower range of predictions for α_D in $^{40,48}\text{Ca}$.

Conclusions. We have extracted the dipole polarizability of ^{40}Ca from a combination of relativistic Coulomb excitation measurement in inelastic proton scattering under very forward angles with total photoabsorption data at high excitation energies. Together with a similar analysis on ^{48}Ca the new data serve as a benchmark test of state-of-the-art theoretical approaches. A representative set of EDFs can describe these data. An improvement is obtained when the EDFs are optimized by adding the dipole polarizability of ^{208}Pb to the calibration dataset. Coupled-cluster computations for the NNLO_{sat} interaction simultaneously describe well the dipole polarizability of ^{40}Ca and ^{48}Ca , as well as the corresponding charge radii and the neutron skin thickness [34]. A nearly linear systematic trend is obtained for other interactions, as in the case of EDF theory. This analysis supports the robustness of current theoretical approaches in the description of α_D and their constraints of symmetry energy parameters discussed, e.g., in Refs. [22,24,52].

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