# Forward doubly-virtual Compton scattering off an unpolarised deuteron in pionless effective field theory 

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#### Abstract

We calculate the forward unpolarised doubly-virtual Compton scattering (VVCS) off the deuteron in the framework of pionless effective field theory, up to next-to-next-to-next-to-leading order (N3LO) for the longitudinal and next-to-leading order (NLO) for the transverse amplitude. The charge elastic form factor of the deuteron, obtained from the residue of the longitudinal VVCS amplitude, is used to extract the value of the single unknown two-nucleon one-photon contact coupling that enters the longitudinal amplitude at N3LO. We also study the lowest spin-independent generalised polarisabilities of the deuteron. The calculated unpolarised VVCS amplitude provides a high-precision model-independent input for a future calculation of the two-photon-exchange correction to the Lamb shift of muonic deuterium.

\section*{CONTENTS} I. Introduction ..... 3 II. Expansions and power-counting for the deuteron VVCS amplitude ..... 4 A. Unpolarised VVCS amplitude and its low-momenta expansion ..... 5 B. EFT expansion and counting ..... 6 C. Lagrangian and coupling constants ..... 9 D. LSZ reduction ..... 13 III. Diagrams and pertinent results, order by order ..... 15 A. Deuteron self energy ..... 15 B. Irreducible four-point function ..... 17 1. LO ..... 17 2. NLO ..... 18 3. NNLO ..... 20 4. N3LO ..... 21 IV. Deuteron charge form factor at N3LO: fitting $l_{1}^{C 0_{S}}$ ..... 24 V. Deuteron generalised polarisabilities ..... 27 VI. Conclusion ..... 31 Acknowledgements ..... 32 References ..... 32


## I. INTRODUCTION

The deuteron is a fundamental source of information about the emerging nuclear force and, in the absence of free neutron targets, is often used to study the neutron structure. Recent advances in muonic-atom spectroscopy, by the CREMA Collaboration at PSI, led to presently the most precise determination of the charge radii of the proton [1, 2], deuteron [3], helium-3 and 4. In case of the deuteron, the isotopic shift measurement also gives an accurate assessment of a subleading nuclear structure contribution - the so-called two-photonexchange (TPE) correction. These accurate measurements provide further challenges for the theoretical description of the low-energy nuclear structure with systematically improvable precision.

While ab initio QCD calculations of these nuclear-structure quantities are still being out of reach, the method of choice for systematic calculations of the nuclear-structure corrections is effective-field theories (EFTs) of the strong interaction. Specifically, we employ the pionless EFT ( $\ddagger$ EFT) [4-11], where the pions are heavy and hence the nucleon-nucleon $(N N)$ interaction is described by contact interactions organized in powers of nucleon threemomentum $P$. This description is limited to momenta well below the pion mass, $P \ll m_{\pi}$, which should be well suited for atomic calculations, where the momenta are of the order of the inverse Bohr radius $\alpha m_{r}$ (with $\alpha$ the fine structure constant and $m_{r}$ the atomic reduced mass). Thus, typical momenta in a light muonic atom, such as muonic hydrogen $(\mu \mathrm{H})$ or deuterium $(\mu \mathrm{D})$, are below 1 MeV . The use of $\nexists \mathrm{EFT}$ is further motivated by its simplicity. The contact interactions lead to separable $N N$ potential, that enables algebraic resummation, resulting in closed analytic expressions for the nuclear force. Furthermore, it is strictly renormalisable (in the EFT sence), gauge invariant and hence exactly fulfills low-energy theorems (LETs) such as the Thomson limit.

The đEFT has already been demonstrated to work very well for low-energy properties of light nuclear systems; in particular, deuteron polarisabilities, the electromagnetic form factors of the deuteron, and the Compton scattering (CS) off the deuteron, see e.g. Refs. [7, 12-14]. Here we shall compute the forward doubly-virtual Compton scattering (VVCS) amplitude, which contains the deuteron structure information on the aforementioned TPE correction. This provides an alternative route to assessing these corrections: the existing calculations employ the nuclear Hamiltonian approach [15-20], or use the approach based on dispersion relations, either data-driven [21], or informed by EFT calculations of the deuteron structure functions [22-24].

The predictive orders of $\not \subset E F T$ for the TPE corrections are exhausted at next-to-next-to-next-to-leading order (N3LO), motivating our choice of calculating at this high order. The counting for the transverse contribution starts one order higher, at the respective N4LO, so the transverse amplitude can be neglected up to N3LO in the calculation of the TPE corrections. This consideration notwithstanding, a calculation of the transverse contribution allows one to study the generalised deuteron polarisabilities, such as the magnetic dipole polarisability and the generalised Baldin sum rule. It also provides input for verifying the
smallness of the transverse contribution in the context of the TPE.
We therefore calculate the longitudinal amplitude to N3LO, and the transverse amplitude up to next-to-leading order (NLO) in the so-called $Z$-expansion scheme [25]. Despite differing only at higher orders, this scheme has shown to converge better than the $\rho$-expansion scheme, in which one chooses to fit the effective range $\rho_{d}$ at NLO, for those observables that are dominated by the long-range properties of the deuteron wave function. This is achieved by fitting the residue of the scattering amplitude at NLO in the $Z$-expansion, thus ensuring the correct reproduction of the long-distance piece of the deuteron wave function starting from NLO.

A recent next-to-next-to-leading order (NNLO) $\not \approx$ EFT result for the longitudinal deuteron structure function [24] (in this connection, see also Refs. [16] and [22]) achieved a precision of about $5 \%$ for the TPE correction. The present N3LO calculation of the longitudinal deuteron VVCS amplitude is expected to further improve the theoretical uncertainty. With the $\not \approx$ EFT expansion parameter $P / m_{\pi} \simeq \gamma / m_{\pi} \simeq 1 / 3$, where $\gamma \simeq 45 \mathrm{MeV}$ is the deuteron binding momentum, one expects to achieve a relative precision of the order of $(1 / 3)^{4} \simeq 1 \%$ - an estimate of the contributions of higher orders in the đEFT expansion, which is the dominant source of uncertainty. This being compatible with other existing results, our work provides an alternative high-precision and model-independent handle on the TPE corrections in $\mu \mathrm{D}$. A detailed study of these corrections, making use of the $\not \approx$ EFT results for the deuteron VVCS, will be presented elsewhere.

The article is organised as follows. Section II concerns the details of the đEFT framework for the calculation of the unpolarised deuteron VVCS amplitude. The results of the calculation are presented, with a detailed description of various contributions, and a discussion of the effect of the $N N$ contact terms entering the calculation, in Sec. III. Section IV considers the deuteron charge form factor and fitting the unknown N3LO contact term, using the deuteron charge radius. A study of the generalised deuteron polarisabilities is presented in Sec. V, with a summary following in Sec. VI.

## II. EXPANSIONS AND POWER-COUNTING FOR THE DEUTERON VVCS AMPLITUDE

Compton scattering off light nuclei - in particular, the proton and deuteron - as a means of accessing information about the nuclear structure, such as nuclear polarisabilities, has been in the focus of experimental and theoretical efforts for several decades; detailed reviews of the proton CS and sum rules can be found in Refs. [26, 27]. At present, the TPE corrections, closely related to the nuclear VVCS, constitute the largest uncertainty in the extraction of the proton and deuteron charge radii from, respectively, $\mu \mathrm{H}[1]$ and $\mu \mathrm{D}$ [28]. Therefore, it is of high interest to systematically calculate and accurately determine the VVCS amplitudes.

Modern studies of proton and deuteron CS rely on the use of EFTs as the main theoretical, rigorous and systematic tool; we refer here to the recent review [29]. For the description of
the nucleons in the pion production energy range, sufficiently far from the confinement scale, $\Lambda \sim 1 \mathrm{GeV}$, the chiral EFT ( $\chi$ EFT) is widely used [19, 20, 22, 30]. On the other hand, đEFT, being appropriate for energies well below the energy scale set by the pion mass, is useful and reliable in the regime where the momentum transfer between the two nucleons (or, broadly speaking, a typical momentum scale of the problem, such as the photon beam momentum) is $P \ll m_{\pi}$, with the expansion parameter set by the ratio $P / m_{\pi} \simeq \gamma / m_{\pi} \simeq 1 / 3$.

In the rest of this section, we briefly recap the essential details of the VVCS amplitude and the $\not \approx$ EFT expansion, set up the counting, and provide the details of the $\not \approx E F T$ Lagrangian relevant to our calculation.

## A. Unpolarised VVCS amplitude and its low-momenta expansion

The unpolarised deuteron VVCS amplitude

$$
\begin{equation*}
T_{f i}=-\mathcal{E}_{\lambda}^{\prime *} \mathcal{E}^{\lambda} \epsilon_{\mu}^{\prime *} \epsilon_{\nu} T^{\mu \nu} \tag{1}
\end{equation*}
$$

where $\mathcal{E}$ and $\mathcal{E}^{\prime}$ are the polarisation vectors of the initial and final deuteron, $\epsilon$ and $\epsilon^{\prime}$ are those of the initial and final photon, is conveniently parametrised by the two scalar amplitudes $T_{1}\left(\nu, Q^{2}\right)$ and $T_{2}\left(\nu, Q^{2}\right)$, entering the VVCS tensor $T^{\mu \nu}$ [31]:

$$
\begin{equation*}
T^{\mu \nu}(p, q)=\left(-g^{\mu \nu}+\frac{q^{\mu} q^{\nu}}{q^{2}}\right) T_{1}\left(\nu, Q^{2}\right)+\frac{1}{M_{d}^{2}}\left(p^{\mu}-\frac{p \cdot q}{q^{2}} q^{\mu}\right)\left(p^{\nu}-\frac{p \cdot q}{q^{2}} q^{\nu}\right) T_{2}\left(\nu, Q^{2}\right) \tag{2}
\end{equation*}
$$

Here, $q$ and $p$ are the photon and deuteron four-momenta, $M_{d}$ is the deuteron mass, $\nu=$ $p \cdot q / M_{d}$ is the laboratory frame energy of the photon, and $Q^{2}=-q^{2}$ its virtuality. While the representation in terms of the amplitudes $T_{1,2}$ is explicitly covariant, the non-relativistic nature of the deuteron allows for a more natural treatment of the deuteron VVCS in terms of the longitudinal and transverse amplitudes,

$$
\begin{equation*}
f_{L}\left(\nu, Q^{2}\right)=-T_{1}\left(\nu, Q^{2}\right)+\left(1+\frac{\nu^{2}}{Q^{2}}\right) T_{2}\left(\nu, Q^{2}\right), \quad f_{T}\left(\nu, Q^{2}\right)=T_{1}\left(\nu, Q^{2}\right) \tag{3}
\end{equation*}
$$

The corresponding decomposition of $T_{f i}$ takes the form

$$
\begin{equation*}
T_{f i}=-\mathcal{E}_{\lambda}^{\prime *} \mathcal{E}^{\lambda}\left[\varepsilon_{0} \varepsilon_{0}^{\prime *} f_{L}\left(\nu, Q^{2}\right)+\left(\varepsilon \cdot \varepsilon^{\prime *}\right) f_{T}\left(\nu, Q^{2}\right)\right] \tag{4}
\end{equation*}
$$

where the modified polarization vector components are given by

$$
\begin{equation*}
\varepsilon_{0}=\left[\epsilon_{0}-\frac{\nu}{|\boldsymbol{q}|}(\boldsymbol{\epsilon} \cdot \hat{\boldsymbol{q}})\right] \frac{|\boldsymbol{q}|}{Q}, \quad \boldsymbol{\varepsilon}=\boldsymbol{\epsilon}-\hat{\boldsymbol{q}}(\boldsymbol{\epsilon} \cdot \hat{\boldsymbol{q}}) \tag{5}
\end{equation*}
$$

with $\boldsymbol{q}$ and $\hat{\boldsymbol{q}}=\boldsymbol{q} /|\boldsymbol{q}|$ the photon three-momentum in the lab frame and its unit vector, and $\left(\epsilon_{0}, \boldsymbol{\epsilon}\right)$ the time and space components of the photon polarisation vector.

Equations (2) and (4) give two equivalent decompositions of the amplitude in terms of gauge invariant tensors, and any other deuteron-spin independent tensors have to vanish in
the sum of a gauge invariant subset of Feynman graphs. This represents an important nontrivial consistency check of our calculation, which is performed without choosing a particular electromagnetic gauge. Note that in what follows we omit the factor $-\mathcal{E}_{\lambda}^{*} \mathcal{E}^{\lambda}=\delta_{\sigma \sigma^{\prime}}$, where $\sigma$ and $\sigma^{\prime}$ are the initial and final deuteron polarisations.

The VVCS amplitudes can be split into the pole (or elastic) and non-pole (or inelastic) parts [31]. The former have a pole at $\nu= \pm Q^{2} /\left(2 M_{d}\right)$ and correspond to the VVCS process going through the unexcited deuteron in the intermediate state; they are parametrised by the deuteron elastic electromagnetic form factors. The latter, on the other hand, admit an expansion in powers of the (small) energies and momenta, and are parametrised by the deuteron (generalised) polarisabilities.

To further arrive at the counting for the VVCS amplitude, we turn to the non-pole part of the amplitude, and first consider its low-energy and low-momenta expansion. The leading terms of that expansion of the non-pole pieces of $f_{L}\left(\nu, Q^{2}\right)$ and $f_{T}\left(\nu, Q^{2}\right)$ are given by the deuteron electric and magnetic dipole polarisabilities $\alpha_{E 1}$ and $\beta_{M 1}$ as [31]

$$
\begin{align*}
& f_{L}\left(\nu, Q^{2}\right)=4 \pi \alpha_{E 1} Q^{2}+\ldots,  \tag{6}\\
& f_{T}\left(\nu, Q^{2}\right)=-\frac{e^{2}}{M_{d}}+4 \pi \beta_{M 1} Q^{2}+4 \pi\left(\alpha_{E 1}+\beta_{M 1}\right) \nu^{2}+\ldots, \tag{7}
\end{align*}
$$

where the dots denote terms at least quartic in $\nu$ and $Q$, and $e$ is the proton charge; the first term in the expansion of $f_{T}\left(\nu, Q^{2}\right)$ - the Thomson term - corresponds to the point-like deuteron. We neglect other terms generated by the expansion of the non-pole pieces of the Born contributions, such as, e.g., the term $e^{2} R_{C}^{2} Q^{2} /(6 M)$, where $R_{C}$ and $M=\left(M_{n}+M_{p}\right) / 2$ are the deuteron charge radius and the average nucleon mass, respectively, in the expansion of $f_{T}$ [32]; see also, e.g., the related discussion in Ref. [33] for the case of a proton. One can show that such terms are all demoted to higher orders in the $đ$ EFT counting than those we consider; in any case, they give numerically negligible contributions to the VVCS amplitude.

## B. EFT expansion and counting

To set up the counting for the deuteron VVCS, we recall that the $\not \subset$ EFT expansion is organised in powers of the ratio $P / m_{\pi}$, where $P$ is the typical momentum scale in the system. Typical energies are counted as $E \propto P^{2}$ : the two-nucleon system is considered as predominantly non-relativistic, and relativistic corrections are taken into account perturbatively. Correspondingly, a nucleon propagator counts as $1 / E \propto P^{-2}$, whereas a loop gives a factor of $P^{5}$ corresponding to an integration over $d^{4} P=\mathrm{d} E \mathrm{~d}^{3} P$. To assign a particular order to a Feynman graph, one counts powers of momenta coming from the interaction vertices, propagators, and loops, assuming that all momenta are of the typical size $\sim P$ (and all energies of the size $\sim P^{2}$ ).

Looking at the leading terms in the $\not \subset E F T$ expansion of $\alpha_{E 1}$ and $\beta_{M 1}[14,25,34]$, one
can see that they are, respectively, $O\left(P^{-4}\right)$ and $O\left(P^{-2}\right)$ :

$$
\begin{align*}
& \alpha_{E 1}=\frac{\alpha M}{32 \pi \gamma^{4}}+\ldots,  \tag{8}\\
& \beta_{M 1}=-\frac{\alpha}{32 M \gamma^{2}}\left[1-\frac{16}{3} \mu_{1}^{2}+\frac{32}{3} \mu_{1}^{2} \frac{\gamma}{\gamma_{s}-\gamma}\right]+\ldots, \tag{9}
\end{align*}
$$

where $\mu_{1}$ is the nucleon isovector magnetic moment (in nucleon magneton units), and $\gamma_{s} \equiv$ $a_{s}^{-1}=O(P)$ is the inverse proton-neutron singlet scattering length. Since $Q^{2}=\boldsymbol{q}^{2}-\nu^{2}=$ $O\left(P^{2}\right)$ and $\nu^{2}=O\left(P^{4}\right)$, we can see, respectively, from Eqs. (6) and (7) that $f_{L}\left(\nu, Q^{2}\right)$ starts at $O\left(P^{-2}\right)$, and $f_{T}\left(\nu, Q^{2}\right)$ starts two orders higher at $O\left(P^{0}\right)$. This derivation applies to the non-pole parts of the amplitudes. However, it is straightforward to deduce that the same counting holds also for the respective pole parts, which is evident from the expressions for the residues of $f_{L}\left(\nu, Q^{2}\right)$ and $f_{T}\left(\nu, Q^{2}\right)$ considered in Sec. IV.

Furthermore, in the TPE correction to the deuterium Lamb shift, one can notice that $f_{T}$ is weighted with another small factor of $O\left(P^{2}\right)$ relative to $f_{L}$ in the integral for the TPE correction [21, 23]. In other words, $f_{T}$ starts to contribute to the TPE correction only at N4LO relative to the leading longitudinal contribution.

To identify the highest order in the $\not \subset E F T$ expansion where one can still hope to obtain a predictive result for the TPE correction, one can notice that the contribution of the polarisabilities of individual nucleons to $f_{L} \propto \alpha_{E 1, N} Q^{2}$ arises at N4LO in the $\not \subset E F T$ counting. This term goes as $Q^{2}$ at large $Q$ and thus leads to a divergent contribution to the TPE correction, which demands an unknown two-nucleon one-lepton contact term at this order to regularise the divergence. Such a contact term would be fitted to the TPE correction (or another two-nucleon one-lepton observable - if such data were available). Consequently, the predictive power of $\not \subset$ EFT for the TPE correction is lost beyond N3LO. In accordance to that, we concentrate on the longitudinal amplitude and calculate $f_{L}\left(\nu, Q^{2}\right)$ up to N3LO, or $O(P)$. Taking the $\not \subset$ EFT expansion parameter as $\gamma / m_{\pi} \sim 1 / 3$, this leads to a naïvely expected relative uncertainty of the calculation of $\sim\left(\gamma / m_{\pi}\right)^{4} \simeq 1 \%$.

This consideration indicates that the transverse amplitude starts to enter the TPE correction one order higher, and only the longitudinal amplitude is needed in case one is solely aiming at an N3LO calculation of the TPE correction. However, we in addition calculate $f_{T}\left(\nu, Q^{2}\right)$ up to $O(P)$, or the respective NLO, in order to precisely quantify the smallness of its contribution to the TPE correction, and also to investigate the (generalised) deuteron magnetic polarisability $\beta_{M 1}\left(Q^{2}\right)$ and the deuteron generalised Baldin sum rule.

To conclude the discussion of the $\not \subset E F T$ counting and expansion, we consider the $N N$ $T$-matrix in the spin-triplet (deuteron) channel, given by

$$
\begin{equation*}
T(k)=-\frac{4 \pi}{M} \frac{1}{-\gamma-i k+\frac{1}{2} \rho_{d}\left(k^{2}+\gamma^{2}\right)+w_{2}\left(k^{2}+\gamma^{2}\right)^{2}+\ldots}, \tag{10}
\end{equation*}
$$

where $k$ is the $N N$ relative momentum, $\rho_{d}$ and $w_{2}$ are the deuteron effective range and shape parameter, and terms of higher orders in $\left(k^{2}+\gamma^{2}\right)$ are not shown explicitly. Ref. [25] notes
that, while the more conventional $\rho$ expansion reproduces the deuteron effective range $\rho_{d}$ at NLO in the $\neq$ EFT expansion:

$$
\begin{equation*}
T(k)=-\frac{4 \pi}{M}[\underbrace{\frac{1}{-\gamma-i k}}_{\text {LO }}+\underbrace{\frac{\gamma \rho_{d}}{-\gamma-i k}+\frac{\gamma \rho_{d}}{2 \gamma}}_{\text {NLO }}+\ldots] \tag{11}
\end{equation*}
$$

one can use an alternative scheme, the $Z$ expansion, choosing instead to reproduce the residue of $T(k)$ at the deuteron pole $k=i \gamma$ at NLO:

$$
\begin{equation*}
T(k)=-\frac{4 \pi}{M}[\underbrace{\frac{1}{-\gamma-i k}}_{\mathrm{LO}}+\underbrace{\frac{Z-1}{-\gamma-i k}+\frac{Z-1}{2 \gamma}}_{\mathrm{NLO}}+\ldots] \tag{12}
\end{equation*}
$$

with the residue $Z$ given by

$$
\begin{equation*}
Z=\frac{1}{1-\gamma \rho_{d}}=1+\gamma \rho_{d}+\left(\gamma \rho_{d}\right)^{2}+\ldots \tag{13}
\end{equation*}
$$

Each scheme introduces a new $O(P)$ small parameter: $\gamma \rho_{d}$, or $Z-1$. The leading-order (LO) $O\left(P^{-1}\right)$ result is the same in both schemes. In fact, they only start to differ at NNLO, i.e., $O(P)$. Furthermore, both schemes ensure that the deuteron pole is located at $k=i \gamma$ at all orders in the expansion, as obtained by re-summation of an infinite chain of diagrams shown in Fig. 1.


FIG. 1. The LO NN T-matrix for the spin-triplet channel. Here and below, big disc (square) vertices denote insertions of a $N N T$-matrix ( $N N$ potential), with the number inside a vertex showing the order of the vertex in the $\not \subset E F T$ counting. The LO $T$-matrix in the spin-singlet ${ }^{1} S_{0}$ channel is obtained analogously.

The residue $Z$ is connected to $A_{S}$, the asymptotic normalisation factor of the deuteron $S$-wave,

$$
\begin{equation*}
\psi(r) \underset{r \rightarrow \infty}{\longrightarrow} \frac{A_{S}}{\sqrt{4 \pi}} \frac{e^{-\gamma r}}{r}=\sqrt{\frac{\gamma Z}{2 \pi}} \frac{e^{-\gamma r}}{r} \tag{14}
\end{equation*}
$$

As argued in Ref. [25], the $Z$ expansion is better suited for quantities that receive mostly long-range contributions and are hence sensitive to the correct description of the long-range tail of the deuteron wave function, one example of such a long-range quantity being the deuteron electric polarisability $\alpha_{E 1}$. We too adopt the Z expansion for our calculation.

## C. Lagrangian and coupling constants

The Lagrangian needed for our calculation is constructed along the usual lines formulated in, e.g, Refs. [4-7, 25, 35, 36], peforming a non-relativistic expansion in the one-nucleon sector and writing out the relevant two-nucleon interactions. The relativistic corrections in both the single-nucleon and two-nucleon sector count as $O\left(P^{2} / M^{2}\right)=m_{\pi}^{2} / M^{2} O\left(P^{2} / m_{\pi}^{2}\right)$ and are therefore more suppressed numerically than suggested by counting powers of momenta. Even when generously counting $m_{\pi} / M \sim P$, as suggested in, e.g., Ref. [36], the relativistic corrections start to appear at N4LO. We therefore neglect them and define $\gamma=\sqrt{-M E_{d}}$, where $E_{d}$ is the deuteron energy relative to the proton and neutron at rest.

We also neglect the isospin violation due to both the proton-neutron mass difference and the isospin-violating terms in the $N N$ interaction, with the caveat that the $N N$ interactions in the singlet channel are fitted to the empirical singlet $p n$ scattering length and effective range and therefore include some isospin-violating effects in that channel.

The one-nucleon Lagrangian needed for the present VVCS calculation reads

$$
\begin{equation*}
\mathcal{L}^{N}=N^{\dagger}\left[i D_{0}+\frac{\boldsymbol{D}^{2}}{2 M}\right] N+\frac{e}{2 M} N^{\dagger} \hat{\mu}(\boldsymbol{\sigma} \cdot \boldsymbol{B}) N+\frac{e}{6} N^{\dagger} \hat{r}_{E}^{2} N(\boldsymbol{\nabla} \cdot \boldsymbol{E}) . \tag{15}
\end{equation*}
$$

The gauge derivatives are defined as

$$
\begin{equation*}
D_{0} N=\left(\partial_{0}+i e \hat{Q} A_{0}\right) N, \quad \boldsymbol{D} N=(\boldsymbol{\nabla}-i e \hat{Q} \boldsymbol{A}) N \tag{16}
\end{equation*}
$$

with the electromagnetic potential $\left(A_{0}, \boldsymbol{A}\right)$ and the nucleon charge operator

$$
\begin{equation*}
\hat{Q}=\frac{1}{2}\left(1+\tau_{3}\right) . \tag{17}
\end{equation*}
$$

The electromagnetic field components are

$$
\begin{equation*}
\boldsymbol{B}=\boldsymbol{\nabla} \times \boldsymbol{A}, \quad \boldsymbol{E}=-\boldsymbol{\nabla} A_{0}-\partial_{0} \boldsymbol{A}, \tag{18}
\end{equation*}
$$

and the nucleon magnetic moment and charge radius operators are defined by

$$
\begin{equation*}
\hat{\mu}=\mu_{0}+\mu_{1} \tau_{3}, \quad \hat{r}_{E}^{2}=r_{0}^{2}+r_{1}^{2} \tau_{3} \tag{19}
\end{equation*}
$$

with $\mu_{0,1}=\left(\mu_{p} \pm \mu_{n}\right) / 2$ and $r_{0,1}^{2}=\left(r_{p}^{2} \pm r_{n}^{2}\right) / 2$ the nucleon isoscalar and isovector magnetic moments (in nuclear magneton units) and charge radii, respectively. The proton charge radius squared appearing here also contains the Darwin-Foldy term $3 /\left(4 M_{p}^{2}\right)$, see, e.g., Ref. [37] for a derivation.

Note that different parts of the minimal charge coupling are of different orders in the \#EFT counting: the longitudinal coupling is $\propto A_{0}=O\left(P^{0}\right)$, whereas the transverse coupling is $\propto \boldsymbol{\nabla} \cdot \boldsymbol{A}=O(P)$, and the seagull term is $\propto \boldsymbol{A}^{2}=O\left(P^{0}\right)$. Furthermore, while $\boldsymbol{B}=O(P)$, $\boldsymbol{E}$ contains two parts that are of different orders, $\boldsymbol{\nabla} A_{0}=O(P)$ and $\partial_{0} \boldsymbol{A}=O\left(P^{2}\right)$. The gauge invariance thus mixes different orders in the $đ$ EFT counting. This is consistent with the transverse amplitude being suppressed in relation to the longitudinal one.

The Lagrangian describing $N N$ interactions in the triplet $S$-wave up to N3LO and in the singlet $S$-wave up to NLO is given by

$$
\begin{align*}
\mathcal{L}_{S}^{N N} & =-C_{0} N^{\dagger} \mathcal{P}_{i} N_{c} N_{c}^{\dagger} \mathcal{P}_{i} N-C_{0}^{s} N^{\dagger} \mathcal{T}_{a} N_{c} N_{c}^{\dagger} \mathcal{T}_{a} N \\
& +\frac{1}{2} C_{2}\left[N^{\dagger} \mathcal{P}_{i} N_{c} N_{c}^{\dagger} \mathcal{O}_{i}^{(2)} N+\text { H.c. }\right]+\frac{1}{2} C_{2}^{s}\left[N^{\dagger} \mathcal{T}_{a} N_{c} N_{c}^{\dagger} \mathcal{O}_{a}^{(2, s)} N+\text { H.c. }\right] \\
& -C_{4} N^{\dagger} \mathcal{O}_{i}^{(2)} N_{c} N_{c}^{\dagger} \mathcal{O}_{i}^{(2)} N-\frac{1}{2} \tilde{C}_{4}\left[N^{\dagger} \mathcal{P}_{i} N_{c} N_{c}^{\dagger} \mathcal{O}_{i}^{(4)} N+\text { H.c. }\right] \\
& +\frac{1}{2} C_{6}\left[N^{\dagger} \mathcal{O}_{i}^{(2)} N_{c} N_{c}^{\dagger} \mathcal{O}_{i}^{(4)} N+\text { H.c. }\right] . \tag{20}
\end{align*}
$$

Here, we defined the charge-conjugated nucleon field as

$$
\begin{equation*}
N_{c}=\tau_{2} \sigma_{2}\left(N^{\dagger}\right)^{T} ; \tag{21}
\end{equation*}
$$

note that

$$
\begin{equation*}
\boldsymbol{D} N_{c}=\left(\boldsymbol{\nabla}+i e \hat{Q}_{c} \boldsymbol{A}\right) N_{c}, \tag{22}
\end{equation*}
$$

with

$$
\begin{equation*}
\hat{Q}_{c}=\tau_{2} Q \tau_{2}=\frac{1}{2}\left(1-\tau_{3}\right) . \tag{23}
\end{equation*}
$$

The spin-triplet-isospin-singlet and spin-singlet-isospin-triplet projectors $\mathcal{P}$ and $\mathcal{T}$ that select the corresponding $N N$ states are defined as

$$
\begin{equation*}
\mathcal{P}_{i}=\frac{1}{\sqrt{8}} \sigma_{i}, \quad \mathcal{T}_{a}=\frac{1}{\sqrt{8}} \tau_{a} \tag{24}
\end{equation*}
$$

with the normalisation

$$
\begin{equation*}
\operatorname{Tr} \mathcal{P}_{i} \mathcal{P}_{j}^{\dagger}=\frac{1}{2} \delta_{i j}, \quad \operatorname{Tr} \mathcal{T}_{a} \mathcal{T}_{b}^{\dagger}=\frac{1}{2} \delta_{a b}, \quad \operatorname{Tr} \mathcal{P}_{i} \mathcal{T}_{b}^{\dagger}=0 \tag{25}
\end{equation*}
$$

The quadratic and quartic Galilean-invariant combinations of the nucleon gauge derivatives and projectors are defined as

$$
\begin{align*}
& \mathcal{O}_{i}^{(2)}=\frac{1}{4}\left[\overleftarrow{\boldsymbol{D}}^{2} \mathcal{P}_{i}-2 \overleftarrow{D}_{j} \mathcal{P}_{i} \vec{D}_{j}+\mathcal{P}_{i} \overrightarrow{\boldsymbol{D}}^{2}\right]  \tag{26}\\
& \mathcal{O}_{i}^{(4)}=\frac{1}{16}\left[\overleftarrow{\boldsymbol{D}}^{4} \mathcal{P}_{i}-4 \overleftarrow{\boldsymbol{D}}^{2} \overleftarrow{D}_{j} \mathcal{P}_{i} \vec{D}_{j}+4 \overleftarrow{D}_{j} \overleftarrow{D}_{k} \mathcal{P}_{i} \vec{D}_{k} \vec{D}_{j}+2 \overleftarrow{\boldsymbol{D}}^{2} \mathcal{P}_{i} \overrightarrow{\boldsymbol{D}}^{2}-4 \overleftarrow{D}_{j} \mathcal{P}_{i} \vec{D}_{j} \overrightarrow{\boldsymbol{D}}^{2}+\mathcal{P}_{i} \overrightarrow{\boldsymbol{D}}^{4}\right] \tag{27}
\end{align*}
$$

for the triplet $N N$ channel, and

$$
\begin{equation*}
\mathcal{O}_{i}^{(2, s)}=\frac{1}{4}\left[\overleftarrow{\boldsymbol{D}}^{2} \mathcal{T}_{i}-2 \overleftarrow{D}_{j} \mathcal{T}_{i} \vec{D}_{j}+\mathcal{T}_{i} \overrightarrow{\boldsymbol{D}}^{2}\right] \tag{28}
\end{equation*}
$$

for the singlet $N N$ channel. The arrows point in the direction of operation of the corresponding derivative. Note that there is a different definition of $\mathcal{O}_{i}^{(4)}$ in literature, e.g., Refs. [35, 36] use

$$
\begin{equation*}
\mathcal{O}_{i}^{(4)}=\frac{1}{16}\left[\overleftarrow{\boldsymbol{D}}^{4} \mathcal{P}_{i}-4 \overleftarrow{\boldsymbol{D}}^{2} \overleftarrow{D}_{j} \mathcal{P}_{i} \vec{D}_{j}+6 \overleftarrow{\boldsymbol{D}}^{2} \mathcal{P}_{i} \overrightarrow{\boldsymbol{D}}^{2}-4 \overleftarrow{D}_{j} \mathcal{P}_{i} \vec{D}_{j} \overrightarrow{\boldsymbol{D}}^{2}+\mathcal{P}_{i} \overrightarrow{\boldsymbol{D}}^{4}\right] \tag{29}
\end{equation*}
$$

The latter expression, however, is potentially problematic, because it does not produce the desirable Galilean-invariant form of the matrix element:

$$
\begin{equation*}
\left\langle\boldsymbol{p}_{1}^{\prime} \boldsymbol{p}_{2}^{\prime}\right| N^{\dagger} \mathcal{P}_{i} N_{c} N_{c}^{\dagger} \mathcal{O}_{i}^{(4)} N\left|\boldsymbol{p}_{1} \boldsymbol{p}_{2}\right\rangle \propto\left(\boldsymbol{p}_{1}-\boldsymbol{p}_{2}\right)^{4} \tag{30}
\end{equation*}
$$

Indeed, it is straightforward to see that

$$
\begin{align*}
\left(\boldsymbol{p}_{1}-\boldsymbol{p}_{2}\right)^{4} & =\left(\boldsymbol{p}_{1}^{2}-2 \boldsymbol{p}_{1} \cdot \boldsymbol{p}_{2}+\boldsymbol{p}_{2}^{2}\right)^{2} \\
& =\boldsymbol{p}_{1}^{2}-4 \boldsymbol{p}_{1}^{2} \boldsymbol{p}_{1} \cdot \boldsymbol{p}_{2}+4\left(\boldsymbol{p}_{1} \cdot \boldsymbol{p}_{2}\right)^{2}+2 \boldsymbol{p}_{1}^{2} \boldsymbol{p}_{2}^{2}-4 \boldsymbol{p}_{1} \cdot \boldsymbol{p}_{2} \boldsymbol{p}_{2}^{2}+\boldsymbol{p}_{2}^{4} \tag{31}
\end{align*}
$$

showing that the operator in Eq. (27) produces the correct contractions of $\boldsymbol{p}_{1}$ and $\boldsymbol{p}_{2}$, whereas Eq. (29) produces an expression that cannot be reduced to $\left(\boldsymbol{p}_{1}-\boldsymbol{p}_{2}\right)^{4}$ in the general case. This issue appears to also affect the sixth-power operator $\mathcal{O}_{i}^{(6)}$ defined in Ref. [36] (this operator starts contributing at N4LO and is therefore not relevant to our calculation). One has to note that Eq. (29) gives the correct result in the center-of-mass frame, and, since it is a relatively high-order operator, the terms that spoil the Galilean invariance are probably rather small in a typical calculation; however, it is generally safer to use the expression given in Eq. (27).

For the $S D$ mixing $N N$ interaction, we consider the assignment where the corresponding coupling is $O\left(P^{-1}\right)$ [7,36], and, consequently, this interaction first appears at $O(P)$ in the $N N$ potential, which corresponds to NNLO. However, the $S D$ mixing term being proportional to a $D$-wave $N N$ operator and there being no other $D$-waves up to N3LO, one needs two insertions of this term in order to receive a contribution to unpolarised VVCS. This demotes the leading $S D$ mixing contribution to $O\left(P^{3}\right)$ in the $N N$ potential, or N4LO, making it irrelevant to the present calculation.

In addition to the $S$-wave interactions, we also include the spin-triplet-isospin-triplet interactions in a $P$-wave, entering at N3LO:

$$
\begin{align*}
\mathcal{L}_{P}^{N N} & =-\frac{1}{4}\left[C_{3_{P_{0}}} \delta_{i a} \delta_{j b}+C_{3_{P_{1}}}\left(\delta_{i j} \delta_{a b}-\delta_{i b} \delta_{j a}\right)+2 C_{3_{P_{2}}}\left(\delta_{i j} \delta_{a b}+\delta_{i b} \delta_{j a}-\frac{2}{3} \delta_{i a} \delta_{j b}\right)\right] \\
& \times N^{\dagger} \mathcal{O}_{i a}^{(1, P)} N_{c} N_{c}^{\dagger} \mathcal{O}_{j b}^{(1, P)} N, \tag{32}
\end{align*}
$$

with

$$
\begin{equation*}
\mathcal{O}_{i a}^{(1, P)}=\overleftarrow{D}_{i} P_{a} \tau_{3}-\tau_{3} P_{a} \vec{D}_{i} \tag{33}
\end{equation*}
$$

To complete the Lagrangian in the two-nucleon sector, one has to include two-nucleon contact interactions with the electromagnetic fields, whose couplings are not fixed by the $N N$ interaction. The contact terms needed for our calculation are

$$
\begin{align*}
\mathcal{L}^{N N \gamma} & =e L_{1}^{M 1_{V}}\left[N^{\dagger} \mathcal{P}_{i} N_{c} N_{c}^{\dagger} \mathcal{T}_{3} N+\text { H.c. }\right] B_{i}-2 i e L_{2}^{M 1 S} \epsilon_{i j k} N^{\dagger} \mathcal{P}_{i} N_{c} N_{c}^{\dagger} \mathcal{P}_{j} N B_{k} \\
& -\frac{e}{2} L_{1}^{E 1_{V}}\left[N^{\dagger} \mathcal{O}_{i j}^{(1, P)} N_{c} N_{c}^{\dagger} \mathcal{P}_{j} N+\text { H.c. }\right] E_{i}+\frac{e}{2} L_{3}^{E 1_{V}}\left[N^{\dagger} \mathcal{O}_{i j}^{(1, P)} N_{c} N_{c}^{\dagger} \mathcal{O}_{j}^{(2)} N+\text { H.c. }\right] E_{i} \\
& +e L_{1}^{C 0_{S}} N^{\dagger} \mathcal{P}_{i} N_{c} N_{c}^{\dagger} \mathcal{P}_{i} N(\boldsymbol{\nabla} \cdot \boldsymbol{E})-\frac{e}{2} L_{3}^{C 0_{S}}\left[N^{\dagger} \mathcal{P}_{i} N_{c} N_{c}^{\dagger} \mathcal{O}_{i}^{(2)} N+\text { H.c. }\right](\boldsymbol{\nabla} \cdot \boldsymbol{E}) . \tag{34}
\end{align*}
$$

The first two magnetic interactions contribute to $f_{T}$ at $O(P)$, or NLO, whereas the four electric contact terms contribute to $f_{L}$ also at $O(P)$, or respective N3LO. Note that the convention in literature is to add to the $L_{2}^{M 11_{S}}$ piece its Hermitian conjugate term, which, however, just gives the factor of 2 that we write out explicitly.

$$
\begin{array}{l|l|l}
\alpha=e^{2} / 4 \pi=1 / 137.036 & \hbar c=197.327 \mathrm{MeV} \mathrm{fm} & m_{\pi}=139.570 \mathrm{MeV} \\
\hline M=938.919 \mathrm{MeV} & \mu_{p}=2.793 & \mu_{n}=-1.913 \\
\hline r_{p}=0.84087 \mathrm{fm} & r_{n}^{2}=-0.1161 \mathrm{fm}^{2} & M_{d}=1875.613 \mathrm{MeV} \\
\hline E_{d}=-2.224575 \mathrm{MeV}[38] & Z=1.6893(30)[38] & w_{2}=0.389 \mathrm{fm}^{3}[35] \\
\hline \gamma_{s}=-8.3208 \mathrm{MeV}[39] & r_{s}=2.750 \mathrm{fm}[39] & C_{{ }_{3} P_{J}}=-1.49 \mathrm{fm}^{4}[35]
\end{array}
$$

TABLE I. Values of parameters entering the calculation of the VVCS amplitude. Unless referred to otherwise, the values are taken from the Particle Data Group listing [40] (with the value of $r_{p}$ corresponding to the $\mu \mathrm{H}$ Lamb shift, and not including the Darwin-Foldy term). See the text for additional information, and also for the values of the two-nucleon electromagnetic couplings.

The coupling constants of the single-nucleon sector are well known, and we show their values in Table I, which also contains the values of the $N N$ parameters used by us. The parameters taken from the Particle Data Group listing [40] are shown rounded, and their uncertainties are negligibly small compared to the projected N3LO precision of $\sim 1 \%$. The empirical values of $E_{d}$ and $Z$ are taken from the recent review [38], with $Z$ calculated from the $S$-wave asymptotic normalisation factor. While the value of $Z$ appears the most important source of uncertainty due to the input parameters, its quoted uncertainty is below $0.2 \%$ and can thus also be neglected in the final error estimate, whose dominant source is the omitted higher-order terms in the $\not \subset$ EFT expansion.

The $S$-wave $N N$ coupling constants are fixed by reproducing the expansion of the $N N$ $T$-matrix in the triplet and singlet channels in $\not \approx E F T$. Throughout our calculation, we use the dimensional regularisation and the power divergence subtraction (PDS) scheme [5, 6] to regularise divergent loop integrals. This introduces a regularisation scale dependence into the calculation, for instance, the loop function that corresponds to a single loop in Fig. 1 in this scheme is

$$
\begin{equation*}
I_{0}(E)=i \int \frac{\mathrm{~d}^{4} l}{(2 \pi)^{4}} \frac{1}{\left[l_{0}-\frac{l^{2}}{2 M}+i 0\right]\left[E-l_{0}-\frac{l^{2}}{2 M}+i 0\right]} \xrightarrow{\mathrm{PDS}}-\frac{M}{4 \pi}(\mu-\sqrt{-M E-i 0}), \tag{35}
\end{equation*}
$$

where $E$ is the relative motion energy of the $N N$ pair, and $\mu$ the regularisation scale. The $N N$ coupling constants also depend on $\mu$ in such a way that the $T$-matrix is $\mu$-independent. Up to the order we are working at, the $N N$ coupling constants are expressed in terms of $\mu$, $\gamma,(Z-1)$, and $w_{2}$ in the triplet channel, and $\mu, \gamma_{s}$, and the singlet effective range $r_{s}$ in the singlet channel. Note that the constants in the triplet channel are expanded in powers of the expansion parameter as well:

$$
\begin{equation*}
C_{0}=C_{0}^{(-1)}+C_{0}^{(0)}+C_{0}^{(1)}+C_{0}^{(2)}+\ldots, \tag{36}
\end{equation*}
$$

and analogously for the other constants. This expansion is needed in order to keep the position of the deuteron pole and the value of the residue unchanged as higher-order corrections are included. We provide the expressions for the constants entering Eq. (20) below for the sake of completeness. The triplet couplings read

$$
\begin{align*}
C_{0}^{(-1)} & =-\frac{4 \pi}{M} \frac{1}{\mu-\gamma}, \quad C_{0}^{(0)}=\frac{2 \pi}{M} \frac{(Z-1) \gamma}{(\mu-\gamma)^{2}}, \quad C_{0}^{(1)}=\frac{\pi}{M} \frac{(Z-1)^{2} \gamma(\gamma-2 \mu)}{(\mu-\gamma)^{3}}, \\
C_{0}^{(2)} & =\frac{\pi}{2 M} \frac{(Z-1)^{3} \gamma(\gamma-2 \mu)^{2}}{(\mu-\gamma)^{4}}+\frac{4 \pi}{M} \frac{w_{2} \gamma^{4}}{(\mu-\gamma)^{2}}, \\
C_{2}^{(-2)} & =\frac{2 \pi}{M} \frac{(Z-1)}{\gamma(\mu-\gamma)^{2}}, \quad C_{2}^{(-1)}=-\frac{2 \pi}{M} \frac{(Z-1)^{2} \mu}{\gamma(\mu-\gamma)^{3}}, \\
C_{2}^{(0)} & =-\frac{\pi}{2 M} \frac{(Z-1)^{3}\left(\gamma^{2}-4 \mu^{2}\right)}{\gamma(\mu-\gamma)^{4}}+\frac{8 \pi}{M} \frac{w_{2} \gamma^{2}}{(\mu-\gamma)^{2}}, \\
C_{4}^{(-3)} & =-\frac{\pi}{M} \frac{(Z-1)^{2}}{\gamma^{2}(\mu-\gamma)^{3}}, \quad S_{4}^{(-2)}=-\frac{\pi}{2 M} \frac{(Z-1)^{3}(\gamma-4 \mu)}{\gamma^{2}(\mu-\gamma)^{4}}+\frac{4 \pi}{M} \frac{w_{2}}{(\mu-\gamma)^{2}}, \\
C_{6}^{(-4)} & =\frac{\pi}{2 M} \frac{(Z-1)^{3}}{\gamma^{3}(\mu-\gamma)^{4}} . \tag{37}
\end{align*}
$$

Here, $S_{4}^{(-2)}=C_{4}^{(-2)}+\tilde{C}_{4}^{(-2)}$ is the only linear combination of $C_{4}^{(-2)}$ and $\tilde{C}_{4}^{(-2)}$ that contributes to $N N$ scattering. We note that for the following consideration, it is convenient to eliminate the coupling $C_{4}^{(-2)}$, expressing it via $S_{4}^{(-2)}$ and $\tilde{C}_{4}^{(-2)}$. The LO, NLO, and NNLO couplings given here coincide with those given in Ref. [25], however, we are not aware of any of the N3LO couplings explicitly appearing in the literature. Finally, the singlet couplings are

$$
\begin{equation*}
C_{0}^{s}=-\frac{4 \pi}{M} \frac{1}{\mu-\gamma_{s}}, \quad C_{2}^{s}=\frac{2 \pi}{M} \frac{r_{s}}{\left(\mu-\gamma_{s}\right)^{2}}, \tag{38}
\end{equation*}
$$

and the $P$-waves couplings contribute in a single linear combination

$$
\begin{equation*}
C_{{ }_{3} P_{J}}=C_{{ }_{3} P_{0}}+2 C_{{ }_{3} P_{1}}+\frac{20}{3} C_{{ }_{3} P_{2}}, \tag{39}
\end{equation*}
$$

whose value as extracted in Ref. [35] from the Nijmegen partial wave analysis [41, 42] is shown in Table I.

The couplings of the contact terms entering $\mathcal{L}^{N N \gamma}$ cannot be inferred from the parameters of the $N N$ scattering amplitude; these contact terms, however, are crucial in compensating the $\mu$ dependence of the VVCS amplitude, and information on their couplings can be obtained from the renormalisation group (RG) equations and from processes involving external electromagnetic probes. We discuss the role played by these contact terms as well as their determination in detail below in Sec. III.

## D. LSZ reduction

In order to calculate the VVCS amplitude, we follow the LSZ reduction procedure along the lines described in, e.g., Ref. [43] for a calculation of the deuteron electromagnetic form
factors. We introduce the coupling of the $N N$ system to an interpolating deuteron field that has the appropriate quantum numbers:

$$
\begin{equation*}
\delta \mathcal{L}^{N N d}=N^{\dagger} \mathcal{P}_{i} N_{c} \mathcal{E}_{i}+\text { H.c. } \tag{40}
\end{equation*}
$$

where $\mathcal{E}_{i}$ are the spatial components of the deuteron polarisation vector (which does not have a zeroth component in the lab frame). The VVCS amplitude is expressed via the sum of all irreducible four-point $(\gamma d \rightarrow \gamma d)$ functions $\mathcal{M}\left(q, p, q^{\prime}, p^{\prime}\right)$ divided by the derivative of the deuteron self-energy $\Sigma(E)$ taken at the deuteron pole:

$$
\begin{equation*}
T_{f i}=\frac{\mathcal{M}\left(q, p, q^{\prime}, p^{\prime}\right)}{\Sigma^{\prime}\left(E_{d}\right)} \tag{41}
\end{equation*}
$$

The irreducibility of $\mathcal{M}$ is defined in a somewhat more complicated way than in a calculation of the deuteron form factors, namely, a diagram is reducible if it contains an insertion of the LO $N N$ potential (or, equivalently, the LO $N N T$-matrix, see Fig. 1) in the triplet channel at energy equal to the deuteron binding energy, i.e., at the deuteron pole. An insertion of this $T$-matrix in an off-shell kinematics is, however, allowed in an irreducible diagram. This is illustrated in Fig. 2. The self energy, in turn, is defined as the sum of all deuteron-deuteron

(a)

(b)

(c)

(d)

FIG. 2. Examples of reducible (a), (b) and irreducible (c), (d) diagrams. The crossed vertex denotes the leading-order $O\left(P^{0}\right) N N d$ coupling.
two-point functions without any insertions of the LO $N N T$-matrix. Note that by virtue of Eq. (41) the normalisation of the interpolating deuteron field in $\delta \mathcal{L}^{N N d}$ is arbitrary. To obtain the order-by-order results, both $\mathcal{M}$ and $\Sigma^{\prime}\left(E_{d}\right)$ have to be expanded, and their ratio has to be expanded after that as well:

$$
\begin{align*}
\mathcal{M}_{L}\left(\nu, Q^{2}\right) & =\mathcal{M}_{L}^{(-3)}+\mathcal{M}_{L}^{(-2)}+\ldots,  \tag{42}\\
\mathcal{M}_{T}\left(\nu, Q^{2}\right) & =\mathcal{M}_{T}^{(-1)}+\mathcal{M}_{T}^{(0)}+\ldots,  \tag{43}\\
\Sigma^{\prime}\left(E_{d}\right) & =\Sigma^{\Sigma^{(-1)}}+\Sigma^{\prime(0)}+\ldots,  \tag{44}\\
f_{L}\left(\nu, Q^{2}\right) & =\underbrace{\frac{\mathcal{M}_{L}^{(-3)}}{\Sigma^{(-1)}}}_{O\left(P^{-2}\right)}+\underbrace{\frac{\mathcal{M}_{L}^{(-2)}}{\Sigma^{((-1)}}-\frac{\mathcal{M}_{L}^{(-3)} \Sigma^{\prime(0)}}{\left[\Sigma^{\prime(-1)}\right]^{2}}}_{O\left(P^{-1}\right)}+\ldots,  \tag{45}\\
f_{T}\left(\nu, Q^{2}\right) & =\underbrace{\frac{\mathcal{M}_{T}^{(-1)}}{\Sigma^{(-1)}}}_{O\left(P^{0}\right)}+\underbrace{\frac{\mathcal{M}_{T}^{(0)}}{\Sigma^{\prime(-1)}}-\frac{\mathcal{M}_{T}^{(-1)} \Sigma^{\prime(0)}}{\left[\Sigma^{\prime(-1)}\right]^{2}}}_{O\left(P^{1}\right)}+\ldots, \tag{46}
\end{align*}
$$

where $\mathcal{M}_{L, T}\left(\nu, Q^{2}\right)$ denote the terms in $\mathcal{M}$ contributing to $f_{L, T}\left(\nu, Q^{2}\right)$.

## III. DIAGRAMS AND PERTINENT RESULTS, ORDER BY ORDER

In this section, we provide details of the order-by-order amplitudes necessary for the extraction of the observables.

## A. Deuteron self energy

To make the graphical representation of the matrix elements relevant for the calculation of the self energy $\Sigma(E)$ and the irreducible four-point function $\mathcal{M}\left(\nu, Q^{2}\right)$ more compact, it is convenient to define certain subgraphs: the order-by-order corrections to the $N N T$-matrix, shown in Fig. 3 and expressed via the dressed potentials that include all possible insertions of the LO NN T-matrix in an intermediate state, shown in Fig. 4; as well as the order-byorder corrections to the $N N d$ vertex, which are shown in Fig. 5. Note that the $T$-matrix always appears in full off-shell kinematics.


FIG. 3. Correction to the $N N T$-matrix vertex at order $k \geq 0$, defined in terms of the dressed potential at that order (denoted by the square with the primed index $k^{\prime}$ ). Only the triplet channel is shown. The dressed potentials at the orders we are considering are shown in Fig. 4.





$+0$
FIG. 4. Dressed potential at $O\left(P^{0}\right), O(P)$, and $O\left(P^{2}\right)$. Only the triplet channel is shown.

With these definitions, one obtains the very compact graphical form for $\Sigma(E)$ shown in Fig. 6. The resulting expression for $\Sigma(E)$ up to N3LO reads




FIG. 5. Corrections to the $N N d$ vertex at $O(P), O\left(P^{2}\right)$, and $O\left(P^{3}\right)$, with the numbers denoting the order of the correction. Not to be confused with the $N N T$ matrices defined in Figs. 1 and 3, which are denoted by bigger discs and have four external legs.


FIG. 6. Graphical expression for $i \Sigma(E)$ up to N3LO.

$$
\begin{align*}
\Sigma(E)=\frac{M}{4 \pi} & {[\mu+i k} \\
& +\frac{(Z-1)\left(\gamma^{2}+k^{2}\right)(k-i \mu)^{2}}{2 \gamma(\gamma-\mu)^{2}} \\
& +\frac{(Z-1)^{2}\left(\gamma^{2}+k^{2}\right)(\mu+i k)^{2}\left(\gamma\left(\gamma^{2}-2 \gamma \mu+2 \mu^{2}\right)+i k^{3}-k^{2}(\gamma-2 \mu)+i \gamma^{2} k\right)}{4 \gamma^{2}(\gamma-\mu)^{4}} \\
& -\frac{(Z-1)^{3}\left(\gamma^{2}+k^{2}\right)(k-i \mu)^{2}\left(-i \gamma\left(\gamma^{2}-2 \gamma \mu+2 \mu^{2}\right)+k^{3}+i k^{2}(\gamma-2 \mu)+\gamma^{2} k\right)^{2}}{8 \gamma^{3}(\gamma-\mu)^{6}} \\
& \left.+\frac{w_{2}\left(\gamma^{2}+k^{2}\right)^{2}(k-i \mu)^{2}}{(\gamma-\mu)^{2}}+\ldots\right], \tag{47}
\end{align*}
$$

where $k=i \sqrt{-M E}$, and the N3LO result occupies the last two lines. Even though $\Sigma(E)$ depends on $\mu$, its derivative at the deuteron pole, $\Sigma^{\prime}\left(E_{d}\right)$, is $\mu$-independent, and the order-by-order expression for it is very compact:

$$
\begin{equation*}
\Sigma^{\prime}\left(E_{d}\right)=\frac{M^{2}}{8 \pi \gamma}\left[1-(Z-1)+(Z-1)^{2}-(Z-1)^{3}+\ldots\right], \tag{48}
\end{equation*}
$$

giving an even simpler expression for the inverse quantity:

$$
\begin{equation*}
\left[\Sigma^{\prime}\left(E_{d}\right)\right]^{-1}=\frac{8 \pi \gamma}{M^{2}}[1+(Z-1)+0+0+\ldots] . \tag{49}
\end{equation*}
$$

As one could expect, the $Z$ factor in the residue is restored at NLO, and there are no corrections at higher orders.

## B. Irreducible four-point function

Here, we present the order-by-order results for the irreducible four-point function $\mathcal{M}$.

1. $L O$




FIG. 7. Diagrams contributing to $\mathcal{M}$ at LO, i.e., $O\left(P^{-3}\right)$ for $\mathcal{M}_{L}$ and $O\left(P^{-1}\right)$ for $\mathcal{M}_{T}$. Dotted vertices are the sum of the charge and the magnetic moment couplings from the single-nucleon Lagrangian; the big grey disc denotes the LO NN T-matrix in the singlet channel. Crossed graphs are not shown.

The diagrams contributing to $i \mathcal{M}$ at LO are shown in Fig. 7, resulting in

$$
\begin{align*}
\mathcal{M}_{L}^{(-3)} & =\frac{e^{2} M^{3}}{\pi} \frac{Q^{2}}{\boldsymbol{q}^{2}}\left[\frac{1}{2 \gamma\left[\boldsymbol{q}^{2}+4\left(\gamma+\lambda_{d}\right)^{2}\right]}-\frac{\phi^{2}\left(\nu, \boldsymbol{q}^{2}\right)}{\boldsymbol{q}^{2}\left(\gamma-\lambda_{d}\right)}\right]+(\nu \rightarrow-\nu),  \tag{50}\\
\mathcal{M}_{T}^{(-1)}=\frac{e^{2} M}{\pi}[ & -\frac{16 \gamma\left(\gamma-\lambda_{d}\right)\left(\gamma+\lambda_{d}\right)^{2}+4 \boldsymbol{q}^{2}\left(2 \gamma^{2}+\gamma \lambda_{d}+\lambda_{d}^{2}\right)-\left(4 \mu_{1}^{2}+4 \mu_{0}^{2}-1\right) \boldsymbol{q}^{4}}{16 \gamma \boldsymbol{q}^{2}\left[\boldsymbol{q}^{2}+4\left(\gamma+\lambda_{d}\right)^{2}\right]} \\
& +\left(\frac{|\boldsymbol{q}|\left(\mu_{0}^{2}-\mu_{1}^{2}\right)}{12 M \nu}+\frac{4 \gamma^{2}-4 \lambda_{d}^{2}+\boldsymbol{q}^{2}}{8|\boldsymbol{q}|^{3}}\right) \phi\left(\nu, \boldsymbol{q}^{2}\right) \\
& \left.-\frac{1}{3}\left(\frac{2 \mu_{0}^{2}}{\gamma-\lambda_{d}}+\frac{\mu_{1}^{2}}{\gamma_{s}-\lambda_{d}}\right) \phi^{2}\left(\nu, \boldsymbol{q}^{2}\right)\right]+(\nu \rightarrow-\nu), \tag{51}
\end{align*}
$$

where the kinematic functions are defined as

$$
\begin{equation*}
\lambda_{d}=\sqrt{\gamma^{2}-M \nu+\frac{\boldsymbol{q}^{2}}{4}}, \quad \phi\left(\nu, \boldsymbol{q}^{2}\right)=\arctan \frac{|\boldsymbol{q}|}{2\left(\gamma+\lambda_{d}\right)} . \tag{52}
\end{equation*}
$$

These expressions (as well as those at higher orders shown below) include the pole parts, with the elastic poles shifted due to the non-relativistic expansion and located at $\nu= \pm \boldsymbol{q}^{2} /(4 M)$. One can also note that all loop integrals at LO are convergent, so no $\mu$ dependence emerges at this step. Another important remark regarding the set of graphs in Fig. 7: together with the respective crossed graphs (not shown in the figure), the contribution of this set to the four-point function (and, correspondingly, to the VVCS amplitude) is gauge-invariant. This pattern is followed by the higher-order contributions presented below, where each
figure shows a set of graphs that, together with the respective crossed counterparts, gives a gauge-invariant result. Note that some of the graphs shown may only serve to recover the gauge invariance and give a zero contribution to $\mathcal{M}$, however, we do not show graphs whose contribution vanishes identically (due to, e.g., isospin or dimensional regularisation factors).

## 2. $N L O$

The NLO contributions to $\mathcal{M}$ come from diagrams shown in Figs. 8 and 9. While the
















FIG. 8. Contributions to $\mathcal{M}$ at NLO, i.e., $O\left(P^{-2}\right)$ for $\mathcal{M}_{L}$ and $O\left(P^{0}\right)$ for $\mathcal{M}_{T}$, due to NLO corrections in the $N N$ interaction. The notation is as in Fig. 7. Crossed graphs are not shown.
contributions to $\mathcal{M}_{L}^{(-2)}$ only come from the graphs in Fig. 8, and their sum is $\mu$ independent, $\mathcal{M}_{T}^{(0)}$ also receives contributions from the two magnetic contact terms in $\mathcal{L}^{N N \gamma}$, corresponding to the graphs in Fig. 9. The renormalisation scale dependence of the coupling constants $L_{1}^{M 1_{V}}$ and $L_{2}^{M 1_{S}}$ has to cancel out the $\mu$ dependence of the total NLO contribution to $\mathcal{M}_{T}$. We find the following RG equations for these two couplings:

$$
\begin{align*}
\mu \frac{\mathrm{d}}{\mathrm{~d} \mu}\left[(\mu-\gamma)\left(\mu-\gamma_{s}\right)\left(L_{1}^{M 1_{V}}-\frac{\mu_{1}}{2}\left\{C_{2}^{(-2)}+C_{2}^{(-2, s)}\right\}\right)\right] & =0  \tag{53}\\
\mu \frac{\mathrm{~d}}{\mathrm{~d} \mu}\left[\frac{L_{2}^{M 1_{S}}}{C_{2}^{(-2)}}\right] & =0 . \tag{54}
\end{align*}
$$

This is consistent with the results obtained in Refs. [14, 35, 36]. The first of these coupling constants, $L_{1}^{M 1_{V}}$, contributes to the deuteron $\beta_{M 1}$. The value of the RG-invariant quantity




FIG. 9. Contributions to $\mathcal{M}$ at NLO due to the magnetic contact terms $L_{1}^{M 1_{V}}$ and $L_{2}^{M 1_{S}}$ (shown as diamonds). The rest of the notation is as in Fig. 7. These graphs give an $O\left(P^{0}\right)$ contribution to $\mathcal{M}_{T}$ only. Crossed graphs are not shown.
in Eq. (53) can be fitted to, e.g., data on $n p \rightarrow d \gamma$; we use the result from Ref. [36] that found, also in the $Z$ expansion,

$$
\begin{equation*}
(\mu-\gamma)\left(\mu-\gamma_{s}\right)\left(L_{1}^{M 1_{V}}-\frac{\mu_{1}}{2}\left\{C_{2}^{(-2)}+C_{2}^{(-2, s)}\right\}\right)=-9.039(27) \mathrm{fm}^{2} \tag{55}
\end{equation*}
$$

The value of $L_{2}^{M 11_{S}}$ is fitted to the deuteron magnetic moment, obtaining [7, 43]

$$
\begin{equation*}
\left.L_{2}^{M 1_{S}}\right|_{\mu=m_{\pi}}=-0.149 \mathrm{fm}^{4} \tag{56}
\end{equation*}
$$

this result is the same at NLO both in the $Z$ and in the $\rho$ expansion. To make the expression for $\mathcal{M}_{T}^{(0)}$ more compact, we introduce dimensionless $\mu$-independent couplings $l_{1}^{M 1_{V}}$ and $l_{2}^{M 1{ }_{S}}$ according to

$$
\begin{align*}
& L_{1}^{M 1_{V}}=\frac{\mu_{1}}{2}\left(C_{2}^{(-2)}+C_{2}^{(-2, s)}\right)+\frac{\pi(Z-1)}{M \gamma} \frac{l_{1}^{M 1_{V}}}{(\mu-\gamma)\left(\mu-\gamma_{s}\right)},  \tag{57}\\
& L_{2}^{M 1_{S}}=l_{2}^{M 1_{S}} C_{2}^{(-2)} \tag{58}
\end{align*}
$$

The resulting expressions for $\mathcal{M}_{L}$ and $\mathcal{M}_{T}$ read

$$
\begin{align*}
& \mathcal{M}_{L}^{(-2)}=\frac{e^{2} M^{3}}{\pi} \frac{Q^{2}}{\boldsymbol{q}^{2}} \frac{(Z-1)}{2 \gamma} \frac{\phi\left(\nu, \boldsymbol{q}^{2}\right)\left[|\boldsymbol{q}|-\left(\gamma+\lambda_{d}\right) \phi\left(\nu, \boldsymbol{q}^{2}\right)\right]}{\boldsymbol{q}^{2}\left(\gamma-\lambda_{d}\right)}+(\nu \rightarrow-\nu)  \tag{59}\\
& \mathcal{M}_{T}^{(0)}=\frac{e^{2} M}{\pi}[ \frac{Z-1}{32 \gamma}-\frac{(Z-1)}{12 \gamma}|\boldsymbol{q}|\left(\frac{\mu_{1} l_{1}^{M 1_{V}}}{\gamma_{s}-\lambda_{d}}-\frac{4 \mu_{0}\left(\mu_{0}-2 l_{2}^{M 1 s}\right)}{\gamma-\lambda_{d}}\right) \phi\left(\nu, \boldsymbol{q}^{2}\right) \\
&\left.+\frac{1}{6}\left(\frac{\mu_{1}^{2} r_{s} \lambda_{d}^{2}}{\left(\gamma_{s}-\lambda_{d}\right)^{2}}-\frac{2(Z-1) \mu_{0}^{2}\left(\gamma+\lambda_{d}\right)}{\gamma\left(\gamma-\lambda_{d}\right)}\right) \phi^{2}\left(\nu, \boldsymbol{q}^{2}\right)\right]+(\nu \rightarrow-\nu) . \tag{60}
\end{align*}
$$

The values of the coupling constants $l_{1}^{M 1_{V}}$ and $l_{2}^{M 11_{S}}$, obtained using the values in Eqs. (55) and (56), are

$$
\begin{equation*}
l_{1}^{M 1_{V}}=-4.596(14), \quad l_{2}^{M 1_{S}}=-8.58 \times 10^{-3} . \tag{61}
\end{equation*}
$$

While $l_{1}^{M 1_{V}}$ could be considered as being of a natural size, $l_{2}^{M 1 s}$ is numerically small due to the LO contribution already coming very close to the empirical value of the deuteron magnetic moment, as pointed out in Refs. [7, 43].

## 3. $N N L O$

From this point on, we only keep track of those interactions that contribute to $f_{L}$, and, as a check, also include contributions to $\mathcal{M}$ that are needed to satisfy the electromagnetic gauge invariance. In practice this means taking into account the minimal coupling, both in the single-nucleon and in the two-nucleon Lagrangian. Incidentally, this allows one to also check that the Thomson term is recovered at NNLO and N3LO. The corresponding diagrams that appear at NNLO are shown in Figs. 10 and 11. At this order, one is still getting a













FIG. 10. Diagrams that contribute to $\mathcal{M}_{L}$ at NNLO, i.e., $O\left(P^{-1}\right)$, due to NNLO terms in the $N N$ interaction. The nucleon-photon vertex is exclusively the minimal coupling. In addition to diagrams that actually contribute to $f_{L}$, we also show those that are necessary in order to keep the electromagnetic gauge invariance. The vertical double dashed lines indicate possible insertions of a LO NN T-matrix in the spin-triplet channel. Crossed graphs are not shown.
$\mu$-independent contribution to $\mathcal{M}_{L}$ from loops with NNLO corrections to $N N$ interactions, Fig. 10. The resulting contribution to the longitudinal part of the four-point function is



FIG. 11. Correction to $\mathcal{M}_{L}$ due to the photon coupling proportional to $\hat{r}_{E}^{2}$, denoted by the black cross vertex, contributing at $O\left(P^{-1}\right)$. The rest of the notation is as in Fig. 10. Crossed graphs are not shown.

$$
\begin{equation*}
\mathcal{M}_{L}^{(-1)}=\frac{e^{2} M^{3}}{\pi} \frac{Q^{2}}{\boldsymbol{q}^{2}} \frac{(Z-1)^{2}}{4 \gamma^{2}}\left[-\frac{1}{4\left(\gamma-\lambda_{d}\right)}-\frac{\phi\left(\nu, \boldsymbol{q}^{2}\right)}{|\boldsymbol{q}|}+\frac{\left(\gamma+\lambda_{d}\right) \phi^{2}\left(\nu, \boldsymbol{q}^{2}\right)}{\boldsymbol{q}^{2}}\right]+(\nu \rightarrow-\nu) \tag{62}
\end{equation*}
$$

In addition, corrections arise due to the nucleon charge radius operator, shown in Fig. 11. Their contribution is

$$
\begin{align*}
\delta \mathcal{M}_{L}^{(-1)}= & -\frac{e^{2} M^{3}}{6 \pi} \frac{Q^{2}}{\boldsymbol{q}^{2}}\left[\frac{\left(r_{0}^{2}+r_{1}^{2}\right) \boldsymbol{q}^{2}}{\gamma\left[\boldsymbol{q}^{2}+4\left(\gamma+\lambda_{d}\right)^{2}\right]}+\frac{\left(r_{0}^{2}-r_{1}^{2}\right)|\boldsymbol{q}| \phi\left(\nu, \boldsymbol{q}^{2}\right)}{M \nu}-\frac{4 r_{0}^{2} \phi^{2}\left(\nu, \boldsymbol{q}^{2}\right)}{\gamma-\lambda_{d}}\right] \\
& +(\nu \rightarrow-\nu) \tag{63}
\end{align*}
$$

## 4. N3LO

Finally, the N3LO contribution to $\mathcal{M}_{L}$ comes from three types of diagrams: those with insertions of the N3LO $N N$ interactions, shown in Fig. 12, diagrams with an insertion of a gauge-invariant electric contact term, shown in Fig. 13, and corrections generated by one insertion of the nucleon charge radii coupling in the NLO graphs, as shown in Fig. 14. The diagrams in Fig. 12 produce a $\mu$-dependent result, and all four electric contact terms are needed in order to render the total result RG invariant. The couplings of these contact terms contribute to $\mathcal{M}_{L}^{(0)}$ in the following $\mu$-independent combinations with the $N N$ couplings:

$$
\begin{array}{ll}
\mu \frac{\mathrm{d}}{\mathrm{~d} \mu}\left[\frac{L_{1}^{E 1_{V}}-\frac{1}{2} M \tilde{C}_{4}^{(-2)}}{C_{0}^{(-1)}}\right]=0, & \mu \frac{\mathrm{~d}}{\mathrm{~d} \mu}\left[\frac{L_{3}^{E 1_{V}}-\frac{1}{2} M C_{6}^{(-4)}}{C_{0}^{(-1)}}\right]=0, \\
\mu \frac{\mathrm{~d}}{\mathrm{~d} \mu}\left[\frac{L_{1}^{C 0_{S}}+\frac{1}{4} M \tilde{C}_{4}^{(-2)}}{\left[C_{0}^{(-1)}\right]^{2}}\right]=0, & \mu \frac{\mathrm{~d}}{\mathrm{~d} \mu}\left[\frac{L_{3}^{C 0_{S}}+\frac{1}{4} M C_{6}^{(-4)}}{\left[C_{0}^{(-1)}\right]^{2}}\right]=0 . \tag{65}
\end{array}
$$

The first two equations have been previously obtained in Ref. [35]; our version, however, contains an additional factor $1 / 2$ in front of the nucleon mass in both equations, at variance with this reference. The second pair of equations is new, to the best of our knowledge.

Considering the $\mu$ running of the quantities entering these RG equations at high momentum scales $\mu \simeq m_{\pi}$, one can conclude that

$$
\begin{equation*}
L_{1}^{E 1_{V}}-\frac{1}{2} M \tilde{C}_{4}^{(-2)}=O\left(P^{-1}\right), \quad L_{3}^{E 1_{V}}-\frac{1}{2} M C_{6}^{(-4)}=O\left(P^{-1}\right) \tag{66}
\end{equation*}
$$















FIG. 12. Diagrams that contribute to $\mathcal{M}_{L}$ at N3LO, i.e., $O\left(P^{0}\right)$, due to N3LO terms in the $N N$ interaction. In addition to diagrams that actually contribute to $f_{L}$, we also show those that are necessary in order to keep the electromagnetic gauge invariance. Grey squares marked by "P" show insertions of the $P$-wave $N N$ interactions. The rest of the notation is as in Fig. 10. Crossed graphs are not shown.

$$
\begin{equation*}
L_{1}^{C 0_{S}}+\frac{1}{4} M \tilde{C}_{4}^{(-2)}=O\left(P^{-2}\right), \quad L_{3}^{C 0_{S}}+\frac{1}{4} M C_{6}^{(-4)}=O\left(P^{-2}\right) \tag{67}
\end{equation*}
$$

The first two combinations, being $O\left(P^{-1}\right)$ instead of the naively expected $O\left(P^{-2}\right)$ and $O\left(P^{-4}\right)$, are thus demoted to at least N4LO and N6LO, respectively. The same happens



FIG. 13. Contributions to $\mathcal{M}_{L}$ at N3LO due to the electric contact terms (shown as crossed diamonds). The rest of the notation is as in Fig. 10. Crossed graphs are not shown.
with the fourth combination, which is $O\left(P^{-2}\right)$ instead of $O\left(P^{-4}\right)$ and is demoted to N5LO. The only combination that gives a contribution at N3LO is the one that involves $L_{1}^{C 0_{S}}$. As we show in Sec. IV, its value can be found from a fit to the deuteron charge form factor.

Note that the cancellations between the contributions of the contact terms and those of the $N N$ couplings are in fact more intricate than given by these RG equations. The $N N$ coupling constants - all apart from $\tilde{C}_{4}^{(-2)}$ - conspire to remove the poles from the N3LO correction to the $N N T$-matrix. The instances of $C_{6}^{(-4)}$ appearing in the RG equations are in fact combinations of all the $N N$ constants appearing at this order, so a statement that $L_{3}^{E 1_{V}}$ and $L_{3}^{C 0_{S}}$ cancel the contribution of $C_{6}^{(-4)}$ might be somewhat imprecise. The constant $\tilde{C}_{4}^{(-2)}$, on the other hand, is singled out from the other $N N$ constants, so its cancellations with the contact terms show a more transparent pattern. In particular, the cancellation is complete at N3LO in the transverse amplitude, consistent with what was shown previously in Refs. [35, 36] (even though our RG equations do not completely coincide with those references).

As before, we write

$$
\begin{equation*}
L_{1}^{C 0_{S}}=-\frac{1}{4} M \tilde{C}_{4}^{(-2)}+\frac{\pi(Z-1)^{3}}{\gamma^{3}(\mu-\gamma)^{2}} l_{1}^{C 0_{S}} \tag{68}
\end{equation*}
$$

getting the following result for the total N3LO contribution of the diagrams in Figs. 12 and 13 :

$$
\begin{align*}
\mathcal{M}_{L}^{(0)} & =\frac{e^{2} M^{3}}{\pi} \frac{Q^{2}}{\boldsymbol{q}^{2}} \frac{(Z-1)^{3}}{\gamma^{3}}\left[\frac{3 \gamma-\lambda_{d}}{2\left(\gamma-\lambda_{d}\right)}+\frac{2 l_{1}^{C 0_{S}} \boldsymbol{q}^{2}+\left(\gamma-\lambda_{d}\right)^{2}}{8|\boldsymbol{q}|\left(\gamma-\lambda_{d}\right)} \phi\left(\nu, \boldsymbol{q}^{2}\right)-\frac{\gamma^{2}-\lambda_{d}^{2}}{8 \gamma^{3} \boldsymbol{q}^{2}} \phi^{2}\left(\nu, \boldsymbol{q}^{2}\right)\right] \\
& -\frac{e^{2} M^{3}}{\pi} \frac{Q^{2}}{\boldsymbol{q}^{2}} \frac{w_{2}\left[|\boldsymbol{q}|-2\left(\gamma+\lambda_{d}\right) \phi\left(\nu, \boldsymbol{q}^{2}\right)\right]^{2}}{4 \boldsymbol{q}^{2}} \\
& +\frac{e^{2} M^{3}}{\pi} \frac{Q^{2}}{\boldsymbol{q}^{2}} \frac{M C_{3 P_{J}}\left[2|\boldsymbol{q}|\left(\gamma-\lambda_{d}\right)-\left(4 \gamma^{2}-4 \lambda_{d}^{2}+\boldsymbol{q}^{2}\right) \phi\left(\nu, \boldsymbol{q}^{2}\right)\right]^{2}}{192 \pi \boldsymbol{q}^{4}}+(\nu \rightarrow-\nu) . \tag{69}
\end{align*}
$$

The nucleon charge radii corrections at this order, coming from the graphs in Fig. 14, yield

$$
\begin{equation*}
\delta \mathcal{M}_{L}^{(0)}=-\frac{e^{2} M^{3}}{3 \pi} \frac{Q^{2}}{\boldsymbol{q}^{2}} \frac{r_{0}^{2}(Z-1)}{\gamma} \frac{\phi\left(\nu, \boldsymbol{q}^{2}\right)\left[|\boldsymbol{q}|-\left(\gamma+\lambda_{d}\right) \phi\left(\nu, \boldsymbol{q}^{2}\right)\right]}{\gamma-\lambda_{d}}+(\nu \rightarrow-\nu) . \tag{70}
\end{equation*}
$$

Having established the order-by-order ingredients for the deuteron VVCS calculation throughout this section, we can now proceed with the extraction of the observables, form factors and generalised polarisabilities, in the following sections.








FIG. 14. Correction to $\mathcal{M}_{L}$ due to the photon coupling proportional to $\hat{r}_{E}^{2}$, contributing at $O\left(P^{0}\right)$. The rest of the notation is as in Figs. 10 and 11. Crossed graphs are not shown.

## IV. DEUTERON CHARGE FORM FACTOR AT N3LO: FITTING $l_{1}^{C 0_{S}}$

In order to extract the deuteron form factors, we use the residues of the pole parts of the VVCS amplitude. Using, e.g., the elastic structure functions [21], one can obtain:

$$
\begin{align*}
& \left.\operatorname{Res} f_{L}\left(\nu, Q^{2}\right)\right|_{\nu=Q^{2} /\left(2 M_{d}\right)}=-e^{2}\left(1+\tau_{d}\right)\left[G_{C}^{2}\left(Q^{2}\right)+\frac{8}{9} \tau_{d} G_{Q}^{2}\left(Q^{2}\right)\right],  \tag{71}\\
& \left.\operatorname{Res} f_{T}\left(\nu, Q^{2}\right)\right|_{\nu=Q^{2} /\left(2 M_{d}\right)}=-\frac{2}{3} e^{2} \tau_{d}\left(1+\tau_{d}\right) G_{M}^{2}\left(Q^{2}\right), \tag{72}
\end{align*}
$$

where $G_{C}\left(Q^{2}\right), G_{M}\left(Q^{2}\right)$, and $G_{Q}\left(Q^{2}\right)$ are the deuteron charge, magnetic, and quadrupole form factors, and $\tau_{d}=Q^{2} /\left(4 M_{d}^{2}\right)$. At the order we are working, $1+\tau_{d}=1$, and $G_{Q}\left(Q^{2}\right)=0$. The shift of the elastic poles, which appear at $\nu= \pm \boldsymbol{q}^{2} /(4 M)$, is also a relativistic correction that can be neglected at this order, allowing one to replace $\boldsymbol{q}^{2} \rightarrow Q^{2}$ in the amplitudes when taking the residues. This results in

$$
\begin{align*}
& \left.\operatorname{Res} f_{L}\left(\nu, Q^{2}\right)\right|_{\nu=Q^{2} /(4 M)}=-e^{2} G_{C}^{2}\left(Q^{2}\right),  \tag{73}\\
& \left.\operatorname{Res} f_{T}\left(\nu, Q^{2}\right)\right|_{\nu=Q^{2} /(4 M)}=-\frac{e^{2}}{24} \frac{Q^{2}}{M^{2}} G_{M}^{2}\left(Q^{2}\right) . \tag{74}
\end{align*}
$$

Evaluating the residues and expanding the square root of $G_{C}^{2}\left(Q^{2}\right)$ and $G_{M}^{2}\left(Q^{2}\right)$ order by order gives

$$
\begin{aligned}
G_{C}\left(Q^{2}\right) & =\frac{4 \gamma}{Q} \arctan \frac{Q}{4 \gamma} \\
& -(Z-1)\left(1-\frac{4 \gamma}{Q} \arctan \frac{Q}{4 \gamma}\right) \\
& -\frac{4}{3} r_{0}^{2} \gamma Q \arctan \frac{Q}{4 \gamma}
\end{aligned}
$$

$$
\begin{align*}
&+\frac{1}{3}(Z-1) r_{0}^{2} Q^{2}\left(1-\frac{4 \gamma}{Q} \arctan \frac{Q}{4 \gamma}\right)-\frac{(Z-1)^{3} l_{1}^{C 0_{S}}}{2 \gamma^{2}} Q^{2},  \tag{75}\\
& \frac{e}{2 M_{d}} G_{M}\left(Q^{2}\right)=\frac{e}{2 M}\left[\left(\mu_{n}+\mu_{p}\right) \frac{4 \gamma}{Q} \arctan \frac{Q}{4 \gamma}\right. \\
&\left.-\left(\mu_{n}+\mu_{p}\right)(Z-1)\left(1-\frac{4 \gamma}{Q} \arctan \frac{Q}{4 \gamma}\right)+\frac{2}{\pi} M L_{2}^{M 1_{S}} \gamma(\mu-\gamma)^{2}\right] \tag{76}
\end{align*}
$$

at N3LO for $G_{C}\left(Q^{2}\right)$ and at NLO for $G_{M}\left(Q^{2}\right)$. The expresion for $G_{M}\left(Q^{2}\right)$ coincides with that obtained in Ref. [7] (where the $\gamma \rho_{d}$ factor in the NLO term has to be replaced by $(Z-1)$ to account for the difference between the $Z$ and $\rho$ expansion schemes). The charge form factor has been previously calculated up to NNLO in the $Z$ expansion in Ref. [25], whose results we also reproduce here.

The single-nucleon contributions can in principle be summed into the isoscalar nucleon form factors $G_{E, M}^{(p+n)}$, resulting in the following compact expressions for the deuteron form factors:

$$
\begin{align*}
& G_{C}\left(Q^{2}\right)=Z G_{E}^{(p+n)}\left(Q^{2}\right) \frac{4 \gamma}{Q} \arctan \frac{Q}{4 \gamma}-\frac{(Z-1)^{3} l_{1}^{C 0_{S}}}{2 \gamma^{2}} Q^{2}  \tag{77}\\
& G_{M}\left(Q^{2}\right)=Z G_{M}^{(p+n)}\left(Q^{2}\right) \frac{4 \gamma}{Q} \arctan \frac{Q}{4 \gamma} \tag{78}
\end{align*}
$$

The order-by-order expression for the deuteron charge radius reads

$$
\begin{align*}
R_{C}^{2} \equiv\left\langle r^{2}\right\rangle_{C}=-\left.6 \frac{\mathrm{~d} G_{C}\left(Q^{2}\right)}{\mathrm{d} Q^{2}}\right|_{Q^{2}=0} & =\frac{1}{8 \gamma^{2}}+\frac{Z-1}{8 \gamma^{2}}+2 r_{0}^{2}+\frac{3(Z-1)^{3}}{\gamma^{2}} l_{1}^{C 0_{S}} \\
& =\left[2.3303+1.6063+0.6241+18.3166 l_{1}^{C 0_{S}}\right] \mathrm{fm}^{2} \tag{79}
\end{align*}
$$

One has to note that the NNLO result $R_{C}^{2}=4.5607 \mathrm{fm}^{2}$ is already very close to the experimentally measured values, for instance, $\mu \mathrm{D}$ spectroscopy gives $R_{C}^{2}=4.5183(33) \mathrm{fm}^{2}$ [44]. Using this value for fitting results in a tiny value of the $l_{1}^{C 0_{S}}$ coupling,

$$
\begin{equation*}
l_{1}^{C 0_{S}}=-2.32(18)(35) \times 10^{-3}, \tag{80}
\end{equation*}
$$

similar to what happens with $l_{2}^{M 1_{S}}$. The uncertainty in the first bracket is calculated using the quoted experimental error of the $\mu \mathrm{D}$ result. Considering other empirical values of $R_{C}^{2}$, e.g., the smaller value $R_{C}^{2}=4.3608 \mathrm{fm}^{2}$ of the empirical parametrisation of Abbott et al. [45], will significantly increase the uncertainty, albeit leaving $l_{1}^{C 0_{S}}$ at a level of at most $10^{-2}$. The value in the second bracket is due to the uncertainty of the value of $Z$, shown in Table I; one can see that the small relative uncertainty of $Z$ is amplified in the small subleading coupling $l_{1}^{C 0}{ }^{C}$.

Figure 15 shows the deuteron charge form factor at the different orders, compared with the result of the recent $\chi$ EFT fit of Ref. [46]. A conservative estimate for an error band at


FIG. 15. Deuteron charge form factor at LO (dash-dot-dotted black), NLO (dash-dotted blue), NNLO (dashed green), and N3LO (solid red, with the band showing the uncertainty due to higherorder terms). The result of the $\chi$ EFT fit [46] is shown by the purple dotted curve.

N3LO due to the higher-order terms could be obtained along the lines suggested in, e.g., Ref. [47]:

$$
\begin{equation*}
\delta G_{C}\left(Q^{2}\right)=\max \left\{\xi^{4} \Delta G_{C}^{\mathrm{LO}}\left(Q^{2}\right), \xi^{3} \Delta G_{C}^{\mathrm{NLO}}\left(Q^{2}\right), \xi^{2} \Delta G_{C}^{\mathrm{NNLO}}\left(Q^{2}\right), \xi \Delta G_{C}^{\mathrm{N} 3 \mathrm{LO}}\left(Q^{2}\right)\right\} \tag{81}
\end{equation*}
$$

where $\xi=\gamma / m_{\pi}$ and $\Delta G_{C}\left(Q^{2}\right)$ are the contributions to $G_{C}\left(Q^{2}\right)$ at the respective orders, with terms $\propto Q^{0}, Q^{2}$ removed to account for the fact that $G_{C}(0)=1$, and $G_{C}^{\prime}(0)=-1 / 6 R_{C}^{2}$ is also fixed at N3LO (the latter up to the small experimental error). The resulting band is rather narrow, its width being about the difference between the $\not \approx$ EFT and $\chi$ EFT results. An analogous band due to the uncertainty of $Z$ would in this scale be narrower than the widths of individual curves. One can see that the shapes of the charge form factor resulting from $\not \subset$ EFT and from $\chi$ EFT agree very well even at values of $Q$ beyond the formal range of validity of $\not$ EFFT $Q \lesssim m_{\pi}$, up to photon virtualities $Q \lesssim 200 \mathrm{MeV}$. This can be attributed to the fact that the slope and, to a lesser extent, the second derivative are by far the only important coefficients in the low-momentum expansion of $G_{C}\left(Q^{2}\right)$ in this range of $Q$, and they are well reproduced at N3LO in $\not \subset E F T$. The recent empirical parametrisations of the charge form factor, such as those derived in Refs. [45, 48] would lie in this figure within a line thickness from the $\chi$ EFT fit (or the N3LO $\not \subset$ EFT result), despite the numerical differences between the form factors (such as, e.g., $G_{C}\left(Q^{2}\right)$ of Abbott et al. having a smaller value of $\left.R_{C}^{2}\right)$. We therefore do not show them here. Note, however, that these superficially small effects can have a rather sizeable influence on the values of the elastic contribution to the $\mu \mathrm{D}$ Lamb shift. One can also note that, since the value of the N3LO contact term $l_{1}^{C 0_{S}}$ affects, through the value of $R_{C}^{2}$, the general shape of the deuteron charge form factor, one can deduce correlations between the value of $R_{C}^{2}$ and the elastic corrections. We shall consider these effects and compare the different variants of the deuteron charge form factor in detail in a dedicated publication.

## V. DEUTERON GENERALISED POLARISABILITIES

Having determined the value of $l_{1}^{C 0_{S}}$, we turn to the (generalised) polarisabilities of the deuteron. We start from the electric and magnetic dipole polarisabilities, the results for which are

$$
\begin{align*}
\alpha_{E 1}= & \frac{\alpha M}{32 \gamma^{4}}\left[1+(Z-1)+0+\frac{M \gamma^{3}}{6 \pi} C_{3 P_{J}}\right] \\
= & {[0.3771+0.2599+0-0.0018] \mathrm{fm}^{3}=0.6353 \mathrm{fm}^{3}, }  \tag{82}\\
\beta_{M 1}= & \frac{\alpha}{32 M \gamma^{2}}\left[-1+\frac{16}{3} \mu_{1}^{2}-\frac{32}{3} \mu_{1}^{2} \frac{\gamma}{\gamma_{s}-\gamma}\right. \\
& \left.\quad+\frac{Z-1}{3}\left(16 \mu_{1}^{2}-3\right)-\frac{32(Z-1)}{3} \mu_{1}\left(\mu_{1}+l_{1}^{M 1}{ }_{V}\right) \frac{\gamma}{\gamma_{s}-\gamma}+\frac{16}{3} \mu_{1}^{2} r_{s} \frac{\gamma^{3}}{\left(\gamma_{s}-\gamma\right)^{2}}\right] \\
= & {[0.0701+0.0003] \mathrm{fm}^{3}=0.0704 \mathrm{fm}^{3} . } \tag{83}
\end{align*}
$$

The expression for $\alpha_{E 1}$ reproduces the N3LO result obtained in Ref. [25] using the $n p \rightarrow \gamma d$ cross section calculated in Ref. [36], however, our $P$-wave contribution is a factor of 2 smaller (which we also cross-checked, reproducing the relevant terms in the $n p \rightarrow d \gamma$ cross section). The numerical value is in agreement with, e.g., the recent evaluation of Ref. [23] that obtained $\alpha_{E 1}=0.626(18) \mathrm{fm}^{3}$ at N3LO in $\chi$ EFT, as well as the calculation of Ref. [19] that used a selection of $\chi$ EFT potentials with various cut-offs, along with the AV18 model potential [49].

The LO expression for $\beta_{M 1}$ reproduces the result of Ref. [14], and the NLO result is new to the best of our knowledge. The numerical value is also in a very good agreement with the N3LO $\chi$ EFT result of Ref. [23], $\beta_{M 1}=0.0715(15) \mathrm{fm}^{3}$; this agreement is remarkable, given that this is a relatively low-order calculation. One can notice that the contribution of the $l_{1}^{M 1_{V}}$ contact term is individually rather sizeable $\simeq-0.06 \mathrm{fm}^{3}$, and that it cancels almost completely with the remaining NLO terms. This cancellation, giving essentially a zero NLO contribution to $\beta_{M 1}$, appears even more surprising. Both these results could be attributed to the procedure used in Ref. [36] to fit the value of $l_{1}^{M 1_{V}}$ to reproduce the $n p \rightarrow d \gamma$ cross section at NLO. Indeed, if this cross section is well described in đEFT, one should also expect a good description of the transverse response function of the deuteron at small non-zero $Q^{2}$. This, by virtue of the sum rule for $\beta_{M 1}$ derived in Ref. [32], should be sufficient in order to reproduce the value of $\beta_{M 1}$ at NLO, explaining at the same time the vanishing NLO contribution. This explanation, however, may imply that the description of the response function has to remain satisfactory up to relatively high energies outside of the validity of $九$ EFT; while this may be the case, investigating it in further detail is outside of the scope of this work.

The generalisation of $\alpha_{E 1}$ and $\beta_{M 1}$ to finite $Q^{2}$ is defined in the usual way,

$$
\begin{equation*}
\alpha_{E 1}\left(Q^{2}\right)=\frac{f_{L}\left(0, Q^{2}\right)}{4 \pi Q^{2}}, \quad \beta_{M 1}\left(Q^{2}\right)=\frac{\bar{f}_{T}\left(0, Q^{2}\right)}{4 \pi Q^{2}} \tag{84}
\end{equation*}
$$

where $f_{L}\left(0, Q^{2}\right)$ is understood as the non-pole part of $f_{L}$, and $\bar{f}_{T}$ stands for the non-pole part of $f_{T}$ with the Thomson term subtracted as well. The resulting curves are shown in Fig. 16. One can see that the patterns shown in the static values repeat in the generalised


FIG. 16. Generalised deuteron polarisabilities $\alpha_{E 1}\left(Q^{2}\right)$ (left panel) and $\beta_{M 1}\left(Q^{2}\right)$ (right panel). The LO, NLO, NNLO, and N3LO results for $\alpha_{E 1}\left(Q^{2}\right)$ in the left panel are coded as in Fig. 15. In the right panel, the LO and NLO results for $\beta_{M 1}\left(Q^{2}\right)$ are shown, respectively, by the black dashed and the red solid curve, with the band showing the estimate of higher-order contributions.
polarisabilities: the bulk of $\alpha_{E 1}\left(Q^{2}\right)$ comes from the LO and NLO contributions (with a small but visible NNLO contribution that is mostly due to the nucleon charge radii corrections and vanishes at $Q=0$ ), whereas the NLO contribution to $\beta_{M 1}\left(Q^{2}\right)$ is small at $Q=0$ as well as at finite virtualities.

The bands that estimate the contribution of higher orders are obtained analogously to Eq. (81), with the obvious modification for the NLO results. The N3LO band on the $\alpha_{E 1}$ curve is almost too narrow to be noticed, while the NLO band on $\beta_{M 1}$ is expectedly much wider. This, in fact, might be an overestimation, especially at low values of $Q$, in view of the agreement between the NLO $\not \approx E F T$ result for $\beta_{M 1}$ and the respective N3LO $\chi$ EFT result. One can expect that the NLO value of $\beta_{M 1}$ is therefore already close to its "true" value and will not change that much at higher orders. An investigation of this issue is, however, also outside of the scope of this work.

It is also interesting to consider two further generalised polarisabilities, namely, the longitudinal polarisability $\alpha_{L}\left(Q^{2}\right)$ and the generalised Baldin sum rule $\left[\alpha_{E 1}+\beta_{M 1}\right]\left(Q^{2}\right)$, defined via the non-pole parts of the amplitudes as

$$
\begin{equation*}
\alpha_{L}\left(Q^{2}\right)=\left.\frac{1}{4 \pi Q^{2}} \frac{\mathrm{~d} f_{L}\left(\nu, Q^{2}\right)}{\mathrm{d} \nu^{2}}\right|_{\nu=0}, \quad\left[\alpha_{E 1}+\beta_{M 1}\right]\left(Q^{2}\right)=\left.\frac{1}{4 \pi} \frac{\mathrm{~d} f_{T}\left(\nu, Q^{2}\right)}{\mathrm{d} \nu^{2}}\right|_{\nu=0} . \tag{85}
\end{equation*}
$$

The $Q^{2}=0$ value of $\alpha_{L}$ is given by

$$
\begin{align*}
\alpha_{L} & =\frac{7 \alpha M^{3}}{768 \gamma^{8}}\left[1+(Z-1)+0+\frac{11 M \gamma^{3}}{126 \pi} C_{3 P_{J}}\right] \\
& =[0.865+0.597+0-0.002] \times 10^{3} \mathrm{fm}^{5}=1.460 \times 10^{3} \mathrm{fm}^{5} \tag{86}
\end{align*}
$$

For the sake of simplicity, this result is obtained by substituting $|\boldsymbol{q}| \rightarrow Q$ in the expressions for $f_{L}$ obtained above and thus neglects relativistic corrections of a relative size of roughly $0.2 \%$. One can see that higher deuteron moments, such as $\alpha_{L}$, are numerically enhanced, unlike what happens in the case of the nucleon, see, e.g, Ref. [50].

The $Q^{2}=0$ value of the Baldin sum rule coincides with the NLO result for $\alpha_{E 1}+\beta_{M 1}$ given above. Strictly speaking, this includes corrections beyond NLO, because $\beta_{M 1}$ starts two orders higher than $\alpha_{E 1}$, however, the explicit gauge invariance allows one to recover $\beta_{M 1}$ by substituting $|\boldsymbol{q}|=\sqrt{Q^{2}+\nu^{2}}$ instead of neglecting $\nu$. Using $\boldsymbol{q}=Q$, on the other hand, changes the value of $\left[\alpha_{E 1}+\beta_{M 1}\right]\left(Q^{2}\right)$ by about $10 \%$ at $Q^{2}=0$ (by dropping the static value of $\beta_{M 1}$ ). The difference quickly decreases with growing $Q$ and becomes negligible already at $Q \simeq 20 \mathrm{MeV}$.


FIG. 17. Deuteron longitudinal polarisability $\alpha_{L}\left(Q^{2}\right)$ (left panel) and the generalised Baldin sum rule $\left[\alpha_{E 1}+\beta_{M 1}\right]\left(Q^{2}\right)$ (right panel). The curves are coded as in the left and right panel of Fig. 16, respectively. The curves for $\left[\alpha_{E 1}+\beta_{M 1}\right]\left(Q^{2}\right)$ are obtained using the exact expression $|\boldsymbol{q}|=\sqrt{\nu^{2}+Q^{2}}$ and reproduce at $Q=0$ the static values of $\alpha_{E 1}+\beta_{M 1}$ at the respective order, see the text for details.

The curves for $\alpha_{L}\left(Q^{2}\right)$ and $\left[\alpha_{E 1}+\beta_{M 1}\right]\left(Q^{2}\right)$ are shown in Fig. 17. One can see that the longitudinal polarisability shares the general features of the previously considered $\alpha_{E 1}\left(Q^{2}\right)$ and $\beta_{M 1}\left(Q^{2}\right)$, with a somewhat quicker falloff with growing $Q$. The generalised Baldin sum rule, on the other hand, demonstrates a sharp increase peaking around $Q=60 \mathrm{MeV}$; this enhancement is due to the magnetic interaction in the singlet channel and is analogous to what has been seen, e.g., in the generalised spin-forward deuteron polarisability $\gamma_{0}\left(Q^{2}\right)$ [51]. The estimate of higher-order corrections to $\alpha_{L}\left(Q^{2}\right)$ and $\left[\alpha_{E 1}+\beta_{M 1}\right]\left(Q^{2}\right)$, shown by the bands, is constructed analogously to, respectively, $\alpha_{E 1}$ and $\beta_{M 1}$. Similarly to those, the N3LO band on the longitudinal polarisability is very narrow, while the NLO band on the Baldin sum rule is significantly wider. In the latter case, one can, again, argue that the effects of higher orders might be overestimated towards $Q=0$ (even though the band is barely visible there in this scale), since we know that the dominant $\alpha_{E 1}$ contribution is well reproduced already at NLO (cf. also the discussion regarding $\beta_{M 1}$ above); however, this is
not any longer the case at larger finite virtualities, where more sizeable contributions could be expected, especially from higher-order magnetic couplings.


FIG. 18. Fourth-order generalised Baldin sum rule. The curves are coded as in the right panel of Fig. 16.

Finally, we consider the fourth-order generalised Baldin sum rule, defined in terms of the transverse amplitude as [27]

$$
\begin{equation*}
\left[\alpha_{E 1, \nu}+\beta_{M 1, \nu}+1 / 12\left(\alpha_{E 2}+\beta_{M 2}\right)\right]\left(Q^{2}\right)=\left.\frac{1}{8 \pi} \frac{\mathrm{~d}^{2} f_{T}\left(\nu, Q^{2}\right)}{\mathrm{d}\left(\nu^{2}\right)^{2}}\right|_{\nu=0} \tag{87}
\end{equation*}
$$

The static limit value of that sum rule gives the corresponding linear combination of the dispersive ( $\alpha_{E 1, \nu}, \beta_{M 1, \nu}$ ) and quadrupole ( $\alpha_{E 2}, \beta_{M 2}$ ) electric and magnetic polarisabilities of the deuteron, defined as in Ref. [52, 53]. ${ }^{1}$ Analogously to the Baldin sum rule considered above, using the exact formula $|\boldsymbol{q}|=\sqrt{\nu^{2}+Q^{2}}$ should allow one to recover the subleading polarisabilities here as well. Expanding the value at $Q=0$ up to N3LO, we get, order-byorder:

$$
\begin{align*}
\alpha_{E 1, \nu}+\beta_{M 1, \nu}+1 / 12\left(\alpha_{E 2}+\beta_{M 2}\right) & =[0.865+0.597+0.345+0.007] \times 10^{3} \mathrm{fm}^{5} \\
& =1.814 \times 10^{3} \mathrm{fm}^{5} \tag{88}
\end{align*}
$$

We refrain from showing here the analytic expression due to its length. The plot of the fourth-order generalised Baldin sum rule is shown in Fig. 18. One can see that it shares the features with the Baldin sum rule, sharply rising from the static limit and peaking at finite values of $Q$, and also getting a rather small NLO contribution at finite virtualities. It has to be noted that such a rapid growth of the subleading coefficients in the expansion of $f_{T}\left(\nu, Q^{2}\right)$ in powers of $\nu^{2}$, i.e., the Baldin sum rule and its fourth-order analogue, at finite $Q$ can be of some concern in the context of a calculation of the TPE correction in $\mu \mathrm{D}$, and it is important to verify that the transverse contribution is small (according to what is given by the counting, and is indeed seen in dispersive calculations [22, 23]).

[^0]
## VI. CONCLUSION

We have calculated the unpolarised deuteron VVCS amplitudes in the framework of $\not \subset E F T$, using the $Z$ expansion. Our results are at N3LO in the $\neq E F T$ expansion for the longitudinal amplitude, and at NLO for the transverse amplitude. We have provided analytic expressions for the VVCS amplitudes - the possibility to do so being one of the advantages of the $\neq$ EFT framework.

Investigating the RG running of the longitudinal amplitude, we show that there is a single unknown one-photon two-nucleon contact term contributing to the amplitude at N3LO, parametrized by the low-energy constant $l_{1}^{C 0_{S}}$. This constant is extracted from a fit to the deuteron charge form factor also at N3LO; the corresponding $\dagger$ EFT result is obtained from the residue of the longitudinal amplitude. We note that two of the RG equations for the N3LO two-nucleon one-photon contact terms obtained by us are at variance with those obtained in Ref. [35].

The deuteron charge and magnetic form factors, calculated, respectively, at N3LO and NLO from the residues of the longitudinal and transverse VVCS amplitudes, recover the previously obtained NNLO result for the charge form factor [25], as well as the NLO result for the magnetic form factor [7]. We furthermore obtain the generalised deuteron polarisabilities, namely, the electric $\alpha_{E 1}\left(Q^{2}\right)$ and magnetic $\beta_{M 1}\left(Q^{2}\right)$ dipole polarisabilities, the longitudinal polarisability $\alpha_{L}\left(Q^{2}\right)$, and the generalised Baldin sum rule $\left[\alpha_{E 1}+\beta_{M 1}\right]\left(Q^{2}\right)$. We also calculate the fourth-order generalised Baldin sum rule. The results for the electric and longitudinal generalised polarisability are at N3LO, while the remaining ones are calculated at NLO. There is a slight discrepancy between our expression for the static electric polarisability of the deutron $\alpha_{E 1}$ and the results obtained previously in Ref. [25]. The numerical values of the static electric and magnetic polarisabilities of the deuteron, $\alpha_{E 1}$ and $\beta_{M 1}$, are in a good agreement with other calculations, in particular, the recent $\chi$ EFT-based evaluation in Ref. [23]. This agreement between the values of $\alpha_{E 1}$ calculated in the two lowenergy effective field theories would be expected, since $\alpha_{E 1}$ is dominated by low-energy (or long-range) properties of the deuteron. On the other hand, the fact that the results for $\beta_{M 1}$ agree so well between the two theories can be reasonably explained by the properties of the procedure used to determine the NLO isovector magnetic contact term in Ref. [36], namely, that it reproduced the cross section in $n p \rightarrow d \gamma$ at low energies (although this remains to be explicitly demonstrated). The results of this work, in particular, the longitudinal amplitude, can be used to benchmark other calculational frameworks, e.g., $\chi$ EFT potentials, especially in settings where reproducing low-energy properties of the deuteron is important, such as the calculation of the TPE corrections in $\mu \mathrm{D}$.

The N3LO $\not \subset$ EFT result for the longitudinal amplitude and the charge form factor provides high-precision model-independent input for a calculation of the TPE correction to the Lamb shift in $\mu \mathrm{D}$. The uncertainty of such a calculation will be dominated by higher-order terms in the $\neq$ EFT expansion, expected to be of the order of $1 \%$ at N3LO. At the same time, the result for $f_{T}\left(\nu, Q^{2}\right)$ can be used to verify the smallness of the transverse contribution to the TPE
correction. This calculation, together with a detailed uncertainty analysis, will be presented in a subsequent publication. We are also investigating the spin-dependent deuteron VVCS amplitudes and the corresponding generalised polarisabilities of the deuteron.

A natural extension of this work would involve an application to unpolarised VVCS in $A=3$ nuclei, where a good description of quantities such as the charge radii has already been obtained in $\not \subset E F T$, see, e.g., Refs. $[54,55]$ for recent results, and Ref. [56] for a review. Given the complications of $\not \approx E F T$ as applied to the three-nucleon sector, pursuing this program would not be straightforward, however, it promises interesting results.

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[^0]:    ${ }^{1}$ The quadrupole polarisabilities (spin independent) are not to be confused with the tensor polarisabilities of the deuteron defined in, e.g., Refs. [14, 34].

