

Antinucleon-nucleon interaction in chiral effective field theory

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ABSTRACT: Results of an exploratory study of the antinucleon-nucleon interaction within chiral effective field theory are reported. The antinucleon-nucleon potential is derived up to next-to-next-to-leading order, based on a modified Weinberg power counting, in close analogy to pertinent studies of the nucleon-nucleon interaction. The low-energy constants associated with the arising contact interactions are fixed by a fit to phase shifts and inelasticities provided by a recently published phase-shift analysis of antiproton-proton scattering data. The overall quality of the achieved description of the antinucleon-nucleon amplitudes is comparable to the one found in case of the nucleon-nucleon interaction at the same order. For most S -waves and several P -waves good agreement with the antinucleon-nucleon phase shifts and inelasticities is obtained up to laboratory energies of around 200 MeV.

KEYWORDS: Chiral Lagrangians, Scattering Amplitudes, QCD

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1 Introduction

The antinucleon-nucleon ($\bar{N}N$) interaction has been studied quite extensively in the past [1–9], not least because of the wealth of data collected at the LEAR facility at CERN, cf. the reviews [10–12]. The majority of those investigations has been performed in the traditional meson-exchange framework where the G -parity transformation is exploited to connect the elastic part of the $\bar{N}N$ interaction with the dynamics in the nucleon-nucleon (NN) system. Annihilation processes are described either by a simple optical potential (which is often assumed to be spin- as well as energy-independent) [1, 2, 5, 7] or in terms of a coupling to a small number of effective two-body annihilation channels [3, 4, 9].

In the last two decades chiral effective field theory (EFT) has become a standard tool in the studies of the NN interaction at low energies. This development was initiated by two seminal papers by Weinberg [13, 14] in which he proposed that EFT and the power-counting rules associated with it should be applied to the NN potential rather than to the reaction amplitude. The reaction amplitude is then obtained from solving a regularized Lippmann-Schwinger equation for the derived interaction potential. His suggestion is based on the observation that diagrams with purely nucleonic intermediate states are strongly enhanced and, therefore, not amenable to a perturbative treatment. However, they can be taken into account and they are actually summed up to infinite order when solving the Lippmann-Schwinger equation. The chiral NN potential contains pion exchanges and a series of contact interactions with an increasing number of derivatives. The latter represent the short-range part of the NN force and are parametrized by low-energy constants (LECs), that need to be fixed by a fit to data. For reviews we refer the reader to the recent refs. [15, 16]. Presently the most refined calculations extend up to next-to-next-to-next-to-leading order ($N^3\text{LO}$) [17, 18] and they yield a rather accurate description of the NN phase shifts up to laboratory energies of 250-300 MeV.

Naturally, the success of chiral EFT in the NN sector provides a strong motivation to apply the same approach also to the $\bar{N}N$ interaction. First and most important for the practical implementation, recently an update of the Nijmegen partial-wave analysis (PWA) of antiproton-proton ($\bar{p}p$) scattering data [19] has been published. For the new PWA [20] the resulting phase shifts and inelasticities are explicitly given and can be readily used for applying the chiral EFT approach to the $\bar{N}N$ interaction in the very same way as it has been done for the NN system.

A further incentive for exploring the feasibility of investigating the $\bar{N}N$ system within chiral EFT comes from the expected increase in interest in the $\bar{N}N$ interaction in the future due to the Facility for Antiproton and Ion Research (FAIR) in Darmstadt whose construction is finally on its way. Among the various project planned at this site is the PANDA experiment [21] which aims to study the interactions between antiprotons and fixed target protons and nuclei in the momentum range of 1.5-15 GeV/c using the high energy storage ring HESR.

Finally, chiral EFT could be a very powerful tool to analyze data from recent measurements of the $\bar{p}p$ invariant mass in the decays of J/ψ , B mesons, etc., and of the reaction $e^+e^- \rightarrow \bar{p}p$. In several of those reactions a near-threshold enhancement in the mass spectrum was found [22–25] and this enhancement could allow one to extract information on the $\bar{p}p$ interaction at very low energies [26–34].

In the present paper we report on results of an exploratory study of the antinucleon-nucleon interaction within chiral EFT. In our application of chiral EFT to the $\bar{N}N$ interaction we follow exactly the approach used by Epelbaum et al. [18, 35, 36] in the NN case. It is consistent with the scheme originally proposed by Weinberg except that one aims for an energy-independent representation of the chiral potential [37]. For the time being we restrict ourselves to an evaluation of the potential up to next-to-next-to-leading order (NNLO). At leading order (LO) the potential is given by one-pion exchange (OPE) and two contact terms without derivatives. At next-to-leading order (NLO) contributions from the leading two-pion exchange (TPE) diagrams as well as seven more contact operators arise. Finally, at NNLO one gets contributions from the subleading TPE with one insertion of dimension two pion-nucleon vertices. Once the potential is established it has to be inserted into a regularized scattering equation in order to obtain the reaction amplitude. For the regularization we follow again closely the procedure adopted by Epelbaum et al. [18, 36] and others [17], in their study of the NN interaction and introduce a momentum-dependent exponential regulator function.

For investigations of the $\bar{N}N$ interaction within EFT based on other schemes see refs. [38, 39], where the Kaplan-Savage-Wise resummation scheme [40] is employed. These authors considered the $\bar{N}N$ interaction up to NLO. There have been also attempts to compute specific $\bar{p}p$ annihilation channels in chiral EFT [41].

The present paper is structured as follows: The effective $\bar{N}N$ potential up to NNLO is described in section 2. We start with a brief review of the underlying power counting and then provide explicit expressions for the contributions from pion exchange and for the contact terms. We also discuss how we treat the annihilation processes. Finally, we introduce the Lippmann-Schwinger equation that we solve and the parameterization of the

S-matrix that we use. In section 3 we indicate our fitting procedure and then we present the results achieved at NLO and at NNLO. Phase shifts and inelasticities for S -, P -, and D - waves, obtained from our EFT interaction, are displayed and compared with those of the $\bar{N}N$ phase-shift analysis. Furthermore, predictions for S -wave scattering lengths are given. A summary of our work and an outlook on future investigations is given in section 4.

2 Chiral potential at next-to-next-to-leading order

The contributions to the NN interaction up to NNLO are described in detail in refs. [18, 35, 36]. The underlying power counting is given by (considering only connected diagrams)

$$\nu = 2L + \sum_i \Delta_i, \quad \Delta_i = d_i + \frac{n_i}{2} - 2 \quad (2.1)$$

where L is the number of loops in the diagram, d_i is the number of derivatives or pion mass insertions, and n_i the number of internal nucleon fields at the vertex i under consideration. The LO potential corresponds to $\nu = 0$ and consists of two four-nucleon contact terms without derivatives and of one-pion exchange. There are no contributions at order $\nu = 1$ due to requirements from parity conservation and time-reversal invariance. At NLO ($\nu = 2$) seven new contact terms (with two derivatives) arise, together with loop contributions from (irreducible) two-pion exchange. Finally, at NNLO ($\nu = 3$) there are additional contributions from two-pion exchange resulting from one insertion of dimension two pion-nucleon vertices, see e.g. ref. [42]. The corresponding diagrams are summarized in figure 1.

The structure of the $\bar{N}N$ interaction is practically identical and, therefore, the potential given in refs. [18, 36] can be adapted straightforwardly for the $\bar{N}N$ case. For the ease of the reader and also for defining our potential uniquely we provide the explicit expressions below.

2.1 Pion exchange

In line with [18] we adopt the following expression for the one-pion exchange potential

$$V_{1\pi}(q) = \left(\frac{g_A}{2F_\pi} \right)^2 \left(1 - \frac{p^2 + p'^2}{2m^2} \right) \boldsymbol{\tau}_1 \cdot \boldsymbol{\tau}_2 \frac{\boldsymbol{\sigma}_1 \cdot \mathbf{q} \boldsymbol{\sigma}_2 \cdot \mathbf{q}}{\mathbf{q}^2 + M_\pi^2}, \quad (2.2)$$

where $\mathbf{q} = \mathbf{p}' - \mathbf{p}$ is the transferred momentum defined in terms of the final ($\mathbf{p}'$) and initial ($\mathbf{p}$) center-of-mass momenta of the baryons (nucleon or antinucleon). Obviously here relativistic $1/m^2$ corrections to the static one-pion exchange potential have been taken into account. As in the work [18] we take the larger value $g_A = 1.29$ instead of $g_A = 1.26$ in order to account for the Goldberger-Treiman discrepancy. This value, together with the used $F_\pi = 92.4 \text{ MeV}$, implies the pion-nucleon coupling constant $g_{NN\pi} = 13.1$ which is consistent with the empirical value obtained from πN and NN data [43, 44] and also with modern determinations utilizing the GMO sum rule [45]. For the nucleon (antinucleon) and pion mass we use the isospin-averaged values $m = 938.918 \text{ MeV}$ and $M_\pi = 138.039 \text{ MeV}$, respectively. Note that the contribution of one-pion exchange to the $\bar{N}N$ interaction is of opposite sign as that in the NN case. This sign difference arises from transforming the

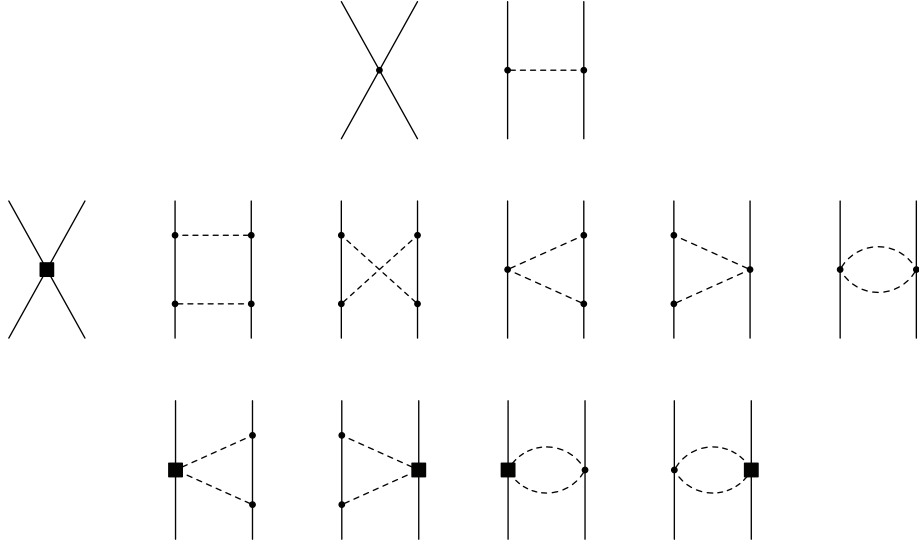


Figure 1. Relevant diagrams up-to-and-including NNLO. Solid and dashed lines denote the antinucleon/nucleon and the pion, respectively. The square symbolizes a contact vertex with two derivatives or a subleading πN vertex. The contributions at LO, NLO, and NNLO are displayed from top to bottom.

$NN\pi$ vertex to the $\bar{N}\bar{N}\pi$ vertex via charge conjugation and a rotation in the isospin space and is commonly referred to as G -parity transformation.

The two-pion exchange potential calculated using spectral function regularization [18] is given at NLO by

$$V_{2\pi}^{(2)}(q) = \boldsymbol{\tau}_1 \cdot \boldsymbol{\tau}_2 V_C^{(2)}(q) + \boldsymbol{\sigma}_1 \cdot \mathbf{q} \boldsymbol{\sigma}_2 \cdot \mathbf{q} V_T^{(2)}(q) + \boldsymbol{\sigma}_1 \cdot \boldsymbol{\sigma}_2 V_S^{(2)}(q), \quad (2.3)$$

where

$$V_C^{(2)}(q) = -\frac{1}{384\pi^2 F_\pi^4} L^{\tilde{\Lambda}}(q) \left\{ 4M_\pi^2(5g_A^4 - 4g_A^2 - 1) + q^2(23g_A^4 - 10g_A^2 - 1) + \frac{48g_A^4 M_\pi^4}{4M_\pi^2 + q^2} \right\},$$

$$V_T^{(2)}(q) = -\frac{1}{q^2} V_S^{(2)}(q) = -\frac{3g_A^4}{64\pi^2 F_\pi^4} L^{\tilde{\Lambda}}(q),$$

and at NNLO by

$$V_{2\pi}^{(3)}(q) = V_C^{(3)}(q) + \boldsymbol{\tau}_1 \cdot \boldsymbol{\tau}_2 \boldsymbol{\sigma}_1 \cdot \mathbf{q} \boldsymbol{\sigma}_2 \cdot \mathbf{q} V_T^{(3)}(q) + \boldsymbol{\tau}_1 \cdot \boldsymbol{\tau}_2 \boldsymbol{\sigma}_1 \cdot \boldsymbol{\sigma}_2 V_S^{(3)}(q), \quad (2.4)$$

with

$$V_C^{(3)}(q) = -\frac{3g_A^2}{16\pi F_\pi^4} \left\{ 2M_\pi^2(2c_1 - c_3) - c_3 q^2 \right\} (2M_\pi^2 + q^2) A^{\tilde{\Lambda}}(q),$$

$$V_T^{(3)}(q) = -\frac{1}{q^2} V_S^{(3)}(q) = -\frac{g_A^2}{32\pi F_\pi^4} c_4 (4M_\pi^2 + q^2) A^{\tilde{\Lambda}}(q).$$

The NLO and NNLO loop functions $L^{\tilde{\Lambda}}(q)$ and $A^{\tilde{\Lambda}}(q)$ are given by

$$L^{\tilde{\Lambda}}(q) = \theta(\tilde{\Lambda} - 2M_\pi) \frac{\omega}{2q} \ln \frac{\tilde{\Lambda}^2 \omega^2 + q^2 s^2 + 2\tilde{\Lambda} q \omega s}{4M_\pi^2(\tilde{\Lambda}^2 + q^2)},$$

$$\omega = \sqrt{q^2 + 4M_\pi^2}, \quad s = \sqrt{\tilde{\Lambda}^2 - 4M_\pi^2}, \quad (2.5)$$

and

$$A^{\tilde{\Lambda}}(q) = \theta(\tilde{\Lambda} - 2M_\pi) \frac{1}{2q} \arctan \frac{q(\tilde{\Lambda} - 2M_\pi)}{q^2 + 2\tilde{\Lambda}M_\pi}. \quad (2.6)$$

For the LECs c_1 and c_4 we adopt the central values from the Q^3 -analysis of the πN system [46]: $c_1 = -0.81 \text{ GeV}^{-1}$, $c_4 = 3.40 \text{ GeV}^{-1}$. For the constant c_3 the value $c_3 = -3.40 \text{ GeV}^{-1}$ is used, which is on the lower side but still consistent with the results from ref. [46]. Note that slightly different values are employed in the $\bar{N}N$ partial-wave analysis [20], namely $c_1 = -0.76 \text{ GeV}^{-1}$, $c_3 = -5.8 \text{ GeV}^{-1}$ and $c_4 = 4.0 \text{ GeV}^{-1}$. These values are also consistent with the recent determination in [47].

2.2 Contact terms

The spin-dependence of the potentials due to the leading order contact terms is given by [48]

$$V_{\bar{N}N}^{(0)} = C_S + C_T \boldsymbol{\sigma}_1 \cdot \boldsymbol{\sigma}_2, \quad (2.7)$$

where the parameters C_S and C_T are low-energy constants (LECs) which need to be determined in a fit to data. At NLO, the spin- and momentum-dependence of the contact terms reads

$$V_{\bar{N}N}^{(2)} = C_1 \mathbf{q}^2 + C_2 \mathbf{k}^2 + (C_3 \mathbf{q}^2 + C_4 \mathbf{k}^2) \boldsymbol{\sigma}_1 \cdot \boldsymbol{\sigma}_2 + \frac{i}{2} C_5 (\boldsymbol{\sigma}_1 + \boldsymbol{\sigma}_2) \cdot (\mathbf{q} \times \mathbf{k})$$

$$+ C_6 (\mathbf{q} \cdot \boldsymbol{\sigma}_1)(\mathbf{q} \cdot \boldsymbol{\sigma}_2) + C_7 (\mathbf{k} \cdot \boldsymbol{\sigma}_1)(\mathbf{k} \cdot \boldsymbol{\sigma}_2), \quad (2.8)$$

where C_i ($i = 1, \dots, 7$) are additional LECs. The average momentum $\mathbf{k}$ is defined by $\mathbf{k} = (\mathbf{p}' + \mathbf{p})/2$. When performing a partial-wave projection, these terms contribute to the two S -wave (1S_0 , 3S_1) potentials, the four P -wave (1P_1 , 3P_0 , 3P_1 , 3P_2) potentials, and the 3S_1 - 3D_1 transition potential in the following way [18]:

$$V(^1S_0) = 4\pi (C_S - 3C_T) + \pi (4C_1 + C_2 - 12C_3 - 3C_4 - 4C_6 - C_7)(p^2 + p'^2)$$

$$= \tilde{C}_{1S_0} + C_{1S_0}(p^2 + p'^2), \quad (2.9)$$

$$V(^3S_1) = 4\pi (C_S + C_T) + \frac{\pi}{3} (12C_1 + 3C_2 + 12C_3 + 3C_4 + 4C_6 + C_7)(p^2 + p'^2)$$

$$= \tilde{C}_{3S_1} + C_{3S_1}(p^2 + p'^2), \quad (2.10)$$

$$V(^1P_1) = \frac{2\pi}{3} (-4C_1 + C_2 + 12C_3 - 3C_4 + 4C_6 - C_7) p p' = C_{1P_1} p p', \quad (2.11)$$

$$V(^3P_1) = \frac{2\pi}{3} (-4C_1 + C_2 - 4C_3 + C_4 + 2C_5 - 8C_6 + 2C_7) p p' = C_{3P_1} p p', \quad (2.12)$$

$$V(^3P_0) = \frac{2\pi}{3} (-4C_1 + C_2 - 4C_3 + C_4 + 4C_5 + 12C_6 - 3C_7) p p' = C_{^3P_0} p p', \quad (2.13)$$

$$V(^3P_2) = \frac{2\pi}{3} (-4C_1 + C_2 - 4C_3 + C_4 - 2C_5) p p' = C_{^3P_2} p p', \quad (2.14)$$

$$V(^3D_1 - ^3S_1) = \frac{2\sqrt{2}\pi}{3} (4C_6 + C_7) p'^2 = C_{^3S_1-^3D_1} p'^2 \equiv C_{\epsilon_1} p'^2, \quad (2.15)$$

$$V(^3S_1 - ^3D_1) = \frac{2\sqrt{2}\pi}{3} (4C_6 + C_7) p^2 = C_{^3S_1-^3D_1} p^2 \equiv C_{\epsilon_1} p^2, \quad (2.16)$$

with $p = |\mathbf{p}|$ and $p' = |\mathbf{p}'|$. There are no additional contact terms at NNLO.

Note that the Pauli principle is absent in case of the $\bar{N}N$ interaction. Accordingly, each partial wave that is allowed by angular momentum conservation occurs in the isospin $I = 0$ and in the $I = 1$ channel. Therefore, there are now twice as many contact terms as in NN .

The main new feature in the $\bar{N}N$ interaction is the presence of annihilation processes. The $\bar{N}N$ system annihilates into a multitude of $n\pi$ channels, where the decay to 4 to 6 pions is dominant in the low-energy region of $\bar{N}N$ scattering [10]. The threshold energy of those channels is in the order of 700 MeV while the $\bar{N}N$ threshold is at 1878 MeV. Therefore, one does not expect that annihilation introduces a new scale into the problem. Accordingly, there should be no need to modify the power counting when going from NN to $\bar{N}N$ because the momenta associated with the annihilation channels should be, in average, much larger than those in the $\bar{N}N$ system itself. This conjecture is supported by the fact that phenomenological models of the $\bar{N}N$ interaction can describe the bulk properties of annihilation very well by simple energy-independent optical potentials of Woods-Saxon or Gaussian type [1, 2, 5, 7]. The ranges associated with those interactions are of the order of 1 fm or less. The above considerations suggest that annihilation processes are primarily tied to short-distance physics and, therefore, can be and should be simply incorporated into the contact terms which anyway are meant to parameterize effectively the short-range part of (elastic) NN and/or $\bar{N}N$ scattering.

Nonetheless we want to emphasize that the above arguments are of pragmatical nature and not fundamental ones. There are definitely annihilation channels that open near the $\bar{N}N$ threshold. Specifically, there are indications that a sizeable part of the annihilation into multipion channels proceeds via two-meson doorway modes like $\bar{N}N \rightarrow \rho\rho \rightarrow 4\pi$ or $\bar{N}N \rightarrow f_2(1270)\omega \rightarrow 5\pi$, and some of those have nominal thresholds close to that of $\bar{N}N$ scattering. On the other hand, according to empirical information the actual branching ratios into individual two-body channels are typically of the order of 1% [7] only and, therefore, they do not have any noticeable impact on the description of the bulk properties of $\bar{N}N$ annihilation. In fact, all the two-body annihilation channels together — as far as they have been measured — yield only about 30% of the total annihilation cross section at the $\bar{N}N$ threshold which is a strong evidence for the dominance of annihilation into 3 or more (uncorrelated) pions.

The study of $\bar{N}N$ scattering in EFT in refs. [38, 39] followed the above arguments and took into account annihilation by simply using complex LECs in eqs. (2.9)–(2.16). However, this prescription has an unpleasant drawback — it does not allow one to impose sensible

unitarity requirements on the resulting scattering amplitude. With unitarity requirements we mean a condition that guarantees that for each partial wave its contribution to the total cross section is larger than its contribution to the integrated elastic cross section. In case of strict two-body unitarity like for NN scattering below the pion production threshold these two quantities are, of course, identical.

Since we want an approach that manifestly fulfils unitarity constraints we treat annihilation in a different way. We start out from the observation that unitarity requires the $\bar{N}N$ annihilation potential to be of the form

$$V_{ann} = \sum_{X=2\pi, 3\pi, \dots} V_{\bar{N}N \rightarrow X} G_X V_{X \rightarrow \bar{N}N} \quad (2.17)$$

where X is the sum over all open annihilation channels, and G_X is the propagator of the intermediate state X . Note that eq. (2.17) is exact under the assumption that there is no interaction in and no transition between the various annihilation channels. Performing an expansion of $V_{\bar{N}N \rightarrow X}$ up to NNLO analogous to the $\bar{N}N$ interaction and evaluating formally the sum and integral in eq. (2.17) yields a contribution from the unitarity cut that can be written as

$$V_{ann}^{L=0} = -i (\tilde{C}_{1S_0}^a + C_{1S_0}^a p^2) (\tilde{C}_{1S_0}^a + C_{1S_0}^a p'^2), \quad V_{ann}^{L=1} = -i (C_\alpha^a)^2 p p', \quad (2.18)$$

where α stands for the 3P_0 , 1P_1 , 3P_1 , and 3P_2 partial waves. For the coupled $^3S_1 - ^3D_1$ partial wave we get

$$\begin{aligned} V_{ann}^{S \rightarrow S} &= -i (\tilde{C}_{3S_1}^a + C_{3S_1}^a p^2) (\tilde{C}_{3S_1}^a + C_{3S_1}^a p'^2), & V_{ann}^{S \rightarrow D} &= -i (\tilde{C}_{3S_1}^a + C_{3S_1}^a p^2) C_{\epsilon_1}^a p'^2, \\ V_{ann}^{D \rightarrow S} &= -i C_{\epsilon_1}^a p^2 (\tilde{C}_{3S_1}^a + C_{3S_1}^a p'^2), & V_{ann}^{D \rightarrow D} &= -i (C_{\epsilon_1}^a)^2 p^2 p'^2. \end{aligned} \quad (2.19)$$

In those expressions the parameters $\tilde{C}^a$ and C^a are real. Thus, for each partial wave we essentially recover the structure of the potential that follows from the contact terms considered above, with the same number of free parameters. However, in eqs. (2.18)–(2.19) the sign of V_{ann} as required by unitarity is already explicitly fixed and does not depend on the sign of the parameters $\tilde{C}^a$ and C^a anymore. Moreover, and most importantly, we see that a term proportional to $p^2 p'^2$ arises in the S waves at NLO and NNLO from unitarity constraints and it has to be included in order to make sure that unitarity is fulfilled at any energy.

Note that, in principle, there is also a contribution from the principal-value part of the integral in eq. (2.17). However, it is real and, therefore, its structure is already accounted for by the standard LECs in eqs. (2.9)–(2.16).

Finally we would like to add that in practice the treatment of annihilation via eqs. (2.18)–(2.19) corresponds to the introduction of an effective two-body annihilation channel with a threshold significantly below the one of $\bar{N}N$ so that the center-of-mass momentum in the annihilation channel is already fairly large and its variation in the low-energy region of $\bar{N}N$ scattering considered by us is negligible.

2.3 Scattering equation

In the actual calculation a partial-wave projection of the interaction potentials is performed, as described in detail in ref. [18]. The reaction amplitudes are obtained from the solution of a relativistic Lippmann-Schwinger (LS) equation:

$$T_{L''L'}(p'', p'; E_q) = V_{L''L'}(p'', p') + \sum_L \int_0^\infty \frac{dp p^2}{(2\pi)^3} V_{L''L}(p'', p) \frac{1}{2E_q - 2E_p + i0^+} T_{LL'}(p, p'; E_q). \quad (2.20)$$

Here, $E_q = \sqrt{m^2 + q^2}$, where q is the on-shell momentum. We adopt here a relativistic scattering equation so that our amplitudes fulfil the relativistic unitarity condition at any order, as done also in the NN sector [16, 18]. On the other hand, relativistic corrections to the potential are calculated order by order, but appear first at next-to-next-to-next-to-leading order ($N^3\text{LO}$) in the Weinberg scheme, see ref. [18].

Like in the NN case we have either uncoupled spin-singlet and triplet waves (where $L'' = L' = L = J$) or coupled partial waves (where $L'', L', L = J - 1, J + 1$). We solve the LS equation in the isospin basis, i.e. for $I = 0$ and $I = 1$ separately, and we compare the resulting phase shifts with those in ref. [20] that are likewise given in the isospin basis. It should be said, however, that for a comparison directly with data a more refined treatment is required. Then one should solve the LS equation in particle basis and consider the coupling between the $\bar{p}p$ and $\bar{n}n$ channels explicitly. In this case one can take into account the mass difference between p ($\bar{p}$) and n ($\bar{n}$) and, thereby, implement the fact that the physical thresholds of the $\bar{p}p$ and $\bar{n}n$ channels are separated by about 2.5 MeV, and also one can add the Coulomb interaction in the $\bar{p}p$ channel. The potential in the LS equation is cut off with a regulator function,

$$f^\Lambda(p', p) = \exp \left[- (p'^6 + p^6) / \Lambda^6 \right], \quad (2.21)$$

in order to remove high-energy components [18]. The cutoff values are chosen in the range $\Lambda = 450 - 600$ MeV at NLO and $\Lambda = 450 - 650$ MeV at NNLO, similar to what was used for chiral NN potentials [18, 36].

The relation between the S - and on-the-energy shell T -matrix is given by

$$S_{LL'}(q) = \delta_{LL'} - \frac{i}{8\pi^2} q E_q T_{LL'}(q). \quad (2.22)$$

The phase shifts in the uncoupled cases can be obtained from the S -matrix via

$$S_{LL} \equiv S_L = e^{2i\delta_L}. \quad (2.23)$$

For the S -matrix in the coupled channels ($J > 0$) we use the so-called Stapp parametrization [49]

$$\begin{pmatrix} S_{J-1 J-1} & S_{J-1 J+1} \\ S_{J+1 J-1} & S_{J+1 J+1} \end{pmatrix} = \begin{pmatrix} \cos 2\epsilon_J e^{2i\delta_{J-1}} & -i \sin 2\epsilon_J e^{i(\delta_{J-1} + \delta_{J+1})} \\ -i \sin 2\epsilon_J e^{i(\delta_{J-1} + \delta_{J+1})} & \cos 2\epsilon_J e^{2i\delta_{J+1}} \end{pmatrix}. \quad (2.24)$$

In case of elastic scattering the phase parameters in eqs. (2.23) and (2.24) are real quantities while in the presence of inelasticities they become complex. Because of that, in the past several generalizations of these formulae have been proposed that still allow one to write the S -matrix in terms of real parameters [20, 50]. We follow here ref. [51] and calculate and present simply the real and imaginary parts of the phase shifts and the mixing parameters obtained via the above parameterization. Note that with this choice the real part of the phase shifts is identical to the phase shifts one obtains from another popular parameterization where the imaginary part is written in terms of an inelasticity parameter η , e.g. for uncoupled partial waves

$$S_L = \eta e^{2i\delta_L}. \quad (2.25)$$

Indeed, for this case $\text{Im } \delta_L = -(\log \eta)/2$ which implies that $\text{Im } \delta_L \geq 0$ since $\eta \leq 1$ because of unitarity. Since our calculation implements unitarity, the optical theorem

$$\text{Im } a_{LL}(q) \geq q \sum_{L'} |a_{LL'}(q)|^2, \quad (2.26)$$

is fulfilled for each partial wave, where $a_{LL'}(q) = (S_{LL'} - \delta_{LL'})/(2iq) = -1/(4\pi)^2 \cdot E_q T_{LL'}(q)$.

For the fitting procedure and for the comparison of our results with those by Zhou and Timmermans we reconstructed the S -matrix based on the phase shifts listed in tables VIII-X in ref. [20] and on the formulae presented in section VII of that paper and then converted them to our convention specified in eqs. (2.23) and (2.24).

3 Results

In the fitting procedure we follow very closely the strategy of Epelbaum et al. in their study of the NN interaction [18, 36]. In particular, we consider the same ranges for the cutoffs, namely for the cutoff in the LS equation values of $\Lambda = 450\text{--}600$ MeV at NLO and $\Lambda = 450\text{--}650$ MeV at NNLO while for the spectral function regularization variations we consider values in the range $\tilde{\Lambda} = 500\text{--}700$ MeV. For any combination of the cutoffs Λ and $\tilde{\Lambda}$, the LECs $C_{S,T}$ and $C_{1\dots 7}$ are fixed from a fit to the $\bar{N}N$ S - and P -waves and the mixing parameter ϵ_1 of ref. [20] for laboratory energies below 125 MeV ($p_{lab} \leq 500$ MeV/c). The numerical values of the LECs are compiled in tables 1 (NLO) and 2 (NNLO) for a selected combination of the cutoffs. The values for $\tilde{C}_{1S_0}$ in the isospin $I = 1$ case found in the fitting procedure turned out to be very small and, therefore, we set them to zero.

Our results are displayed and compared with the $\bar{N}N$ PWA [20] in figures 2–6. The bands represent the variation of the obtained phase shifts and mixing parameters with the cutoff. Those variations can be viewed as an estimate for the theoretical uncertainty. Thus, in principle for the same variation of the cutoff those bands should become narrower and narrower when one goes to higher order. However, as argued in ref. [36], in practice one has to be careful in the interpretation of the bands, specifically for the transition from NLO to NNLO. Since the same number of contact terms are present in the interactions at NLO and NNLO one rather should expect variations of similar magnitude. In particular, for

reasons discussed in [36] the cutoff variation underestimates the uncertainty for the NLO results. In any case one has to keep in mind that, following ref. [36], we use a larger cutoff region at NNLO than for the NLO case.

Let us now discuss the individual partial waves. Results for the 1S_0 channel can be found in the upper part of figure 2. Obviously, the phase shift for isospin $I = 0$ (we use here the spectral notation $^{(2I+1)(2S+1)}L_J$) is very well described up to fairly high energies — even at NLO — and likewise the inelasticity, presented in terms of the imaginary part of the phase shift. Moreover, the dependence on the cutoff is very small. In the $I = 1$ channel the situation is rather different. Here we observe a sizeable cutoff dependence of the results for energy above 150 MeV. This has to do with the fact that the PWA suggests a resonance-like behavior of the phase in this region. Since this resonance lies in an energy region where we expect our results to show increasing uncertainties, based on the experience from the NN case [36], it is not surprising that it is difficult to reproduce this structure quantitatively. Nevertheless, there is a visible improvement when going from NLO to NNLO and at the latter order the empirical phase shifts already lie within the error bands of theory.

We want to emphasize that this improvement is entirely due to inclusion of the sub-leading two-pion exchange potential, since as already stressed above no new contact terms arise at NNLO and thus the number of adjustable parameters is the same at NLO and NNLO. Also, it should be said that the NLO result, shown here up to $T_{lab} = 200$ MeV, exhibits a similar trend like the one for NNLO at higher energies, i.e. the phases reach a maximum and then become more negative again.

The situation for the 3P_0 partial wave is similar, see figure 2 (lower part). Also here the $I = 0$ phase shifts are well reproduced while in the $I = 1$ case there is an even larger cutoff dependence than in the 3S_0 . Obviously also the $^{33}P_0$ amplitude of the PWA [20] exhibits a resonance-like behavior. Its reproduction requires a potential that is repulsive at large separations of the antinucleon and nucleon but becomes attractive for short distances. Since there is only a single LEC up to NNLO for P waves, the magnitude and range of such an attraction cannot be adequately accounted for. For improvements one has to wait for a N^3 LO calculation.

Results for the 1P_1 and 3P_1 partial waves are shown in figure 3. In general, the description improves when going from NLO to NNLO. Specifically for the two 1P_1 channels and the $^{33}P_1$ the results at NNLO agree with those of the PWA within the uncertainty bands for energies up to 150 MeV and often even up to 250 MeV. An exception is the $^{13}P_1$ partial wave where the phase shift can only be described up to 50 MeV or so. Similar to the $^{33}P_0$, the PWA yields a negative phase at low energies which tends towards positive values at larger energies [20] and one encounters the same difficulty as discussed above.

In figure 4 one can find our results for the coupled 3S_1 – 3D_1 partial wave. Here the S -wave phase shifts (and also the inelasticity) are satisfactorily described over the whole energy range considered with uncertainties comparable to those observed for the NN interaction [36]. There is a larger cutoff dependence in the D waves and the mixing parameter ϵ_1 , specifically for $I = 0$. However, one has to keep in mind that there is no LEC up to NNLO for the D waves. The $^{33}D_1$ exhibits the trend of turning from negative to positive values at higher energies which cannot be described in an NNLO calculation, as discussed above.

LEC	{450, 500}	{600, 500}	{450, 700}	{600, 700}	
$I = 0$	$\tilde{C}_{1S_0}$	-0.151	-0.267	-0.151	-0.273
	C_{1S_0}	0.455	0.436	0.454	0.426
	$\tilde{C}_{1S_0}^a$	0.270	0.232	0.232	0.177
	$C_{1S_0}^a$	-0.915	-0.277	-0.905	-0.206
	C_{3P_0}	1.150	1.453	1.398	1.724
	$C_{3P_0}^a$	0.769	0.478	0.754	0.455
$I = 1$	$\tilde{C}_{1S_0}$	0	0	0	0
	C_{1S_0}	0.446	0.692	0.449	0.675
	$\tilde{C}_{1S_0}^a$	1.329	2.108	1.460	2.202
	$C_{1S_0}^a$	-1.118	-0.369	-1.214	-0.498
	C_{3P_0}	-0.357	-0.074	-0.321	0.041
	$C_{3P_0}^a$	0.501	0.232	0.498	0.222
$I = 0$	C_{1P_1}	0.384	-0.015	0.394	0.020
	$C_{1P_1}^a$	0.711	0.714	0.709	0.705
	C_{3P_1}	-0.374	-0.235	-0.296	-0.146
	$C_{3P_1}^a$	0.381	0.190	0.378	0.194
	$\tilde{C}_{3S_1}$	-0.132	-0.083	-0.122	-0.075
	C_{3S_1}	-0.497	-0.623	-0.731	-0.853
	$\tilde{C}_{3S_1}^a$	0.334	0.325	0.319	0.301
	$C_{3S_1}^a$	0.221	-0.573	0.325	-0.438
	C_{ϵ_1}	0.496	0.520	0.557	0.585
	$C_{\epsilon_1}^a$	-0.599	-0.218	-0.653	-0.290
$I = 1$	C_{1P_1}	-0.623	-0.735	-0.659	-0.858
	$C_{1P_1}^a$	0.682	0.544	0.688	0.573
	C_{3P_1}	-0.180	-0.373	-0.201	-0.443
	$C_{3P_1}^a$	0.716	0.628	0.719	0.645
	$\tilde{C}_{3S_1}$	-0.089	-0.120	-0.087	-0.122
	C_{3S_1}	0.698	0.148	0.707	0.188
	$\tilde{C}_{3S_1}^a$	0.399	0.210	0.398	0.224
	$C_{3S_1}^a$	0.164	0.665	0.124	0.602
	C_{ϵ_1}	0.245	0.182	0.279	0.237
	$C_{\epsilon_1}^a$	0.015	0.111	-0.019	-0.046
$I = 0$	C_{3P_2}	0.225	0.466	0.363	0.630
	$C_{3P_2}^a$	0.674	0.428	0.661	0.410
$I = 1$	C_{3P_2}	-0.362	-0.268	-0.361	-0.266
	$C_{3P_2}^a$	0.528	0.350	0.529	0.351

Table 1. The LECs at NLO for the different cutoff combinations $\{\Lambda [\text{MeV}], \tilde{\Lambda} [\text{MeV}]\}$. The values of the $\tilde{C}_i$ are in unit of 10^4 GeV^{-2} and the C_i in 10^4 GeV^{-4} . The parameters related to annihilation, $\tilde{C}_i^a$ and C_i^a (see eqs. (2.18)–(2.19)), are in units of 10^2 GeV^{-1} and 10^2 GeV^{-3} , respectively.

LEC	{450, 500}	{650, 500}	{450, 700}	{650, 700}	
$I = 0$	$\tilde{C}_{1S_0}$	-0.140	-0.278	-0.141	-0.299
	C_{1S_0}	0.456	0.459	0.456	0.463
	$\tilde{C}_{1S_0}^a$	0.208	0.247	0.155	0.219
	$C_{1S_0}^a$	-1.063	-0.337	-1.045	-0.233
	C_{3P_0}	0.031	0.310	-0.444	-0.217
	$C_{3P_0}^a$	0.796	0.492	0.828	0.556
$I = 1$	$\tilde{C}_{1S_0}$	0.025	0.095	0.052	-0.011
	C_{1S_0}	0.453	0.213	0.450	0.189
	$\tilde{C}_{1S_0}^a$	1.884	2.483	2.129	3.847
	$C_{1S_0}^a$	-1.733	-2.778	-2.566	-4.474
	C_{3P_0}	-0.535	-0.117	-0.531	-0.116
	$C_{3P_0}^a$	0.514	0.182	0.517	0.182
$I = 0$	C_{1P_1}	0.400	-0.113	0.438	-0.069
	$C_{1P_1}^a$	0.722	0.637	0.721	0.634
	C_{3P_1}	-0.521	-0.339	-0.596	-0.432
	$C_{3P_1}^a$	0.417	0.168	0.421	0.175
	$\tilde{C}_{3S_1}$	-0.162	-0.100	-0.183	-0.103
	C_{3S_1}	0.353	0.204	0.728	0.526
	$\tilde{C}_{3S_1}^a$	0.364	0.371	0.397	0.415
	$C_{3S_1}^a$	0.087	-0.841	-0.117	-1.125
	C_{ϵ_1}	0.205	0.236	0.062	0.106
	$C_{\epsilon_1}^a$	-0.485	-0.002	-0.362	0.167
$I = 1$	C_{1P_1}	-1.013	-1.294	-1.349	-1.869
	$C_{1P_1}^a$	0.711	0.535	0.775	0.668
	C_{3P_1}	-0.530	-0.902	-0.794	-1.356
	$C_{3P_1}^a$	0.742	0.630	0.788	0.735
	$\tilde{C}_{3S_1}$	-0.067	-0.143	-0.044	-0.125
	C_{3S_1}	1.150	0.764	1.325	1.235
	$\tilde{C}_{3S_1}^a$	0.413	0.282	0.411	0.402
	$C_{3S_1}^a$	-0.336	0.211	-0.896	-0.441
	C_{ϵ_1}	0.320	0.287	0.376	0.383
	$C_{\epsilon_1}^a$	-0.065	0.021	-0.182	-0.162
$I = 0$	C_{3P_2}	-0.300	-0.120	-0.518	-0.399
	$C_{3P_2}^a$	0.707	0.402	0.731	0.443
$I = 1$	C_{3P_2}	-0.648	-0.558	-0.821	-0.782
	$C_{3P_2}^a$	0.544	0.329	0.565	0.377

Table 2. The LECs at NNLO for the different cutoff combinations $\{\Lambda [\text{MeV}], \tilde{\Lambda} [\text{MeV}]\}$. The values of the $\tilde{C}_i$ are in unit of 10^4 GeV^{-2} and the C_i in 10^4 GeV^{-4} . The parameters related to annihilation, $\tilde{C}_i^a$ and C_i^a (see eqs. (2.18)–(2.19)), are in units of 10^2 GeV^{-1} and 10^2 GeV^{-3} , respectively.

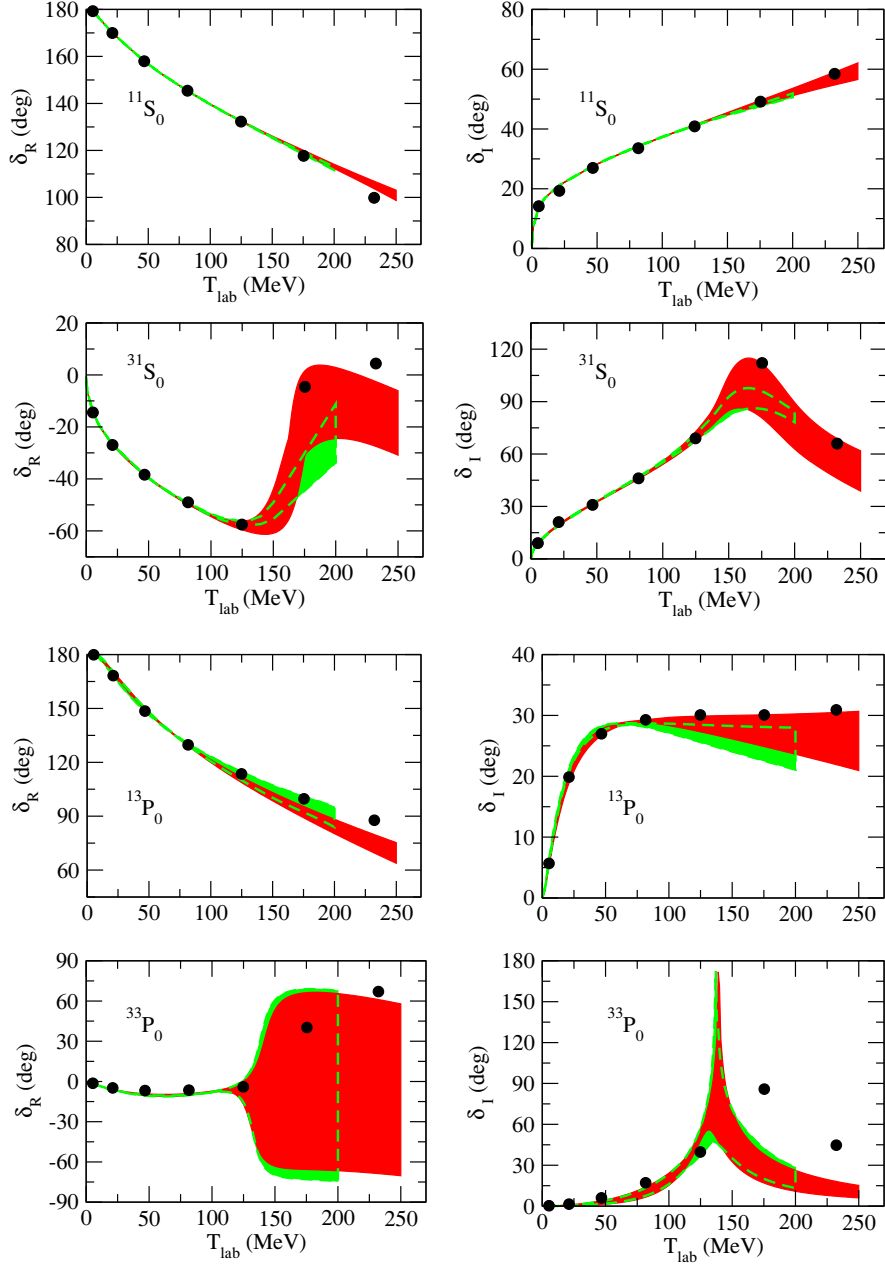


Figure 2. Real and imaginary parts of the phase shift in the 1S_0 and 3P_0 partial waves. The red/dark band shows the chiral EFT results up to NNLO for variations of the cutoff in the range $\Lambda = 450\text{--}650$ MeV in the Lippmann-Schwinger equation, while the green/light band are results to NLO for $\Lambda = 450\text{--}600$ MeV. The cutoff in the pion loops is varied independently in the range $\tilde{\Lambda} = 500\text{--}700$ MeV. The solid circles represent the solution of the PWA of ref. [20].

The situation in the $^3P_2\text{--}^3F_2$ channel is displayed in figure 5. In general our results agree with those of the PWA up to about 200 MeV within the uncertainty. Stronger deviations are visible again for those phases which show a resonance-like behavior like, e.g., the $^{13}P_2$.

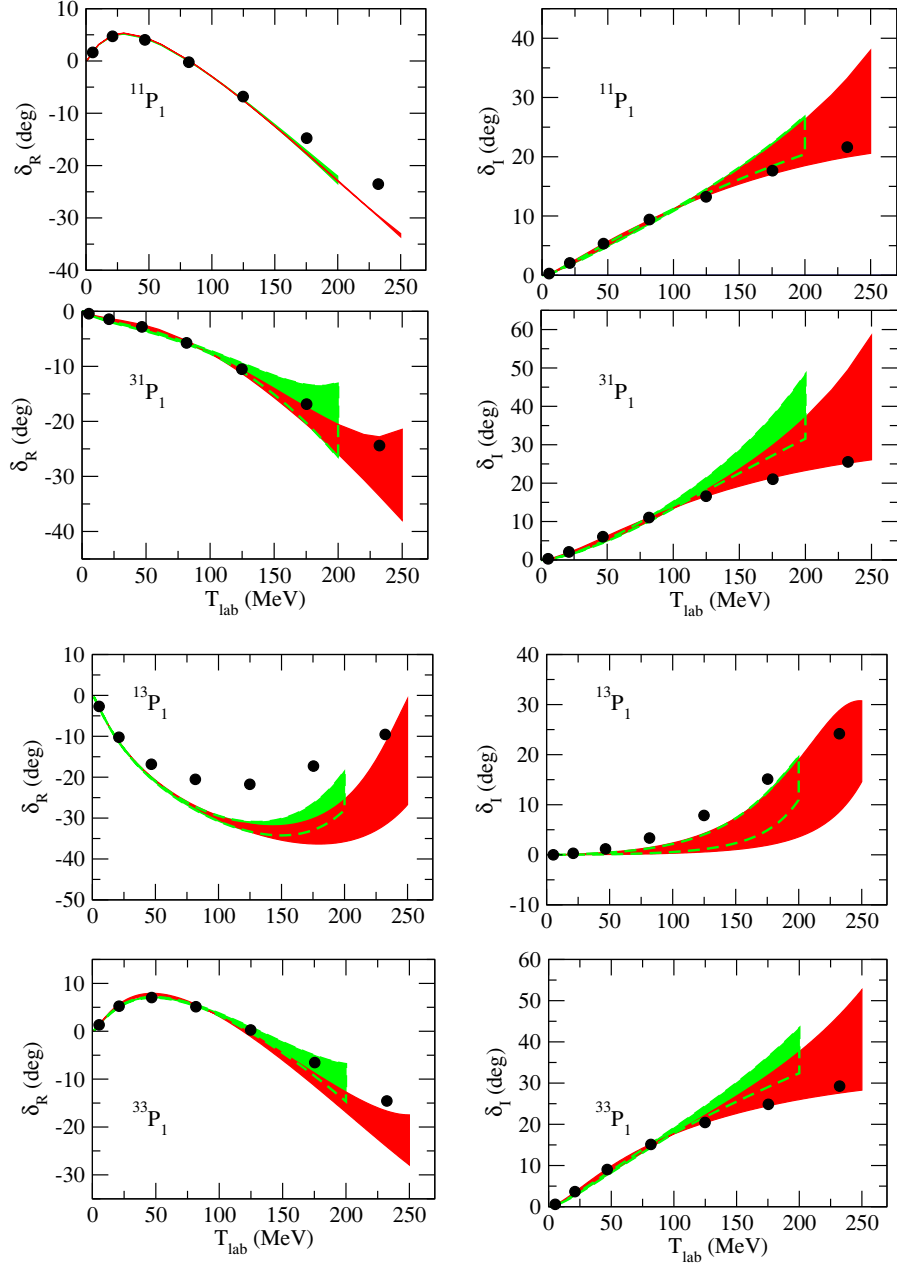


Figure 3. Real and imaginary parts of the phase shift in the 1P_1 and 3P_1 partial waves. The red/dark band shows the chiral EFT results up to NNLO for variations of the cutoff in the range $\Lambda = 450\text{--}650$ MeV in the Lippmann-Schwinger equation, while the green/light band are results to NLO for $\Lambda = 450\text{--}600$ MeV. The cutoff in the pion loops is varied independently in the range $\tilde{\Lambda} = 500\text{--}700$ MeV. The solid circles represent the solution of the PWA of ref. [20].

At last, in figure 6 the 1D_2 and 3D_2 phase shifts are presented. There are no LECs in those partial waves up to NNLO and, thus, our results are genuine predictions. The potential consists only of one- and two-pion exchange and, consequently, there is no contribution to annihilation. Thus, $\delta_I \equiv 0$ and we do not show this quantity.

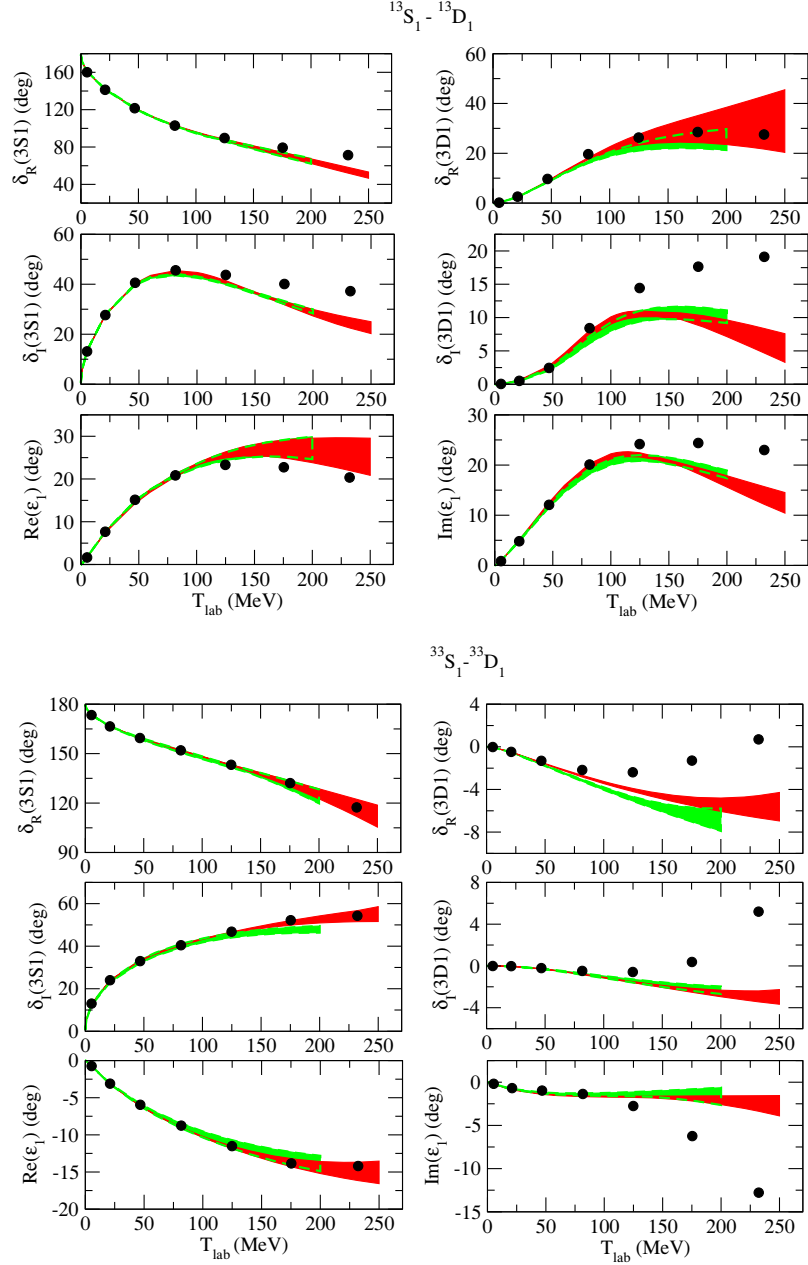


Figure 4. Real and imaginary parts of the phase shift in the 3S_1 - 3D_1 partial wave. The red/dark band shows the chiral EFT results up to NNLO for variations of the cutoff in the range $\Lambda = 450$ – 650 MeV in the Lippmann-Schwinger equation, while the green/light band are results to NLO for $\Lambda = 450$ – 600 MeV. The cutoff in the pion loops is varied independently in the range $\tilde{\Lambda} = 500$ – 700 MeV. The solid circles represent the solution of the PWA of ref. [20].

Results for the scattering lengths (for 1S_0 and 3S_1) and for scattering volumes (for the P waves) are summarized in table 3. These are complex numbers because of the presence of annihilation. The effective range parameters implied directly by the PWA of [20] are not provided in that reference. Thus, the lowest energy that enters our fitting procedure

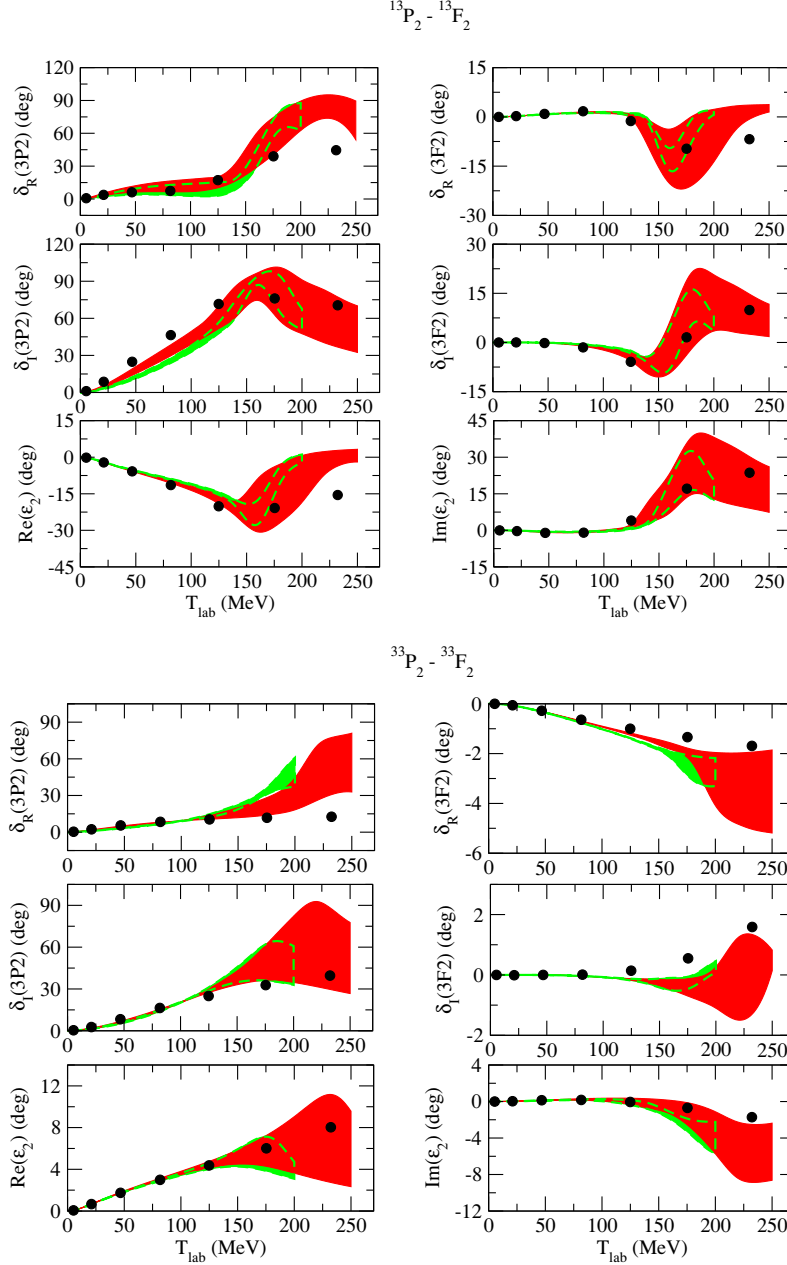


Figure 5. Real and imaginary parts of the phase shift in the 3P_2 - 3F_2 partial wave. The red/dark band shows the chiral EFT results up to NNLO for variations of the cutoff in the range $\Lambda = 450$ – 650 MeV in the Lippmann-Schwinger equation, while the green/light band are results to NLO for $\Lambda = 450$ – 600 MeV. The cutoff in the pion loops is varied independently in the range $\tilde{\Lambda} = 500$ – 700 MeV. The solid circles represent the solution of the PWA of ref. [20].

concerns the phase shifts at $p_{lab} = 100 \text{ MeV}/c$ which corresponds to $T_{lab} = 5.3 \text{ MeV}$. In view of that one can consider our values as predictions of chiral EFT. As one can see in table 3 we get practically the same results at NLO and at NNLO and, moreover, there is very little cutoff dependence. Actually, in case of $\text{Re } a_{1S_0}$ in the $I = 0$ channel there is no variation in the first two digits and, therefore, only a single number is given.

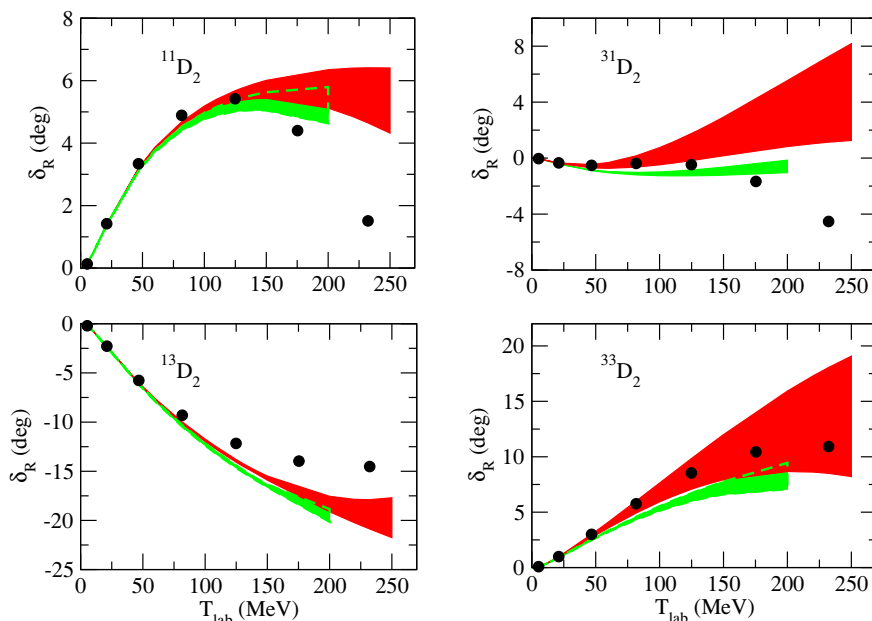


Figure 6. Real part of the phase shift in the 1D_2 and 3D_2 partial waves. The red/dark band shows the chiral EFT results up to NNLO for variations of the cutoff in the range $\Lambda = 450\text{--}650$ MeV in the Lippmann-Schwinger equation, while the green/light band are results to NLO for $\Lambda = 450\text{--}600$ MeV. The cutoff in the pion loops is varied independently in the range $\tilde{\Lambda} = 500\text{--}700$ MeV. The solid circles represent the solution of the PWA of ref. [20].

Table 3 contains also scattering lengths and volumes predicted by the most refined meson-exchange potential developed by the Jülich group, namely model D published in [7]. It is interesting to see that at least for the S waves the results are very similar not only on a qualitative level but in most cases even on a quantitative level. One has to keep in mind that there are no data that would allow one to fix the relative magnitude of the singlet- and triplet- contributions near threshold. Moreover, the Jülich $\bar{N}N$ potential was only fitted to integrated cross sections. Differential cross sections or polarization data were not considered.

There is some experimental information that puts constraints on these scattering lengths. Measurements of the level shifts and widths of antiproton-proton allow one to deduce values for the spin-averaged $\bar{p}p$ scattering lengths via the Deser-Trueman formula. Corresponding results taken from ref. [52] are listed in table 4. In that reference one can also find values for the imaginary part of the scattering lengths that are inferred from measurements of the $(\bar{n}p$ and $\bar{p}p)$ annihilation cross section. A comparison directly with the measured level shifts and widths [53–56] is provided in table 5 where now the Deser-Trueman formula was applied to the theory results.

As far as we know, this experimental evidence was not taken into account in the PWA [20]. Nonetheless, for completeness we provide the predictions based on our EFT interaction. One should be cautious, however, in comparing our results with the experimental numbers. As said above, our calculations are performed in the isospin basis so that $a_{\bar{p}p}$ is simply given by $(a_{I=0} + a_{I=1})/2$. It is known that the presence of the Coulomb force

		I=0	I=1
1S_0	NLO	$-0.21 - i(1.20 \cdots 1.21)$	$(1.03 \cdots 1.04) - i(0.56 \cdots 0.58)$
	NNLO	$-0.21 - i(1.21 \cdots 1.22)$	$(1.02 \cdots 1.04) - i(0.57 \cdots 0.61)$
	model D	$-0.23 - i1.01$	$0.99 - i0.58$
3S_1	NLO	$(1.34 \cdots 1.37) - i(0.88 \cdots 0.90)$	$(0.43 \cdots 0.44) - i(0.87 \cdots 0.90)$
	NNLO	$(1.37 \cdots 1.38) - i(0.86 \cdots 0.88)$	$(0.43 \cdots 0.44) - i(0.91 \cdots 0.92)$
	model D	$1.55 - i1.45$	$0.33 - i0.96$
3P_0	NLO	$-(3.55 \cdots 4.32) - i(7.35 \cdots 8.45)$	$(2.42 \cdots 2.47) - i(0.03 \cdots 0.10)$
	NNLO	$-(3.08 \cdots 3.78) - i(6.93 \cdots 7.55)$	$(2.35 \cdots 2.42) - i(0.03 \cdots 0.12)$
	model D	$-7.40 - i3.21$	$2.50 - i1.23$
1P_1	NLO	$-(2.84 \cdots 2.86) - i(0.24 \cdots 0.29)$	$(0.89 \cdots 0.92) - i(0.19 \cdots 0.20)$
	NNLO	$-(2.87 \cdots 2.89) - i(0.25 \cdots 0.31)$	$(0.78 \cdots 0.86) - i(0.20 \cdots 0.29)$
	model D	$-3.26 - i0.50$	$0.45 - i0.55$
3P_1	NLO	$(4.80 \cdots 4.82) - i(0.00 \cdots 0.02)$	$-(1.95 \cdots 1.97) - i(0.37 \cdots 0.40)$
	NNLO	$(4.76 \cdots 4.77) - i(0.00 \cdots 0.02)$	$-(2.02 \cdots 2.09) - i(0.39 \cdots 0.52)$
	model D	$4.87 - i0.06$	$-2.05 - i1.37$
3P_2	NLO	$-(0.31 \cdots 0.42) - i(0.27 \cdots 0.51)$	$-(0.20 \cdots 0.21) - i(0.16 \cdots 0.21)$
	NNLO	$-(0.45 \cdots 0.78) - i(0.47 \cdots 0.65)$	$-(0.28 \cdots 0.37) - i(0.18 \cdots 0.25)$
	model D	$-0.14 - i1.27$	$-0.37 - i0.50$

Table 3. Scattering lengths (in fm) for the S -waves and scattering volumes (in fm³) for the P -waves in the isospin $I = 0$ and $I = 1$ channels. Results based on the NLO and NNLO potentials are given and compared with the predictions of the Jülich $\bar{N}N$ model D [7].

	chiral EFT	model D	Experiment
$\bar{a}_{S, \bar{p}p}$	NLO	$0.80 - i1.10$	(0.95 ± 0.02) $-i(0.73 \pm 0.03)$
	NNLO		
	model D		
$\text{Im } \bar{a}_{S, I=1}$	NLO	-0.86	(-0.83 ± 0.07)
	NNLO		
$\text{Im } \bar{a}_{S, I=0}$	NLO	-1.34	(-0.63 ± 0.08)
	NNLO		
$\bar{a}_{P, \bar{p}p}$	NLO	$-0.31 - i0.87$	-0.61 ± 0.81 $-i(0.77 \pm 0.06)$
	NNLO		
	model D		

Table 4. Spin-averaged scattering lengths for S -wave ($\bar{a}_S$; in fm) and scattering volumes for P -wave ($\bar{a}_P$; in fm³). Results based on the NLO and NNLO potentials are given and compared with the predictions of the Jülich $\bar{N}N$ model D [7]. The experimental information is taken from ref. [52].

in $\bar{p}p$ and the p - n mass difference lead to changes of the S -wave scattering lengths in the order of 0.1 fm [57] and, therefore, one should not take quantitative differences too serious. Note also that additional assumptions have to be made in order to deduce the splitting of the 1S_0 and 3S_1 level shifts from the experiment [52, 58].

		ΔE (eV)	Γ (eV)
1S_0	NLO	$-(306 \cdots 361)$	$(1528 \cdots 1553)$
	NNLO	$-(302 \cdots 361)$	$(1545 \cdots 1589)$
	model D	-330	1380
	Experiment [53]	-740 ± 150	1600 ± 400
	[54]	-440 ± 75	1200 ± 250
3S_1	NLO	$-(768 \cdots 786)$	$(1519 \cdots 1562)$
	NNLO	$-(781 \cdots 790)$	$(1537 \cdots 1563)$
	model D	-816	2092
	Experiment [54]	-785 ± 35	940 ± 80
	[55]	-850 ± 42	770 ± 150
		ΔE (meV)	Γ (meV)
3P_0	NLO	$+(13 \cdots 23)$	$(159 \cdots 232)$
	NNLO	$+(8 \cdots 17)$	$(173 \cdots 186)$
	model D	$+60$	109
	Experiment [56]	$+139 \pm 28$	120 ± 25

Table 5. Hadronic shifts and broadenings in hyperfine states of $\bar{p}H$. Results based on the NLO and NNLO potentials are given and compared with the predictions of the Jülich $\bar{N}N$ model D [7]. The experimental information is taken from refs. [53–56].

Finally, let us discuss $\bar{N}N$ bound states. Several of the phase shifts tabulated in ref. [20] start at 180° at $T_{lab} = 0$ MeV, namely $^{11}S_0$, $^{13}P_0$, $^{13}S_1$, and $^{33}S_1$, which according to the standard convention based on the Levinson theorem signals the presence of a bound state. Therefore, we performed a search for possible bound states generated by our EFT interaction where we restricted ourselves to energies not too far from the $\bar{N}N$ threshold. We did not find any near-threshold poles in the $^{11}S_0$ and $^{33}S_1$ – $^{33}D_1$ partial waves. In case of the $^{13}S_1$ – $^{13}D_1$ interaction there is a pole which corresponds to a “binding” energy of $E_B = +(5.6 \cdots 7.7) - i(49.2 \cdots 60.5)$ MeV, depending on the cutoffs $\{\Lambda, \tilde{\Lambda}\}$, at NLO and $E_B = +(4.8 \cdots 21.3) - i(60.6 \cdots 74.9)$ MeV at NNLO. The positive sign of the real part of E_B indicates that the poles we found are actually located above the $\bar{N}N$ threshold. But they move below the threshold when we switch off the imaginary part of the potential and that is the reason why we refer to them as bound states. To be precise these are unstable bound states in the terminology of ref. [59]. Note that those poles lie on the physical sheet and, therefore, do not correspond to resonances. Evidently, the width of the state, $\Gamma = -2\text{Im } E_B$, is rather large. There is also a pole in the $^{13}P_0$ partial wave. It corresponds to a binding energy of $E_B = (-1.1 \cdots +1.9) - i(17.8 \cdots 22.4)$ MeV at NLO and $E_B = -(3.7 \cdots 0.2) - i(22.0 \cdots 26.4)$ MeV at NNLO. In this context we want to mention that bound states and also resonances have been likewise found in other studies of the $\bar{N}N$ interaction, see refs. [8, 9] for recent examples.

4 Summary and outlook

In this paper we presented an exploratory study of the $\bar{N}N$ interaction in a chiral effective field theory approach based on a modified Weinberg power counting, analogous to the NN

case in [18, 36]. The $\bar{N}N$ potential has been evaluated up to NNLO in the perturbative expansion and the arising low-energy constants have been fixed by a fit to the phase shifts and inelasticities provided by a recently published phase-shift analysis of $\bar{p}p$ scattering data [20]. It turned out that the overall quality of the description of the $\bar{N}N$ amplitudes that can be achieved at NNLO is comparable to the one found in case of the NN interaction at the same order [36]. Specifically, for the S -waves ($^{11}S_0$, $^{13}S_1$, $^{33}S_1$) nice agreement with the phase shifts and inelasticities of [20] has been obtained up to laboratory energies of about 200 MeV, i.e. over almost the whole energy region considered. The same is also the case for many of the P -waves. Thus, we conclude that the chiral EFT approach, applied successfully in refs. [17, 18] to the NN interaction and in refs. [60, 61] to the hyperon-nucleon interaction, is very well suited for studies of the $\bar{N}N$ interaction too.

Of course, there are also some visible deficiencies in our results. They occur primarily in those partial waves where the partial-wave analysis of [20] suggests the presence of (presumably strongly inelastic) resonances at energies around $T_{lab} \approx 200 - 250$ MeV. It is not surprising that structures in this energy region cannot be reproduced reliably within our NNLO calculation. Clearly, here an extension of our investigation to N³LO is necessary for improving the description of the $\bar{N}N$ interaction. Therefore, we plan to extend our study to N³LO in the future. At this stage it will become sensible to perform the calculation in particle basis so that the Coulomb interaction in the $\bar{p}p$ system can be taken into account rigorously, and to compute observables and compare them directly with scattering data for $\bar{p}p$ elastic scattering and for the charge-exchange reaction $\bar{p}p \rightarrow \bar{n}n$. Annihilation processes that occur predominantly at short distances reduce the magnitude of the S -wave amplitudes so that higher partial waves start to become important at much lower energies as compared to what one knows from the NN interaction. Thus, without a realistic description of higher partial waves, and particularly of the D -waves, it is not meaningful to confront the amplitudes resulting from our NNLO interaction directly with $\bar{N}N$ data and, therefore, we have refrained from doing so in the present work.

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