The Problem of Renormalization of Chiral Nuclear Forces

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Ever since quantum field theory was first applied to the derivation of nuclear forces in the mid-twentieth century, the renormalization of pion exchange with realistic couplings has presented a challenge. The implementation of effective field theories (EFTs) in the 1990s promised a solution to this problem but unexpected obstacles were encountered. The response of the nuclear community has been to focus on “chiral potentials” with regulators chosen to produce a good description of data. Meanwhile, a successful EFT without explicit pion exchange—Pionless EFT—has been formulated where renormalization is achieved order by order in a systematic expansion of low-energy nuclear observables. I describe how lessons from Pionless EFT are being applied to the construction of a properly renormalized Chiral EFT.

Keywords: effective field theory, renormalization, nuclear physics, singular potentials, pion exchange

1. INTRODUCTION

In the aftermath of the solution of the “problem of infinities” in Quantum Electrodynamics (QED), an intense quest set in to renormalize nuclear forces, where pion exchange replaced the photon exchange responsible for atomic forces. (For an early example, see reference [1].) It was quickly understood that the only relativistic pion-nucleon coupling that is renormalizable is pseudoscalar [2]. However, pseudoscalar coupling differs from pseudovector coupling by a large nucleon-pair term, which was found to be in conflict with pion phenomenology [3]. For the favored pseudovector coupling, the description of two-nucleon data depended sensitively on the high-momentum (or short-distance) cutoff (see, for example, reference [4]). Efforts moved toward the investigation of various prescriptions for handling short-range effects, including specific cocktails of (usually single-)heavier-meson exchange, form factors with ad hoc shapes, and/or boundary conditions at some finite distance. Nuclear theory acquired an increasingly phenomenological character. Typically, the non-relativistic Schrödinger equation was solved with a two-nucleon (2N) potential including one-pion exchange, some approximation to two-pion exchange, and a more or less arbitrary short-range form, with sufficiently many parameters to fit data to the desired accuracy. The end result was that potentials including quite different physics could produce very good parameterizations of 2N data up to around the pion-production threshold, while typically underpredicting three- and more-nucleon binding by more than 10%. A serious difficulty was to infer a satisfactory form of three-nucleon (3N) forces and, for reactions, 2N currents. Reference [5] recounts some of this history.

In contrast, by the mid-1970s renormalizable quantum field theories had won the day in particle physics, leading to the formulation of Quantum Chromodynamics (QCD) as the theory of strong interactions. Out of the attempts to make predictions for QCD at low energies and to understand how the Standard Model (SM) can arise from a more fundamental theory, the concept of effective field theory (EFT) was born [6]. An EFT comprises all the interactions among relevant degrees of freedom at the low-energy scale. The hope has been that some physical phenomenon can be described in terms of a low-energy EFT, with the high-energy physics encoded by the coupling constants of the EFT. The standard procedure is to make a systematic expansion of the low-energy observables in terms of powers of a non-dimensional energy scale parameter, $\Lambda$. This parameter is often related to a high-energy scale parameter, $\Lambda'$, by $\Lambda \ll \Lambda'$. The goal is to extract the high-energy physics encoded in the EFT by systematically improving the description of the low-energy observables as $\Lambda' \to \infty$.
freedom that are allowed by symmetries, including an arbitrary number of fields and derivatives. For predictions, contributions to observables must be ordered according to their expected size. This “power counting” allows for an \textit{a priori} error estimate from neglected higher-order contributions. At each order in the expansion, only a finite number of “low-energy constants” (LECs)—the interaction strengths—appear. In a consistent power counting, they are sufficient to ensure that any dependence on the regulator can be made arbitrarily small by taking the cutoff large. Thus, EFTs are renormalizable in the modern sense that at each order a finite number of parameters generate results for observables that are independent of details of the arbitrary regularization procedure.

A successful EFT, Chiral Perturbation Theory (ChPT), was developed in the 1980s to handle interactions among pions and one nucleon below the characteristic QCD scale \(M_{\text{QCD}} \sim 1 \text{ GeV}\) [7, 8]. Requiring renormalization in a perturbative expansion, a consistent power counting was developed based on “naive dimensional analysis” (NDA) [9]. Taking the typical external momentum in a reaction to be of the order of the pion mass, \(Q \sim m_\pi \ll M_{\text{QCD}}\), observables are expanded in a series of powers of \(Q/M_{\text{QCD}}\) times calculable functions of \(Q/m_\pi\). When Weinberg remarked [10, 11] that ChPT, now generalized as “Chiral EFT” (ChEFT), could be used to derive nuclear forces, he identified an infrared enhancement in nuclear amplitudes by the nucleon mass \(m_N = O(M_{\text{QCD}})\), which can lead to the failure of perturbation theory—a good thing since nuclei are bound states and resonances. He proposed that the ChPT power counting could still be applied to the nuclear potential, defined as the sum of diagrams lacking an explicit enhancement. Then, the Lippmann-Schwinger equation, or equivalently the Schrödinger equation, would be solved with a truncated “chiral potential.”

The potential defined by Weinberg contains pion exchange diagrams where all LECs are fixed, at least in principle, from ChPT. But it also includes shorter-range interactions with LECs that can only be determined in nuclear systems. Implicit in Weinberg’s proposal was that the short-range LECs would not contain an \textit{implicit} enhancement. This would be the case if the solution of the dynamical equation does not generate cutoff dependence beyond that which can be compensated by the LECs already present up to that order according to NDA.

Whether this assumption is true was not immediately clear. NDA says that the potential at leading order (LO) consists of two non-derivative, chirally symmetric contact interactions together with one-pion exchange (OPE). More-pion exchange should come at higher orders together with more-derivative contact interactions. Non-perturbative pion exchange prevents an analytical solution even at the 2N level. The first numerical solution of a chiral potential in the 2N system [12, 13] tested renormalizability of the amplitude: a variation from 0.5 to 1 GeV in the cutoff of a local Gaussian regulator seemed to be compensated by a refitting of the LECs at hand. However, the fitting procedure was cumbersome as an over-complete set of interactions was used and the local regulator mixed different partial waves, limiting the range of cutoffs that could be explored. Since then a large variety of chiral potentials have been developed (for reviews, see for example references [14, 15]).

A landmark was a 2N potential [16] that was perceived to match the accuracy of phenomenological potentials (for a recent comparison between chiral 2N potentials and data, see reference [17]). Chiral potentials have become the favorite input to “\textit{ab initio}” methods, which provide numerically controlled solutions of the Schrödinger equation for multi-nucleon systems.

Unfortunately, pretty early on the first evidence appeared [18] that Weinberg’s prescription does not provide amplitudes, and thus observables, that are renormalized order by order. In the 2N \(1S_0\) channel at LO, a semi-analytical argument shows that there remains a logarithmic dependence on the cutoff proportional to the average quark mass. The only way to eliminate it, at least with a momentum- or coordinate-space cutoff, is to include at LO a non-derivative, chirally \textit{breaking} contact interaction, which according to NDA should appear two orders down the expansion, that is, at next-to-next-to-leading order (N\textsuperscript{2}LO)\textsuperscript{1}. More dramatically, it was later shown [20, 21] that oscillatory cutoff dependence appears at LO in waves where OPE is attractive, singular, and accounted for non-perturbatively. A chirally symmetric LEC is needed for renormalization in each wave, but again NDA assigns those in partial waves beyond \(S\) to higher orders. Similar problems afflict processes with external probes [22].

As I describe in section 3, the origin of these problems is the renormalization of attractive singular potentials [23, 24]. NDA might fail because exact solutions of the Schrödinger equation depend on the cutoff differently than perturbative solutions. The LECs needed for the renormalization of the amplitude are enhanced by implicit powers of \(M_{\text{QCD}}\).

How to account for this? In response to the renormalization failure of Weinberg’s power counting a simpler nuclear EFT [25–27] was developed in the late 1990s. In this “Pionless EFT” pions are integrated out and only contact interactions remain. The effects of loops in the Lippmann-Schwinger equation are much easier to see, including the \(m_N\) enhancement and a further enhancement of \(4\pi\) that was not pointed out by Weinberg. The lessons of Pionless EFT for ChEFT are summarized in section 2.

The first attempt to fix power counting using the insights from Pionless EFT was initiated [28, 29] at the same time as the main elements of the power counting of Pionless EFT were being understood. Valid for sufficiently small values of the pion mass and external momenta, this version of ChEFT treats pion exchange in perturbation theory, removing the renormalization problems mentioned above. Unfortunately, in the 2N system at physical pion mass one cannot go in this way to momenta much beyond those of Pionless EFT [30]. The alternative is \textit{partly} perturbative pions: OPE is iterated only in the low partial waves where it is sufficiently strong, together with the contact interactions whose LECs are necessary

\textsuperscript{1}A note on notation: It has become usual in the nuclear community to refer to a subleading chiral potential of order \(n \geq 2\) as \(N^{n-1}\text{LO}\); because with Weinberg’s power counting the parity- and time-reversal-invariant potential of order \(n = 1\) vanishes [19]. However, this usage is too provincial to accommodate experience with other observables and power countings in ChEFT or other EFTs. Here, a correction of order \(n\) in the expansion is denoted as \(N^n\text{LO}\), whether it is non-zero or not.
for LO renormalization [20]. All subleading pion exchanges, together with the remaining contact interactions, are treated in perturbation theory [31]. This approach is discussed in section 4, including what little has been done to confront it with data.

Section 5 offers the conclusion that this approach solves the renormalization woes of nuclear forces while accounting for the long-range interactions from pion exchange systematically. Although they differ in detail from the field-theoretical renormalization described below, renormalization-group analyses of the Schrödinger equation [22, 32–34] support this picture. How it can meet the accuracy requirements of the nuclear community remains to be seen. My emphasis here is on the internal consistency of ChPT. I expand on the renormalization issues summarized in reference [35], but I refer the reader to the latter for a more complete review of ChPT and its relation to other nuclear EFTs.

2. SAY WHAT?

As reviewed in reference [35], defining the nuclear potential as the sum of "irreducible" diagrams without the \( m_N \) infrared (IR) enhancement does indeed ensure that the cutoff-independent parts of pion-exchange diagrams can be ordered according to ChPT power counting. These components of the pion-exchange potentials are in general non-analytic functions of momenta and pion mass that can be calculated in terms of pion-baryon interactions.

The ChPT power counting is designed for processes where the typical external momentum is comparable to the pion mass, \( Q \sim m_\pi \). A (relativistic) pion propagator scales as \( Q^{-2} \). In contrast, a nucleon is heavy compared to \( Q \) and thus non-relativistic. Moreover, energies and three-momenta being comparable, nucleon recoil is suppressed by one power of \( Q/m_N \equiv O(Q/M_{QCD}) \)—that is, the nucleon is static, its propagator scaling as \( Q^{-1} \). Because the Delta-nucleon mass difference is (at physical quark masses) only about twice the pion mass, a Delta propagator scales in the same way. In integrals from the loops that make up the potential one picks poles from the pion propagators, typically resulting in factors of \( (4\pi)^{-2} \). They combine with factors of the pion decay constant \( f_\pi \approx 92 \text{ MeV} \) from the pion-baryon interactions to produce inverse factors of \( 4\pi f_\pi = O(M_{QCD}) \). The power counting explicitly relies on an estimate, NDA [9], of the factors of \( 4\pi \) that distinguish between \( f_\pi \) and the breakdown scale \( M_{QCD} \), which appears in interactions with derivatives and powers of the pion mass. In summary, the ChPT rules (in momentum space) are:

\[
\begin{align*}
\text{(pion) loop integral} & \sim (4\pi)^{-2} Q^4, \\
\text{baryon, pion propagator} & \sim Q^{-1}, Q^{-2}, \\
\text{vertex} & \sim Q^4 f_\pi^{2-b-f} M_{QCD}^{2-d-f/2},
\end{align*}
\]

where \( d, b, \) and \( f \) are the numbers of derivatives/pion masses, pion fields, and baryon fields, respectively, in an interaction.

The expected size of any diagram can be found using the identities \( I = L - 1 + \sum_i V_i \) and \( 2I + E = \sum_i V_i (b_i + f_i) \) involving the number of loops \( L \), internal (external) lines \( I \) (\( E \)), and vertices \( (V_i) \) having a set of values \( d = d_i, b = b_i, \) and \( f = f_i \). In particular,

\[
2N \text{ potential} \sim 4\pi m_N^{-1} M_{NN}^{-1} \left( Q M_{QCD}^{-1} \right)^\mu ,
\]

where [28, 29]

\[
M_{NN} \equiv \frac{16\pi f_\pi^2}{k_\Lambda m_N} = O(f_\pi)
\]

in terms of the pion-nucleon axial-vector coupling \( g_A \approx 1.27 \) and [11]

\[
\mu \equiv 2L + \sum_i V_i (d_i + f_i/2 - 2).
\]

Because every additional loop (without increase in the number of derivatives/pion masses at vertices) leads to a relative factor \( O(Q^2/M_{QCD}^2) \), one gets the well-known ordering where p-pion exchange starts at \( \mu = 2(p-1) \). Note that the NLO correction vanishes due to parity and time-reversal symmetries [19].

This power counting applies to diagrams that make up the long-range potential. Yet physics, as opposed to metaphysics, is about observables. The meaning of Equation (4) is that it indirectly orders the contributions to amplitudes. For the direct link, we need to consider as well "reducible" diagrams where intermediate states contain only nucleons. One picks poles from the non-relativistic nucleon propagators, for which energies are of the order of recoil—in those diagrams, one cannot approximate nucleons as static. (This of course has nothing to do with relativistic corrections, as sometimes misstated in the literature.) These poles lead not only to an \( m_N \) enhancement [10, 11], but typically also to different powers of \( (4\pi)^{-1} \). Experience with Pionless EFT [35, 36], where these are all the loops one needs to deal with, shows that the factors associated with reducible loops are

\[
\text{nucleon propagator} \sim m_N Q^{-2},
\]

\[
\text{reducible loop integral} \sim (4\pi m_N)^{-1} Q^5.
\]

When one inserts the order-\( \mu \) potential into a 2N diagram, we need one extra reducible loop with two nucleon propagators (compare Figures 1A, B), leading to a relative factor \( O(Q^2/M_{QCD}^2) \). This amount to an IR enhancement of \( 4\pi m_N/Q \) over the factor that arises from Equations (1) and (2). As a consequence, the series in the LO potential fails to converge for \( Q \sim M_{NN} \). This is what makes ChEFT different for \( A \geq 2 \) nucleons compared to ChPT for \( A \leq 1 \).

The factor of \( 4\pi \) in the IR enhancement had not been recognized before Pionless EFT was developed, but it is important to understand the failure of perturbation theory for pions. The exact solution of the LO potential for \( Q \sim M_{NN} \) can give rise to a binding energy per nucleon

\[
\frac{B_A}{A} \sim \frac{M_{NN}^2}{M_{QCD}} \sim \frac{f_\pi}{4\pi} \sim 10 \text{ MeV}.
\]

This is somewhat larger than observed for light nuclei, indicating a certain amount of fine tuning in the 2N interactions. But it is on
To estimate the respective contributions to the AN amplitude, one can first consider the LO ($\mu = 0$), 2N potential: to produce a connected diagram, we need at least $A - 1$ 2N interactions linked by $A - 2$ propagators. Next, one insertion of a subleading $aN$ potential between two LO amplitudes comes with $A + a - 2$ propagators and $A + a$ loops. Another insertion of the same subleading potential takes $a$ additional propagators and $a - 1$ additional loops, and so on. The rules (7), (8) imply that an $aN$ potential of index $\mu$ gives, at $Q \sim M_{NN}$,

$$A\!N \text{ amplitude } \sim (4\pi)^{A-1} m_N^{-1} M_{NN}^{a-1} Q^{2-a} \left( QM_{\text{QCD}}^{-1} \right)^{\nu}, \quad (11)$$

where

$$\nu \equiv \mu + a - 2 \quad (12)$$

and $n$ is the order in perturbation theory. While $\nu$ is the perturbative cost of one insertion of a subleading potential characterized by $\mu$ (6) and $a$, $n$ insertions cost $\nu n$ as indicated by the power of $Q/M_{\text{QCD}}$ in Equation (11). The presence of $a - 2$ [instead of $2(a-2)$] in $\nu$ reflects the suppression by $(4\pi)^{-1}$ [instead of $(4\pi)^{-2}$] in more-nucleon forces. A sample of pion-range diagrams that contributes at various values of $\nu$ is shown in Figure 2 (see reference [35] for more details).

The $n$ in Equation (11) encodes the perturbative character of any subleading interaction. A common fallacy is that the mere definition of a potential means that the corresponding dynamical (Lippmann-Schwinger or Schrödinger) equation must be solved exactly. On the contrary, if there is a sense in which a subleading potential can be treated non-perturbatively, then it should also be possible to include it in distorted-wave perturbation theory, where the distortion is caused by the LO potential. If that is not the case, then at least part of that "subleading" potential is not subleading. Such a consistency test is almost completely ignored in the community. The one exception I am aware of is reference [42], where it is shown that this test is not met by most available chiral potentials.

"But surely," you might be reasoning, "a subleading potential can be treated non-perturbatively." That is certainly the case for a regular subleading potential, but not necessarily for a singular potential, for which neither the perturbative series nor the exact solution of the dynamical equation are well-defined without (potentially distinct) counterterms. So far I have been glossing over the cutoff dependence that usually arises in loops and is, of course, present in the LECs. A regulator is nothing...
FIGURE 2 | Sample of pion-range diagrams in the aN nuclear potential ordered according to the expected size of their contributions to the amplitude (Equation 12).

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but a way to split short-range physics between loops and LECs. If we increase a momentum cutoff \( \Lambda \) (or decrease a coordinate cutoff \( R \sim \Lambda^{-1} \)), we account, correctly or incorrectly, for more short-range physics through the loops of the Lippmann-Schwinger equation. As long as \( \Lambda \gtrsim M_{QCD} \), we can compensate by changing the LECs present at the same order, without increasing the relative truncation error of \( O(Q/M_{QCD}) \). The crucial point is that only the combination of the two effects matter, and physics enters through the fitting of as many observables as LECs—observables which are either calculated in the underlying theory (when we speak of “matching” the EFT to the underlying theory) or measured experimentally. This process of renormalization is essential for amplitudes to be free of detailed assumptions about short-range physics, and in general only the sum of all contributions at a given order—loops and LECs ensuring renormalization—can be said to be perturbative or not.

If all we needed was to eliminate the cutoff-dependent parts of pion exchange in the potential, the LECs for the job would be given by NDA, by construction [9]. It is crucial to realize, though, that reducible loops introduce further cutoff dependence, which we need eliminate as well. The potential itself has to depend on the cutoff so that observables do not. The LECs that renormalize this part of the \( A \gtrsim 2 \) problem will not in general satisfy NDA. We examine this aspect of renormalization next.

3. RENORMALIZATION OF SINGULAR POTENTIALS

The difficulty we face is that EFT potentials are singular and, because of additional derivatives and loops, they get more and more singular as the order of the EFT expansion increases. Singularities are apparent already in the LO (\( \mu = 0, a = 2 \)) pion-range potential, OPE: labeling the two nucleons 1 and 2,

\[
V_{OPE}(\vec{r}) = \frac{\tau_1 \cdot \tau_2}{m_N M_{NN}} \left[ \frac{e^{-m_\pi r}}{r^5} \left( 1 + m_\pi r + \frac{m_\pi^2 r^2}{3} \right) S_{12}(\vec{r}) + \left( \frac{m_\pi^2 e^{-m_\pi r}}{r} - 4\pi \delta(\vec{r}) \right) \frac{\vec{\sigma}_1 \cdot \vec{\sigma}_2}{3} \right],
\]

where \( \vec{r} = r \hat{r} \) is the relative position, \( \vec{\sigma}_i (\tau_i) \) is the spin (isospin) Pauli matrix for nucleon \( i \), and

\[
S_{12}(\vec{r}) = 3 \vec{\sigma}_1 \cdot \hat{r} \vec{\sigma}_2 \cdot \hat{r} - \vec{\sigma}_1 \cdot \vec{\sigma}_2
\]

is the spin-tensor operator. While the delta function contributes only to \( S \) waves, the tensor potential is non-vanishing for total
spin \( s = 1 \) and can mix waves with orbital angular momentum \( l = j \pm 1 \). It is attractive in some uncoupled waves like \( ^3P_0 \) and \( ^3D_2 \), and in one of the eigenchannels of each coupled wave. The regular Yukawa potential is attractive in isovector (isoscalar) channels for \( s = 0 \) (\( s = 1 \)). More-pion exchange leads to more singular terms, \( p \)-pion exchange containing for example terms \( \alpha^{-2(p+1)} \) in addition to delta functions and their derivatives.

For \( Q \approx M_{\text{NN}} \) OPE is expected to be non-perturbative by the argument of the previous section. It has been known for a long time (see e.g., the review [44]) that attractive singular potentials, treated exactly, do not fully determine the solution of the Schrödinger equation [44]. This is a manifestation that renormalization of a singular potential requires contact terms that naturally exist in EFT [23, 24]. In contrast, pion-range corrections to OPE are expected to be perturbative according to the power counting embodied in Equations (11) and (12). From an EFT perspective, additional contact interactions are needed to make these corrections well-defined [31].

The issue I address in this section is how many, and which, contact interactions must be present for the renormalization of specific singular potentials. For simplicity, I consider central potentials; we return to the nuclear potential in section 4.

### 3.1. Non-perturbative Renormalization

Renormalization is usually discussed at the level of loops in Feynman diagrams for the Lippmann-Schwinger equation in momentum space, but it can also be formulated in terms of the Schrödinger equation in coordinate space. In the latter, which is more familiar to many, renormalization deals with distance scales on the order of those where the EFT breaks down, which I will call \( R_{\text{und}} \). The fall off of the potential at much larger distances is not important, as it affects instead the near-threshold behavior. For definiteness, let us take a central two-body potential

\[
V_L(r) = -\frac{\alpha}{2\mu r^n}
\]

in the center-of-mass frame, where \( \mu \) is the reduced mass, \( \alpha \) is a constant with mass dimension 2 \( - n \), and \( n > 0 \) is an integer. The long-range potential is characterized by an intrinsic distance scale \( r_0 = |\alpha|^{1/(n-2)} \). For \( n = 2 \) the action is scale invariant.

In the radial Schrödinger equation the potential is supplemented by the centrifugal barrier with orbital angular momentum \( l, l(l+1)/(2\mu r^2) \). The uncertainty principle implies the kinetic term scales similarly, as \( 1/(2\mu r^2) \). For \( 0 < n < 2 \) the potential is relatively small at small distances and the corresponding behavior of the wavefunction is determined by \( l \): we find ourselves in the familiar situation where one solution, labeled regular, behaves as \( r^l \) for small \( r \), while the other, labeled irregular and discarded, as \( r^{-l-1} \). In contrast, for \( n = 2 \) and \( |\alpha| \) is sufficiently large, or for \( n \geq 3 \), \( V_L(r) \) dominates at small distances. If \( \alpha < 0 \), the strong repulsion prevents any short-range approach; one can again keep just the regular solution, from which the scattering amplitude can be calculated. But when the potential is attractive, \( \alpha > 0 \), observables are sensitive to short-distance physics and renormalization is needed.

To see this in detail, consider first \( n \geq 3 \) at zero energy. For \( r \lesssim [(l+1)]^{-1/(n-2)} r_0 \), where \( V_L(r) \) dominates, the Schrödinger equation becomes an ordinary Bessel equation, and the solution is a combination of spherical Bessel functions. Both solutions are equally irregular as \( r \to 0 \) [44]. One can write the wavefunction in the \( l \) wave at small distances as

\[
\psi_l(r) \propto r^{l/4-1} \cos \left( \sqrt{\alpha} r^{3/2} / n - 1 + \phi_l \right) + \ldots ,
\]

where \( \phi_l \) is a phase that determines the relative importance of the two irregular solutions and is not fixed by the long-range potential \( V_L \). This is in strong contrast with the repulsive case, where the solutions are regular and irregular modified Bessel functions, which respectively decrease and increase exponentially as \( r \) decreases.

The case \( n = 2 \) is borderline singular, the character of the solution depending on the relative size of \( \alpha \) and a combination of \( l(l+1) \) with a number \( O(1) \) coming from the kinetic repulsion. It turns out that the critical value is \( \alpha_l = (l+1/2)^2 \). For \( l \geq l_a = \sqrt{\alpha} - 1/2 \), repulsion wins; one solution is more singular than the other and can again be discarded [45]. For \( l < l_a \) the situation is similar to \( n \geq 3 \): Equation (16) holds with \( \sqrt{\alpha} r^{3-n/2}((n/2) - 1) \to \alpha - \alpha_1 \ln(r/r_0) \), where \( r_0 \) is an arbitrary dimensionful parameter and \( \phi_l = \phi_l(r_0) \). This is an example of an anomaly [46, 47] where the scale invariance of the classical system is broken by the renormalization of the quantum system.

Equation (16) is the quantum version of the “fall to the center” in a classical singular potential [45, 48]. The phases \( \phi_l \) determine the asymptotic behavior of the wavefunction, from which the zero-energy scattering amplitude is extracted. For example, the \( S \)-wave scattering length is well-defined for a pure \( n \geq 4 \) potential [48] and given for \( n = 4 \) by

\[
a_0 = \sqrt{\alpha} \tan \phi_0 .
\]

If one imposes a particular value on \( \psi_l(R) \) at a chosen distance \( R \)—for example, that the wavefunction \( \psi_l(R) = 0 \)—the phases are fixed. However, a different value of \( R \) leads to different phases. In EFT, this arbitrariness is replaced by the values of LECs. The minimal set of contact interactions is determined by demanding renormalizability.

### 3.1.1. S Wave

Let us look into the \( S \) wave first. Choosing a sharp cutoff in coordinate space at \( R \), we replace the potential (15) by [23]

\[
V(r) = V_S(R) \theta(R-r) + V_L(r) \theta(r-R) .
\]

The depth \( V_S(R) \) of the spherical well is related to the LEC \( C_0 \) of a contact interaction,

\[
C_0 \delta(r) = \frac{C_0}{4\pi r^2} \delta(r) \to \frac{3C_0}{4\pi R^2} \theta(R-r) \equiv V_S(R) \theta(R-r) \ \text{at} \ R \to R ,
\]

A solution of the Schrödinger equation for the augmented potential requires the matching of the logarithmic derivatives of outside and regular spherical-well wavefunctions at \( r = R \),

\[
\sqrt{-2\mu R^2 V_S(R)} \cot \sqrt{-2\mu R^2 V_S(R)} = \frac{\partial}{\partial r} \ln \left( r\psi_0(r) \right) \bigg|_{r=R} .
\]
Having renormalized zero-energy scattering, an important question is whether the problem is well-defined also at finite energy $E \equiv k^2/(2\mu)$. That this is the case can be shown [23] with the WKB approximation, which applies to the region where the wavelength is small compared to the characteristic distance over which the potential varies appreciably. For distances where $|V_L(r)| \gg E$, one recovers Equation (16) for the wavefunction, up to energy-dependent corrections that are determined by Equation (16) itself. In the absence of a short-range interaction, decrease in $R$ would lead to the repeated appearance of low-energy bound states due to the unstoppable growth in attraction, a phenomenon reflected in the never-ending oscillations of the wavefunction [48]. With $V_S(R)$ preventing this collapse and ensuring the description of one low-energy datum, bound states can accrete only from negative energies, converging to finite values as $R$ decreases. How many of the bound states are within the region of validity of the EFT depends, of course, on the scales in the problem: the very low-energy spectrum will be affected by the long-distance tail of the potential while states with binding energies $\gtrsim (2\mu R^2)^{-1}$ are irrelevant for the distances of interest. For $n = 2$ and $\alpha > \alpha_0$, which is equivalent [57] to the three-boson system with short-range interactions at unitarity, the bound states form a geometric tower ("Efimov states" [58]) that signals the remaining discrete scale invariance stemming from the limit cycle in the contact interaction [59, 60]. While the existence of the tower is a consequence of the symmetry, its position is fixed by the LEC. It is remarkable that it is the proper renormalization of the EFT that underlies the "Efimov physics" intensely explored with cold atoms [61].

A particularly simple example of a singular potential is the delta function itself. In this case the external potential vanishes and the external zero-energy wavefunction is replaced by

$$\psi_0(r) \propto r^{-1} \left(1 - \frac{r}{a_0} + \ldots\right),$$

where $a_0$ determines the ratio between irregular and regular solutions and is nothing but the scattering length. The solution for Equation (20) can be written explicitly,

$$V_S(R) = -\frac{1}{2\mu R^2} \left[\left(1 + 2m\right)\frac{\pi^2}{4} + \frac{2R}{a_0} + \ldots\right],$$

where $m$ is an integer. It is apparent how a cutoff-dependent $C_0(R) \propto R$ softens the delta function. The scattering length enters in the smaller $R^3$ term. Of course, a similar result is obtained for a momentum cutoff $\Lambda \sim R^{-1}$ [27].

A subtlety arises when a regular potential with $n = 1$ in Equation (15) is present together with the delta function, as is the case for OPE. By itself, the long-range potential needs no regularization; with the delta function, a new cutoff dependence emerges in the irregular solution [24, 62]:

$$\psi_0(r) \propto r^{-1} \left\{1 - r \left[\frac{1}{a_0} + \alpha \left(\ln \frac{r}{R_*} - 1\right) + \ldots\right]\right\},$$

When $n = 2$ and $\alpha \leq \alpha_0$, or $n = 1$, we can solve this equation with $V_S(R) = 0$ if the admixture of the most singular external solution tends to zero as $R \to 0$. Thus the amplitude is renormalized properly without a contact interaction as long as we retain only the least singular wavefunction behavior, the prescription offered in reference [45].

For $n = 2$ and $\alpha > \alpha_0$, or for $n \geq 3$, because the two external solutions differ only by a phase, the contact interaction is necessary. Substituting the wavefunction (16) into Equation (20), yields a transcendental equation linking $\phi_0$ to $V_S(R)$ [23]. Two approximate solutions are

$$\sqrt{-2\mu R^2 V_S(R)} \simeq m\pi \left\{1 - \left[1 - 1 - \frac{n}{4} + \sqrt{\alpha} R^{1-n/2}\right.\right.$$

$$\left.\tan\left(\frac{2\sqrt{\alpha}}{n-2} R^{1-n/2} + \phi_0\right)\right\}^{-1},$$

when the right-hand side of Equation (20) is large, and

$$\sqrt{-2\mu R^2 V_S(R)} \simeq \frac{(1 + 2m)\pi}{2} - \frac{2}{(1 + 2m)\pi} \left[n - \sqrt{\alpha} R^{1-n/2}\right.\right.$$

$$\left.\tan\left(\frac{2\sqrt{\alpha}}{n-2} R^{1-n/2} + \phi_0\right)\right],$$

when it is small, where in both cases $m$ is an integer. Now one can keep the scattering amplitude at zero energy fixed at its experimental value by adjusting $2\mu R^2 V_S(R)$, which displays an periodic dependence on a power of the cutoff [23, 24, 49–54]. For $n = 2$, the dependence is periodic in $\ln R$, characteristic of a limit cycle and a remaining discrete scale invariance. (For discussions of limit cycles, see references [55, 56].) The $n \geq 3$ oscillation indicates a generalized limit cycle. The case $n = 4$ is displayed in Figure 3 [23].
where \( \bar{v}_0 \) and \( R_* \) are length scales that enter the zero-energy scattering amplitude. Instead of Equation (24),

\[
V_S(R) = -\frac{1}{2\mu R^2} \left[ (1 + 2m^2) \frac{\pi^2}{4} + 2R \left( \frac{1}{\bar{v}_0} + \alpha \ln \frac{R}{R_*} + \ldots \right) \right].
\]

(26)

The main difference is the appearance of the \( \ln R \) with a coefficient \( \propto \alpha \).

In both these cases, where the outside potential is not singular, it is easy to see that the amplitude at finite energy is well-defined. The energy enters both internal and external wavefunctions as \( (kr)^2 \) and can only affect the depth of the spherical well by a term of \( O(R^0) \), an effect that disappears as \( R \) decreases. The multiple branches in Equations (24) and (26) are a consequence of the fact that a spherical well can have multiple bound states. The zero-energy amplitude is essentially determined by the shallowest state, and we can choose different well depths to place any one state at the desired position. Deeper states have energies \( \propto \sqrt{r}(< R) \) and do not disappear as \( r \to 0 \). A long-range singular potential leads to the same conclusion. For the argument with a delta-shell regularization, see reference \([63]\).

In contrast, when the external potential is attractive and singular with \( n \geq 3 \),

\[
R_l(R) = l + 1 - \frac{n}{4} + \sqrt{\alpha} R^{1-n/2} \tan \left( \frac{\sqrt{\alpha}}{n/2 - 1} R^{1-n/2} + \phi_1 \right).
\]

(31)

Matching in the S wave makes \( \phi_1 \) \( R \)-independent. Since \( 2\mu R^2 V_S(R) \) is approximately cutoff independent as can be seen from either of the two approximate solutions (21) and (22), Equation (28) gives

\[
R_l(R) = 3 - \Delta_1(R),
\]

(32)

where \( \Delta_1(R < r_0) \) is finite. Comparison with Equation (31) then shows that \( \phi_1 \propto R^{1-n/2} \). Continuing to larger \( l \) we find

\[
\phi_l(R \ll r_0) = -\frac{\sqrt{\alpha}}{n/2 - 1} R^{1-n/2}.
\]

(33)

The phases are thus angular-momentum and energy independent \([63]\) in this limit, \textit{but cutoff dependent} \([48]\).

What is needed for renormalization is a single contact interaction with a minimum number of derivatives in each wave, with LECs \( C_{2l}' \). The interaction is non-local, for example for \( l = 1 \),

\[
\frac{C_{2l}'}{4\pi R^2} \left( \delta \frac{\partial}{\partial r} \right) \frac{\partial}{\partial r} \bigg|_{r_0=0} \to C_{2l}'(R) \frac{2}{\pi R^2} \left( \frac{\partial \theta(R - r) - \delta(r - R)}{r} \right) \frac{\partial}{\partial r} \bigg|_{r_0=0},
\]

(34)

where \( C_{2l}'(R) \) is determined so as to keep the phase \( \phi_l \), and thus one \( P \)-wave low-energy datum, fixed. The contact interactions are all determined by the underlying interactions, but without additional dynamical assumptions we do not know how they relate to each other. Model independence requires we keep them free.

3.1.3. Implications

Much of the above had been understood without EFT. The use of boundary conditions, for example, goes back at least to the work of Breit \([64]\). In EFT, a boundary condition corresponds to a specific regulator. At the two-body level, in the \( S \) wave we have simply traded the dependence in \( R \) by that of \( V_S(R) \). Renormalization means that, as far as observables are concerned, the regulator choice is irrelevant (within the error of the truncation); only the unobservable cutoff dependence of the
LECs depends on the regulator. What matters is that a LEC encodes one parameter. The LO EFT in coordinate space is in the spirit of atomic Quantum-Defect Theory, where the interaction of far-away electrons with an ionic core or molecule is solved for exactly and a few parameters (“defects”) account for short-range interactions [65].

The model independence of the EFT is manifest in the fact that the same two-body contact interactions that renormalize the two-body problem contribute to other processes. For example, the three-boson system was considered in reference [54], where binding energies and the particle-dimer scattering length were calculated. Convergence was observed in a range of cutoffs, with asymptotic values representing model-independent predictions. The role of $D$ and higher waves in these results was, however, not discussed.

The contact interactions can also be seen as providing a self-adjoint extension of the Hamiltonian. As stressed in reference [66], the so-called deficiency index for a singular potential is $(\infty, \infty)$, i.e., an infinite number of parameters—the phases $\phi_l$ in Equation (16) for all values of $l$—are needed to determine the self-adjoint extension uniquely. In the EFT this translates into the existence of an infinite number of contact interactions, one with the minimal number of derivatives for each wave (of course, the EFT contains also contact interactions with an arbitrary number of derivatives).

While mathematically the problem looks hopeless, on physical grounds this is clearly a red herring. As remarked in reference [20], increasing $l$ strengthens the centrifugal barrier and shrinks the distances $r \lesssim [(l+1)^{-1/(n-2)}]r_0$ where the attractive $n \geq 3$ potential takes over. The distance of closest approach at momentum $k$ can be estimated from the point where the energy is comparable to the centrifugal barrier, or $r \gtrsim [(l+1)^{1/2}k^{-1}]$. For $k \approx M_{\text{und}}$, the breakdown scale, we are only interested in distances $r \approx R_{\text{und}} \sim [(l+1)^{1/2}M_{\text{und}}^{-1}]$. We might then expect that only in waves with $l \lesssim l_\alpha$ does a singular potential need to be treated exactly and Equation (16) apply, where [32]

$$l_\alpha(l_\alpha + 1) \sim \frac{r_0}{R_{\text{und}}}.$$  

A more precise semi-analytical estimate comes from the investigation of the critical strength $\alpha$ where a Bessel series solution of the Schrödinger equation exhibits a square-root branch point characteristic of non-perturbative behavior. For $n = 3$ [32], it is described pretty well for large $l$ by the estimates above. For $n = 2$, consideration of the first two orders in the perturbative expansion suggests $l_\alpha = (\pi |\alpha| - 2)/4$ [31]. An attractive singular potential defined with a step function at $l_\alpha$ has a finite deficiency index $(l_\alpha l_\alpha)$.

The situation is different in the case of $n = 1$. The potential is larger than both centrifugal barrier and kinetic repulsion for $r \gtrsim n^2(l/r_0$, where $n(l)$ is $O(1)$ for $l = 0$ and grows as $l$ for large $l$. Balance among these terms leads to bound states of sizes $r_n \sim 2n^2r_0$ and binding energies $E_n \sim \alpha^2/(8\mu n^2)$. (Taking as an example the Coulomb interaction, where $\alpha = 2\mu \alpha_e$ in terms of the fine-structure constant $\alpha_e$, we get the proper result $B \sim \mu \alpha_e^2/(2n^2)$ if we interpret $n$ as the principal quantum number.)

These estimates are in any case affected by the long-range tail of the potential, which we are not considering in this section. But at distances $R_{\text{und}} \lesssim r \lesssim r_0$, we expect $l_\alpha \approx 1$: while the $S$ wave might be non-perturbative and perhaps require a short-range potential (26) to generate a bound state at the observed location, higher waves should be perturbative.

### 3.2. Perturbative Corrections

EFT provides a framework where we can systematically incorporate corrections to the leading interactions, which can be checked with the method developed in reference [67]. We pair subleading long-range interactions with the subleading short-range interactions needed for renormalization order by order. As stressed in reference [68], renormalization at a given order contains clues about the relative importance of higher corrections. Just as a negative power of $R$ indicates at least one missing LEC, so positive powers of $R$ point to the order before at least one new LEC should appear. If the error in an observable not used in the fit of LECs at N$^\text{LO}$ (with some integer $i$) scales as a positive power of the coordinate cutoff, say $R^j$, then we may expect that corrections appear at N$^{i+1}$LO, where $j \leq x$ is an integer (not necessarily the largest integer). This constraint comes from the demand that the regulator error should not exceed the truncation error when $R \lesssim R_{\text{und}}$. (It does not exclude the presence of a LEC at a lower order than that estimated by the cutoff dependence, corresponding to boundary conditions of the RG equation [22].) We will see examples below.

The next renormalization challenge arises from the more-singular corrections to the long-range potential. An almost automatic reflex is to simply add the correction to the LO potential, as Weinberg prescribed, and solve the Schrödinger equation. For a regular potential, adding a regular correction that is small everywhere can be done in perturbation theory, but it can also be done by solving the Schrödinger equation exactly. For a more-singular correction, however, the perturbing potential will be larger than the LO potential at sufficiently small $r$. One risks destroying the systematic character of the EFT unless one keeps $R$ relatively large. Whether this risk materializes needs to be checked explicitly. As we will see, renormalization requires distorted-wave perturbation theory around the LO solution [20, 31]. Implications for nuclear interactions are discussed in section 4.1.

#### 3.2.1. Distorted-Wave Perturbation

A pedagogical toy model that nicely illustrates the need for perturbation theory on singular corrections was presented in reference [69]. The model consists of two separable, regular potentials, one of range $m_i^{-1}$, the other of range $m_i^{-1} \ll m_i^{-1}$. Because the potentials are separable, exact answers can be found for the effective-range parameters. The potential parameters are fine-tuned so that each potential separately produces a natural scattering length, that is, $a_0 \sim m_i^{-1}$ ($a_0 \sim m_i^{-1}$) in the absence of the short-range (long-range) potential. Next, the short-range potential is expanded in powers of $k/m_i$, creating a series of singular interactions. While for $k \sim m_i$ the long-range potential is non-perturbative, the singular corrections should be treated in distorted-wave perturbation theory. Lo and behold, the results up
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counting in Pionless EFT \[V_1(R) = 0 \] in Equation (18). As discussed above, the energy dependence first affects the matching between internal and external wavefunctions at relative $O(k^2 R^2)$. The ratio of irregular and regular solutions, which determines $k \cot \delta_0(k)$ where $\delta_0(k)$ is the S-wave phase shift, starts at $O(R)$. Thus, at LO

$$k \cot \delta_0(k) = -\frac{1}{\delta_0} \left(1 + O(Ra_0k^2)\right),$$

which means that the fractional error in \(\delta_0\) is

$$\frac{\Delta \delta_0(k)}{\delta_0(k)} = O(Ra_0k^2).$$

For example, the effective range \(r_0 \sim R\). This again can be easily obtained with a momentum regulator [27]. In ChEFT, where away from the chiral limit the delta function is accompanied in the singlet S wave by the Yukawa potential, the situation is not substantially different [24]. Aside the $O(aR \ln R)$ dependence in Equation (26), the argument does not change and Equation (37) still holds with $\delta_0 \to \delta_0$. Despite the presence of pions, the error is still $\ll R$. It can be removed in first-order perturbation theory by a two-derivative contact interaction

$$\delta V_S = C_2 \left[\nabla^2 \delta(\vec{r}) + 2 \vec{V} \delta(\vec{r}) \cdot \vec{V} + 2 \delta(\vec{r}) \nabla^2 \right],$$

whose LEC $C_2(R) \propto R^2$ fixes $r_0 \sim R_{\text{und}}$. For $R \ll R_{\text{und}}$, this contact interaction is an NLO correction to the LO interaction with LEC $C_0$. This is in fact one of the elements in the power counting in Pionless EFT [35]. Note that, if we were to impose that $C_2/C_0$ scaled with $R_{\text{und}}^2$ as implied by NDA, we would obtain an effective range that scaled the same way, in contrast to what one obtains for typical short-range potentials [27]. Once again, renormalization automatically enforces a general property of short-range interactions.

But what if we solved the Schrödinger equation exactly following Weinberg’s prescription? In the simpler case without a long-range potential, it has been shown explicitly [72–74] that this can be done in a renormalized way only if $r_0 \lesssim R$, which is arbitrarily small. In other words, the two-derivative contact interaction is non-perturbatively renormalizable only if the theory satisfies a “Wigner bound” [75] $r_0 \geq 0$. In contrast, when the two-derivative contact interaction is treated in perturbation theory, at second order and higher, which contain loops involving two or more powers of $C_2$-four- and higher-derivative contact interactions appear to guarantee renormalization. When we resum the two-derivative contact interaction we generate diagrams with an arbitrary number of loops, but lack the counterterms to remove the cutoff dependence. A calculator committed to exact solutions might be tempted to eschew renormalization (and thus model independence) and live with a relatively large $R$. Still, such stubbornness in resumming what needs no resummation might be rewarded by results that are worse than those of the perturbative expansion. An example is provided by a calculation [76] of the S-wave scattering phase shifts for a harmonically trapped unitary system, where the regulator was implemented in the form of a maximum number of shells. One can see explicitly how in first-order perturbation theory the derivatives in Equation (38) give a contribution to the NLO energy which is proportional to the LO energy, apart from a shift in the LO LEC. The result of resumming the NLO interaction is not only cutoff dependent but also gives rise to a larger violation of unitarity than even NLO.

Note that one can introduce an auxiliary “dimeron” field in the EFT Lagrangian [77] whose kinetic term provides an energy-dependent correction to the potential. Exploiting the redundancy of interactions in the enlarged Lagrangian, one can remove the momentum-dependent corrections (38). Renormalization changes with an energy-dependent potential and, in particular, a resummation does not restrict $r_0$. However, unless there is evidence for $r_0 \gg R_{\text{und}}$, this is still an NLO correction and the resummation does not affect observables up to higher-order terms [27].

Resummation of subleading interactions can lead to an even more paradoxical situation. The problem is that subleading singular potentials are not in general attractive in all the same waves as OPE. If the corrections are iterated together with OPE, the cutoff behavior of the amplitude will change completely: channels that required a counterterm at LO may not require, or even tolerate, one at subleading order [24]. Take a wave where the LO potential is singular with a power $n$ and attractive, thus requiring a counterterm, but the subleading potential is repulsive (strength $a'$) with a power $n' > n$. The exact solution of the Schrödinger equation for the sum of the external potentials is now dominated at short distances by the irregular solution of the subleading potential, which grows exponentially as $r$ decreases. Matching to the short-range potential $V_S$ will force a non-vanishing irregular solution, which in turn leads to an exponentially increasing dependence of the fractional phase shift error in $R$, $\propto R^{1+n'/2} \exp[2\sqrt{-a'a''} R^{1-n'/2}/(n'/2 - 1)]$ [24]. The only solution is to remove the LO LEC at subleading order! There is hardly a way to keep the systematic expansion of the EFT.

Another toy model [78] illustrates this paradox. This time the underlying potential consists of a repulsive $r^{-3}$ component associated with a mass $m_3$ together with an attractive $r^{-3}$ from a heavier $m_S \gg m_3$, as well as less singular terms. Its exact S-wave results are compared to those of a potential consisting of the repulsive $r^{-3}$ potential plus a delta-function interaction. Parameters are chosen so that the repulsive potential is non-perturbative. Despite the fact that the phase shifts of the repulsive component are well-defined by themselves, reference [78] includes the delta function non-perturbatively, fixing it to reproduce the scattering length of the underlying potential. For
The phase shifts are in reasonably good agreement with those of the underlying potential. However, agreement deteriorates as \( R \) decreases. Disregarding conceptual differences in renormalization of attractive and repulsive singular potentials \([23, 24]\), reference \([78]\) concludes that cutoff dependence cannot be removed in general, rather than in the particular case of resuming the subleading delta function. In response, reference \([79]\) included the \( 2n \)-derivative delta functions, which account for the short-range potential, at \( N^{2(n+1)} \)LO in perturbation theory. Calculations up to \( N^3 \)LO show convergence to the exact phase shifts up to at least \( k \sim 2m_l \) without significant restriction on \( R \). (Reference \([80]\) nevertheless points to some ambiguity in the values of the NLO phase shifts, apparently implying that it is sufficient reason to abandon renormalization.)

Thus the singular nature of the potentials that we want to treat in an EFT expansion of the amplitude requires the use of perturbation theory on corrections, as implied by the power counting of section 2. This in fact ensures small changes in amplitudes after renormalization \([31]\). But then one might wonder to which extent the singular nature of the LO potential affects the order of the corrections. As we have seen, when the only singular part of the LO potential is a delta function, the first correction comes at NLO. When the outside potential is singular and attractive, the situation is different. For an LO singular attraction, one finds \([24]\) that after fixing the phase \( \phi_0 \) the S-wave phase shifts scale as

\[
\frac{\Delta \delta_0(k)}{\delta_0(k)} \propto R^{1+n/2}.
\]

(39)

This means that corrections are expected at (or before) \( N^2 \)LO for \( n = 2, 3, N^3 \)LO for \( n = 4, 5, \ldots \). It is unclear why the results reported in reference \([54]\) indicate higher sensitivity to \( R \) than given by Equation (39).

Now, the power counting for nuclear interactions in section 2 says that at \( N^2 \)LO there are corrections to the long-range potential with an \( r^{-(n+2)} \) singularity. The additional singularity can be removed in first-order perturbation theory by additional contact interactions with two derivatives. This can be shown relatively simply in a toy model where a \( \pm r^{-4} \) potential is added to an \( n = 2 \) attractive LO potential \([31]\). The analysis was carried out in momentum space with a sharp cutoff \( \Lambda \). At \( N^2 \)LO, where the \( \pm r^{-4} \) potential is considered as a first-order perturbation, two forms of additional, oscillating cutoff dependence appear: one proportional to \( \Lambda^2 \), reflecting the stronger singularity of the perturbing potential, the other proportional to \( k^2 \). In the S-wave, a two-derivative potential \([38]\) is sufficient, together with an \( N^2 \)LO shift in the \( \phi_0 \) of Equation (19), to remove the two additional divergences. This argument can presumably be continued at higher orders and repeated for \( l \geq 1 \) waves by considering interactions of type \([34]\) with two more derivatives. One tentatively concludes that NDA holds in distorted-wave perturbation once it has been corrected at LO.

### 3.2.2. Simple Perturbation

In partial waves \( l \geq l_\ell \), where the LO potential is perturbative and particles are free in zeroth approximation, corrections are included in simple perturbation theory. The first task in this case is to quantify the angular-momentum suppression for the long-range potentials so as to establish the orders they come in. The second need is to find the orders the associated contact interactions appear at.

For the \( \mu = 0 \) long-range potential, rules \((7)\) and \((8)\) indicate that a contact interaction is needed for renormalization at \( n \)th order in perturbation theory, where \( n \geq 2l + 1 \). This is consistent with the inference from the residual cutoff dependence of the non-derivative contact interaction. As we saw in section 3.1.2, \( l \)-wave scattering “lengths” \( a_l \) are induced through matching at finite \( R \). Just as for the S-wave effective range, they can be made arbitrarily small by taking \( R \to 0 \). However, the higher power of \( R, R^{2l+1} \), suggests that contact interactions in higher waves enter in perturbation theory at \( N^{2l+1} \)LO or lower, another element of Pionless EFT power counting \([35]\).

The increased singularity of subleading potentials asks for counterterms at lower orders in perturbation theory. The first-order perturbative correction due to subleading potentials involving pion loops is renormalized with LECs assigned by NDA. Making further general statements about the order contact interactions are needed is cumbersome without an explicit angular-momentum suppression factor.

If one were to solve the Schrödinger equation exactly in one of these waves, renormalization would require a LEC, which then determines the asymptotic properties of the wavefunction. The tail of the non-perturbative wavefunction can be reproduced with increasing accuracy as the order of perturbation theory increases \([23]\). Being a series in \( \alpha \), the perturbative expansion cannot reproduce the oscillations found in Equation (16), which are tied to the non-analytic dependence \( \sqrt{r} \). This is no problem because, by definition of \( l_\ell \), these oscillations take place at distances smaller than those probed by the EFT. Their effects can be “averaged out” and appear through contact interactions at subleading orders. If one wants to save all the perturbative work by sticking to a non-perturbative solution, one loses some predictive power at LO but, because it is a single LEC (in one wave), this is perhaps acceptable. Alternatively, one could simply not include the LEC if \( l \) is sufficiently high for oscillations to happen below \( R \), which might be limited in numerical calculations anyway. In this case \( R \) is in the region where perturbation theory works and the result will be relatively insensitive to \( R \). Unnecessary iteration in high waves is thus relatively harmless, other than obscuring the systematic EFT expansion.

### 4. Renormalization of Chiral EFT

By this point in the manuscript it should be clear how to proceed with ChEFT in the nuclear sector. The power counting of ChPT is based on NDA, which comes from demanding that the EFT expansion be renormalized order by order so as to ensure model independence. In the more general ChEFT we continue to insist on model independence, but now LO is non-perturbative. The results of the previous section apply to pion-exchange potentials, where the spin-isospin factors and
4.1. Partly Perturbative Pions

The simple power counting of Equations (7) and (8) does not capture factors of $l^{-1}$, just as it misses other dimensionless factors. More realistically, OPE in the radial Schrödinger equation is an expansion in $Q/M_{NN}^{(l)}$, where $M_{NN}^{(0,l)} \sim M_{NN}$ but $M_{NN}^{(l)}$ increases with $l$ depending in general also on the spin $s$. Once $M_{NN}^{(l,0)} \sim M_{QCD}$, OPE is perturbative. What do we know about $M_{NN}^{(l,0)}$ and $f_{\pi}^{(l)}$?

The bold suggestion was made in references [28, 29] that $f_{\pi}^{(l)} \approx 0$, so that pion exchange would be amenable to perturbation theory in all waves. The estimate in Equation (5) assumed NDA for the one-nucleon quantities $m_N = O(M_{QCD})$, $f_\pi = O(M_{QCD}/4\pi)$, and $g_A = O(1)$, plus neglected any dimensionless factors. Numerically, $M_{NN} \approx 290$ MeV. What if the various spin/isospin factors and other numbers floating around, each of $O(1)$, conspire to make OPE more perturbative, so that $M_{NN}$ is effectively comparable to $M_{QCD}$?

In that case, at LO ChEFT would be formally the same as Pionless EFT [35, 36], where the binding of light nuclei rests on the shoulders of non-derivative 2N and 3N contact interactions [81, 82]. But because pions are explicit, the range of validity of the EFT is enlarged—at least near the chiral limit where integrating out pions becomes a very restrictive condition. An attractive feature of this proposal is that it could potentially explain why Pionless EFT works better than expected, for example, for binding energies [35].

This proposal also neatly solves the renormalization issues of the last section. OPE is now an NLO effect of relative $O(Q/M_{NN})$, so no problems with its singularity emerge. Being perturbative, it brings NLO cutoff dependence only to $S$ waves. Because at LO the external potential vanishes, Equation (36) requires at NLO one chirally symmetric two-derivative contact interaction in each $S$ wave. Then $Q \sim m_\pi$ implies the concomitant presence of a chiral-symmetry-breaking non-derivative interaction with LEC proportional to the quark masses, $m_\pi^2 D_2$. In the background of an LO wavefunction of the type (23), OPE generates an $m_\pi^2 \ln \Lambda$ cutoff dependence which can be absorbed in $D_2$. The 2N amplitude is renormalized and in good agreement [28, 29, 83] with the Nijmegen partial-wave analysis (PWA) [84] up to $Q \sim m_\pi$.

Alas, calculations at $O(Q^2/M_{NN}^2)$ have shown [30, 85] that in the low, spin-triplet partial waves, where the OPE tensor force is attractive, the expansion fails for $Q \sim 100$ MeV. In partial waves with $l = j \geq 1$, where counterterms are needed only at a very large number of loops $L \geq 2l$, the breakdown of perturbation theory was estimated in the chiral limit to be at a critical momentum [86]

$$p_{cr} \approx \frac{\beta}{\sqrt{27(2(-1)^l + 1)}} M_{NN}.$$  \hspace{1cm} (40)

If we impose $p_{cr} \sim M_{QCD}$, we get $\beta \approx 2.5$. The radius of convergence of the perturbative series is not as large in waves with $l = j \pm 1$. In both cases the first few orders were found [86] not to be representative of the large-order convergence. For low partial waves counterterms enter already at low orders. When they were assigned arbitrary but natural values, all waves except $3S_1$, $3D_1$, $3P_0$, and perhaps $3P_1$ were found to converge up to $p_{cr} \approx M_{NN}$. An example of failure, $3P_0$, is given in Figure 4 [87], where OPE is NLO, $n$-iterated OPE N$^3$LO, leading two-pion exchange (TPE) N$^3$LO, and subleading TPE N$^4$LO. The LECs are assumed to be given by NDA instead of being introduced only at the order where they are first needed for renormalization. These signs of the breakdown of perturbative pions are consistent with an expansion in $Q/M_{NN}^{(l)}$ with $M_{NN}^{(l)} \sim f_\pi$ as indicated by NDA.

It seems inevitable that pions must be treated non-perturbatively in the low partial waves if we want to go beyond Pionless EFT at physical quark masses. Still, based on the general arguments of section 3.1.3 we expect pions to be perturbative for sufficiently high partial waves. The $n = 3$ tensor force, for which $r_0 \sim M_{NN}^{-1}$, does not vanish for spin $s = 1$. Equation (35) with $B_{\text{rand}} \sim [(l+1)]^{1/2} M_{QCD}^{-1}$ provides an estimate $\beta_{\text{cr}} \approx 2$ for the critical angular momentum in attractive triplet waves. This conclusion is made firmer by a generalization to the tensor potential of the analysis of the onset of square-root branch points in the Bessel series solution of the Schrödinger equation [32]. Given that the strength of OPE is fixed by $M_{NN}$, it translates
into an upper bound on the critical momentum $p_{cr}$, including repulsive waves. The results, listed in Table 1, are obtained in the chiral limit; a realistic pion mass could affect the smaller values by factors of $O(1)$ but is not expected to be important for the larger values. They indicate that OPE in $^3\!S_1-^3\!D_1$ and $^3\!P_0$ likely fails to converge already before $M_{NN}$. In contrast, OPE in high waves, such as $F$ and higher, converges beyond $M_{QCD}$. The gray zone is the $D$ and $P$ waves other than $^3\!P_0$. Given the low values of $p_{cr}$ on the scale set by $M_{QCD}$, one might conclude that $l_{cr}^{(1)} \approx 3$. An analysis of spin-triplet phase shifts where OPE and TPE are removed in distorted-wave perturbation [89] supports this conclusion.

A different but closely related estimate for $l_{cr}^{(1)}$ comes from the cutoff values where the first bound state crosses threshold in the absence of contact interactions. The very early work on ChEFT and much of its phenomenological improvements, which continue to this day, have used Weinberg's prescription. Unfortunately this prescription assigns to triplet waves a single non-derivative contact interaction at LO, which is incapable to determine more than one phase in a model-independent way. In particular, for a separable regulator the contact interaction contributes only to the S wave. Spurious low-energy bound states can be kept at bay at LO in the $^3\!S_1-^3\!D_1$ coupled channel [62, 90–93], but only in this channel [20, 21]. In triplet waves where OPE is repulsive there is no need for counterterms at LO [20, 94], but without them bound states repeatedly cross threshold in attractive waves and lead to wild variations in the phase shifts at energies within the realm of ChEFT [20, 21, 93, 95]. With a super-Gaussian separable regulator, bound states first emerge at, roughly, $\Lambda \sim 0.5, 1, 2, 4$, and 6 GeV in respectively $^3\!S_1-^3\!D_1$, $^3\!P_0$, $^3\!D_2$, $^3\!P_2-^3\!F_2$, and $^3\!D_3-^3\!G_3$ channels [20, 93]. Except for $^3\!D_3-^3\!G_3$, this sequence is similar to that of the attractive channels in Table 1. The lowest two channels would display shallow states when $\Lambda \sim M_{QCD}$, indicating that OPE is non-perturbative, while the higher waves are less clear—numerical experimentation suggested [20] their effects were not negligible, which can be understood from the results of reference [32].

Perhaps even more seriously, in Weinberg’s scheme more-pion exchange and other contact interactions, which should be treated perturbatively, are not. This leads to the pathologies discussed in section 3.2. Indeed, renormalization problems have been reported [21, 96–100] within Weinberg’s prescription also for higher-order potentials. These renormalization failures prevent taking a momentum-space cutoff at the breakdown scale $M_{QCD}$ or higher. A “physical cutoff” $\Lambda_{phys} \leq 1$ GeV, before $^3\!P_0$ would develop a bound state [20], is needed, and results are sensitive to the choice of regulator. No wonder then that much effort in phenomenology with chiral potentials has been dedicated to finding the “best” regulator. The limitation to small cutoffs leads to startling dependence on what should be equivalent forms of interactions in the Lagrangian, see for example reference [101].

One concludes that, while it seems well-established that to handle triplet waves beyond $M_{NN}$ pions are non-perturbative in at least $^3\!S_1-^3\!D_1$ and $^3\!P_0$, there is some uncertainty as to the partial wave up to which this is so. Part of the uncertainty comes from the presence of LECs in lowest orders of the amplitude, which require a closer comparison with data (section 4.2). What is clear is that there is an angular-momentum suppression. The perturbative expressions in reference [86] suggest

$$M_{NN}^{(1)} \sim \tilde{P} M_{NN},$$

apart from an overall suppression of $\tilde{P}$. In contrast, the analyses of reference [32] leads to $\tilde{P} \sim [ll+l+1]^{1/2}$.

Singlet channels are somewhat simpler, but not devoid of subtleties. Since the tensor force vanishes for $s = 0$, OPE has $n = 1$ and $r_0 \sim M_{NN}/m_0^2$. The general argument from section 3.1.3 indicates that only in the S wave should we expect non-perturbative effects, $|l_{cr}^{(0)}| = 1$. In higher waves, the OPE potential dominates over kinetic and centrifugal repulsion only at large distances, and there the exponential fall-off of OPE leads to further suppression.

The perturbative convergence of the $l \geq 1$ channels was studied in reference [102]. This is particularly easy because the Yukawa potential is well-defined for an arbitrary number of loops. The phase shifts are seen to converge quickly already for $^1\!P_1$, and faster as $l$ increases. The suppression factor $M_{NN}^{(l0)}$ can be estimated from the critical strength $M_{NN}^{(l0)}$ needed to generate a zero-energy bound state in the corresponding $l$ wave, shown in Table 2. There are two sequences of channels that alternate because of the factor of $-3$ in the ratio between isospin singlet and triplet: if we multiply the isosinglet entries in Table 2 the results form a single monotonous sequence. Assuming $Q \sim m_\pi$, we find that in each sequence increasing $l$ by 2 roughly suppresses OPE by one order in the expansion, starting with $^1\!P_1$ at NLO and $^1\!D_2$ at N$^2$LO. Moreover,

$$M_{NN}^{(l0)} \sim [ll+l+1] M_{NN},$$

in the isosinglet waves, with a factor 3 larger in isotriplets. If one insists on the full solution for the Yukawa potential in higher partial waves, there are no renormalization problems [20, 94], as the potential is regular. In the S wave, however, interference with the delta function leads to an unexpected violation of NDA. As first noticed in reference [18] and confirmed many times since—for example, references [62, 86],

| Partial wave | $p_{cr}$/MeV |
|--------------|--------------|
| $^3\!S_1-^3\!D_1$ | 66 |
| $^3\!P_0$ | 182 |
| $^3\!P_1$ | 365 |
| $^3\!P_2-^3\!F_2$ | 470 |
| $^3\!D_2$ | 403 |
| $^3\!D_3-^3\!G_3$ | 382 |
| $^3\!F_3$ | 2860 |
| $^3\!F_4-^3\!H_4$ | 2330 |
| $^3\!G_4$ | 1870 |
A non-derivative, chiral-symmetry-breaking contact interaction with LEC $m_\pi^2 D_{2(0)}$ if OPE is treated non-perturbatively in the $^1S_0$ channel. This LEC is $D_{2(0)} \sim C_{0(0)}/M_{NN}^2$ on the basis of NDA, and thus $N^2LO$. Renormalization of non-perturbative OPE instead requires $D_{2(0)} \sim C_{0(0)}/M_{NN}^2$ [18].

One chirally symmetric contact interaction with the minimum number of derivatives for each wave where attractive tensor OPE is iterated. The most dramatic effect is in $^3P_0$, where a contact interaction $C_{2(1)} \tilde{p} \cdot \tilde{p}$ with $C_{2(1)} \sim C_{0(1)}/M_{NN}^2$ is needed [20]. NDA would give instead $C_{2(1)} \sim C_{0(1)}/M_{QCD}^2$. The two-order enhancement comes from the running of pion exchange, and similar enhancements apply for the LECs in other attractive, singular waves where OPE is non-perturbative.

These counterterms are schematically displayed in Table 3, assuming $f_\pi^1 = 3$.

Results can be found in references [20, 62, 93, 95] for cutoff values as high as 10 GeV in super-Gaussian separable regulators. In comparison with the Nijmegen PWA, one finds:

- In the $^3S_1$-$^3D_1$ coupled channels, where Weinberg's prescription is consistent with renormalization, phase shifts come out well with one fitted LEC. Results improve for $\Lambda \gtrsim M_{QCD}$; even the mixing angle, which is somewhat overpredicted with a small $\Lambda \sim 500$ MeV, agrees with the Nijmegen PWA to within 1° up to a laboratory energy $E_{lab} \simeq 200$ MeV for $\Lambda \gtrsim 4$ GeV. When the scattering length is used to fix the LEC, the deuteron binding energy is $B_{20}^D \sim 2.0$ MeV, which is essentially the same as for lower cutoffs [106].

- For low uncoupled, attractive triplet channels ($^3P_0$, $^3D_2$) iterating pions with one fitted LEC works equally well. As an example, Figure 5 [20] shows $^3P_0$, which comes out much better than in Weinberg's prescription with $\Lambda \sim 500$ MeV (compare this also with Figure 4 where pions are treated perturbatively). The vanishing of the amplitude beyond $E_{lab} \simeq 200$ MeV can be described, because attraction from OPE is compensated by the contact interaction. Again, agreement improves with increasing cutoff.

- For low coupled triplet channels ($^3P_2$, $^3F_3$, $^3D_3$, $^3G_3$)—see Figure 5 [20] again for an example—iterated pions with the associated LEC do not improve significantly over Weinberg's prescription with $\Lambda \sim 500$ MeV. While $^3D_3$ is much better, changing from repulsion to attraction, $^3P_2$ goes from underprediction to considerable overprediction.

- In triplet channels without free parameters ($^3P_1$, $^3F_3$, $^3F_3$, $^3H_4$, $^3G_4$) iterated pions tend to work well, whether they are expected to be perturbative or not. In these channels results are the same as in Weinberg's prescription; there is not much change as $\Lambda \gtrsim M_{QCD}$.

- In $I \geq 1$ singlet channels ($^1P_1$, $^1D_2$, $^1F_3$, $^1G_4$), iterated pions undershoot data except in $^1F_3$. Again results essentially agree with Weinberg's prescription at small $\Lambda \sim 500$ MeV.

- In $^1S_0$, the phase shifts resemble those of Pionless EFT, where after the fast rise due to the existence of a virtual state, they remain essentially flat as $E_{lab}$ increases. Weinberg's
TABLE 3 | Schematic momentum dependence of the lowest-order contact interactions in the 2N system up to D waves, according to references [20, 31, 62, 68, 104, 105].

|       | \(^1S_0\) | \(^3S_1\) | \(\epsilon_1\) | \(^3P_0, ^3P_2\) | \(^1P_1, ^3P_1\) | \(\epsilon_2\) | \(^3D_2, ^3D_3\) |
|-------|----------|----------|----------------|----------------|----------------|----------------|----------------|
| LO    | 1        | \(p^2 + p^2\) | \(p^2\) | \(p^2(p^2 + p^2)\) | \(p^2p^2\) | \(p^2(p^2 + p^2)\) |
| NLO   | \(p^4 + p^4\) | \(p^2\) | \(p^2\) | \(p^2p^2\) | \(p^2(p^2 + p^2)\) | \(p^2(p^2 + p^2)\) |
| N^2LO | \(p^2 + p^2\) | \(p^2\) | \(p^2\) | \(p^2p^2\) | \(p^2(p^2 + p^2)\) | \(p^2(p^2 + p^2)\) |

FIGURE 5 | Two-nucleon \(^3P_0\) and \(^3P_2\)-\(^3F_2\) phase shifts (\(\delta\)) and mixing angle (\(\epsilon_2\)) as functions of the laboratory energy \(T_L\). The LO results (solid lines) at a cutoff \(\Lambda = 3.94\) GeV are compared with the Nijmegen PWA [84] (dashed lines). Reprinted figure with permission from reference [20]. Copyright (2005) by the American Physical Society.

prescription applies, and renormalization allows us to increase the cutoff beyond \(M_{QCD}\), but agreement with the Nijmegen PWA deteriorates as we do so.

Thus, a renormalized approach where the regulator is unimportant gives a qualitative guide to 2N data at LO, which is slightly better than Weinberg’s prescription with specific regulators and small momentum-cutoff parameters. It has been shown recently [107] that, with a non-separable regulator, a specific combination of the four possible spin-isospin non-derivative contact interactions that yields only one \(^3S_1\)-\(^3D_1\) bound state simultaneously prevents bound states in other channels. While this is not true for an arbitrary regulator, it does allow to extend LO results with Weinberg’s prescription to higher cutoff values, in general improving agreement with the Nijmegen PWA. However, results are not clearly better than the renormalized approach, particularly in the \(^3P_0\) channel which lacks the repulsion to produce the amplitude zero.

In addition to simple perturbative corrections in higher partial waves, one needs to account in subleading orders for potential corrections via distorted-wave perturbation theory in the lowest partial waves. The residual \(\Lambda^{-1}\) dependence of the LO amplitude means that at NLO—relative \(\mathcal{O}(Q/M_{QCD})\)—there is also:

- A two-derivative, chirally symmetric contact interaction with LEC \(C_{2(0)}\) in the \(^1S_0\) channel. In order to render cutoff effects on the effective range no larger than \(N^2\)LO, \(C_{2(0)} \sim C_{0(2)}/(M_{\text{NN}}M_{QCD})\) [68]. NDA gives instead \(C_{2(0)} \sim C_{0(2)}/M_{QCD}^4\), or \(N^2\)LO (confusingly denoted NLO in the nuclear community), which produces a short-range contribution to the effective range smaller than pion’s by two powers of the expansion parameter. Yet, only about half of the \(^1S_0\) effective range comes from OPE.

The cutoff dependence in other channels is milder, in agreement with the discussion of section 3.2. The NLO interaction is shown in the second line of Table 3. At NLO in the amplitude, the NLO interaction should be included in first order in the distorted-wave expansion.

At higher orders, corrections to the long-range potential enter according to the power counting of section 2. Barring unforeseen renormalization issues, at \(\mathcal{O}(Q^4/M_{QCD}^4)\) we need
to include LECs with up to µ derivatives more than the LECs appearing at LO [31], except in the $^1S_0$ channel where the Yukawa/delta-function interference takes place. The momentum structures of the LECs up to $N^3LO$ are shown in Table 3, again under the assumption $J^{(1)} = 3$. They are:

- In each triplet channel where attractive OPE is iterated at LO ($^3S_1$-$^3D_1$, $^5P_0$, etc.), a contact interaction with two derivatives more than the contact interaction at LO [104, 105]. While for $^3S_1$-$^3D_1$ this coincides with NDA, for other channels NDA would say these contact interactions only appear at $N^3LO$ or higher.
- Contact interactions with two derivatives [68] for singlet ($^1P_1$) and triplet $P$ waves where OPE is repulsive ($^3P_1$). This is the NDA scaling.
- Four- and six-derivative contact interactions in the $^1S_0$ channel at $N^2LO$ and $N^3LO$, respectively [68]. Again, NDA would have these contact interactions at $N^4LO$ or higher.

Up to $N^3LO$ in the amplitude, their contributions are included in first order in the distorted-wave expansion. Meanwhile, the NLO interaction must be included in second and third orders, either by itself or with one $N^2LO$ interaction.

The phase shifts have been calculated up to $N^3LO$ along these lines in references [68, 104, 105], together with Deltaless TPE:

- In the $^3S_1$-$^3D_1$ coupled channels, where LO already yielded very good results at LO, results improve only marginally at $N^2LO$.
- In $^3P_0$, which was also relatively well-described at LO, results improve quite a bit around the maximum phase shift at $N^2LO$.

Not much improvement, if any, is seen at $N^3LO$. Results from reference [104] are shown in Figure 6, to be compared with LO in Figure 5. Other uncoupled, attractive triplet channels ($^3D_2$ etc.) were not calculated.

- The coupled $^3P_2$-$^3F_2$ wave with OPE iterated at LO shows no real improvement at $N^2LO$, and only mildly better agreement with the Nijmegen PWA at $N^3LO$. No results are available for higher coupled triplet channels ($^3D_3$-$^3G_3$ etc.).
- In $^1P_1$, which works well at LO with no free parameter, results deteriorate at $N^2LO$. Higher repulsive triplet channels ($^3F_3$ etc.) were not considered.
- In $^3P_1$, agreement with the Nijmegen PWA improves at $N^2LO$, although results are very sensitive to the pion-nucleon parameters that enter the $\mu = 3$ TPE. Higher singlet partial waves were not studied.
- The $^1S_0$ phase shift improves considerably at NLO but is still not very close to the Nijmegen PWA. $N^2LO$ improves further, but the zero of the amplitude is still poorly described.

Overall, there is some improvement at $N^2LO$ but not much at $N^3LO$. This is perhaps an indication that a better description of the pion-nucleon subamplitude with an explicit Delta isobar is needed.

Note that subleading corrections have also been calculated in references [108, 109] with a slightly different accounting of higher orders. For example, TPE is taken to start three orders higher than OPE, which is contrary to the power counting of section 2 and difficult to conciliate with the power counting used in ChPT. Still, results are generically not much different from those described above. A third power-counting variant has been proposed [32] with similar features. It has not been tested in detail, perhaps because no clear prescription is given for handling the LO cutoff dependence in a channel like $^3P_0$ where a counterterm is assigned relative $O(Q^{1/2}/M_{QCD}^{1/2})$. Reference [67] discusses these alternatives.

The main phenomenological shortcomings of the renormalized approach are $^3P_1$, $^3P_2$ and singlet partial waves. For most of these channels, subsequent work indicates OPE might be perturbative. Equation (42) shows that OPE should be included in $^1P_1$ at NLO, in $^1D_2$ at $N^2LO$, and so on. On the basis of NDA, contact interactions with the minimal number of derivatives are expected at respectively $N^2LO$, $N^4LO$, and so on. Under the assumption that the angular-momentum suppression of TPE is the same as OPE, reference [87] provided evidence that the perturbative expansion converges for singlet waves up to $k \approx 300$ MeV and $N^4LO$ without explicit Delta isobars. Reference [87] goes further by showing that under NDA for the LECs also triplet waves converge in the same range, except for $^3P_0$ and possibly $^3D_3$. For illustration, results for the $^3P_2$-$^3F_2$ coupled channels are shown in Figure 7 [87], which should be compared to Figure 5 where OPE was treated non-perturbatively at LO. The maximum momentum $k \approx 300$ MeV seems tied to the absence of an explicit Delta isobar [87] but no similar calculation is available in Deltaful ChEFT. Earlier studies [110–112], which indicated that pions are perturbative in high waves, sometimes included Deltas but did not take into account the IR enhancement in iterated pion exchange. Clearly a more comprehensive study of higher orders with Deltas is needed to confront this renormalized approach with phenomenology.
The situation is particularly unsatisfactory in the \(^1S_0\) channel, where LO—same as in Weinberg’s prescription at fixed pion mass—is far off, just as in Pionless EFT [35]. In particular, the Nijmegen PWA displays a zero at a relatively low momentum \(k_0 \sim 340\, \text{MeV}\), which is absent at LO. It is possible that the inclusion of an explicit Delta isobar (separated in mass from the nucleon by \(\sim 300\, \text{MeV}\)) improves the convergence in this region, as a large part of the central potential moves from \(N^2\)LO to \(N^3\)LO. However, the expansion will in any case converge at best very slowly for \(k > k_0\), as all subleading orders have to conspire to cancel against LO. Since numerically \(k_0 \sim M_{NN}\), only for a fully perturbative-pion approach is this of no concern. Note that also \(^3S_1\) and \(^3P_0\) have amplitude zeros at relatively low energies, but in both cases they arise at LO from the combination of non-perturbative OPE and contact interactions need for renormalization.

The \(^1S_0\) channel is special also for the presence of an unnaturally shallow virtual state that requires a fine-tuning of the short-range interaction. It is the interference between the non-derivative contact interaction and the Yukawa potential that causes a violation of NDA in this channel. It also leads to the piling up of higher-order counterterms seen in Table 3. Given the uniqueness of this channel, it is perhaps not surprising that power counting might require refinement. In reference [113] it was shown that short-range interactions show strong energy dependence. To ameliorate the expansion in \(^1S_0\), it was suggested in references [62, 114] that the chirally symmetric two-derivative interaction with LEC \(C_{2(0)}\) should be promoted from NLO to LO, following an earlier suggestions for Pionless EFT [115] and ChEFT with purely perturbative pions [116]. To avoid the Wigner bound, this is done through a dibaryon field [77] whose kinetic term is taken to be LO together with its residual mass. This promotion induces promotions at higher orders of the contact interactions with more derivatives. Results of course improve at LO and further at NLO, but not at \(N^2\)LO, in particular near \(k_0\).

In reference [117] it was then proposed—similarly to an earlier attempt [118]—that the zero be included at LO by a combination of dibaryon field and contact interaction (or alternatively two dibaryon fields, the kinetic term of one of which is higher order). Again this induces the promotion of contact interactions with more derivatives at higher orders. Phase shifts come out great at LO and essentially on the nose at NLO, even beyond \(k_0\) (see Figure 8). Unfortunately these reorganizations of the expansion produce energy-dependent potentials at LO, which complicate few-body calculations.

A further proposed reorganization of ChEFT arises from treating selected relativistic corrections, which are small for the momenta of interest, as LO—see, for example, reference [120]. A modified nucleon propagator ensures less dependence on the regulator, but a \(^3P_0\) LEC still has to be promoted compared to NDA, as in the purely non-relativistic context [20]. By resumming higher-order terms into LO whether they are relativistic corrections or not, one can soften the large-momentum behavior of loops and alter the cutoff dependence. This is no different than picking a regulator, which effectively includes an infinite number of higher-derivative interactions. Results then depend on the corresponding cutoff parameter \(\Lambda\). Renormalization exchanges this dependence for the minimal number of parameters allowed without dynamical assumptions.
Achieving cutoff independence with a resummation of a selected interaction merely replaces $\Lambda$ by the mass parameter characterizing this interaction, call it $M$. If $M \ll M_{\text{QCD}}$ is inferred from data, this resummation is justified because the interaction is not of higher order. However, when resumming relativistic corrections $M \gtrsim m_N$: it corresponds to one fixed cutoff value and convergence cannot be used to demote interactions that are needed for renormalization without resummation. As long as no LECs are promoted or demoted, a resummation of higher-order corrections is safe. There is growing interest in the development of a covariant version of ChEFT, which could perhaps be used as input to relativistic formulations of nuclear physics [121, 122].

4.3. More Nucleons

There is not much known about renormalized ChEFT beyond 2N. The power counting of section 2 shows that the 3N force is expected to start at NLO from two-pion exchange when Delta isobars are included explicitly, and at N^2LO when they are not. The crucial issue is whether shorter-range interactions are enhanced as in the 2N system. Such an enhancement does take place in Pionless EFT [81] and it has been suggested for ChEFT on phenomenological grounds in reference [123].

In calculations for more than two nucleons in the renormalized approach, one needs to truncate the LO 2N potential for $l \lesssim l_{ct}^{(0)}$, which is reminiscent of the truncation in total 2N angular momentum typically invoked in solutions of the Faddeev and Faddeev-Yakubovski equations for 3N and 4N systems with phenomenological potentials. As we have seen the optimal values for $l_{ct}^{(0)}$ are uncertain and the $l$ dependence of $M_{NN}^{(l)}$ is not fully determined. Of course, as in the 2N system, subleading orders should be treated in distorted-wave perturbation theory.

Existing calculations are limited to the 3N system and took $l_{ct}^{(0,1)} = 3$. At LO [20, 93] and, without explicit Deltas, also at NLO [93], observables converge as the cutoff increases to at least 10 GeV without 3N forces (see Figure 9) [93]. The triton binding energy is $B_3^{\text{LO}} \approx 4$ MeV and $B_3^{\text{NLO}} \approx 6$ MeV, quite different from results for a low cutoff in Weinberg’s prescription, $\approx 11$ MeV ($\approx 6.5$ MeV) at LO (N^2LO) [106]. Results were shown not to change significantly when waves beyond $l_{ct} = 3$ were included. Conversely, if it turns out that $l_{ct}^{(0,1)} < 3$, results might change quantitatively, but qualitative statements should stand. In particular, one concludes there is no renormalization justification in ChEFT to take the non-derivative 3N contact interaction as LO. Most likely the same conclusion holds for higher-body forces, but no calculations have been carried out.

The tendency for underbinding at LO seen in the deuteron and triton seems to persist for symmetric nuclear matter. In a cutoff-converged Brueckner pair approximation [124], nuclear matter was found to saturate, but with significant underbinding. This is in contrast to Weinberg’s prescription, where Deltaless [125] or Deltaful [126] potentials of $O(1)$ and $O(Q^2/M_{\text{QCD}}^2)$ do not yield saturation within the EFT domain. Yet higher potentials do lead to saturation with this prescription [125–127]. Although usually presented as a success, the emperor has no clothes: it means that, if nuclear matter is within the regime of ChEFT, interactions that are formally of higher order according to NDA must actually be LO to balance against other LO interactions. Presumably it is the extra repulsion from $^3P_0$ in a renormalized approach that saturates nuclear matter. It is not clear how saturation in Chiral EFT would relate, if it can be related at all, to the proposal of reference [128] where saturation arises from the 3N parameter that appears at LO in Pionless EFT. What is clear is, more EFT calculations beyond the 2N system are sorely needed!

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**Figure 9** | Triton binding energy $E_{3\text{H}}$ and doublet neutron-deuteron scattering length $2a_{nd}$ as functions of the cutoff $\Lambda$. Results at LO (solid lines) and NLO (dashed and dotted lines) for various $2N$ fitting procedures are compared with experiment (horizontal red lines). Reprinted figure with permission from reference [93]. Copyright (2019) by the American Physical Society.
5. CONCLUSION

The longstanding problem of renormalization of chiral nuclear forces has been solved at the 2N and 3N levels. Perhaps not surprising in hindsight, this solution is a middle ground between Weinberg’s original prescription and Kaplan, Savage, and Wise’s suggestion of fully perturbative pions. One-pion exchange is iterated in lower waves together with the necessary contact interactions, while all corrections are included in distorted-wave perturbation theory.

That is not to say that the best solution has been found. Issues remain regarding exactly how strong the angular-momentum suppression is and where the non-perturbative/perturbative boundary lies. Whether the ordering of few-body forces holds remains an open question, although the first step in grounding it on a renormalized approach has been made [42]. Fortunately, there is still plenty to learn.

AUTHOR CONTRIBUTIONS

The author confirms being the sole contributor of this work and has approved it for publication.

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Conflict of Interest: The author declares that the research was conducted in the absence of any commercial or financial relationships that could be construed as a potential conflict of interest.

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