

Covariant spectator theory of np scattering: Effective range expansions and relativistic deuteron wave functions

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We present the effective range expansions for the 1S_0 and 3S_1 scattering phase shifts, and the relativistic deuteron wave functions that accompany our recent high precision fits (with $\chi^2/N_{\text{data}} \simeq 1$) to the 2007 world np data below 350 MeV. The wave functions are expanded in a series of analytical functions (with the correct asymptotic behavior at both large and small arguments) that can be Fourier-transformed from momentum to coordinate space and are convenient to use in any application. A FORTRAN subroutine to compute these wave functions can be obtained from the authors.

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I. INTRODUCTION

This article presents the effective range expansions and relativistic deuteron wave functions that accompany the recent high-precision fits to the 2007 world np data (containing 3788 data) below 350 MeV [1]. These fits, obtained using the covariant spectator theory (CST) [2,3], resulted in two new one-boson-exchange models of the np interaction, both of which also reproduce the experimental triton binding energy without introducing additional irreducible three-nucleon forces [4]. One model (WJC-1) has 27 parameters and fits with a $\chi^2/N_{\text{data}} = 1.06$. The other model (WJC-2) has only 15 parameters and fits with a $\chi^2/N_{\text{data}} = 1.12$.

The main body of the article is divided into three sections. In Sec. II the sixth-order effective range expansions for the 1S_0 and 3S_1 phase shifts are given; it is shown that these expansions provide an excellent description of the phase shifts up to 50 MeV laboratory kinetic energy. Then, in Sec. III, the relativistic deuteron wave functions that automatically emerge from these fits are discussed. The deuteron binding energy was constrained during the fits, so both models have the correct binding energy.

Section III begins with a review of the definitions of the wave functions and a description of how they are related to the relativistic dnp vertex functions defined in the CST. It is shown how the CST wave functions are decomposed into four independent amplitudes and how these can be identified with the familiar u (S state) and w (D state) components, plus two P -state components of purely relativistic origin [5,6]. Because these P states are associated with the virtual antinucleon degrees of freedom they have positive parity. One P -wave component has a symmetric total spin triplet structure (denoted v_t) and the other an antisymmetric total spin singlet structure (denoted v_s). The antisymmetric v_s wave function would be zero if both nu-

cleons had the same energy, but because one is on-shell and the other off-shell, the relative energy is $E_p - \frac{1}{2}M_d$ (where E_p is the total relativistic energy of a nucleon with three-momentum \mathbf{p} and M_d is the deuteron mass), allowing for the existence of this odd state. The normalization of the wave functions is discussed [7]. It turns out that the probability of the WJC-1 P -state components is only about 0.3% while the WJC-2 P -state components are much smaller (with a combined probability of less than 0.02%). These components are retained because they are *required by the manifest covariance of the CST* and are comparable to the larger wave functions in the momentum region of 400 to 600 MeV. Even though they are not present in some other relativistic approaches [8], they have been found previously [9] to make surprisingly important contributions to the deuteron form factors. The wave functions are expanded in a series of analytic functions that can be conveniently Fourier-transformed to coordinate space, giving analytical wave functions in both momentum and coordinate space convenient to use in any application. The section concludes with a discussion of the asymptotic D/S ratio. Calculations of the magnetic and quadrupole moments require the evaluation of interaction currents, which is beyond the scope of this article and is deferred to future work.

Our results are summarized in the conclusion, Sec. IV. There are also three appendices. Appendix A gives details of the construction and evaluation of the kernels omitted from Ref. I [1]. Appendix B discusses the method used to evaluate the angular integrals, complicated by the presence of a cusp in the region of integration. Finally, Appendix C gives more details about the definitions, extraction, and normalization of deuteron wave functions.

II. EFFECTIVE RANGE EXPANSIONS

The four-term effective range expansion for the 1S_0 and 3S_1 phases that we use is

$$k \cot \delta = -\frac{1}{a} + r_0 k^2 \left\{ \frac{1}{2} + (r_0 k)^2 [P + Q(r_0 k)^2] \right\}, \quad (2.1)$$

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TABLE I. The effective range parameters defined in Eq. (2.1) and the fractional errors Δ at 1 and 5 MeV for each model. The fractional error is defined to be $\Delta = (\cot \delta - \cot \delta_{\text{eff}}) / \cot \delta$, where δ_{eff} is the phase predicted by Eq. (2.1) and δ the phase emerging from solutions of the scattering equation.

Parameter	1S_0		3S_1	
	WJC-1	WJC-2	WJC-1	WJC-2
a (fm)	-23.7494	-23.7496	5.4295	5.4342
r_0 (fm)	2.6261	2.6623	1.7601	1.7666
P	0.0007	-0.0005	0.0172	0.0206
Q	0.0033	0.0032	0.0060	0.0058
Δ (1 MeV)	0.0044	0.0038	-0.0011	-0.0009
Δ (5 MeV)	0.0033	0.0030	-0.0012	-0.0012

where a is the scattering length, r_0 the effective range, and P and Q the dimensionless third- and fourth-order parameters. Here

$$k^2 = \frac{1}{2}mE_{\text{lab}}, \quad (2.2)$$

where k is the magnitude of the center-of-mass momentum of the nucleons and E_{lab} the laboratory kinetic energy and we used an average nucleon mass $m = 938.9$ MeV and $\hbar c = 197.3288$ MeV fm = 1. Recall that Eq. (2.2) holds for both relativistic and nonrelativistic kinematics, allowing us to compare our results directly with nonrelativistic calculations.

To fix these parameters we evaluated the phase shifts at four energies: 0.0001, 10, 25, and 50 MeV. The first energy of 0.0001 MeV is so low that it effectively fixes the scattering length (and hence the total cross section at the np threshold). The use of these four energies resulted in a small error between the phases calculated at 1 and 5 MeV, and those predicted using (2.1). The effective range parameters and the errors at 1 and 5 MeV are summarized in Table I.

The effective range expansion is surprisingly accurate over a wide energy range. As shown in Fig. 1, the expansions are qualitatively accurate up to E_{lab} of about 150 MeV (the small errors shown in Table I are completely invisible on the figure). In the fitting we used the expansions instead of the phase shifts to calculate the S -wave contributions to the observables of all energies below 50 MeV. In doing the final minimizations we constrained both the deuteron binding energy and the 1S_0 phase shift at 0.0001 MeV (which we fixed at $\delta = 1.4937^\circ$). This was a very effective way to maintain an accurate fit to the low-energy cross-section data.

III. DEUTERON WAVE FUNCTIONS

The CST deuteron wave functions have been developed and defined in a number of references. The earliest reference is the work of Buck and Gross [5]; more recent references include Ref. [10] (referred to as Ref. II) and Ref. [11], where many practical details for how to use the wave functions in practical calculations are developed. While all of these references use the same basic definitions, details of how to interpret and use the wave functions have improved with time. In this section we first present a brief review of the necessary definitions (relating them to the earlier references) and then present the numerical results for the wave functions of models WJC-1 and WJC-2.

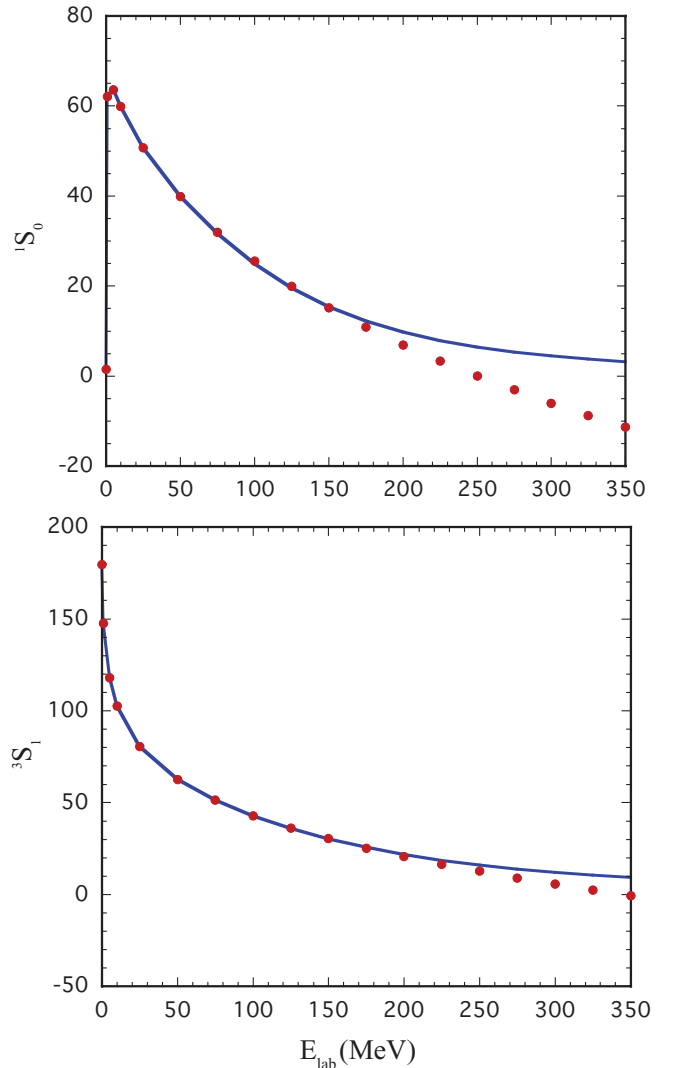


FIG. 1. (Color online) The effective range expansions for the S waves (smooth line) compared to calculated phase shifts (solid dots). The phases are in degrees.

A. Definitions of the vertex function and the covariant wave function

The equation satisfied by the np -scattering amplitude is [in the notation of Eq. (2.3) of Ref. I]

$$M_{12}(p, p'; P) = \bar{V}_{12}(p, p'; P) - \int_{k_1} \bar{V}_{12}(p, k; P) G_2(k_2) M_{12}(k, p'; P), \quad (3.1)$$

where P is the conserved total four-momentum, and $p, p',$ and k are relative four-momenta related to the momenta of particles 1 and 2 by $p_1 = \frac{1}{2}P + p$, $p_2 = \frac{1}{2}P - p$. $E_{k_1} = \sqrt{m^2 + \mathbf{k}_1^2}$ is the energy of the on-shell particle 1, and the covariant integral is

$$\int_k \equiv \int \frac{d^3k}{(2\pi)^3} \frac{m}{E_k}. \quad (3.2)$$

Note that these covariant operators can be written either in terms of the independent momenta $\{P, p, p'\}$ or $\{P, p_1, p'\}$. The scattering amplitude,

$$\begin{aligned} M_{12}(p, p'; P) &\equiv M_{\lambda_1 \lambda'_1, \beta \beta'}(p, p'; P) \\ &= \bar{u}_\alpha(\mathbf{p}_1, \lambda_1) \mathcal{M}_{\alpha \alpha', \beta \beta'}(p, p'; P) u_{\alpha'}(\mathbf{p}'_1, \lambda'_1) \end{aligned} \quad (3.3)$$

is the matrix element of the Feynman scattering amplitude \mathcal{M} between positive-energy Dirac spinors of particle 1. The definition of the nucleon spinors $u(\mathbf{p}_1, \lambda_1)$ (with λ_1 the helicity of the nucleon) is given in Eq. (A3); the on-shell spinor has four-momentum p_1 with three-momentum component \mathbf{p}_1 , and $p_1^2 = m^2$. The propagator for the off-shell particle 2 is

$$G_2(k_2) \equiv G_{\beta \beta'}(k_2) = \frac{(m + \not{k}_2)_{\beta \beta'}}{m^2 - k_2^2 - i\epsilon} h^2(k_2) \quad (3.4)$$

with $k_2 = P - k_1$, $k_1^2 = m^2$, and $h(k_2)$ [denoted $H(k_2)$ in Ref. I] the form factor of the off-shell nucleon (related to its self-energy), normalized to unity when $k_2^2 = m^2$

$$h(k_2) = \left[\frac{(\Lambda_N^2 - m^2)^2}{(\Lambda_N^2 - m^2)^2 + (m^2 - k_2^2)^2} \right]^2. \quad (3.5)$$

See Appendix C for further discussion of the nucleon form factor h and the role it plays in the deuteron wave functions.

Just below threshold, the np scattering amplitude has a pole at the deuteron mass, $P^2 = M_d^2$. Near the pole this amplitude, in the CST, can be written

$$\begin{aligned} M_{\lambda_1 \lambda'_1, \beta \beta'}(p, p'; P) &= R_{\lambda_1 \lambda'_1, \beta \beta'}(p, p'; P) \\ &\quad - \sum_{\lambda_d} \frac{\mathcal{G}_{\lambda_1 \beta}(p, P, \lambda_d) \bar{\mathcal{G}}_{\lambda'_1 \beta'}(p', P, \lambda_d)}{M_d^2 - P^2 - i\epsilon}, \end{aligned} \quad (3.6)$$

where λ_1, λ'_1 are the helicities of the outgoing and incoming on-shell particle 1, β, β' are the Dirac indices of the off-shell particle 2, λ_d is the helicity of the deuteron, \mathcal{G} is the *contracted* dnp vertex functions describing the coupling of the deuteron to the neutron and proton, and R is a remainder function finite at the pole. The equation for \mathcal{G} can be found by substituting (3.6) into the scattering Eq. (3.1) and demanding that it hold at the pole (see Refs. [3,7] and II), giving

$$\begin{aligned} \mathcal{G}_{\lambda_1 \beta}(p, P, \lambda_d) \\ = - \int_{k_1} \bar{V}_{\lambda_1 \lambda'_1, \beta \alpha}(p, k; P) G_{\alpha \beta'}(k_2) \bar{\mathcal{G}}_{\lambda'_1 \beta'}(k, P, \lambda_d). \end{aligned} \quad (3.7)$$

The contracted vertex function is a Lorentz scalar product of a Dirac matrix element and the covariant polarization vector of the deuteron, $\xi_\mu(\lambda_d)$:

$$\mathcal{G}_{\lambda_1 \beta}(p, P, \lambda_d) \equiv \mathcal{O}_{\lambda_1 \beta}^\mu(p, P) \xi_\mu(\lambda_d), \quad (3.8)$$

where \mathcal{O} is the “uncontracted” dnp vertex function for particle 2 off-shell [cf. Eq. (2.27) of Ref. II]

$$\mathcal{O}_{\lambda_1 \beta}^\mu(p, P) \equiv [\Gamma^\mu(p, P) C]_{\beta \beta'} \bar{u}_{\beta'}^T(\mathbf{p}_1, \lambda_1), \quad (3.9)$$

with C the Dirac charge conjugation matrix. To simplify the language, \mathcal{G} will always be called the contracted vertex

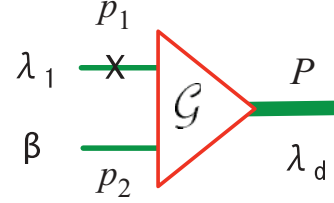


FIG. 2. (Color online) Diagrammatic representation of the vertex function (3.8). The on-shell particle is labeled by the \times , so $p_1^2 = m^2$, and the relative four-momentum is $p = \frac{1}{2}(p_1 - p_2)$. The deuteron is also on-shell, with $P^2 = M_d^2$.

function, and the functions \mathcal{O} (or sometimes Γ) will be referred to simply as the “vertex function.” The contracted vertex function is illustrated diagrammatically in Fig. 2.

The relativistic deuteron wave functions were defined previously in Eq. (2.32) of Ref. II and Eq. (2.29) of Ref. [11]:

$$\psi_{\lambda_1 \beta, \lambda_d}(p, P) = N_d G_{\beta \beta'}(p_2) \mathcal{G}_{\lambda_1 \beta'}(p, P, \lambda_d). \quad (3.10)$$

Here N_d is a normalization constant, chosen to be

$$N_d = \frac{1}{\sqrt{(2\pi)^3 2M_d}}. \quad (3.11)$$

As discussed in Ref. [11], it is the vertex function that enters directly into any Feynman diagram involving an incoming or outgoing deuteron state, so if a Feynman amplitude is expressed in terms of the wave function instead of the vertex function, one must be careful to divide by the normalization constant (3.11). The reason for this choice of normalization constant will be discussed further below.

The normalization of the wave function (3.10) is discussed in detail in Refs. [3,7], II, and Appendix C. The exact result, in the notation of Eq. (3.7), is

$$\begin{aligned} \delta_{\lambda_d \lambda'_d} &= \frac{1}{2M_d N_d^2} \left\{ \int_{p_1} \psi_{\lambda_1 \beta, \lambda_d}^\dagger(p, P) \gamma_{\beta \beta'}^0 \psi_{\lambda_1 \beta', \lambda'_d}(p, P) \right. \\ &\quad - \int_{p_1} \int_{p'_1} \psi_{\lambda_1 \beta, \lambda_d}^\dagger(p, P) \Delta \bar{V}_{\lambda_1 \lambda'_1, \beta \beta'}(p, p'; P) \\ &\quad \left. \times \psi_{\lambda_1 \beta, \lambda'_d}(p', P) \right\}, \end{aligned} \quad (3.12)$$

where summation over repeated indices is implied and

$$\Delta \bar{V}_{\lambda_1 \lambda'_1, \beta \beta'}(p, p'; P) = \frac{\partial}{\partial M_d} \bar{V}_{\lambda_1 \lambda'_1, \beta \beta'}(p, p'; P), \quad (3.13)$$

with the partial derivative holding p_1 and p'_1 constant. This normalization integral will be simplified and related to the usual nonrelativistic normalization in Appendix C 3 below.

While it is often convenient to use the fully covariant expression (3.10), the physical content of the wave function can be displayed if we expand the off-shell propagator into positive- and negative-energy pieces. (Note that an alternative expansion given in Eq. (A21) of Ref. [11] is sometimes more convenient for applications.) Working in the rest frame, where we choose $\mathbf{p}_1 = -\mathbf{p}_2 = \mathbf{p}$, the propagator can be

decomposed

$$\begin{aligned} G_{\alpha\beta}(p_2) &= \frac{m}{E_p} \sum_{\lambda_2} \left\{ \frac{u_{\alpha}(-\mathbf{p}, \lambda_2) \bar{u}_{\beta}(-\mathbf{p}, \lambda_2)}{2E_p - M_d} - \frac{v_{\alpha}(\mathbf{p}, \lambda_2) \bar{v}_{\beta}(\mathbf{p}, \lambda_2)}{M_d} \right\}, \\ &= \frac{m}{E_p} \sum_{\lambda_2, \rho_2} G^{\rho_2}(p) u_{\alpha}^{\rho_2}(-\mathbf{p}, \lambda_2) \bar{u}_{\beta}^{\rho_2}(-\mathbf{p}, \lambda_2), \end{aligned} \quad (3.14)$$

where ρ_2 is the ρ spin of the off-shell particle 2 [ρ spin is a shorthand notation for positive-energy ($\rho = +$) or negative-energy ($\rho = -$) Dirac subspaces; see Eq. (A3)]. The ρ -spin propagators $G^{\rho}(p) = G^{\pm}(p)$ are

$$G^{+}(p) = \frac{1}{2E_p - M_d} \quad G^{-}(p) = -\frac{1}{M_d}, \quad (3.15)$$

and u and v are the positive- and negative-energy spinors of particle 2, with $u^{\rho} = u^{\pm}$ and $u^{+}(-\mathbf{p}, \lambda_2) = u(-\mathbf{p}, \lambda_2)$ and $u^{-}(-\mathbf{p}, \lambda_2) = v(\mathbf{p}, \lambda_2)$ [see Eq. (A3)]. Using this notation the wave function can be conveniently written

$$\psi_{\lambda_1\alpha, \lambda_d}(p, P) = \sum_{\rho_2, \lambda_2} \psi_{\lambda_1\lambda_2, \lambda_d}^{\rho_2}(\mathbf{p}) u_{\alpha}^{\rho_2}(-\mathbf{p}, \lambda_2), \quad (3.16)$$

where the wave functions ψ^{ρ_2} are

$$\psi_{\lambda_1\lambda_2, \lambda_d}^{\rho_2}(\mathbf{p}) = N_d \frac{m}{E_p} G^{\rho_2}(p) \Gamma_{\lambda_1\lambda_2, \lambda_d}^{\rho_2}(\mathbf{p}). \quad (3.17)$$

Suppressing the Dirac indices in the matrix elements gives

$$\begin{aligned} \Gamma_{\lambda_1\lambda_2, \lambda_d}^{\rho_2}(\mathbf{p}) &= \bar{u}_{\beta}^{\rho_2}(-\mathbf{p}, \lambda_2) \mathcal{G}_{\lambda_1\beta}(p, P, \lambda_d) \\ &= \bar{u}^{\rho_2}(-\mathbf{p}, \lambda_2) \Gamma^{\mu}(p, P) C \bar{u}^T(\mathbf{p}, \lambda_1) \xi_{\mu}(\lambda_d). \end{aligned} \quad (3.18)$$

These definitions agree with Eq. (2.67) of Ref. II. Note that $\Gamma_{\lambda_1\lambda_2, \lambda_d}^{\rho_2}(\mathbf{p})$ contains no additional factors of m/E ; these are written explicitly in (3.17), and will be identified whenever they appear below.

B. Partial-wave CST equations for the deuteron vertex functions

Equation (3.1), and the companion Eq. (3.7) for the bound state are solved by expanding the amplitudes into partial waves. The connection between the wave functions in momentum space and their partial waves will be briefly outlined in this subsection; for further details see Appendix E of Ref. I and, for the deuteron channel, Appendix A.

The first step is to extract the $J = 1$ partial waves from the momentum-space amplitudes $\Gamma_{\lambda_1\lambda_2, \lambda_d}^{\rho_2}(\mathbf{p})$ defined in Eq. (3.18). These amplitudes have a simple angular dependence. If $\mathbf{p} = p\{\sin\theta \cos\phi, \sin\theta \sin\phi, \cos\theta\}$ (where, from here up to Sec. III G below and again in Appendix C, we will use the notation $p = |\mathbf{p}|$ in order to avoid confusion with the four-momentum p), the form of these functions is

$$\Gamma_{\lambda_1\lambda_2, \lambda_d}^{\rho_2}(\mathbf{p}) = \sqrt{\frac{3}{4\pi}} D_{\lambda_d\lambda}^{1*}(\phi, \theta, 0) \Gamma_{\lambda_1\lambda_2, \lambda_d}^{\rho_2}(p), \quad (3.19)$$

where $\lambda = \lambda_1 - \lambda_2$, $\lambda_d = M_J$ is the polarization of the deuteron, and D^1 are the $J = 1$ rotation matrices. [This can be obtained from Eq. (E29) of Ref. I by recalling that

the scattering amplitude at the deuteron pole is proportional to the product of vertex functions for the incoming and outgoing nucleons.] In order to keep our notation succinct, we suppress the usual index J on the partial-wave vertex function $\Gamma_{\lambda_1\lambda_2, \lambda_d}^{J\rho_2}(\mathbf{p})$. An argument \mathbf{p} instead of \mathbf{p} identifies it unambiguously as a partial-wave vertex function. If we choose \mathbf{p} to lie in the xz plane ($\phi = 0$), we can use (3.19) to find the connection between $\Gamma_{\lambda_1\lambda_2, \lambda_d}^{\rho_2}(\mathbf{p})$ and $\Gamma_{\lambda_1\lambda_2, \lambda_d}^{\rho_2}(p)$. Using the normalization condition for the d^J 's

$$1 = \frac{1}{2}(2J + 1) \int_0^{\pi} \sin\theta d\theta [d_{\lambda'\lambda}^J(\theta)]^2, \quad (3.20)$$

the relation is

$$\Gamma_{\lambda_1\lambda_2, \lambda_d}^{\rho_2}(\mathbf{p}) = \sqrt{3\pi} \int_0^{\pi} \sin\theta d\theta d_{\lambda_d, \lambda}^1(\theta) \Gamma_{\lambda_1\lambda_2, \lambda_d}^{\rho_2}(p), \quad (3.21)$$

and the deuteron wave functions corresponding to (3.17) are

$$\psi_{\lambda_1\lambda_2, \lambda_d}^{\rho_2}(p) = N_d \frac{m}{E_p} G^{\rho_2}(p) \Gamma_{\lambda_1\lambda_2, \lambda_d}^{\rho_2}(p). \quad (3.22)$$

The next step is to use the properties of the partial-wave amplitudes under parity and particle interchange (derived in Refs. I and II) to show that there are only four independent vertex functions. We begin by restoring reference to the ρ spin of particle 1 and the relative energy $p_0 = (p_{10} - p_{20})/2$ (previously suppressed for convenience), which transforms the partial-wave vertex functions (3.21) into

$$\Gamma_{\lambda_1\lambda_2, \lambda_d}^{\rho_2}(\mathbf{p}) \rightarrow \Gamma_{\lambda_1\lambda_2, \lambda_d}^{+\rho_2}(\mathbf{p}, p_0), \quad (3.23)$$

where, in the center-of-mass frame with particle 1 on shell, $p_0 = E_p - M_d/2 > 0$. Under parity (\mathcal{P}) and particle interchange (\mathcal{P}_{12}), it was shown in Appendix E of Ref. I that these partial-wave amplitudes (for $J = 1$) satisfy the transformation properties

$$\begin{aligned} \mathcal{P} \Gamma_{\lambda_1\lambda_2, \lambda_d}^{+\rho_2}(\mathbf{p}, p_0) &= \rho_2 \Gamma_{-\lambda_1-\lambda_2, \lambda_d}^{+\rho_2}(\mathbf{p}, p_0) \\ \mathcal{P}_{12} \Gamma_{\lambda_1\lambda_2, \lambda_d}^{+\rho_2}(\mathbf{p}, p_0) &= \Gamma_{\lambda_2\lambda_1, \lambda_d}^{\rho_2+}(\mathbf{p}, -p_0). \end{aligned} \quad (3.24)$$

The parity relation shows that only two (of the four possible) helicity states are independent (which we chose to be $\lambda_1 = +$; the states with $\lambda_1 = -$ can be obtained using parity) and the exchange relation shows that the exchange amplitudes (with particle 2 on-shell and *negative* relative energy) can be obtained from the direct amplitudes (with particle 1 on-shell and *positive* relative energy). The detailed construction of the deuteron bound-state equation is reviewed in Appendix C 1, where it is shown how these symmetry properties are derived and used to reduce the initial set of coupled equations to only four independent ones.

Using these results, the vertex functions (3.21) can be organized into the column vector

$$|\Gamma_{\lambda_d}(p)\rangle = N_d \frac{m}{E_p} \begin{pmatrix} \Gamma_{++}^{+}(\mathbf{p}) \\ \Gamma_{+-}^{-}(\mathbf{p}) \\ \Gamma_{-+}^{+}(\mathbf{p}) \\ \Gamma_{--}^{-}(\mathbf{p}) \end{pmatrix}, \quad (3.25)$$

where, in the matrix representations, we return to the notation p to represent the momentum dependence of the vertex function. With this notation, the original Eq. (3.7) can now be written

as four independent equations for these vertex functions. In a convenient matrix form the equations are

$$|\Gamma_{\lambda_d}(p)\rangle = - \int_0^\infty k^2 dk \mathbf{V}_d(p, k) \mathbf{g}(k) |\Gamma_{\lambda_d}(k)\rangle, \quad (3.26)$$

where $\mathbf{V}_d(p, k)$ (the kernel at the deuteron pole) and $\mathbf{g}(k)$ are 4×4 matrices. The equations are independent of the projection of the total angular momentum, which for the deuteron is its helicity ($m_J = \lambda_d$). With the factors of m/E explicitly included in the definition of the column vector (3.25) and in the kernel $\mathbf{V}_d(p, k)$, the propagator is the 4×4 diagonal matrix:

$$\mathbf{g}(k) = \begin{pmatrix} G^+(k) & 0 & 0 & 0 \\ 0 & G^-(k) & 0 & 0 \\ 0 & 0 & G^+(k) & 0 \\ 0 & 0 & 0 & G^-(k) \end{pmatrix}, \quad (3.27)$$

with $G^\rho(k)$ defined in Eq. (3.15), and the kernel is the 4×4 matrix

$$\mathbf{V}_d(p, k) = \begin{pmatrix} v_9^{++} & v_1^{+-} & v_{11}^{++} & v_3^{+-} \\ v_9^{-+} & v_1^{--} & v_{11}^{-+} & v_3^{--} \\ v_{12}^{++} & v_8^{+-} & v_{10}^{++} & v_6^{+-} \\ v_{16}^{-+} & v_4^{--} & v_{14}^{-+} & v_2^{--} \end{pmatrix}, \quad (3.28)$$

where the $v_i^{\rho_1 \rho_2}$ were previously defined in Eq. (E45) of Ref. I and also in Eq. (C6).

C. Decomposition into the S , D , and P states

The Dirac operator Γ^μ [from Eq. (3.9)] can be expanded in terms of four scalar invariant functions, F , G , H , and I as given in Eq. (33) of Ref. [5] and Eq. (2.28) of Ref. II,

$$\Gamma^\mu(p, P) = F\gamma^\mu + \frac{G}{m}p^\mu - \frac{(m - \not{p})}{m} \left(H\gamma^\mu + \frac{I}{m}p^\mu \right). \quad (3.29)$$

These functions differ by some factors of 2 from those originally introduced by Blankenbecler and Cook [6]. Since the deuteron is on mass-shell, so $P^2 = W^2 = M_d^2$, they are functions of p^2 only.

Using the explicit definition of the spinors given in Appendix A and substituting the expansion (3.29) into the Dirac matrix elements (3.18), the wave functions (3.17) can be reduced to two-component matrix elements, with four independent scalar wave functions u , w , v_t , and v_s written as linear combinations of the four invariant functions F , G , H , and I [5]:

$$\begin{aligned} \psi_{\lambda_1 \lambda_2, \lambda_d}^+(\mathbf{p}) &= \frac{1}{\sqrt{4\pi}} \chi_{-\lambda_2}^\dagger \left[u(\mathbf{p}) \boldsymbol{\sigma} \cdot \boldsymbol{\xi}_{\lambda_d} + \frac{w(\mathbf{p})}{\sqrt{2}} \right. \\ &\quad \left. \times (3 \hat{\mathbf{p}} \cdot \boldsymbol{\xi}_{\lambda_d} \boldsymbol{\sigma} \cdot \hat{\mathbf{p}} - \boldsymbol{\sigma} \cdot \boldsymbol{\xi}_{\lambda_d}) \right] \frac{i\sigma_2}{\sqrt{2}} \chi_{\lambda_1} \end{aligned} \quad (3.30)$$

$$\begin{aligned} \psi_{\lambda_1 \lambda_2, \lambda_d}^-(\mathbf{p}) &= \sqrt{\frac{3}{4\pi}} \chi_{-\lambda_2}^\dagger \left[v_s(\mathbf{p}) \hat{\mathbf{p}} \cdot \boldsymbol{\xi}_{\lambda_d} - \frac{v_t(\mathbf{p})}{\sqrt{2}} \right. \\ &\quad \left. \times (\boldsymbol{\sigma} \cdot \hat{\mathbf{p}} \boldsymbol{\sigma} \cdot \boldsymbol{\xi}_{\lambda_d} - \hat{\mathbf{p}} \cdot \boldsymbol{\xi}_{\lambda_d}) \right] \frac{i\sigma_2}{\sqrt{2}} \chi_{\lambda_1}. \end{aligned}$$

For explicit formulas connecting the F , G , H , and I to u , w , v_t , and v_s , see Eqs. (45) and (46) of Ref. [5] [note that the normalization factor (3.11) sets the scale of the connection]. Here $\hat{\mathbf{p}} = \{\sin \theta, 0, \cos \theta\}$ lies in the xz plane. For a deuteron at rest with its polarization vector defined along the z direction, the three-vector components of the polarization are

$$\xi_0 = \begin{pmatrix} 0 \\ 0 \\ 1 \end{pmatrix}, \quad \xi_\pm = \mp \frac{1}{\sqrt{2}} \begin{pmatrix} 1 \\ \pm i \\ 0 \end{pmatrix}. \quad (3.31)$$

The nucleon spinors can be either helicity spinors, or spinors with spin projections along the (fixed) z direction. In the latter case, we must replace $\lambda_1 \rightarrow s_1$ and $-\lambda_2 \rightarrow s_2$ (since, when $\theta = 0$, the helicity and spin projection of particle 2 are in opposite directions). In either case (by inspection) the wave functions are real.

The wave functions u , w , v_t , and v_s have a convenient physical interpretation that follows from the structure of the two-component matrix elements that multiply them. This follows most directly if the nucleon spin projections are fixed in the z direction. In this case the $\rho = +$ wave functions is symmetric in the spins,

$$(\psi_{s_1 s_2, \lambda_d}^+(\mathbf{p}))^* = \psi_{s_1 s_2, \lambda_d}^+(\mathbf{p}) = \psi_{s_2 s_1, \lambda_d}^+(\mathbf{p}), \quad (3.32)$$

showing that the spin state is a spin-one triplet. The $\rho = -$ term includes a symmetric piece (the spin triplet v_t) and an antisymmetric piece (corresponding to a spin singlet state, denoted by the s subscript on v_s).

In the CST, wave functions in coordinate space are *defined* to be the Fourier transforms of the momentum-space wave functions (3.30) [5]

$$\psi_{s_1 s_2, \lambda_d}^\rho(\mathbf{r}) = \frac{1}{(2\pi)^{3/2}} \int d^3 p e^{i\mathbf{p}\cdot\mathbf{r}} \psi_{s_1 s_2, \lambda_d}^\rho(\mathbf{p}). \quad (3.33)$$

Using the expansion of the plane wave

$$e^{i\mathbf{p}\cdot\mathbf{r}} = 4\pi \sum_{\ell} \sum_{m=-\ell}^{\ell} i^\ell j_\ell(pr) Y_{\ell m}^*(\hat{\mathbf{p}}) Y_{\ell m}(\hat{\mathbf{r}}) \quad (3.34)$$

and noting that the coefficient of $w(\mathbf{p})$ involves a linear combination of the Y_{2m} spherical harmonics

$$3 \hat{\mathbf{p}}^i \hat{\mathbf{p}}^j - \delta^{ij} = \sum_m c_m^{ij} Y_{2m}(\hat{\mathbf{p}}) \quad (3.35)$$

(where the c_{ij} are constants), it follows that the coordinate space spin representation of the deuteron wave functions is

$$\begin{aligned} \mathbf{r} \psi_{s_1 s_2, \lambda_d}^+(\mathbf{r}) &= \frac{1}{\sqrt{4\pi}} \chi_{s_2}^\dagger \left[u(\mathbf{r}) \boldsymbol{\sigma} \cdot \boldsymbol{\xi}_{\lambda_d} - \frac{w(\mathbf{r})}{\sqrt{2}} \right. \\ &\quad \left. \times (3 \hat{\mathbf{r}} \cdot \boldsymbol{\xi}_{\lambda_d} \boldsymbol{\sigma} \cdot \hat{\mathbf{r}} - \boldsymbol{\sigma} \cdot \boldsymbol{\xi}_{\lambda_d}) \right] \frac{i\sigma_2}{\sqrt{2}} \chi_{s_1} \end{aligned} \quad (3.36)$$

$$\begin{aligned} \mathbf{r} \psi_{s_1 s_2, \lambda_d}^-(\mathbf{r}) &= i \sqrt{\frac{3}{4\pi}} \chi_{s_2}^\dagger \left[v_s(\mathbf{r}) \hat{\mathbf{r}} \cdot \boldsymbol{\xi}_{\lambda_d} - \frac{v_t(\mathbf{r})}{\sqrt{2}} \right. \\ &\quad \left. \times (\boldsymbol{\sigma} \cdot \hat{\mathbf{r}} \boldsymbol{\sigma} \cdot \boldsymbol{\xi}_{\lambda_d} - \hat{\mathbf{r}} \cdot \boldsymbol{\xi}_{\lambda_d}) \right] \frac{i\sigma_2}{\sqrt{2}} \chi_{s_1}, \end{aligned}$$

where, denoting the typical wave function by z_ℓ (so that $z_0 = u$, $z_2 = w$, and $z_1 = v_t$ or v_s), the momentum- and position-space wave functions are related by the spherical Bessel transforms

$$\begin{aligned} z_\ell(\mathbf{p}) &= \sqrt{\frac{2}{\pi}} \int_0^\infty r dr j_\ell(\mathbf{p}r) z_\ell(r) \\ \frac{z_\ell(r)}{r} &= \sqrt{\frac{2}{\pi}} \int_0^\infty p^2 dp j_\ell(\mathbf{p}r) z_\ell(\mathbf{p}), \end{aligned} \quad (3.37)$$

where j_ℓ is the spherical Bessel function of order ℓ with the convenient recursion relation

$$j_\ell(z) = z^\ell \left(-\frac{1}{z} \frac{d}{dz} \right)^\ell \frac{\sin z}{z}. \quad (3.38)$$

Note the appearance of the minus sign multiplying $w(r)$ and the factor of i multiplying both of the P -state terms. These come from the factor of i^ℓ in the plane-wave expansion and can be easily overlooked.

D. Normalization of the wave functions

In Appendix C3 the normalization condition (3.12) is reduced to the following simple form

$$1 = \int_0^\infty p^2 dp [u^2(\mathbf{p}) + w^2(\mathbf{p}) + v_t^2(\mathbf{p}) + v_s^2(\mathbf{p})] + \left\langle \frac{dV}{dM_d} \right\rangle. \quad (3.39)$$

The derivative term is negative and of the order of a few percentages (see Table II).

The origin of the derivative term can be understood in two rather different ways. First, it can be derived from the requirement that a nonlinear version of the equation, used to derive the unitarity relation for positive energies, also holds at the deuteron pole [3,7]. From this point of view, the normalization condition ensures that the strength of the pole is not altered by repeated interactions near the pole and shows that the normalization is a consequence of the equation itself. But the normalization condition also follows from the requirement of current conservation, where it is seen to give precisely the correct factor to ensure that the conservation of charge is an automatic consequence of the correct normalization of the deuteron wave functions [7].

The second interpretation shows, indirectly, that interaction current contributions to the deuteron charge (which give

rise to the derivative term) are necessarily of the order of a few percentages. Hence, the calculation of any electromagnetic property of the deuteron (including the magnetic and quadrupole moments) that does *not* include interaction currents can be expected *ab initio* to be in error by a few percentages. Since the famous discrepancy in the quadrupole moment is about 5%, it is quite possible that a careful CST calculation (including interaction currents) could explain it, but this calculation is, unfortunately, beyond the scope of this article. For this reason, we will not report deuteron moments here.

If these wave functions are used in any calculation that neglects interaction currents, it is probably a better approximation to use what we will refer to as the *scaled* normalization condition

$$1 = \int_0^\infty p^2 dp [u^2(\mathbf{p}) + w^2(\mathbf{p}) + v_t^2(\mathbf{p}) + v_s^2(\mathbf{p})]. \quad (3.40)$$

This is the normalization condition used in the earliest treatments of the CST deuteron wave functions [5], and is also appropriate for electromagnetic calculations using the relativistic impulse approximation (RIA) [9,11].

E. Asymptotic behavior of the wave functions

Using the CST equations for the bound state it is possible to predict the asymptotic behavior of the wave functions. The derivation and discussion are given in Appendix C. The result is:

$$u(\mathbf{p}) \rightarrow \frac{1}{p^5} \quad w(\mathbf{p}) \rightarrow \frac{1}{p^5} \quad v_t(\mathbf{p}) \rightarrow \frac{1}{p^4} \quad v_s(\mathbf{p}) \rightarrow \frac{1}{p^4}. \quad (3.41)$$

Before the convergence of the angular integrals was improved (as discussed in Appendix B) the numerical solutions had power law behaviors which differed from the integers predicted in (3.41) by about 0.1 to 0.2. With the newly converged angular integrals, the actual solutions exhibit the correct behavior.

The odd powers for u and w (and even powers for the v 's) require special consideration when analytic representations of the wave functions are constructed in the following sections.

F. Numerical results for the wave functions

The solutions for the wave functions are tabulated in Tables III and IV, with the high-momentum behavior shown graphically in Fig. 3. The figure shows that the asymptotic estimates (3.41) hold for momenta larger than about 10 GeV and that in this large-momentum region the small P -state wave functions are larger than the dominant S - and D -state wave functions. The wave functions of model WJC-1 satisfy the approximate relations $u \simeq w$, and the D -state wave functions of the two models are also comparable, but the S -state wave function of model WJC-2 has a zero at large momentum and is considerably smaller than its counterpart. The P -state wave functions for both models are comparable to within a factor of 2, with $v_t \sim -v_s$.

TABLE II. The deuteron probabilities (in percentages). Both the exact and scaled probabilities are shown.

Probability	WJC-1		WJC-2	
	Exact	Scaled	Exact	Scaled
P_s	97.3876	92.3330	95.7607	93.5985
P_d	7.7452	7.3432	6.5301	6.3827
P_{v_t}	0.1180	0.1119	0.0103	0.0101
P_{v_s}	0.2234	0.2118	0.0090	0.0088
$\sum P$	105.4743	100.0000	102.3101	100.0000
$\langle V \rangle$	-5.4743	-	-2.3101	-
Total	100.0000	-	100.0000	-

TABLE III. Model WJC-1 momentum-space wave functions $z_\ell(p)$, defined in Eq. (3.30). The momenta are in GeV and the wave functions are in units of $(\text{GeV})^{-3/2}$. These wave functions are normalized to 105.4743%. The original value of w at the lowest momentum point (-0.690345×10^{-2}), which resulted from the imprecise cancellation of two very large terms, can be replaced by the number in bold obtained using p^2 scaling. In this and the following table we use the notation $0.527313(-03) = 0.527313 \times 10^{-3}$.

p	u	w	v_t	v_s
0.527313 (-03)	0.146227 (+03)	0.495991 (-03)	-0.623843 (-03)	-0.358658 (-03)
0.277683 (-02)	0.145699 (+03)	0.137542 (-01)	-0.326777 (-02)	-0.192465 (-02)
0.681763 (-02)	0.142975 (+03)	0.813298 (-01)	-0.802292 (-02)	-0.471289 (-02)
0.126402 (-01)	0.135554 (+03)	0.264888	-0.148513 (-01)	-0.871988 (-02)
0.202305 (-01)	0.121587 (+03)	0.607884 (+00)	-0.236810 (-01)	-0.139049 (-01)
0.295712 (-01)	0.101834 (+03)	0.108549 (+01)	-0.343738 (-01)	-0.201942 (-01)
0.406428 (-01)	0.798054 (+02)	0.160146 (+01)	-0.466979 (-01)	-0.274694 (-01)
0.534249 (-01)	0.594035 (+02)	0.204959 (+01)	-0.603055 (-01)	-0.355616 (-01)
0.678972 (-01)	0.428005 (+02)	0.236963 (+01)	-0.747258 (-01)	-0.442522 (-01)
0.840413 (-01)	0.303155 (+02)	0.255133 (+01)	-0.893826 (-01)	-0.532839 (-01)
0.101842	0.213130 (+02)	0.261197 (+01)	-0.103638	-0.623822 (-01)
0.121288	0.149401 (+02)	0.257733 (+01)	-0.116855	-0.712774 (-01)
0.142375	0.104515 (+02)	0.247282 (+01)	-0.128471	-0.796948 (-01)
0.165107	0.728541 (+01)	0.232096 (+01)	-0.137990	-0.876522 (-01)
0.189495	0.504321 (+01)	0.213982 (+01)	-0.145164	-0.948577 (-01)
0.215565	0.344843 (+01)	0.194353 (+01)	-0.149827	-0.101215
0.243352	0.231128 (+01)	0.174361 (+01)	-0.151964	-0.106954
0.272908	0.149976 (+01)	0.154713 (+01)	-0.151653	-0.111870
0.304301	0.922698	0.135976 (+01)	-0.149087	-0.115989
0.337619	0.515394	0.118496 (+01)	-0.144410	-0.119420
0.372971	0.231152	0.102511 (+01)	-0.137607	-0.122164
0.410488	0.350502 (-01)	0.881942	-0.128414	-0.123858
0.450334	-0.994438 (-01)	0.755982	-0.116562	-0.124047
0.492699	-0.189481	0.645099	-0.102598	-0.121827
0.537814	-0.242417	0.544069	-0.878148 (-01)	-0.116934
0.585950	-0.262050	0.448848	-0.732091 (-01)	-0.109318
0.637426	-0.255246	0.360144	-0.592921 (-01)	-0.995041 (-01)
0.692620	-0.230996	0.280430	-0.465163 (-01)	-0.878985 (-01)
0.751979	-0.196819	0.211737	-0.352080 (-01)	-0.752255 (-01)
0.816027	-0.159320	0.154625	-0.255953 (-01)	-0.621228 (-01)
0.885387	-0.122585	0.109110	-0.177552 (-01)	-0.494042 (-01)
0.960798	-0.900194 (-01)	0.742130 (-01)	-0.116628 (-01)	-0.377210 (-01)
0.104314 (+01)	-0.631224 (-01)	0.485750 (-01)	-0.718102 (-02)	-0.276183 (-01)
0.113346 (+01)	-0.421787 (-01)	0.305968 (-01)	-0.408644 (-02)	-0.193771 (-01)
0.123304 (+01)	-0.270876 (-01)	0.185189 (-01)	-0.208407 (-02)	-0.130319 (-01)
0.134340 (+01)	-0.166413 (-01)	0.107639 (-01)	-0.902849 (-03)	-0.841563 (-02)
0.146641 (+01)	-0.978419 (-02)	0.600927 (-02)	-0.275271 (-03)	-0.522831 (-02)
0.160437 (+01)	-0.557178 (-02)	0.322545 (-02)	0.208268 (-04)	-0.313370 (-02)
0.176010 (+01)	-0.306684 (-02)	0.165657 (-02)	0.125739 (-03)	-0.181763 (-02)
0.193715 (+01)	-0.161854 (-02)	0.809542 (-03)	0.135471 (-03)	-0.102257 (-02)
0.213998 (+01)	-0.824465 (-03)	0.375303 (-03)	0.110722 (-03)	-0.559203 (-03)
0.237429 (+01)	-0.413458 (-03)	0.166244 (-03)	0.808655 (-04)	-0.297772 (-03)
0.264741 (+01)	-0.199638 (-03)	0.674221 (-04)	0.532186 (-04)	-0.154652 (-03)
0.296895 (+01)	-0.923325 (-04)	0.240784 (-04)	0.325043 (-04)	-0.783631 (-04)
0.335166 (+01)	-0.410173 (-04)	0.695348 (-05)	0.187892 (-04)	-0.386958 (-04)
0.381276 (+01)	-0.175134 (-04)	0.111592 (-05)	0.103540 (-04)	-0.185709 (-04)
0.437596 (+01)	-0.741984 (-05)	-0.231960 (-06)	0.551835 (-05)	-0.861492 (-05)
0.507462 (+01)	-0.295514 (-05)	-0.452993 (-06)	0.276257 (-05)	-0.386774 (-05)
0.595690 (+01)	-0.110492 (-05)	-0.326876 (-06)	0.131169 (-05)	-0.167290 (-05)
0.709434 (+01)	-0.387509 (-06)	-0.173383 (-06)	0.590317 (-06)	-0.691226 (-06)
0.859696 (+01)	-0.126207 (-06)	-0.755500 (-07)	0.249574 (-06)	-0.269811 (-06)
0.106411 (+02)	-0.375810 (-07)	-0.277414 (-07)	0.976773 (-07)	-0.979770 (-07)
0.135237 (+02)	-0.999262 (-08)	-0.854233 (-08)	0.346150 (-07)	-0.323958 (-07)
0.177768 (+02)	-0.228855 (-08)	-0.215203 (-08)	0.107312 (-07)	-0.945155 (-08)

TABLE III. (Continued.)

p	u	w	v_t	v_s
0.244325 (+02)	-0.428701 (-09)	-0.419613 (-09)	0.272700 (-08)	-0.230841 (-08)
0.357133 (+02)	-0.637757 (-10)	-0.539018 (-10)	0.510267 (-09)	-0.433381 (-09)
0.571587 (+02)	-0.579367 (-11)	-0.486092 (-11)	0.700683 (-10)	-0.590489 (-10)
0.105975 (+03)	-0.256335 (-12)	-0.211897 (-12)	0.562191 (-11)	-0.463011 (-11)
0.260189 (+03)	-0.281688 (-14)	-0.229780 (-14)	0.150575 (-12)	-0.121213 (-12)
0.137016 (+04)	-0.682890 (-18)	-0.553955 (-18)	0.193220 (-15)	-0.153367 (-15)

The probabilities for each of the components of the wave function are summarized in Table II. This table reports both the exact probabilities (3.39) and the renormalized probabilities (3.40).

G. Fitted wave functions

It is convenient to fit the deuteron wave functions to a series of simple functions that can be analytically integrated to obtain the wave functions in r space and to interpolate for any value of the momentum. This subsection will describe in detail how this is done.

1. Scaling the wave functions

The first step in the fitting process is to scale out the rapid dependence of each wave function on the momentum, so it may be studied on a linear plot. Constructing these scaling functions took some care, but in the end we found that the

functions

$$\begin{aligned}
 u_{\text{scale}}(p) &= \frac{N}{(m_{s1}^2 + p^2)^{3/2}} \left[\frac{1}{(\alpha_0^2 + p^2)} + \frac{R}{(m_{s1}^2 + p^2)} \right] \\
 w_{\text{scale}}(p) &= \frac{Np^2}{(m_{s1}^2 + p^2)^{5/2}} \left[\frac{1}{(\alpha_0^2 + p^2)} + \frac{R}{(m_{s1}^2 + p^2)} \right] \\
 v_{\text{scale}}(p) &= Np \left[\frac{1}{(m_{s0}^2 + p^2)^{5/2}} + \frac{R}{(m_{s1}^2 + p^2)^{5/2}} \right] \quad (3.42)
 \end{aligned}$$

(where the functional form of v_{scale} is used to scale both v_t and v_s , and we return to denoting the magnitude of the three-momentum by p [instead of p , as we did in the previous sections]) work very well. Note that these scaling functions have the desired (and observed) p^ℓ behavior at small p and

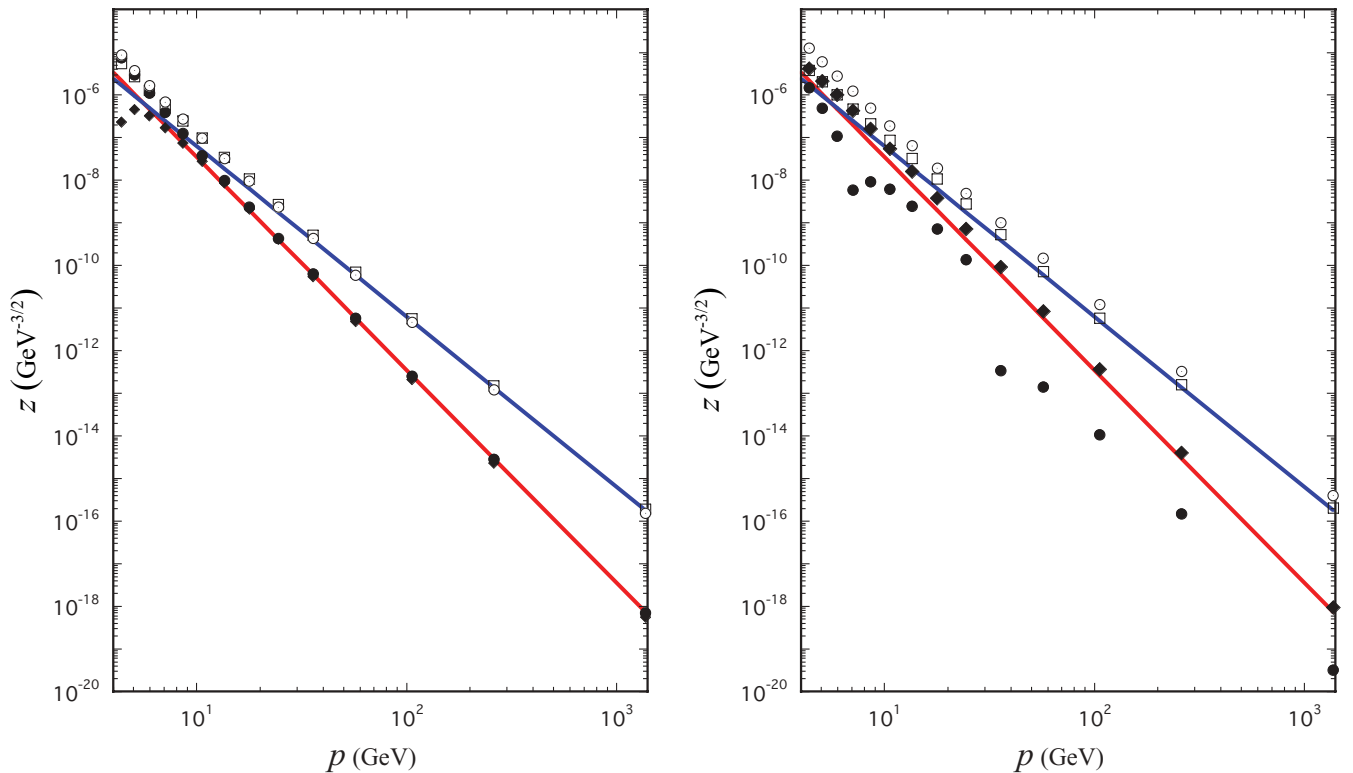


FIG. 3. (Color online) Log-log plot of the momentum-space wave functions $z_\ell(p)$, defined in Eq. (3.30). (Left) WJC-1 wave functions tabulated in Table III, compared to the asymptotic limits given in Eq. (3.41). Solid circles: u_0 ; solid diamonds: w_2 ; open squares: v_t ; open circles: v_s . The solid lines are proportional to p^{-5} and p^{-4} . (Right) WJC-2 wave functions tabulated in Table IV. The solid lines are proportional to p^{-5} and p^{-4} , and are the same as those shown in the left panel.

TABLE IV. Model WJC-2 momentum-space wave functions $z_\ell(p)$, defined in Eq. (3.30). The momenta are in GeV and the wave functions are in units of $(\text{GeV})^{-3/2}$. These wave functions are normalized to 102.3101%. The original value of w at the lowest momentum point (-0.743127×10^{-2}), which resulted from the imprecise cancellation of two very large terms, can be replaced by the number in bold obtained using p^2 scaling. As in Table III, we use numbers in parentheses to denote power-of-10 factors.

p	u	w	v_t	v_s
0.527313 (-03)	0.146161 (+03)	0.497127 (-03)	0.673653 (-04)	0.139978 (-04)
0.277683 (-02)	0.145632 (+03)	0.137857 (-01)	0.372102 (-03)	0.368387 (-04)
0.681763 (-02)	0.142908 (+03)	0.815168 (-01)	0.908775 (-03)	0.100679 (-03)
0.126402 (-01)	0.135483 (+03)	0.265495	0.168223 (-02)	0.192525 (-03)
0.202305 (-01)	0.121510 (+03)	0.609254	0.268867 (-02)	0.316519 (-03)
0.295712 (-01)	0.101747 (+03)	0.108784 (+01)	0.392312 (-02)	0.478611 (-03)
0.406428 (-01)	0.797033 (+02)	0.160461 (+01)	0.537988 (-02)	0.685611 (-03)
0.534249 (-01)	0.592857 (+02)	0.205290 (+01)	0.705316 (-02)	0.942644 (-03)
0.678972 (-01)	0.426693 (+02)	0.237212 (+01)	0.893675 (-02)	0.125069 (-02)
0.840413 (-01)	0.301750 (+02)	0.255185 (+01)	0.110216 (-01)	0.160505 (-02)
0.101842	0.211677 (+02)	0.260934 (+01)	0.132917 (-01)	0.199444 (-02)
0.121288	0.147943 (+02)	0.257044 (+01)	0.157173 (-01)	0.240633 (-02)
0.142375	0.103090 (+02)	0.246071 (+01)	0.182443 (-01)	0.286055 (-02)
0.165107	0.714947 (+01)	0.230262 (+01)	0.208563 (-01)	0.313824 (-02)
0.189495	0.491677 (+01)	0.211456 (+01)	0.234179 (-01)	0.337366 (-02)
0.215565	0.333407 (+01)	0.191087 (+01)	0.258522 (-01)	0.353615 (-02)
0.243352	0.221125 (+01)	0.170294 (+01)	0.280836 (-01)	0.333024 (-02)
0.272908	0.141604 (+01)	0.149824 (+01)	0.300543 (-01)	0.284267 (-02)
0.304301	0.856967	0.130257 (+01)	0.317477 (-01)	0.198415 (-02)
0.337619	0.468983	0.111930 (+01)	0.332767 (-01)	0.574512 (-03)
0.372971	0.205168	0.950678	0.348229 (-01)	-0.146236 (-02)
0.410488	0.303889 (-01)	0.798954	0.363889 (-01)	-0.389499 (-02)
0.450334	-0.828321 (-01)	0.665687	0.373820 (-01)	-0.658983 (-02)
0.492699	-0.154049	0.551266	0.364870 (-01)	-0.910693 (-02)
0.537814	-0.193834	0.452057	0.328528 (-01)	-0.113767 (-01)
0.585950	-0.206848	0.363715	0.273257 (-01)	-0.131172 (-01)
0.637426	-0.198813	0.284931	0.211844 (-01)	-0.143788 (-01)
0.692620	-0.177269	0.216739	0.152062 (-01)	-0.150297 (-01)
0.751979	-0.148546	0.159772	0.980525 (-02)	-0.151078 (-01)
0.816027	-0.118123	0.113878	0.537109 (-02)	-0.145747 (-01)
0.885387	-0.890513 (-01)	0.783167 (-01)	0.202329 (-02)	-0.134838 (-01)
0.960798	-0.639660 (-01)	0.518825 (-01)	-0.137523 (-03)	-0.119400 (-01)
0.104314 (+01)	-0.437558 (-01)	0.329899 (-01)	-0.128366 (-02)	-0.101160 (-01)
0.113346 (+01)	-0.283328 (-01)	0.201385 (-01)	-0.167665 (-02)	-0.815911 (-02)
0.123304 (+01)	-0.176467 (-01)	0.117937 (-01)	-0.152450 (-02)	-0.629427 (-02)
0.134340 (+01)	-0.104205 (-01)	0.658279 (-02)	-0.117834 (-02)	-0.462869 (-02)
0.146641 (+01)	-0.581003 (-02)	0.349531 (-02)	-0.802520 (-03)	-0.324538 (-02)
0.160437 (+01)	-0.315563 (-02)	0.176854 (-02)	-0.460812 (-03)	-0.218651 (-02)
0.176010 (+01)	-0.164951 (-02)	0.831201 (-03)	-0.227186 (-03)	-0.141637 (-02)
0.193715 (+01)	-0.800130 (-03)	0.349702 (-03)	-0.970097 (-04)	-0.879905 (-03)
0.213998 (+01)	-0.366387 (-03)	0.123929 (-03)	-0.288285 (-04)	-0.527587 (-03)
0.237429 (+01)	-0.175371 (-03)	0.333862 (-04)	0.684399 (-05)	-0.308052 (-03)
0.264741 (+01)	-0.775657 (-04)	-0.229350 (-05)	0.151418 (-04)	-0.173993 (-03)
0.296895 (+01)	-0.303299 (-04)	-0.119914 (-04)	0.132282 (-04)	-0.949967 (-04)
0.335166 (+01)	-0.104213 (-04)	-0.111922 (-04)	0.932916 (-05)	-0.502335 (-04)
0.381276 (+01)	-0.309564 (-05)	-0.763178 (-05)	0.592149 (-05)	-0.257338 (-04)
0.437596 (+01)	-0.144689 (-05)	-0.413223 (-05)	0.374297 (-05)	-0.128285 (-04)
0.507462 (+01)	-0.488220 (-06)	-0.212835 (-05)	0.203145 (-05)	-0.613856 (-05)
0.595690 (+01)	-0.107614 (-06)	-0.100778 (-05)	0.101279 (-05)	-0.280394 (-05)
0.709434 (+01)	-0.588336 (-08)	-0.430027 (-06)	0.476792 (-06)	-0.121466 (-05)
0.859696 (+01)	0.931326 (-08)	-0.163878 (-06)	0.211653 (-06)	-0.493916 (-06)
0.106411 (+02)	0.613471 (-08)	-0.550345 (-07)	0.873897 (-07)	-0.185900 (-06)
0.135237 (+02)	0.246297 (-08)	-0.159505 (-07)	0.327267 (-07)	-0.635345 (-07)
0.177768 (+02)	0.718045 (-09)	-0.386029 (-08)	0.106738 (-07)	-0.191935 (-07)
0.244325 (+02)	0.135807 (-09)	-0.732832 (-09)	0.281290 (-08)	-0.493279 (-08)
0.357133 (+02)	0.340558 (-12)	-0.932273 (-10)	0.532783 (-09)	-0.101591 (-08)

TABLE IV. (Continued.)

p	u	w	v_t	v_s
0.571587 (+02)	-0.140690 (-12)	-0.832796 (-11)	0.721192 (-10)	-0.146914 (-09)
0.105975 (+03)	-0.105532 (-13)	-0.362792 (-12)	0.579471 (-11)	-0.119175 (-10)
0.260189 (+03)	-0.149527 (-15)	-0.394161 (-14)	0.157413 (-12)	-0.317980 (-12)
0.137016 (+04)	-0.319205 (-19)	-0.952160 (-18)	0.204258 (-15)	-0.406280 (-15)

the correct asymptotic p dependence for all the states. The parameter α_0 is determined from the asymptotic relation

$$2\sqrt{m^2 + \alpha_0^2} - 2m = \epsilon, \quad (3.43)$$

where $\epsilon = 2.2246$ MeV is the deuteron binding energy. This gives

$$\alpha_0 = \sqrt{m\epsilon + \frac{1}{4}\epsilon^2} = 45.7159 \text{ MeV}, \quad (3.44)$$

which differs slightly from the nonrelativistic $\alpha_0 = \sqrt{m\epsilon}$, as has been emphasized in the literature [8]. This parameter is fixed by the deuteron binding energy, but since the P states vanish outside of the range of the potential, their leading mass, $m_{s,0}$ is treated as a free parameter. This and the other parameters used in the scale functions are given in Tables V and VI.

2. Fitting the momentum-space wave functions

In order to obtain simple and accurate fits to the wave functions, we found it convenient to use two types of expansion functions. For all but the high-momentum tail, we use functions that go asymptotically like an *even* power of p for the S and D states (in practice, p^{-6}), and an *odd* power of p for the P states (p^{-5}). These functions go to zero precisely one power of p faster than the observed asymptotic behavior and are very conveniently transformed to coordinate space giving a superposition of simple exponentials. However, the asymptotic tail cannot be accurately described by such functions, and it is best to use one special function with the correct asymptotic behavior in order to describe the wave functions at very large p . This function has a Fourier transform that can be written in terms of the modified Bessel functions of the second kind, $K_n(z)$.

For the first set of functions we choose

$$G_\ell^i(p) = \sqrt{\frac{2}{\pi}} \frac{p^\ell m_i^2 M_i^{2n_\ell - \ell}}{(m_i^2 + p^2)(M_i^2 + p^2)^{n_\ell}}, \quad (3.45)$$

where the factor of $\sqrt{2/\pi}$ is introduced for convenience, ℓ is the angular momentum of the state, $n_\ell = 2$ for the S and P states, and $n_\ell = 3$ for the D state. Note that, near $p = 0$, the

TABLE V. The parameters used in the scaling functions (3.42) for model WJC-1.

	N (GeV ²)	R	$m_{s,1}$ (GeV)	$m_{s,0}$ (GeV)
u	0.4775	-0.9856	1.160	-
w	0.5131	-0.9942	0.685	-
v_t	-0.03024	-0.9772	1.600	0.488
v_s	-0.1038	-0.9942	0.840	0.633

expansion functions have the normalization

$$G_\ell^i(0) \rightarrow \sqrt{\frac{2}{\pi}} \left(\frac{p}{M_i}\right)^\ell. \quad (3.46)$$

The tail wave function, chosen to have the correct falloff as $p \rightarrow \infty$ and the same normalization at $p = 0$, is

$$G_\ell^n(p) = \sqrt{\frac{2}{\pi}} \frac{p^\ell M_n^{2n_\ell + 1 - \ell}}{(M_n^2 + p^2)^{n_\ell + \frac{1}{2}}}. \quad (3.47)$$

Denoting the typical wave function by z_ℓ (as we did above) the full momentum-space wave functions are expanded in terms of G_ℓ

$$z_\ell(p) = \sum_{i=1}^n b_i^z G_\ell^i(p). \quad (3.48)$$

As this notation implies, the last function in the sum (when $i = n$) is the ‘‘tail’’ function (3.47), while the first $n - 1$ are of the type (3.45). Since all of the functions have the same normalization at $p = 0$, the relative size of the expansion coefficients is a measure of the relative size of each function (at least at small momenta).

The masses m_i and M_i (for $i = 1$ to $n - 1$) used in each of the wave functions (3.45) were defined by the relations

$$m_i = \alpha_\ell + (i - 1)m_x \quad M_i = m_i + m_x \quad (3.49)$$

where the ‘‘step’’ mass m_x was chosen to depend on the number of functions n used in the expansion, so

$$m_x = \frac{M_0}{n - 1}. \quad (3.50)$$

Hence, as n increases, m_x decreases, giving a finer ‘‘grid’’ of mass scales. It was found that the precise value of M_0 was not critical, except that fitting the complicated structure of $u(2)$ at large p [to save writing in this section, we will sometimes use the notation $z(i)$ to denote the generic wave function for model WJC- i , with $i = 1, 2$] required a large value of M_0 which we chose to be 10 GeV. For all other wave functions we chose $M_0 = 1.2$ GeV. The leading mass α_ℓ determines the asymptotic behavior of the wave function at large distances in coordinate

TABLE VI. The parameters used in the scaling functions (3.42) for model WJC-2.

	N (GeV ²)	R	$m_{s,1}$ (GeV)	$m_{s,0}$ (GeV)
u	0.3060	-0.9978	1.000	-
w	0.3474	-0.9805	0.632	-
v_t	0.02700	-0.9610	0.700	0.551
v_s	-0.2356	-0.9671	1.054	0.977

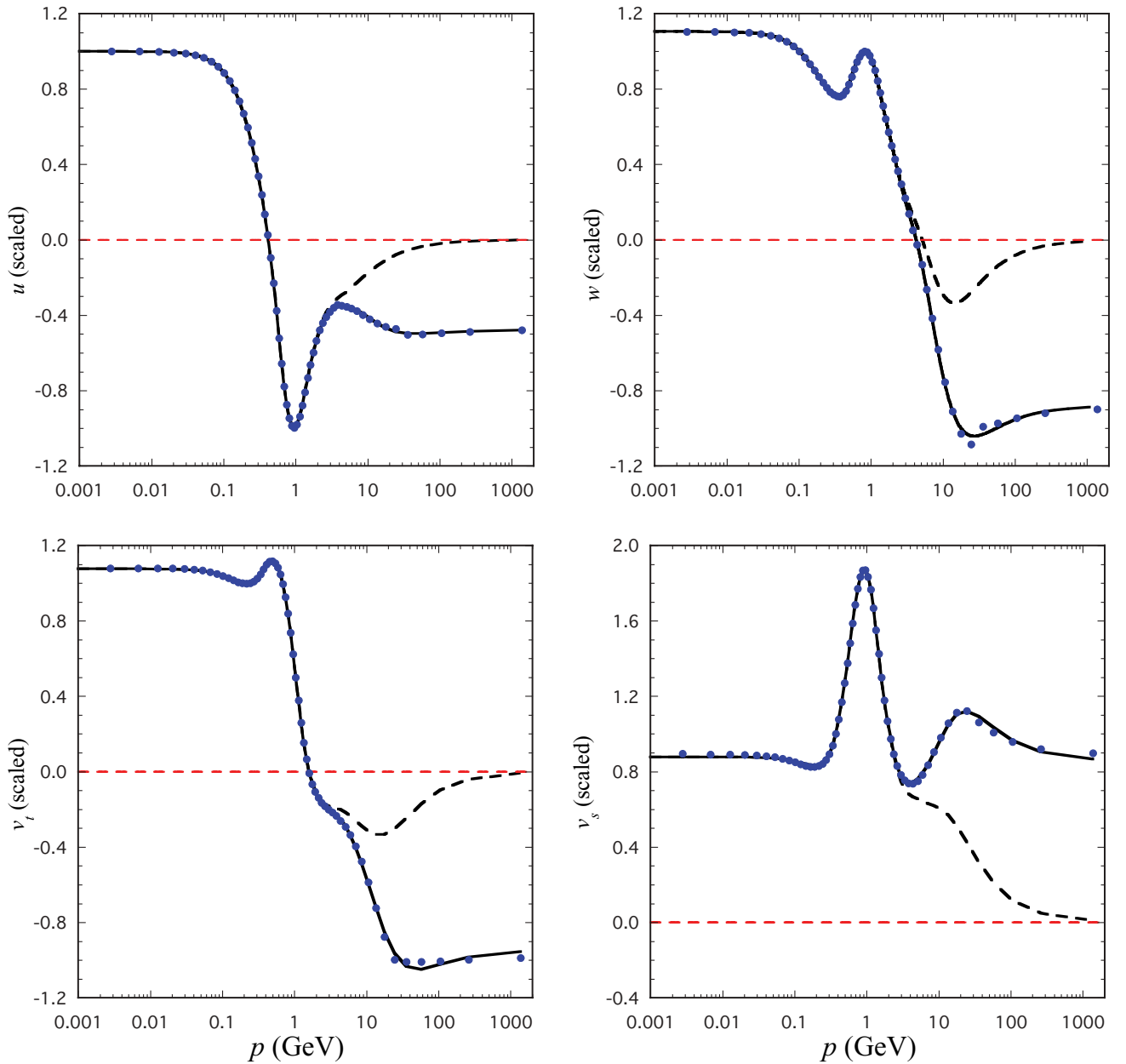


FIG. 4. (Color online) Ratio of the raw wave functions for model WJC-1 (taken from Table III with the probabilities “scaled” to 100% as outlined in Table II) to the scale functions from Eq. (3.42). In each panel the dots are the tabulated wave functions for 60 Gauss points, the solid curve the fit, and the dashed curve the fit *without* the tail wave function of Eq. (3.47).

space, and was therefore chosen to be α_0 of Eq. (3.44) for the S and D states and m_{s0} from Tables V and VI for the P states. The mass M_n used in the tail wave functions was fixed at 2 GeV for all z_i except $u(2)$, where it was fixed at 11 GeV. Using these choices, excellent fits to the momentum-space wave functions were found. These are shown in Figs. 4 and 5. The expansion parameters are given in Tables VII and VIII.

Note that the coefficients of the tail wave functions are very small. The tail makes a negligible contribution at low momenta but is very important for the description of the high-momentum components (for $p \gtrsim 3$ GeV). Remember that the figures show *ratios*; the actual values of the wave functions are quite small above $p \gtrsim 3$ GeV, as already shown in Fig. 3. In any case it will be possible to study the sensitivity of any observable to

the very high momentum components simply by setting this last coefficient to zero.

The components $z(i)$ are compared in Fig. 6, which shows the wave functions for momenta up to 1 GeV, and in Fig. 7, which gives an exploded view of the wave functions from 500 MeV to 1 GeV, a region where all of the components might be important, depending on whether the larger u and w components interfere in a given matrix element.

In general we observe that the P -state components are very small, particularly so for WJC-2. However, in the region shown in Fig. 7 all of the WJC-1 components are larger than the WJC-2 ones, and it is possible that, in some observables, P -state components might compensate for the differences between the larger S - and D -state components.

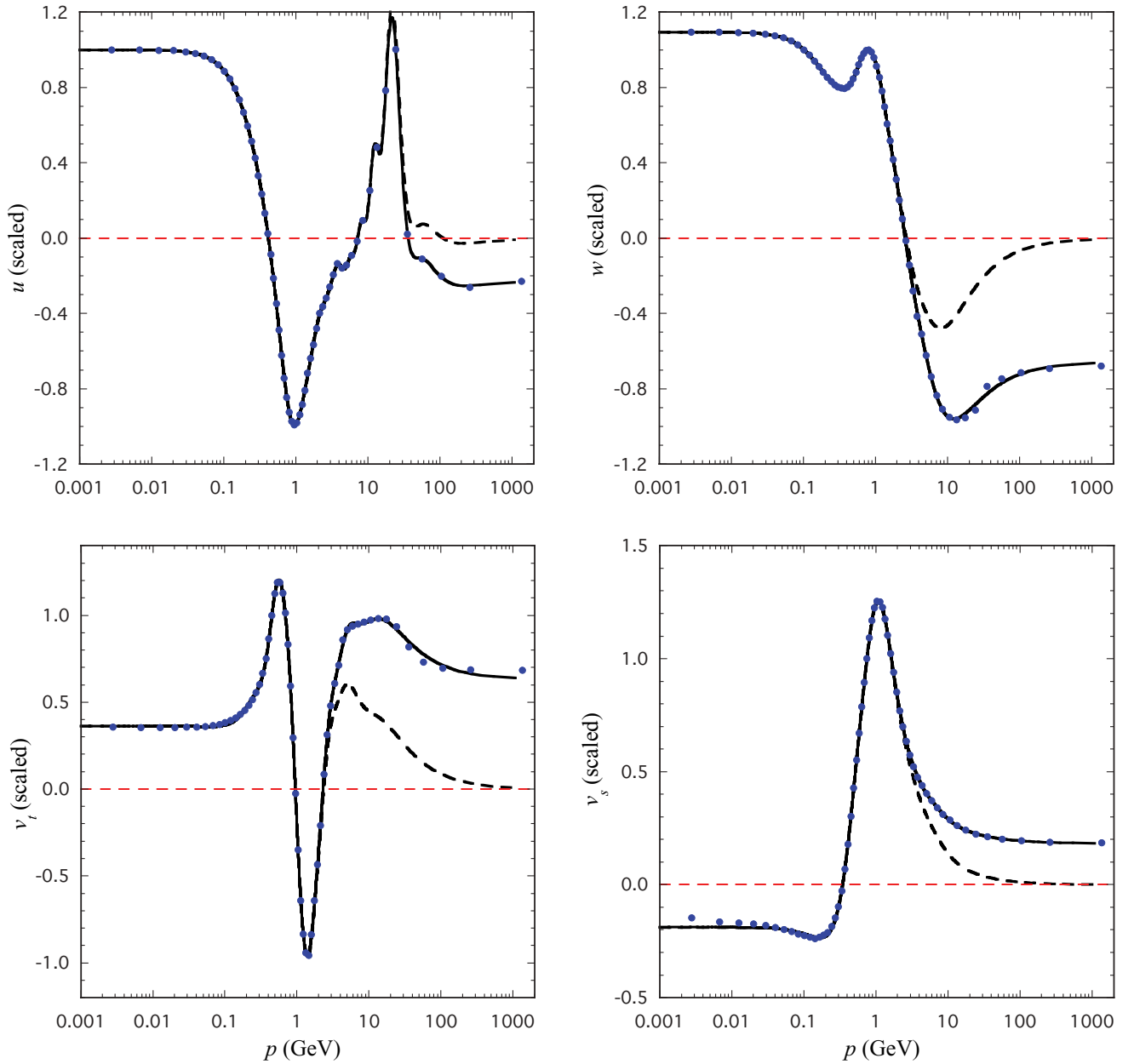


FIG. 5. (Color online) Ratio of the raw wave functions for model WJC-2 (taken from Table IV with the probabilities “scaled” to 100% as outlined in Table II) to the scale functions from Eq. (3.42). In each panel the dots are the tabulated wave functions for 60 Gauss points, the solid curve the fit, and the dashed curve the fit *without* the tail wave function of Eq. (3.47). The origin of the many small wiggles in u is not clear, but u is *very* small in this region, as illustrated in Fig. 3. The first v_s point is probably inaccurate because of numerical cancellations.

3. Transformations to coordinate space

The coordinate space wave functions are constructed from the spherical Bessel transforms (3.37). The details are given in Appendix C 4. The coordinate space expansion functions for the terms $i < n$ are

$$G_0^i(r) = A_i \left\{ e^{-z_i} - e^{-Z_i} \left[1 + \frac{1}{2} Z_i (1 - R_i^2) \right] \right\}$$

$$G_1^i(r) = A_i \left\{ R_i e^{-z_i} \left[1 + \frac{1}{z_i} \right] - e^{-Z_i} \left[1 + \frac{1}{z_i} + \frac{1}{2} Z_i (1 - R_i^2) \right] \right\},$$

$$G_2^i(r) = B_i \left\{ R_i^2 e^{-z_i} \left[1 + \frac{3}{z_i} + \frac{3}{z_i^2} \right] - e^{-Z_i} \left[1 + \frac{3}{Z_i} + \frac{3}{Z_i^2} + \frac{1}{2} (1 - R_i^2) (1 + Z_i) + \frac{1}{8} Z_i^2 (1 - R_i^2)^2 \right] \right\}, \quad (3.51)$$

where

$$z_i = m_i r, \quad Z_i = M_i r, \quad R_i = \frac{m_i}{M_i}, \quad (3.52)$$

$$A_i = \frac{m_i^2 M_i^4}{(M_i^2 - m_i^2)^2}, \quad B_i = A_i \frac{M_i^2}{M_i^2 - m_i^2}.$$

TABLE VII. The expansion parameters for model WJC-1 wave functions. First line is the total number of terms n in the sum (3.48), second is the step mass m_x (in MeV), the next lines are the coefficients b_i^z , $\{i = 1, n-1\}$, with the last line the coefficient b_n^z , written in the notation $125.(-6) = 125 \times 10^{-6}$, all in units of $\text{GeV}^{-3/2}$. The “tail” mass $M_n = 2$ GeV in all cases. We have given the b_i^z coefficients to three decimal places, sufficient to reproduce the fits to better than 1% accuracy; if greater accuracy is needed use the routine supplied by the authors.

u	w	v_t	v_s
17	16	12	12
75.00	80.00	109.09	109.09
134.963	23.813	-27.120	-78.763
52.871	32.709	233.155	688.895
-217.709	-111.381	-998.893	-2910.298
1876.699	844.861	2679.013	7690.084
-11369.449	-4376.965	-4894.348	-13861.596
49427.176	16489.297	6277.553	17571.825
-156695.247	-45122.587	-5674.481	-15720.926
369322.468	91061.493	3550.952	9747.293
-655367.189	-136672.748	-1467.429	-3994.747
879178.453	152411.033	360.796	975.086
-887103.150	-124633.670	-40.010	-107.489
662786.733	72617.712	0.000	0.000
-355720.583	-28549.974	0.000	0.000
129738.381	6790.953	0.000	0.000
-28805.739	-738.436	0.000	0.000
2939.797	0.000	0.000	0.000
-125.0(-6)	-100.0(-6)	498.0(-7)	-394.0(-7)

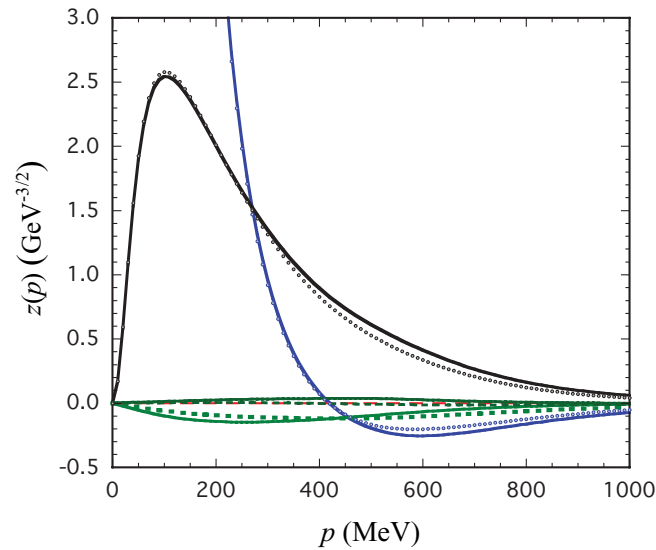


FIG. 6. (Color online) Momentum-space wave functions for both models. The largest two wave functions are u (off-scale at small momenta) and w (zero at small momenta), with WJC-1 wave functions solid and WJC-2 dotted. The P -state wave functions are very small, with both v_t wave functions solid ($v_t(1)$ is negative and $v_t(2)$ positive) and both v_s wave functions dotted (with $v_s(1)$ negative and $v_s(2)$ very near zero). Note that the S -state wave function has a zero around 400 MeV and that above this momentum the P -state wave functions for WJC-1 are no longer completely negligible, as illustrated in Fig. 7.

TABLE VIII. The expansion parameters for each of the model WJC-2 wave functions. For an explanation, see the caption to Table VII. Here the “tail” mass $M_n = 11$ GeV for u and 2 GeV for all other wave functions.

u	w	v_t	v_s
30	18	21	18
344.83	70.59	60.00	70.59
181.569	19.228	33002.947	-74492.403
-0.817	30.600	-580810.178	1001357.141
11.779	-107.606	4678729.666	-6147239.139
-64.782	897.969	-22669425.846	22510592.546
247.517	-5434.481	72755203.760	-53617329.732
-772.915	25662.551	-159377601.760	83750225.006
1948.950	-93771.705	232305154.041	-77842609.919
-3813.276	265664.025	-193963168.786	20071103.929
5397.722	-582644.240	15154785.470	48069105.480
-4637.491	986160.259	160238368.984	-61037252.628
533.832	-1280768.137	-139567635.494	8837981.631
3806.187	1262888.436	-52643827.848	50533251.186
-3119.875	-927821.391	170318699.570	-65739754.705
-2058.321	491712.394	-69098137.722	43130588.287
3773.554	-177560.697	-117233319.092	-16799889.802
1092.507	39095.262	192985378.690	3717624.361
-3919.342	-3959.947	-138452165.069	-363261.256
-736.912	0.000	56855870.564	0.000
4018.336	0.000	-13049893.573	0.000
612.635	0.000	1310791.818	0.000
-4213.343	0.000	0.000	0.000
-260.753	0.000	0.000	0.000
4581.603	0.000	0.000	0.000
-1171.113	0.000	0.000	0.000
-4559.802	0.000	0.000	0.000
5566.714	0.000	0.000	0.000
-2984.961	0.000	0.000	0.000
814.407	0.000	0.000	0.000
-92.509	0.000	0.000	0.000
-118.0(-11)	-172.0(-6)	514.0(-7)	-110.0(-6)

The last terms with $i = n$ are

$$\begin{aligned}
 G_0^n(r) &= \frac{2}{3\pi} M_n^2 Z_n^2 K_1(Z_n) \\
 G_1^n(r) &= \frac{2}{3\pi} M_n^2 Z_n^2 K_0(Z_n) \\
 G_2^n(r) &= \frac{2}{15\pi} M_n^2 Z_n^3 K_0(Z_n),
 \end{aligned} \tag{3.53}$$

where $K_n(z)$ are the modified Bessel functions of the second kind (see Appendix C).

At large r the functions $G_\ell^i(i < n)$ have the asymptotic behavior expected for solutions of the Schrödinger equation with orbital angular momentum ℓ . The tail functions fall off like exponentials multiplied by a fractional power of r but, because of their large mass (short range), do not contribute to the overall asymptotic behavior of the wave functions in coordinate space. At small r the functions $G_\ell^i(i < n)$ have the expected $r^{\ell+1} \sim Z_i^{\ell+1}$ behavior. However, the tail functions contribute some nonanalytic logarithmic behavior at small r ,

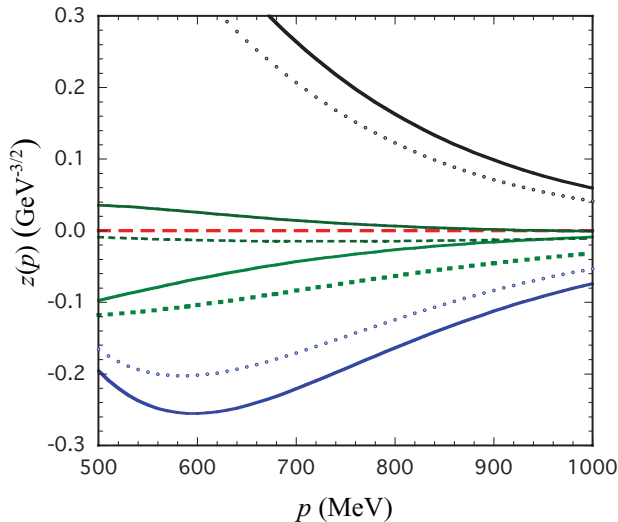


FIG. 7. (Color online) Enlarged view of the momentum-space wave functions for both models. The curves are labeled as in Fig. 6, with a dashed zero reference line. The large positive states are w ; the large negative ones u , with WJC-1 (solid) and WJC-2 (dotted). Both of the v_t wave functions are solid, and both v_s wave functions are dotted, with WJC-1 negative in both cases and WJC-2 both close to zero.

as described in Appendix C. As it turns out, this behavior is too small to be seen at the level of 1%.

4. Wave functions in coordinate space

The coordinate-space wave functions are shown in Figs. 8 and 9. Note that the P -state wave functions vanish beyond 2 fm and that they are larger for WJC-1, as already seen in momentum space. Particularly notable is the zero in $u(1)$ near

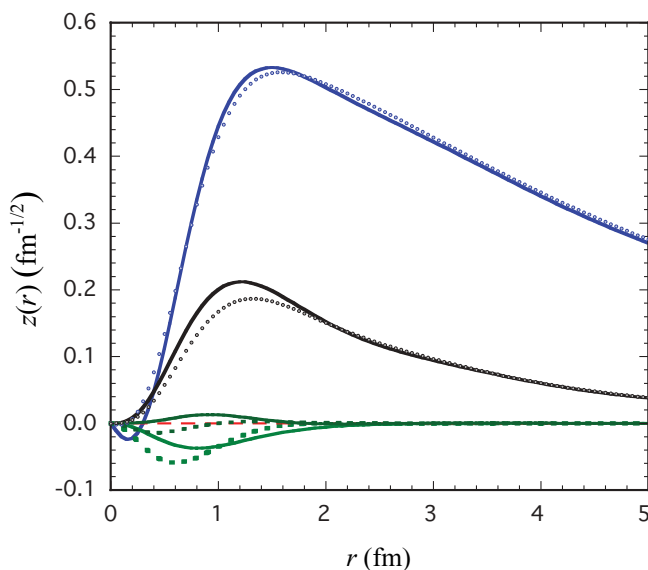


FIG. 8. (Color online) The coordinate space wave functions for both models. The curves are labeled as in Fig. 6. The largest two wave functions are the familiar u and w wave functions, with WJC-1 solid and WJC-2 dotted. The v_t wave functions are both solid, with WJC-1 negative and WJC-2 small and positive. The v_s wave functions are both dotted, with WJC-1 negative and WJC-2 close to zero.

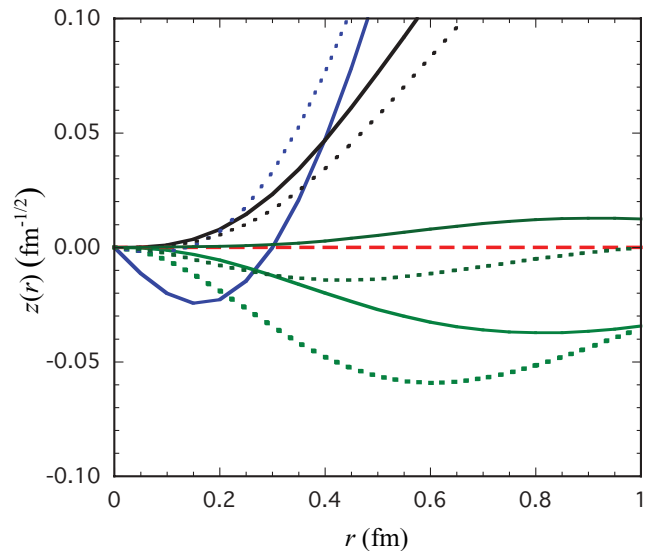


FIG. 9. (Color online) Expanded view of the coordinate space wave functions for both models at small r . The curves are labeled as in Fig. 6.

0.35 fm, leading to a dip at very small r . This dip is not an artifact of the fits; it is related to the deeper dip in the momentum space $u(1)$ near 600 MeV, clearly seen in Figs. 6 and 7. A similar (but much smaller and not visible in the figures) dip is also present in $u(2)$ inside of 0.15 fm. The fact that the dip in WJC-1 is larger may be a consequence of its larger P -state components.

Before any physical conclusions can be drawn from the differences in the wave functions for WJC-1 and WJC-2, observables must be calculated from these models. Electromagnetic observables (such as the quadrupole moment) will be sensitive to interaction currents, which will differ in each case, and might very well compensate for any difference arising from the wave functions. However, these calculations, which will be a subject for a future article, are beyond the scope of this work.

5. Asymptotic normalization and the D/S ratio

One observable that can be calculated immediately is the ratio of the asymptotic normalization constants for the S - and D -state wave functions. The results for the two models, together with the asymptotic normalization A_s , are summarized in Table IX. These quantities can be extracted from the expansions (3.51):

$$A_s = b_1^u A_1, \quad A_d = b_1^w B_1 R_1^2, \quad \eta = A_d/A_s. \quad (3.54)$$

Note that η is more accurately determined than A_s and that the values of η predicted by the two models are in agreement with each other (within errors) and also agree (to less than 2 standard deviations) with the results of Ref. [12]. The agreement is even better with most other experimental and theoretical results for η quoted in Ref. [12], which are all situated in the range between 0.0259 and 0.0272.

IV. CONCLUSIONS

In this article we present (Table I) effective range expansions for the 1S_0 and 3S_1 phase shifts and relativistic deuteron

TABLE IX. The asymptotic normalization A_s and D/S ratio η for the two models. Results are shown for variations of n around the central (best) values of 17 and 16 for WJC-1 and 30 and 18 for WJC-2, with errors based on the fluctuations. These may be compared to the recently determined $\eta = 0.0256(4)$ [12].

n	WJC-1		WJC-2	
	η	A_s	η	A_s
$n_0 - 2$	0.02625	0.86585	0.02635	0.87471
$n_0 - 1$	0.02623	0.86837	0.02639	0.87693
n_0	0.02619	0.86416	0.02630	0.87768
$n_0 + 1$	0.02623	0.86604	0.02643	0.87829
$n_0 + 2$	0.02622	0.86436	0.02614	0.87879
$n_0(\text{error})$	0.02619(4)	0.864(2)	0.0263(1)	0.8777(15)

wave functions based on the precision models WJC-1 and WJC-2 of Ref. I.

Since the effective range expansions emerge directly from the new high-precision phase shift analysis described in Ref. I, they can be regarded as an up-to-date, precision determination of the effective range parameters that should constrain any modern theory of the nuclear force, such as effective (or chiral effective) field theories. It does not matter that they were determined using the CST. Note that throughout this work we assumed equal masses for the proton and neutron; if their mass difference were taken into account we could expect corrections of about 0.07%.

The covariant deuteron wave functions presented here have four components: two that have a nonrelativistic analog and two (the P states) of purely relativistic origin. Convenient analytic representations of these wave functions are presented, with expansion coefficients given in Tables VII and VIII. These expansions (and subroutines that are available from the authors) make it easy and convenient to use these wave functions with any calculation. When these wave functions are used in a covariant theory with one nucleon off-shell they provide a precise description of the nonperturbative interactions that lead to the deuteron bound state. However, even in this case a completely consistent description requires that we include other contributions, such as final state interactions or interaction currents, that are also generated by the one-boson-exchange dynamics. That these effects cannot be ignored, even in the low-energy or momentum limit, is indicated by the size of the derivative of the kernel that contributes to the normalization [recall Eq. (3.39)]. This is about -5% for WJC-1 and -2% for WJC-2 and could, of course, be much larger at higher energies for some observables. In particular, the calculation of any electromagnetic property of the deuteron (including the magnetic and quadrupole moments) that does *not* include interaction currents can be expected *ab initio* to be in error by a few percentages. It is therefore quite possible that a careful CST calculation of the quadrupole moment could resolve the current $\sim 5\%$ discrepancy between theory and experiment, but this calculation must wait until the exchange currents have been accurately calculated.

For this reason it is also unclear how to incorporate the wave functions given in this article into a calculation which uses

dynamics different from the CST (nonrelativistic or another form of relativistic dynamics, including light-front). In such cases we are inclined to suggest that the best procedure is to renormalize the wave functions as shown in Eq. (3.40), which is also why our numerical wave functions are presented here in this normalization.

One observable that does *not* depend on the normalization of the wave functions is the asymptotic D/S ratio. Our results (Table IX) determine this ratio to an accuracy about 10 times smaller than the experimental accuracy and are in good agreement with measured values. Still, it would be interesting to review the theory that has gone into the analysis of the experiments that determine the D/S ratio and to see if there are any relativistic corrections previously overlooked.

ACKNOWLEDGMENTS

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APPENDIX A: CONSTRUCTION OF THE KERNEL

In this appendix we present details, omitted from Ref. I [1], of the construction of the NN kernel directly from the sum of Feynman amplitudes that define it. This method differs completely from that used in Ref. II [10]; it is simpler and more transparent, allowing changes in the kernel to be made more easily.

1. Review of definitions from Ref. I

In Ref. I the kernels are defined to be

$$\begin{aligned}
 & \mathbf{V}_{\lambda_1 \lambda_2, \lambda'_1 \lambda'_2}^{J \rho_1 \rho_2, \rho'_1 \rho'_2}(\delta_{p_0}, \delta_S) \\
 &= \frac{1}{2} \left\{ V_{\text{dir } \lambda_1 \lambda_2, \lambda'_1 \lambda'_2}^{J \rho_1 \rho_2, \rho'_1 \rho'_2}(p_0) + \delta_S V_{\text{dir } \lambda_1 \lambda_2, -\lambda'_1 -\lambda'_2}^{J \rho_1 \rho_2, \rho'_1 \rho'_2}(p_0) \right. \\
 & \quad \left. + \delta_{p_0} V_{\text{dir } \lambda_1 \lambda_2, \lambda'_1 \lambda'_2}^{J \rho_2 \rho_1, \rho'_1 \rho'_2}(-p_0) + \delta_{p_0} \delta_S V_{\text{dir } \lambda_1 \lambda_2, -\lambda'_1 -\lambda'_2}^{J \rho_2 \rho_1, \rho'_1 \rho'_2}(-p_0) \right\}, \tag{A1}
 \end{aligned}$$

where $\delta_{p_0} = (-1)^I \eta(\rho_T \delta_S)^\lambda$ (with $\rho_T = \rho_1 \rho_2 \rho'_1 \rho'_2$, $\lambda = \lambda_1 - \lambda_2$, and I the isospin) is the phase of \mathbf{V} under $p_0 \rightarrow -p_0$ and

$\rho_1 \leftrightarrow \rho_2$, and $\delta_S = \delta_P \rho'_1 \rho'_2 \eta$ (with $\delta_P = \pm 1$ the parity), and in both expressions $\eta = (-1)^{J-1}$ is a ubiquitous phase. *These linear combinations are very similar to (but differ from) those used previously in Ref. II.*

If particle 1 is on-shell, $\rho_1 = \rho'_1 = +$. Using the parity relation,

$$V_{\text{dir}\lambda_1\lambda_2,\lambda'_1\lambda'_2}^{J\rho_1\rho_2,\rho'_1\rho'_2}(\pm p_0) = \rho_T V_{\text{dir}\rightarrow\lambda_1\rightarrow\lambda_2,\rightarrow\lambda'_1\rightarrow\lambda'_2}^{J\rho_1\rho_2,\rho'_1\rho'_2}(\pm p_0), \quad (\text{A2})$$

the amplitudes (A1) can always be organized so that $\lambda_1 = \lambda'_1 = +1/2$, leaving the helicities of particle 2 and the phases δ_S and δ_{ρ_0} unconstrained. Hence there are $2^4 = 16$ independent kernels for each J, ρ_2 and ρ'_2 . Our particular choice of independent kernels is denoted v_i and defined in Table XI of Ref. I. Their behavior under parity and interchange symmetry for each combination of $\{\rho_2, \rho'_2\}$ is given in Table XII of Ref. I.

The evaluation of the matrix elements V^J is simplified by the fact that one can write the nucleon spinors as direct products of two two-component spinors, one of which contains all angle dependence, while the other describes the dependence on momentum and ρ spin. Recall that the nucleon spinors [from Eq. (E1) of Ref. I] are:

$$u_1^\rho(\mathbf{p}, \lambda) = N_\rho(p\lambda) \otimes \chi_\lambda(\theta) = \begin{cases} u(\mathbf{p}, \lambda) & \rho = + \\ v(-\mathbf{p}, \lambda) & \rho = - \end{cases} \quad (\text{A3})$$

$$u_2^\rho(\mathbf{p}, \lambda) = N_\rho(p\lambda) \otimes \chi_{-\lambda}(\theta) = \begin{cases} u(-\mathbf{p}, -\lambda) & \rho = + \\ v(\mathbf{p}, -\lambda) & \rho = - \end{cases},$$

where the ρ -space spinors are

$$N_+(p\lambda) = \begin{pmatrix} \cosh \frac{1}{2}\zeta \\ 2\lambda \sinh \frac{1}{2}\zeta \end{pmatrix} \quad N_-(p\lambda) = \begin{pmatrix} -2\lambda \sinh \frac{1}{2}\zeta \\ \cosh \frac{1}{2}\zeta \end{pmatrix}, \quad (\text{A4})$$

with $p = |\mathbf{p}|$ and $\tanh \zeta = p/E_p$. Note that, at large p ,

$$\cosh \frac{1}{2}\zeta = \sqrt{\frac{E_p + m}{2m}} \rightarrow \sqrt{\frac{p}{2m}}$$

$$\sinh \frac{1}{2}\zeta = \frac{p}{\sqrt{2m(E_p + m)}} \rightarrow \sqrt{\frac{p}{2m}}. \quad (\text{A5})$$

For momenta limited to the $\hat{x}\hat{z}$ plane, the spin 1/2 spinors are

$$\chi_{1/2}(\theta) = R_y(\theta) \begin{pmatrix} 1 \\ 0 \end{pmatrix} = \begin{pmatrix} \cos \frac{1}{2}\theta \\ \sin \frac{1}{2}\theta \end{pmatrix} \quad (\text{A6})$$

$$\chi_{-1/2}(\theta) = R_y(\theta) \begin{pmatrix} 0 \\ 1 \end{pmatrix} = \begin{pmatrix} -\sin \frac{1}{2}\theta \\ \cos \frac{1}{2}\theta \end{pmatrix}.$$

2. Extracting the angular integrals

Each term in the Feynman meson-exchange operator can be written in the generic form

$$\mathcal{V}_b(\theta, p_0) = \sum_{ij} g_i s_j \mathcal{D}_b(\theta, p_0) [\mathcal{O}^i(p_0) \otimes \mathcal{S}^j]_1^b \times [\mathcal{O}^i(p_0) \otimes \mathcal{S}^j]_2^b, \quad (\text{A7})$$

where \mathcal{O}^i are operators in the ρ -spin space, \mathcal{S}^j operators in the spin-space, g_i and s_j are constants, and the sums are over all operators needed to describe the meson exchange interaction. Here \mathcal{D}_b is the scalar part of the propagator

$$\mathcal{D}_b(\theta, p_0) = \frac{f_b(\Lambda_b, q)}{m_b^2 + |q(\theta, p_0)|^2}, \quad (\text{A8})$$

with $q(\theta, p_0)$ the four-momentum transferred by the exchanged meson and f_b the meson form factor. The operator $\mathcal{O}(p_0)$ acts in the 2×2 ρ -spin space, and \mathcal{S} operates in the 2×2 spin space, so $\mathcal{O} \otimes \mathcal{S}$ is a 4×4 matrix. Using this notation, the matrix elements of \mathcal{V} reduce to (suppressing the index b for simplicity)

$$V_{\text{dir}\lambda_1\lambda_2,\lambda'_1\lambda'_2}^{\rho_1\rho_2,\rho'_1\rho'_2}(\theta, \pm p_0) = \mathcal{O}_{\lambda_1\lambda_2,\lambda'_1\lambda'_2,\pm}^{\rho_1\rho'_1,\rho_2\rho'_2} \langle \lambda_1\lambda_2 | \lambda'_1\lambda'_2 \rangle \mathcal{D}(\theta, \pm p_0), \quad (\text{A9})$$

where

$$\mathcal{O}_{\lambda_1\lambda_2,\lambda'_1\lambda'_2,\pm}^{\rho_1\rho'_1,\rho_2\rho'_2} = \sum_i g_i [\bar{N}_{\rho_1}(p\lambda_1) \mathcal{O}^i(\pm p_0) N_{\rho'_1}(p'\lambda'_1)] \times [\bar{N}_{\rho_2}(p\lambda_2) \mathcal{O}^i(\pm p_0) N_{\rho'_2}(p'\lambda'_2)] \quad (\text{A10})$$

$$\langle \lambda_1\lambda_2 | \lambda'_1\lambda'_2 \rangle = \sum_j s_j [\chi_{\lambda_1}^\dagger(\theta) \mathcal{S}^j \chi_{\lambda'_1}(0)] [\chi_{-\lambda_2}^\dagger(\theta) \mathcal{S}^j \chi_{-\lambda'_2}(0)],$$

with $\bar{N} = N^\dagger \tau_3$ the Dirac conjugation in ρ -spin space, and use is made of the fact that the partial-wave amplitude may be calculated by aligning the initial momentum \mathbf{p}' in the $+z$ direction. Note that the sign of p_0 in the operators $\mathcal{O}(\pm p_0)$ is captured in the last subscript of the matrix elements \mathcal{O} , that the superscripts for \mathcal{O} group together the ρ -spin indices for particle 1, followed by the ρ -spin indices for particle 2, and that the matrix elements of \mathcal{S} do not depend on the ρ -spin of the states. We keep λ_1 unspecified in these formulas even though only the case $\lambda_1 = +$ needs to be considered. The linear combinations (A1) restrict $\lambda'_1 = +$, but the individual terms in the sum use both signs of λ'_1 , so it cannot be restricted in (A10). Recalling that the partial-wave projections are of the form

$$V_{\text{dir}\lambda_1\lambda_2,\lambda'_1\lambda'_2}^{J\rho_1\rho_2,\rho'_1\rho'_2}(p_0) = 2\pi \int_0^\pi \sin \theta d\theta d_{\lambda'\lambda}^J(\theta) \times V_{\text{dir}\lambda_1\lambda_2,\lambda'_1\lambda'_2}^{J\rho_1\rho_2,\rho'_1\rho'_2}(p, p'; P), \quad (\text{A11})$$

we see that each of the matrix elements $\langle \lambda_1\lambda_2 | \lambda'_1\lambda'_2 \rangle$ will be multiplied by the rotation function $d_{\lambda'\lambda}^J(\theta)$, where $\lambda = \lambda_1 - \lambda_2$ and $\lambda' = \lambda'_1 - \lambda'_2$. It turns out that the ρ -spin matrix elements generate an additional factor of $z = \cos \theta$, so matrix elements with an additional factor of z are needed; these are handled by multiplying each of the matrix elements by z^i , where $i = 0$ or 1. The 16 integrals we require are defined in Table X, with explicit expressions for these 16 matrix elements in terms of five independent integrals summarized in Table XI. The five independent integrals over Legendre polynomials and their derivatives are given in Table XII, with convenient identities showing how to reduce five other integrals to these five given in Table XIII.

The 16 generic integrals (for each value of ρ_1) defined in Tables XI and XII of Ref. I can then be written as products of the 16 angular-dependent terms given in Table X and Dirac

TABLE X. The 16 matrix elements $A_{\lambda_1 \lambda_2, \lambda'_1 \lambda'_2}^i = z^i \langle +\lambda_2 | \lambda'_1 \lambda'_2 \rangle d_{\lambda/\lambda}^J$ for $\lambda_1 = +$ and $i = 0$ or 1. Here $\lambda = \pm \frac{1}{2}$ is written as $\lambda = \pm$ and $z = \cos \theta$.

$A_{++++}^i = z^i \langle ++ ++ \rangle d_{00}^J = z^i \frac{1}{2}(1+z) d_{00}^J$	$A_{+---}^i = z^i \langle ++ -- \rangle d_{00}^J = -z^i \frac{1}{2}(1-z) d_{00}^J$
$A_{+-++}^i = z^i \langle +- ++ \rangle d_{01}^J = z^i \frac{1}{2} \sin \theta d_{01}^J$	$A_{+-- -}^i = z^i \langle +- -- \rangle d_{01}^J = z^i \frac{1}{2} \sin \theta d_{01}^J$
$A_{+++-}^i = z^i \langle ++ +- \rangle d_{10}^J = -z^i \frac{1}{2} \sin \theta d_{10}^J$	$A_{+--+}^i = z^i \langle ++ +- \rangle d_{-1,0}^J = z^i \frac{1}{2} \sin \theta d_{-1,0}^J$
$A_{+--+}^i = z^i \langle +- +- \rangle d_{11}^J = z^i \frac{1}{2}(1+z) d_{11}^J$	$A_{+--+}^i = z^i \langle +- +- \rangle d_{-1,1}^J = z^i \frac{1}{2}(1-z) d_{-1,1}^J$

matrix elements (defined below) as follows:

$$\begin{aligned}
v_5^{\rho_2 \rho'_2} &= \sum_i \left\{ \int_+ (\mathcal{O}_{++++}^{\rho_1+, \rho_2 \rho'_2, i} A_{++++}^i - \mathcal{O}_{+-+ -}^{\rho_1+, \rho_2 \rho'_2, i} A_{+-+ -}^i) \right. \\
&\quad \left. \pm \int_- (\mathcal{O}_{+++-}^{\rho_2+, \rho_1 \rho'_2} A_{+++-}^i - \mathcal{O}_{++- -}^{\rho_2+, \rho_1 \rho'_2} A_{++- -}^i) \right\} \\
v_6^{\rho_2 \rho'_2} &= \sum_i \left\{ \int_+ (\mathcal{O}_{+-+ -}^{\rho_1+, \rho_2 \rho'_2, i} A_{+-+ -}^i - \mathcal{O}_{+-+ -}^{\rho_1+, \rho_2 \rho'_2, i} A_{+-+ -}^i) \right. \\
&\quad \left. \mp \int_- (\mathcal{O}_{+++-}^{\rho_2+, \rho_1 \rho'_2} A_{+++-}^i - \mathcal{O}_{++- -}^{\rho_2+, \rho_1 \rho'_2} A_{++- -}^i) \right\} \\
v_7^{\rho_2 \rho'_2} &= \sum_i \left\{ \int_+ (\mathcal{O}_{++++}^{\rho_1+, \rho_2 \rho'_2, i} A_{++++}^i - \mathcal{O}_{+-+ -}^{\rho_1+, \rho_2 \rho'_2, i} A_{+-+ -}^i) \right. \\
&\quad \left. \pm \int_- (\mathcal{O}_{+++-}^{\rho_2+, \rho_1 \rho'_2} A_{+++-}^i - \mathcal{O}_{++- -}^{\rho_2+, \rho_1 \rho'_2} A_{++- -}^i) \right\} \\
v_8^{\rho_2 \rho'_2} &= \sum_i \left\{ \int_+ (\mathcal{O}_{+-+ -}^{\rho_1+, \rho_2 \rho'_2, i} A_{+-+ -}^i - \mathcal{O}_{+-+ -}^{\rho_1+, \rho_2 \rho'_2, i} A_{+-+ -}^i) \right. \\
&\quad \left. \mp \int_- (\mathcal{O}_{+++-}^{\rho_2+, \rho_1 \rho'_2} A_{+++-}^i - \mathcal{O}_{++- -}^{\rho_2+, \rho_1 \rho'_2} A_{++- -}^i) \right\} \\
v_9^{\rho_2 \rho'_2} &= \sum_i \left\{ \int_+ (\mathcal{O}_{++++}^{\rho_1+, \rho_2 \rho'_2, i} A_{++++}^i + \mathcal{O}_{+-+ -}^{\rho_1+, \rho_2 \rho'_2, i} A_{+-+ -}^i) \right. \\
&\quad \left. \pm \int_- (\mathcal{O}_{+++-}^{\rho_2+, \rho_1 \rho'_2} A_{+++-}^i + \mathcal{O}_{++- -}^{\rho_2+, \rho_1 \rho'_2} A_{++- -}^i) \right\} \\
v_{10}^{\rho_2 \rho'_2} &= \sum_i \left\{ \int_+ (\mathcal{O}_{+-+ -}^{\rho_1+, \rho_2 \rho'_2, i} A_{+-+ -}^i + \mathcal{O}_{+-+ -}^{\rho_1+, \rho_2 \rho'_2, i} A_{+-+ -}^i) \right. \\
&\quad \left. \pm \int_- (\mathcal{O}_{+++-}^{\rho_2+, \rho_1 \rho'_2} A_{+++-}^i + \mathcal{O}_{++- -}^{\rho_2+, \rho_1 \rho'_2} A_{++- -}^i) \right\}
\end{aligned}$$

$$\begin{aligned}
v_{11}^{\rho_2 \rho'_2} &= \sum_i \left\{ \int_+ (\mathcal{O}_{++++}^{\rho_1+, \rho_2 \rho'_2, i} A_{++++}^i + \mathcal{O}_{+-+ -}^{\rho_1+, \rho_2 \rho'_2, i} A_{+-+ -}^i) \right. \\
&\quad \left. \pm \int_- (\mathcal{O}_{+++-}^{\rho_2+, \rho_1 \rho'_2} A_{+++-}^i + \mathcal{O}_{++- -}^{\rho_2+, \rho_1 \rho'_2} A_{++- -}^i) \right\} \\
v_{12}^{\rho_2 \rho'_2} &= \sum_i \left\{ \int_+ (\mathcal{O}_{+-+ -}^{\rho_1+, \rho_2 \rho'_2, i} A_{+-+ -}^i + \mathcal{O}_{+-+ -}^{\rho_1+, \rho_2 \rho'_2, i} A_{+-+ -}^i) \right. \\
&\quad \left. \pm \int_- (\mathcal{O}_{+++-}^{\rho_2+, \rho_1 \rho'_2} A_{+++-}^i + \mathcal{O}_{++- -}^{\rho_2+, \rho_1 \rho'_2} A_{++- -}^i) \right\}, \tag{A12}
\end{aligned}$$

where the $+$ and $-$ integrals include the propagator

$$\int_{\pm} = \pi \int_0^{\pi} \sin \theta d\theta \mathcal{D}(\theta, \pm p_0) \tag{A13}$$

and, in each case, the upper sign goes with the upper index (of the two lower indices) and the lower sign with the lower index (of the two lower indices), so that, for example, v_1 is symmetric under $p_0 \rightarrow -p_0$ and v_5 is antisymmetric. The 16 ρ -spin matrix elements that enter into these expressions, all independent of θ , result from the fact that the general ρ -spin matrix elements \mathcal{O} given in Eq. (A10) are linear in z and can therefore be separated into two terms using the notation

$$\mathcal{O}_{\lambda_1 \lambda_2, \lambda'_1 \lambda'_2, \pm}^{\rho_1 \rho'_1, \rho_2 \rho'_2} = \mathcal{O}_{\lambda_1 \lambda_2, \lambda'_1 \lambda'_2, \pm}^{\rho_1 \rho'_1, \rho_2 \rho'_2, 0} + z \mathcal{O}_{\lambda_1 \lambda_2, \lambda'_1 \lambda'_2, \pm}^{\rho_1 \rho'_1, \rho_2 \rho'_2, 1}. \tag{A14}$$

The \mathcal{O}^1 's arise only from the exchange of vector mesons; all of the \mathcal{O} 's arising from the exchange of each meson will be computed below.

3. Particle momenta

In some applications there is the possibility that *both* particles in the final state could be off-shell (the possibility that

TABLE XI. Matrix elements of Table X expressed in terms of the angular functions from Table XII and a_6 , a_7 , a_8 , and a_9 from Table XIII. Note that $d_{-1,0}^J = -d_{1,0}^J$ and $d_{01}^J = -d_{10}^J$. Here $a_i = a_i^J(\theta)$. The integrals a_4 and a_5 are needed only for the A^1 and hence only for vector mesons. The matrix elements $\mathcal{A}_{ij,k}$ have $i, j = \{1, 4\}$ according to the code $1 = ++$, $2 = +-$, $3 = -+$, $4 = --$, and $k = 1$ for A^0 and $k = 2$ for A^1 .

$A_{++++}^0 = \mathcal{A}_{11,1} = \frac{1}{2}(1+z) d_{00}^J = \frac{1}{2}(a_1 + a_2)$	$A_{+---}^0 = \mathcal{A}_{14,1} = -\frac{1}{2}(1-z) d_{00}^J = -\frac{1}{2}(a_1 - a_2)$
$A_{+-++}^1 = \mathcal{A}_{11,2} = \frac{1}{2} z (1+z) d_{00}^J = \frac{1}{2}(a_2 + a_4)$	$A_{+-- -}^1 = \mathcal{A}_{14,2} = -\frac{1}{2} z (1-z) d_{00}^J = -\frac{1}{2}(a_2 - a_4)$
$A_{+-+ -}^0 = \mathcal{A}_{21,1} = \frac{1}{2} \sin \theta d_{01}^J = -\frac{1}{2} a_7$	$A_{+-- -}^0 = \mathcal{A}_{24,1} = \frac{1}{2} \sin \theta d_{01}^J = -\frac{1}{2} a_7$
$A_{+-+ -}^1 = \mathcal{A}_{21,2} = \frac{1}{2} z \sin \theta d_{01}^J = -\frac{1}{2} a_8$	$A_{+-- -}^1 = \mathcal{A}_{24,2} = \frac{1}{2} z \sin \theta d_{01}^J = -\frac{1}{2} a_8$
$A_{+++-}^0 = \mathcal{A}_{12,1} = -\frac{1}{2} \sin \theta d_{10}^J = -\frac{1}{2} a_7$	$A_{+--+}^0 = \mathcal{A}_{13,1} = \frac{1}{2} \sin \theta d_{-1,0}^J = -\frac{1}{2} a_7$
$A_{+++-}^1 = \mathcal{A}_{12,2} = -\frac{1}{2} z \sin \theta d_{10}^J = -\frac{1}{2} a_8$	$A_{+--+}^1 = \mathcal{A}_{13,2} = \frac{1}{2} z \sin \theta d_{-1,0}^J = -\frac{1}{2} a_8$
$A_{+--+}^0 = \mathcal{A}_{22,1} = \frac{1}{2}(1+z) d_{11}^J = \frac{1}{2}(a_6 + a_1)$	$A_{+--+}^0 = \mathcal{A}_{23,1} = \frac{1}{2}(1-z) d_{-1,1}^J = \frac{1}{2}(a_6 - a_1)$
$A_{+--+}^1 = \mathcal{A}_{22,2} = \frac{1}{2} z (1+z) d_{11}^J = \frac{1}{2}(a_9 + a_2)$	$A_{+--+}^1 = \mathcal{A}_{23,2} = \frac{1}{2} z (1-z) d_{-1,1}^J = \frac{1}{2}(a_9 - a_2)$

TABLE XII. The five independent angular integrals, $a_i^J(\theta)$, needed for this calculation. Four more frequently used linear combinations of these, denoted a_j , with $j = \{6, 9\}$, are given in Table XIII.

$a_1^J(\theta)$	$P_J(z)$
$a_2^J(\theta)$	$zP_J(z)$
$a_3^J(\theta)$	$P_{J-1}(z)$
$a_4^J(\theta)$	$z^2P_J(z)$
$a_5^J(\theta)$	$zP_{J-1}(z)$

both particles in the initial state are off-shell is not considered at this time). To allow for this possibility, we will write the final-state momenta in the center-of-mass system of the two-nucleon system in the general form

$$p_1 = \left\{ \frac{1}{2}W + p_0, \mathbf{p} \right\} = \left\{ x_0 E_p + (1 - x_0) \frac{1}{2}W, \mathbf{p} \right\} \quad (\text{A15})$$

$$p_2 = \left\{ \frac{1}{2}W - p_0, -\mathbf{p} \right\} = \left\{ -x_0 E_p + (1 + x_0) \frac{1}{2}W, -\mathbf{p} \right\}$$

with

$$p_0 = x_0(E_p - \frac{1}{2}W), \quad (\text{A16})$$

where x_0 is a dimensionless number varying between $-\infty$ and ∞ and W is the total two-body energy. Note that when $x_0 = 1$, particle 1 is on-shell (with energy E_p), while when $x_0 = -1$, particle 2 is on-shell. Hence changing the sign of x_0 is a convenient way to interchange the energies of particles 1 and 2 in the final state, and we may construct the $\pm p_0$ combinations of Eq. (A1) merely by changing the sign of x_0 . (A similar variable, y_0 , could be used for the initial state.)

4. Single particle ρ -spin matrix elements

We now turn our attention to the form of the ρ -spin matrix elements $\mathcal{O}_{\lambda_1 \lambda_2, \lambda'_1 \lambda'_2}^{\rho_1 \rho'_1, \rho_2 \rho'_2, i}$. In all but the simplest cases, these are best calculated by first constructing matrix elements on *each* nucleon line and then multiplying these together to get the total \mathcal{O} . The one-nucleon matrix elements are evaluated numerically by matrix multiplication, best described separately for each type of meson exchange. The total ρ -spin matrix elements for each meson will be summarized in the next subsection.

TABLE XIII. Useful identities involving the d functions. Here $P_j = P_j(z)$, $a_i = a_i^J(\theta)$, and the new combinations, a_6, a_7, a_8 , and a_9 are defined.

$\frac{1}{2} [(1+z)d_{11}^J - (1-z)d_{-1,1}^J] = d_{00}^J = P_J = a_1$
$\frac{1}{2} [(1+z)d_{11}^J + (1-z)d_{-1,1}^J] = a_6$
$= \frac{1}{J+1} [JzP_J + P_{J-1}] = \frac{1}{J+1} [Ja_2 + a_3]$
$\frac{1}{2} z [(1+z)d_{11}^J + (1-z)d_{-1,1}^J] = a_9$
$= \frac{1}{J+1} [Jz^2P_J + zP_{J-1}] = \frac{1}{J+1} [Ja_4 + a_5]$
$\sin \theta d_{10}^J = \sqrt{\frac{J}{J+1}} [zP_J - P_{J-1}]$
$= \sqrt{\frac{J}{J+1}} [a_2 - a_3] \equiv a_7$
$z \sin \theta d_{10}^J = \sqrt{\frac{J}{J+1}} [a_4 - a_5] \equiv a_8$

a. Scalar mesons

The calculation of the *on-shell* scalar exchange is very straightforward, with

$$\mathcal{O}_{\lambda_1 \lambda_2, \lambda'_1 \lambda'_2, \pm}^{\rho_1 \rho'_1, \rho_2 \rho'_2, 0, s} = -[\bar{N}_{\rho_1}(p\lambda_1) \mathbf{1} N_{\rho'_1}(p'\lambda'_1)] \times [\bar{N}_{\rho_2}(p\lambda_2) \mathbf{1} N_{\rho'_2}(p'\lambda'_2)] \quad (\text{A17})$$

$$\mathcal{O}_{\lambda_1 \lambda_2, \lambda'_1 \lambda'_2, \pm}^{\rho_1 \rho'_1, \rho_2 \rho'_2, 1, s} = 0,$$

where $\mathbf{1}$ is the unit matrix in 2×2 space, and the additional superscript s labels these \mathcal{O} 's as the contributions from scalar mesons. It is convenient to calculate the general matrix element

$$\mathbf{1}_i \equiv [\bar{N}_{\rho_i}(p\lambda_i) \mathbf{1} N_{\rho'_i}(p'\lambda'_i)], \quad (\text{A18})$$

and then construct all of the products (A17), avoiding as much algebra as possible. This also allows the off-shell matrix elements to be calculated without any extra work. In what follows we will generalize the notation of (A18):

$$\mathcal{O}_i \equiv [\bar{N}_{\rho_i}(p\lambda_i) \mathcal{O} N_{\rho'_i}(p'\lambda'_i)], \quad (\text{A19})$$

where \mathcal{O} is any 2×2 operator in the ρ -spin space. Explicitly, the Dirac matrices are written as a direct product of two 2×2 matrices (where the first matrix operates in the 2×2 ρ -spin space and the second in the spin space):

$$\mathbf{1} = \mathbf{1} \otimes \mathbf{1} \quad \gamma^0 = \tau^3 \otimes \mathbf{1} \quad \gamma^5 = \tau^1 \otimes \mathbf{1} \quad \gamma^i = i\tau^2 \otimes \sigma^i. \quad (\text{A20})$$

Hence, the ρ -spin matrices in the 2×2 spin independent part of the Dirac space are the familiar Pauli matrices.

We frequently encounter off-shell couplings, which in every case give factors of

$$(m - \not{p}) u^\rho(\mathbf{p}, \lambda) = (m + \boldsymbol{\gamma} \cdot \mathbf{p} - \gamma^0 p_0) u_i^\rho(\mathbf{p}, \lambda) = (\rho E_p - p_0) \gamma^0 u^\rho(\mathbf{p}, \lambda), \quad (\text{A21})$$

where we have used the Dirac equation for the (*always*) on-shell spinor $u^\rho(\mathbf{p}, \lambda)$ and recalled that $u^-(\mathbf{p}, \lambda) = v(-\mathbf{p}, \lambda)$ so its Dirac equation is

$$(m + \not{p}) v(-\mathbf{p}, \lambda) = (m + \gamma^0 E_p + \boldsymbol{\gamma} \cdot \mathbf{p}) v(-\mathbf{p}, \lambda) = 0.$$

This use of the Dirac equations allows us to replace the angular dependent $\boldsymbol{\gamma} \cdot \mathbf{p}$ term by an angular independent factor. (Similar steps work off-shell couplings in the final state.) Off-shell couplings therefore do not involve any new angular integrals, but they do require evaluation of a new 2×2 Dirac matrix element. Since a more general scalar Feynman operator is

$$\Lambda_s(p, p') = g_s \mathbf{1} - v_s \left[\frac{m - \not{p}}{2m} + \frac{m - \not{p}'}{2m} \right] \quad (\text{A22})$$

the *spin-independent* part of the matrix element in the general scalar case is then

$$\begin{aligned} \Lambda_{si}(p_i, p'_i) &\equiv \bar{u}^{\rho_i}(\mathbf{p}, \lambda_i) \Lambda_s(p_i, p'_i) u^{\rho'_i}(\mathbf{p}', \lambda'_i) \\ &= g_s \mathbf{1}_i - \frac{v_s}{2m} \tau_i^3 [\rho_i E_p - p_{i0} + \rho'_i E_{p'} - p'_{i0}] \\ &\equiv R_{\lambda_i \lambda'_i, s}^{\rho_i \rho'_i}(\pm p_0, \pm p'_0), \end{aligned} \quad (\text{A23})$$

where, in the last line, we *define* the first of the one-particle ρ -spin matrix elements needed in this calculation. The

TABLE XIV. Matrix elements $\langle \rho, \lambda | \mathcal{O}_n | \rho', \lambda' \rangle = \bar{N}_\rho(p\lambda) \mathcal{O}_n N_{\rho'}(p'\lambda') \equiv D_{\lambda\lambda',n}^{\rho\rho'}$ for various operators \mathcal{O}_n , where $n = \{1, 4\}$ with the correspondence $\tau^1 \rightarrow 1, i\tau^2 \rightarrow 2, \tau^3 \rightarrow 3$, and $\mathbf{1} \rightarrow 4$. The two-component ρ -spin spinors N_ρ were defined in Eq. (A4). In this table $c = \cosh \frac{1}{2}\zeta$, $c' = \cosh \frac{1}{2}\zeta'$, $s = 2\lambda \sinh \frac{1}{2}\zeta$, and $s' = 2\lambda' \sinh \frac{1}{2}\zeta'$, where ζ was defined below Eq. (A4).

$\langle +, \lambda \mathbf{1} +, \lambda' \rangle = -\langle -, \lambda \mathbf{1} -, \lambda' \rangle = cc' - ss'$	$\langle +, \lambda \mathbf{1} -, \lambda' \rangle = \langle -, \lambda \mathbf{1} +, \lambda' \rangle = -(cs' + sc')$
$\langle +, \lambda \tau^3 +, \lambda' \rangle = \langle -, \lambda \tau^3 -, \lambda' \rangle = cc' + ss'$	$\langle +, \lambda \tau^3 -, \lambda' \rangle = -\langle -, \lambda \tau^3 +, \lambda' \rangle = -(cs' - sc')$
$\langle +, \lambda \tau^1 +, \lambda' \rangle = \langle -, \lambda \tau^1 \lambda -, \lambda' \rangle = cs' - sc'$	$\langle +, \lambda \tau^1 -, \lambda' \rangle = -\langle -, \lambda \tau^1 +, \lambda' \rangle = cc' + ss'$
$\langle +, \lambda i\tau^2 +, \lambda' \rangle = -\langle -, \lambda i\tau^2 -, \lambda' \rangle = cs' + sc'$	$\langle +, \lambda i\tau^2 -, \lambda' \rangle = \langle -, \lambda i\tau^2 \lambda +, \lambda' \rangle = cc' - ss'$

definition of R_s restores explicit reference to the helicities λ_i and λ'_i implicitly contained in the matrix elements $\mathbf{1}_i$ and τ_i^3 , and replaces p_{i0} by $\pm p_{i0}$, with $p_{10} = W/2 + p_0$ and $p_{20} = W/2 - p_0$ (so that sign of p_0 is positive for particle 1 and negative for particle 2). We emphasize that the spin dependent angular part is unchanged by the off-shell couplings.

We encounter the off shell factors numerous times, so it is convenient to denote

$$\Delta E_i \equiv \rho_i E_p - p_{i0} \quad \Delta E'_i \equiv \rho'_i E_{p'} - p'_{i0}. \quad (\text{A24})$$

However, note that

$$\Delta E_1 \Delta E_2 = \Delta E'_1 \Delta E'_2 = 0 \quad (\text{A25})$$

because one of the two particles is always in shell.

b. Pseudoscalar mesons

The pseudoscalar Feynman operator is

$$\Lambda_p(p, p') = g_p \gamma^5 - g_p (1 - \lambda_p) \times \left\{ \left[\frac{m - \not{p}}{2m} \right] \gamma^5 + \gamma^5 \left[\frac{m - \not{p}'}{2m} \right] \right\}. \quad (\text{A26})$$

Using the same arguments, the pseudoscalar matrix elements become

$$\Lambda_{pi}(p_i, p'_i) = g_p \left\{ \tau_i^1 - \frac{(1 - \lambda_p)}{2m} i \tau_i^2 [\Delta E_i - \Delta E'_i] \right\} \equiv R_{\lambda_i \lambda'_i, p}^{\rho_i \rho'_i}(\pm p_0, \pm p'_0), \quad (\text{A27})$$

where we introduced the new matrix element R_p . This requires evaluation of the matrix elements of τ^1 and $i\tau^2$, but the spin-dependent matrix elements are the same as in the scalar case. For the evaluation of the matrix elements, see Table XIV.

c. Vector mesons

The vector matrix elements introduce new spin-dependent factors, and the terms \mathcal{O}^1 linear in z .

The first step in the reduction of the vector-meson exchange terms is to reduce the Pauli interaction term using the generalized Gordon decomposition:

$$\frac{i\sigma^{\mu\nu}(p - p')_\nu}{2m} = \gamma^\mu - \frac{(p + p')^\mu}{2m} - \left(\frac{m - \not{p}}{2m} \right) \gamma^\mu - \gamma^\mu \left(\frac{m - \not{p}'}{2m} \right). \quad (\text{A28})$$

Combining this result with the general definition of the vector Feynman operator

$$\Lambda_v^\mu(p, p') = g_v \left\{ \gamma^\mu + v_\nu \left(\left[\frac{m - \not{p}}{2m} \right] \gamma^\mu + \gamma^\mu \left[\frac{m - \not{p}'}{2m} \right] \right) + \frac{\kappa_v}{2m} i\sigma^{\mu\nu}(p - p')_\nu \right\} \quad (\text{A29})$$

gives

$$\Lambda_v^\mu(p, p') = g_v \left\{ \gamma^\mu (1 + \kappa_v) - \frac{\kappa_v}{2m} (p + p')^\mu + (v_\nu - \kappa_\nu) \left(\left[\frac{m - \not{p}}{2m} \right] \gamma^\mu + \gamma^\mu \left[\frac{m - \not{p}'}{2m} \right] \right) \right\} \rightarrow g_v \left\{ \gamma^\mu (1 + \kappa_v) - \frac{\kappa_v}{2m} (p + p')^\mu + \frac{(v_\nu - \kappa_\nu)}{2m} [\Delta E \gamma^0 \gamma^\mu + \gamma^\mu \gamma^0 \Delta E'] \right\}, \quad (\text{A30})$$

where the last line anticipates later use of the Dirac equation to reduce the operators $m - \not{p}$.

These two vertex operators are contracted with the spin-one meson propagator, which gives

$$\Lambda_{v1}^\mu(p_1, p'_1) \Lambda_{v2}^\nu(p_2, p'_2) \Delta_{\mu\nu} = \Lambda_{v1}^0(p_1, p'_1) \Lambda_{v2}^0(p_2, p'_2) - \Lambda_{v1}^j(p_1, p'_1) \Lambda_{v2}^j(p_2, p'_2) + \frac{\eta}{m_v^2} \left\{ \Lambda_{v1}^\mu(p_1, p'_1) (p_1 - p'_1)_\mu \Lambda_{v2}^\nu(p_2, p'_2) (p_2 - p'_2)_\nu \right\} \quad (\text{A31})$$

where the factor $\eta = 1$ is included solely to keep track of the effect of the $q^\mu q^\nu / m_v^2$ in the meson propagator (note that the sign follows from $-q^\mu q^\nu = (p_1 - p'_1)^\mu (p_2 - p'_2)^\nu$). To simplify, note that

$$B_i \equiv \Lambda_{vi}^\nu(p_i, p'_i) (p_i - p'_i)_\nu = g_v \left\{ \mathbf{1}_i v_\nu \frac{(p_i^2 - p_i'^2)}{2m} - \tau_i^3 (1 + v_\nu) (\Delta E_i - \Delta E'_i) \right\} \equiv R_{\lambda_i \lambda'_i, v2}^{\rho_i \rho'_i}(\pm p_0, \pm p'_0) \quad (\text{A32})$$

$$\Lambda_{vi}^0(p_i, p'_i) = g_v \left\{ \tau_i^3 (1 + \kappa_v) + \mathbf{1}_i \left[\frac{v_\nu}{2m} (\Delta E_i + \Delta E'_i) - \frac{\kappa_v}{2m} (\rho_i E_p + \rho'_i E_{p'}) \right] \right\} \equiv R_{\lambda_i \lambda'_i, v1}^{\rho_i \rho'_i}(\pm p_0, \pm p'_0),$$

where we introduced the new matrix elements R_{v1} and R_{v2} .

The three-vector part introduces the spin-dependent operators σ^i and angle-dependent terms from the factor $(p + p')^j$. These are first separated into two terms,

$$\Lambda_{vi}^j(p_i, p'_i) = A_i \times \sigma_i^j - \mathbf{1}_i \frac{g_v \kappa_v}{2m} (p_i + p'_i)^j, \quad (\text{A33})$$

where we remind the reader that the subscript i labels the particle number (1 or 2), and A_i is the spin-independent coefficient of the spin-dependent operator σ^j , and the spin-dependent part of the second term is the identity, suppressed in the expression. The coefficient A_i is

$$A_i = g_v \left\{ i \tau_i^2 (1 + \kappa_v) + \tau_i^1 \frac{(\nu_v - \kappa_v)}{2m} (\Delta E_i - \Delta E'_i) \right\} \\ \equiv R_{\lambda_i \lambda'_i, \nu_3}^{\rho_i \rho'_i} (\pm p_0, \pm p'_0), \quad (\text{A34})$$

which defines the matrix element R_{ν_3} .

In calculating the operator $\Lambda_{v1}^j(p_1, p'_1) \Lambda_{v2}^j(p_2, p'_2)$ we encounter the cross terms

$$\boldsymbol{\sigma}_1 \cdot (\mathbf{p}_2 + \mathbf{p}'_2) = -\boldsymbol{\sigma}_1 \cdot (\mathbf{p}_1 + \mathbf{p}'_1) = -2(\lambda_1 p + \lambda'_1 p'), \quad (\text{A35})$$

where the evaluation has been carried out in the center-of-mass system and use was made of the fact that the helicity eigenstates satisfy the eigenvalue conditions

$$\boldsymbol{\sigma} \cdot \mathbf{p} |\lambda\rangle = 2\lambda p |\lambda\rangle. \quad (\text{A36})$$

(For incoming particle 2, for example, $|\frac{1}{2}\rangle$ is a state with spin down in the \hat{z} direction and \mathbf{p}'_2 is in the $-\hat{z}$ direction, so $\boldsymbol{\sigma} \cdot \mathbf{p}'_2 |\frac{1}{2}\rangle = -p\sigma^3 |\frac{1}{2}\rangle = p |\frac{1}{2}\rangle$.) Finally, note that

$$(\mathbf{p}_1 + \mathbf{p}'_1) \cdot (\mathbf{p}_2 + \mathbf{p}'_2) = -(\mathbf{p}_1 + \mathbf{p}'_1)^2 \\ = -(\mathbf{p}^2 + \mathbf{p}'^2 + 2pp'z), \quad (\text{A37})$$

where we use \mathbf{p}^2 , p^2 , and p to denote the square of the three-vector, the four-vector, and the magnitude of the three-vector, respectively. Combining (A33)–(A37) gives the following:

$$\Lambda_{v1}^j(p_1, p'_1) \Lambda_{v2}^j(p_2, p'_2) \\ = A_1 A_2 \langle \lambda_1 \lambda_2 | \boldsymbol{\sigma}_1 \cdot \boldsymbol{\sigma}_2 | \lambda'_1 \lambda'_2 \rangle \\ + \left\{ \frac{g_v \kappa_v}{m} [A_1 \mathbf{1}_2 (\lambda_1 p + \lambda'_1 p') + A_2 \mathbf{1}_1 (\lambda_2 p + \lambda'_2 p')] \right. \\ \left. - \left(\frac{g_v \kappa_v}{2m} \right)^2 \mathbf{1}_1 \mathbf{1}_2 (\mathbf{p}^2 + \mathbf{p}'^2 + 2pp'z) \right\} \langle \lambda_1 \lambda_2 | \lambda'_1 \lambda'_2 \rangle, \quad (\text{A38})$$

where, for clarity, we have included the term $\langle \lambda_1 \lambda_2 | \lambda'_1 \lambda'_2 \rangle$ that multiplies the terms in the $\{\}$; this term is not part of the ρ -spin matrix element and will be removed once the effect of the operator $\boldsymbol{\sigma}_1 \cdot \boldsymbol{\sigma}_2$ has been expressed using the relation

$$\langle \lambda_1 \lambda_2 | \boldsymbol{\sigma}_1 \cdot \boldsymbol{\sigma}_2 | \lambda'_1 \lambda'_2 \rangle \rightarrow \mathcal{T} \langle \lambda_1 \lambda_2 | \lambda'_1 \lambda'_2 \rangle, \quad (\text{A39})$$

where \mathcal{T} will be determined shortly. The extra $z = \cos \theta$ dependence in the last term of (A38) is the origin of the special linear z -dependent terms described in Eq. (A14) above. In our model, only vector-meson exchanges contribute such terms.

To determine \mathcal{T} correctly note that the helicity states $|+\rangle$ or $|-\rangle$ correspond to the *spin* states $|\uparrow\rangle$ or $|\downarrow\rangle$, respectively, and hence are spin triplet states with $T = 1$. The same argument works for both initial or final states. This leaves

only the combinations with equal helicities in *both* the initial and final states: $\lambda_1 = \lambda_2 = \pm$ and $\lambda'_1 = \lambda'_2 = \pm$. Looking at Eq. (A12), the only potentials affected are v_1 , v_5 , v_9 , and v_{13} , and using the results

$$\boldsymbol{\sigma}_1 \cdot \boldsymbol{\sigma}_2 |++\rangle = -|++\rangle + 2|--\rangle \\ \boldsymbol{\sigma}_1 \cdot \boldsymbol{\sigma}_2 |--\rangle = -|--\rangle + 2|++\rangle \quad (\text{A40})$$

the \mathcal{T} term in these kernels leads to the decomposition

$$v_i^{\rho_2 \rho'_2} \sim (\mathcal{O}_{++}^{++}, \rho_2 \rho'_2, i} \mathcal{T} A_{++}^i \pm \mathcal{O}_{++}^{+-}, \rho_2 \rho'_2, i} \mathcal{T} A_{++}^i) \\ = \frac{1}{2} (\mathcal{O}_{++}^{++}, \rho_2 \rho'_2, i} \pm \mathcal{O}_{++}^{+-}, \rho_2 \rho'_2, i}) (A_{++}^i + A_{++}^i) \\ + \frac{1}{2} S (\mathcal{O}_{++}^{++}, \rho_2 \rho'_2, i} \mp \mathcal{O}_{++}^{+-}, \rho_2 \rho'_2, i}) (A_{++}^i - A_{++}^i), \quad (\text{A41})$$

where $S = -3$. If the \mathcal{T} operator is absent, the same expression holds, but with $S = 1$, so this decomposition can be used for both cases. With this relation it is not necessary to explicitly calculate the matrix elements $\langle \lambda_1 \lambda_2 | \boldsymbol{\sigma}_1 \cdot \boldsymbol{\sigma}_2 | \lambda'_1 \lambda'_2 \rangle$.

d. Axial vector mesons

Only the simplest possible coupling was used for the axial vector mesons, namely

$$\Lambda_{ai}^\mu(p_i, p'_i) = g_a \gamma_i^5 \gamma_i^\mu, \quad (\text{A42})$$

and this was contracted with the $g_{\mu\nu}$ tensor, giving

$$\Lambda_{a1}^\mu(p_1, p'_1) \Lambda_{a2}^\nu(p_2, p'_2) g_{\mu\nu} = \Lambda_{a1}^0 \Lambda_{a2}^0 \langle \lambda_1 \lambda_2 | \lambda'_1 \lambda'_2 \rangle \\ - C_1 C_2 \langle \lambda_1 \lambda_2 | \boldsymbol{\sigma}_1 \cdot \boldsymbol{\sigma}_2 | \lambda'_1 \lambda'_2 \rangle, \quad (\text{A43})$$

with

$$\Lambda_{ai}^0 = -i g_a \tau_i^2 \equiv R_{\lambda_i \lambda'_i, a1}^{\rho_i \rho'_i} (\pm p_0, \pm p'_0) \\ C_i = -g_a \tau_i^3 \equiv R_{\lambda_i \lambda'_i, a2}^{\rho_i \rho'_i} (\pm p_0, \pm p'_0), \quad (\text{A44})$$

which defines the last two ρ -spin matrix elements, R_{a1} and R_{a2} .

5. Two-particle ρ -spin matrix elements

The two-particle ρ -spin matrix elements are the ones defined in Eq. (A14). They are constructed from the helicity matrix elements $D_{\lambda \lambda', n}^{\rho \rho'}$ given in Table XIV, and the one-nucleon matrix elements R_b defined in Eqs. (A23), (A27), (A32), (A34), and (A44) above. In constructing the matrix elements, note that $p_0 \rightarrow -p_0$ corresponds to the interchange of $p_{10} \leftrightarrow p_{20}$, and similarly for p'_0 . The results, with all indices included, are as follows:

scalar, from Eq. (A23):

$$\mathcal{O}_{\lambda_1 \lambda_2, \lambda'_1 \lambda'_2, \pm}^{\rho_1 \rho'_1, \rho_2 \rho'_2, 0, s} = -R_{\lambda_1 \lambda'_1, s}^{\rho_1 \rho'_1} (\pm p_0, p'_0) R_{\lambda_2 \lambda'_2, s}^{\rho_2 \rho'_2} (\mp p_0, -p'_0) \\ \mathcal{O}_{\lambda_1 \lambda_2, \lambda'_1 \lambda'_2, \pm}^{\rho_1 \rho'_1, \rho_2 \rho'_2, 1, s} = 0 \quad (\text{A45})$$

pseudoscalar, from Eq. (A27):

$$\mathcal{O}_{\lambda_1 \lambda_2, \lambda'_1 \lambda'_2, \pm}^{\rho_1 \rho'_1, \rho_2 \rho'_2, 0, p} = R_{\lambda_1 \lambda'_1, p}^{\rho_1 \rho'_1} (\pm p_0, p'_0) R_{\lambda_2 \lambda'_2, p}^{\rho_2 \rho'_2} (\mp p_0, -p'_0) \\ \mathcal{O}_{\lambda_1 \lambda_2, \lambda'_1 \lambda'_2, \pm}^{\rho_1 \rho'_1, \rho_2 \rho'_2, 1, p} = 0 \quad (\text{A46})$$

vector, from Eqs. (A32), (A34), and (A38):

$$\begin{aligned} \mathcal{O}_{\lambda_1\lambda_2,\lambda'_1\lambda'_2,\pm}^{\rho_1\rho'_1,\rho_2\rho'_2,0,v} &= R_{\lambda_1\lambda'_1,v_1}^{\rho_1\rho'_1}(\pm p_0, p'_0)R_{\lambda_2\lambda'_2,v_1}^{\rho_2\rho'_2}(\mp p_0, -p'_0) \\ &+ \frac{\eta}{m^2}R_{\lambda_1\lambda'_1,v_2}^{\rho_1\rho'_1}(\pm p_0, p'_0)R_{\lambda_2\lambda'_2,v_2}^{\rho_2\rho'_2}(\mp p_0, -p'_0) \\ &- \mathcal{T}R_{\lambda_1\lambda'_1,v_3}^{\rho_1\rho'_1}(\pm p_0, p'_0)R_{\lambda_2\lambda'_2,v_3}^{\rho_2\rho'_2}(\mp p_0, -p'_0) \\ &+ \left(\frac{g_v\kappa_v}{2m}\right)^2(\mathbf{p}^2 + \mathbf{p}'^2)D_{\lambda_1\lambda'_1,4}^{\rho_1\rho'_1}D_{\lambda_2\lambda'_2,4}^{\rho_2\rho'_2} \\ &- \frac{g_v\kappa_v}{m}[(\lambda_1 p + \lambda'_1 p')R_{\lambda_1\lambda'_1,v_3}^{\rho_1\rho'_1}(\pm p_0, p'_0)D_{\lambda_2\lambda'_2,4}^{\rho_2\rho'_2} \\ &+ (\lambda_2 p + \lambda'_2 p')R_{\lambda_2\lambda'_2,v_3}^{\rho_2\rho'_2}(\mp p_0, -p'_0)D_{\lambda_1\lambda'_1,4}^{\rho_1\rho'_1}] \\ \mathcal{O}_{\lambda_1\lambda_2,\lambda'_1\lambda'_2,\pm}^{\rho_1\rho'_1,\rho_2\rho'_2,1,v} &= 2pp' \left(\frac{g_v\kappa_v}{2m}\right)^2 D_{\lambda_1\lambda'_1,4}^{\rho_1\rho'_1}D_{\lambda_2\lambda'_2,4}^{\rho_2\rho'_2} \end{aligned} \quad (\text{A47})$$

(in these expressions we used the fact that D_4 does not depend on p_0), and

axial vector, from Eqs. (A43) and (A44):

$$\begin{aligned} \mathcal{O}_{\lambda_1\lambda_2,\lambda'_1\lambda'_2,\pm}^{\rho_1\rho'_1,\rho_2\rho'_2,0,a} &= g_b^2 \{ D_{\lambda_1\lambda'_1,2}^{\rho_1\rho'_1}D_{\lambda_2\lambda'_2,2}^{\rho_2\rho'_2} - \mathcal{T}D_{\lambda_1\lambda'_1,3}^{\rho_1\rho'_1}D_{\lambda_2\lambda'_2,3}^{\rho_2\rho'_2} \} \\ \mathcal{O}_{\lambda_1\lambda_2,\lambda'_1\lambda'_2,\pm}^{\rho_1\rho'_1,\rho_2\rho'_2,1,a} &= 0 \end{aligned} \quad (\text{A48})$$

Once again we note that the D 's do not depend on p_0 .

Treatment of the photon. For the photon we assume that $F_{1n} = 0$ and that all other form factors are equal: $F_{2n} = F_{1p} = F_{2p} = F_D$. The photon is an isovector-vector exchange, obtained using the following couplings (with $\alpha \simeq 0.007297$):

$$\begin{aligned} g^2 &\rightarrow 0 \quad \text{no } e^2 \text{ term} \\ g^2\kappa &\rightarrow e^2\kappa_n \rightarrow -\alpha(1.913) = -0.01396 \equiv G_1 \\ g^2\kappa^2 &\rightarrow e^2\kappa_p\kappa_n \rightarrow -\alpha(1.913)(1.793) = -0.02503 \equiv G_2. \end{aligned} \quad (\text{A49})$$

These couplings require a special construction and cannot easily be incorporated into the general vector formulas above.

As a first step, define artificial couplings designed to reproduce the ratio $G_2/G_1 = \kappa_{\text{eff}} = \kappa_p = 1.7930$ and $g_{\text{eff}}^2 = G_1/\kappa_p = G_1^2/G_2 = -0.00779$. The upshot of this is to use the proton anomalous moment for κ and a fictitious charge of $g_{\text{eff}}^2 = -0.00779$ so $g_{\text{eff}}^2\kappa_p = G_1$ and $g_{\text{eff}}^2\kappa_p^2 = G_2$ are reproduced. Then, we will factor out the coupling g_{eff} from the R 's so it occurs as an overall multiplicative factor with the correct (negative) sign.

Next, note that all terms in the photon exchange must involve at least one factor of κ , so $B \sim R_{v_2}$ does not contribute. The R_{v_1} and R_{v_3} squared terms must have the g_{eff}^2 terms subtracted. Noting that $v = 0$ for the photon, and *redefining the photon R 's so $g_v = 1$ and $\kappa_v = \kappa_p$* , this gives

Photon:

$$\begin{aligned} \mathcal{O}_{\lambda_1\lambda_2,\lambda'_1\lambda'_2,\pm}^{\rho_1\rho'_1,\rho_2\rho'_2,0,\gamma} &= g_{\text{eff}}^2 (R_{\lambda_1\lambda'_1,\gamma_1}^{\rho_1\rho'_1}(\pm p_0, p'_0)R_{\lambda_2\lambda'_2,\gamma_1}^{\rho_2\rho'_2}(\mp p_0, -p'_0) \\ &- D_{\lambda_1\lambda'_1,3}^{\rho_1\rho'_1}D_{\lambda_2\lambda'_2,3}^{\rho_2\rho'_2}) - \mathcal{T}g_{\text{eff}}^2 (R_{\lambda_1\lambda'_1,\gamma_3}^{\rho_1\rho'_1}(\pm p_0, p'_0) \\ &\times R_{\lambda_2\lambda'_2,\gamma_3}^{\rho_2\rho'_2}(\mp p_0, -p'_0) - D_{\lambda_1\lambda'_1,2}^{\rho_1\rho'_1}D_{\lambda_2\lambda'_2,2}^{\rho_2\rho'_2}) \\ &+ G_2 \frac{(\mathbf{p}^2 + \mathbf{p}'^2)}{4m^2} D_{\lambda_1\lambda'_1,4}^{\rho_1\rho'_1}D_{\lambda_2\lambda'_2,4}^{\rho_2\rho'_2} \end{aligned}$$

$$\begin{aligned} &- \frac{G_1}{m} [(\lambda_1 p + \lambda'_1 p')R_{\lambda_1\lambda'_1,\gamma_3}^{\rho_1\rho'_1}(\pm p_0, p'_0)D_{\lambda_2\lambda'_2,4}^{\rho_2\rho'_2} \\ &+ (\lambda_2 p + \lambda'_2 p')R_{\lambda_2\lambda'_2,\gamma_3}^{\rho_2\rho'_2}(\mp p_0, -p'_0)D_{\lambda_1\lambda'_1,4}^{\rho_1\rho'_1}] \end{aligned}$$

$$\mathcal{O}_{\lambda_1\lambda_2,\lambda'_1\lambda'_2,\pm}^{\rho_1\rho'_1,\rho_2\rho'_2,1,\gamma} = G_2 \frac{pp'}{2m^2} D_{\lambda_1\lambda'_1,4}^{\rho_1\rho'_1}D_{\lambda_2\lambda'_2,4}^{\rho_2\rho'_2}, \quad (\text{A50})$$

where the redefined R 's are [recalling Eqs. (A32) for R_{v_1} and (A34) for R_{v_3} with $v = 0$ and $g_v = 1$]

$$\begin{aligned} R_{\lambda\lambda',\gamma_1}^{\rho\rho'}(x_0, y_0) &= (1 + \kappa_p)D_{\lambda_i\lambda'_i,3}^{\rho_i\rho'_i} - \frac{\kappa_p}{2m}D_{\lambda\lambda',4}^{\rho\rho'}(\rho E_p + \rho' E_{p'}) \\ R_{\lambda\lambda',\gamma_3}^{\rho\rho'}(x_0, y_0) &= (1 + \kappa_p)D_{\lambda\lambda',2}^{\rho\rho'} \\ &- \frac{\kappa_p}{2m}D_{\lambda\lambda',1}^{\rho\rho'}[\Delta E_\rho(x_0) - \Delta E'_{\rho'}(y_0)]. \end{aligned} \quad (\text{A51})$$

6. Isospin factor

For the exchange of isovector mesons, one must evaluate the factor $\tau_1 \cdot \tau_2$. This operator preserves the isospin symmetry of the state; therefore its value is the same for each of the 16 kernels (or 4 when $J = 0$) that make up the coupled array describing the singlet (S), triplet (T), or coupled (C) channels: namely it must equal unity on isovector np states and -3 on isoscalar np states. To determine the correct value, look at the leading amplitudes (those which are symmetric under the change in sign of p_0 and hence survive when both particles are on shell) and note that, for these amplitudes, the Pauli principle requires that $L + S + I$ be odd. For singlet and triplet states, $J = L$ while for coupled states $J = L \pm 1$. Hence we have the identification:

$$\begin{aligned} \text{singlet and coupled: } &\begin{cases} J \text{ even} &\rightarrow I = 1 \\ J \text{ odd} &\rightarrow I = 0 \end{cases} \\ \text{triplet: } &\begin{cases} J \text{ even} &\rightarrow I = 0 \\ J \text{ odd} &\rightarrow I = 1. \end{cases} \end{aligned} \quad (\text{A52})$$

These are summarized by the following:

$$\begin{aligned} \text{singlet and coupled: } &\tau_1 \cdot \tau_2 = -1 + 2(-1)^J \\ \text{triplet: } &\tau_1 \cdot \tau_2 = -1 - 2(-1)^J. \end{aligned} \quad (\text{A53})$$

To determine the correct isospin factor for each $v_i^{\rho_1\rho'_1}$ the channel to which it contributes must first be determined by examining the matrices Eqs. (E43)–(E48) of Ref. I. Inspection shows that the kernels v_1, v_3, v_9 , and v_{11} do not contribute to the triplet channels, while kernels v_5, v_7, v_{13} , and v_{15} contribute *only* to triplet channels. The rest contribute to either, depending on their ρ spin. In particular, contributions to the triplet channels come from (i) the $++$ sector of v_6 and v_8 , (ii) the $+-$ sector of v_{10} and v_{12} , (iii) the $-+$ sector of v_2 and v_4 , and (iv) the $--$ sector of v_{14} and v_{16} , with the other sectors of these kernels contributing to either singlet or coupled.

7. Charge symmetry breaking induced by pion exchange

In order to include important violations of charge symmetry and to better describe the important pion exchange, both models treat the charged π^\pm and the neutral π^0 as independent

mesons, with their masses fixed to their experimental values. Model WJC-1 allows the couplings of the pions to be varied independently while model WJC-2 assumes that the couplings (written in terms of g and not f) of both are the same.

In addition to the couplings and masses, the correct isospin factor for the exchange of each pion must be worked out. There are two ways to do this. The operator $\tau_1 \cdot \tau_2$ can be decomposed into

$$\tau_1 \cdot \tau_2 = \tau_1^3 \tau_2^3 + 2(\tau_1^+ \tau_2^- + \tau_1^- \tau_2^+), \quad (\text{A54})$$

where $\tau^\pm = (\tau^1 \pm \tau^2)/2$ are the raising and lowering operators (normalized to 1 and not $\sqrt{2}$ as is more commonly done) and the first term arises from π^0 exchange and the second from π^\pm exchange. Noting that

$$\begin{aligned} \tau_1^3 \tau_2^3 |pn\rangle &= -|pn\rangle \\ \tau_1^3 \tau_2^3 |np\rangle &= -|np\rangle \\ 2(\tau_1^+ \tau_2^- + \tau_1^- \tau_2^+) |pn\rangle &= 2|np\rangle \\ 2(\tau_1^+ \tau_2^- + \tau_1^- \tau_2^+) |np\rangle &= 2|pn\rangle \end{aligned} \quad (\text{A55})$$

we see that

$$\begin{aligned} \tau_1 \cdot \tau_2 |_{\pi^0} &\rightarrow -1 \\ \tau_1 \cdot \tau_2 |_{\pi^\pm} &\rightarrow \begin{cases} 2 & I = 1 \text{ states} \\ -2 & I = 0 \text{ states.} \end{cases} \end{aligned} \quad (\text{A56})$$

These factors can also be obtained by using the fact that isospin invariance gives couplings for $\{p \rightarrow p\pi^0, n \rightarrow n\pi^0, p \rightarrow n\pi^+, n \rightarrow p\pi^-\}$ in the ratios $\{1, -1, \sqrt{2}, \sqrt{2}\}$.

The implication of (A56) for π^\pm couplings is to replace the isospin factors from (A53) by

$$\tau_1 \cdot \tau_2 |_{\pi^\pm} \rightarrow \begin{cases} 2(-1)^J & \text{S and C states} \\ -2(-1)^J & \text{T states.} \end{cases} \quad (\text{A57})$$

APPENDIX B: EVALUATION OF ANGULAR INTEGRALS

1. Angular integrals

The angular integrals needed for this problem are defined to be

$$\mathcal{A}_{\lambda_1 \lambda_2, \lambda'_1 \lambda'_2}^{i, a, \pm} = \int_{-1}^1 dz A_{\lambda_1 \lambda_2, \lambda'_1 \lambda'_2}^i \left[\theta, x_0 \left(E_p - \frac{1}{2} W \right) \right] \mathcal{D}_b \quad (\text{B1})$$

where \mathcal{D}_b is the dressed propagator (including form factor) for meson b , defined in Eq. (A8), and the four-momentum transfer for the direct (+) and alternating (−) terms is

$$\begin{aligned} q^2(x_0) &= \left[x_0 \left(E_p - \frac{1}{2} W \right) - y_0 \left(E_{p'} - \frac{1}{2} W \right) \right]^2 \\ &\quad - p^2 - p'^2 + 2pp'z, \end{aligned} \quad (\text{B2})$$

where in this article the initial state always has particle 1 on-shell, so $y_0 = 1$. These integrals are assembled from the matrix elements of Table X and the functions of Table XIII, leading to the identities of Table XI. All of the integrals \mathcal{A} are recalculated for each J , p , x_0 , p' , and y_0 , so these arguments are not included in the list of subscripts denoting the dependencies of \mathcal{A} .

2. Mapping for the direct terms

For the direct term the meson propagator is strongly peaked in the forward direction ($z = 1$). For the recent NN fits, the meson propagators have the form

$$\mathcal{D}_b(z) \equiv F(z) = \frac{m_b^2}{m_b^2 + A|z_0 - z|} \left[\frac{\Lambda_b^2}{\Lambda_b^2 + A|z_0 - z|} \right]^4, \quad (\text{B3})$$

where m_b and Λ_b are the meson mass and form factor mass of meson b , respectively, and

$$\begin{aligned} A &= 2pp' \\ z_0 &= \frac{p^2 + p'^2 - (E_p - E_{p'})^2}{2pp'} \equiv 1 + \xi, \end{aligned} \quad (\text{B4})$$

with $\xi = 0$ if the incoming and outgoing momenta (p and p') are equal in magnitude, and $\xi > 0$ when $p \neq p'$. [Since $z_0 \geq 1$, the absolute value in (B3) can be ignored.] At the largest momenta used in the numerical solutions, $A \simeq 1800 \text{ GeV}^2$, compared to (for example) the pion mass squared $m_\pi^2 \simeq 2 \times 10^{-4} \text{ GeV}^2$. Under these conditions, numerical evaluation of the integral, with its narrow and large forward peak is difficult.

To carry out the integration numerically, we apply a Gauss-Legendre quadrature rule to discretize the integration variable z into a grid of Gauss points distributed between -1 and $+1$. However, it is convenient to map the integration variable $z \rightarrow y$, where y will vary slowly in the region where $F(z)$ is large. One possible function that will accomplish this, and still remain bounded between -1 and 1 is

$$z(y) = \frac{\tanh[y/(b_0 - 1)]}{\tanh[1/(b_0 - 1)]}. \quad (\text{B5})$$

When b_0 is close to 1, this function spreads out the intervals near $z = \pm 1$, so more Gauss points may be placed in the region of the forward peak. The mapping function, for four different values of b_0 , is shown in Fig. 10. Using these mappings, the propagator becomes a function of the integration variable y

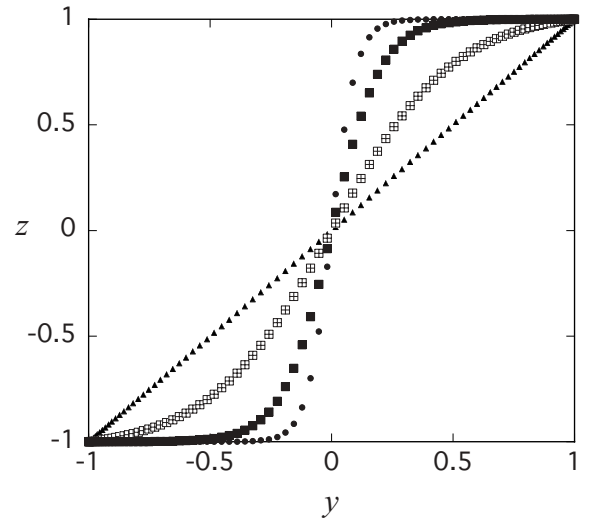


FIG. 10. Plot of the mapping function $z(y)$ defined in Eq. (B5). The four curves, each with 90 Gauss points, correspond to $b_0 = 1.1$ (small solid triangles), $b_0 = 1.5$ (open boxes), $b_0 = 1.2$ (solid boxes), and $b_0 = 1.1$ (small solid circles).

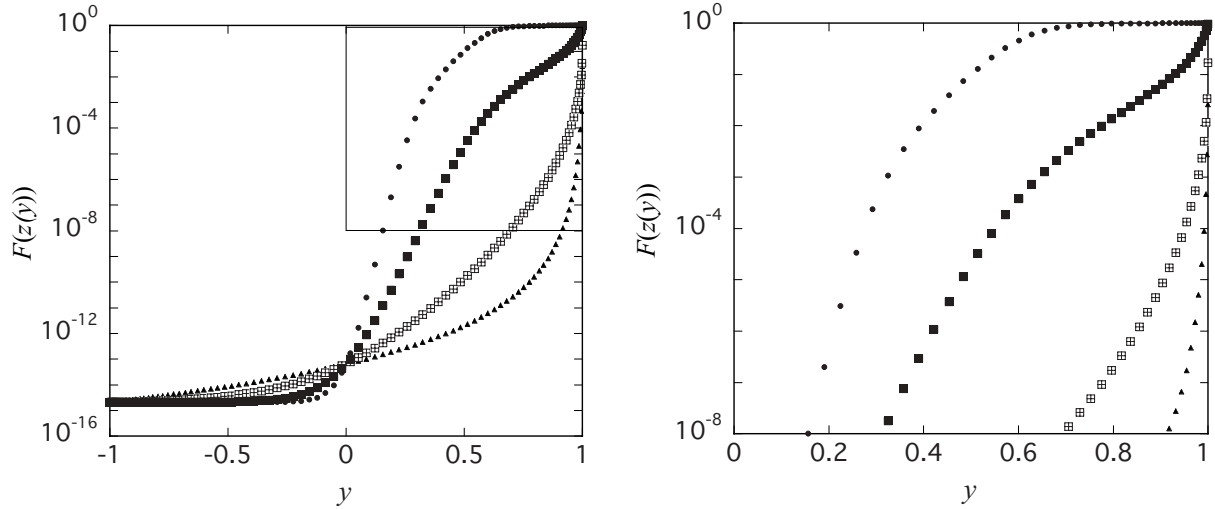


FIG. 11. Plot of the meson propagator F (with $m_b = 0.14$, $\Lambda_b = 4$, $z_0 = 1$, and $A = 1800$) as a function of the mapped variable y defined in Eq. (B5). The four curves, each with 90 Gauss points, correspond to $b_0 = 10$ (small solid triangles), $b_0 = 1.5$ (open boxes), $b_0 = 1.2$ (solid boxes), and $b_0 = 1.1$ (small solid circles). The right panel covers the region shown in the upper right corner of the left panel.

(instead of the physical variable z), and the integrand, as a function of the mapped variable y , is shown in Fig. 11 (for the same four choices of b). The angular integral is transformed to

$$\int_{-1}^1 dz F(z) = \int_{-1}^1 dy \frac{dz}{dy} F[z(y)]. \quad (\text{B6})$$

Clearly, without any mapping ($b_0 > 10$) the forward peak is very poorly sampled. If we take $b_0 = 1 + \epsilon$, with ϵ small, but larger than zero (to avoid the “singularity” at $b_0 = 1$), we are able to evaluate the integrals accurately. In practice, we found that $b_0 = 1.2$ (one of the cases shown in the figures) gave a good sampling of the forward peak and the integration converged rapidly.

3. Mapping for the exchange terms

The exchange propagator has the same form as Eq. (B3), except now

$$z_0 = \frac{p^2 + p'^2 - (E_p + E_{p'} - W)^2}{2pp'} \rightarrow -1 + \frac{2W}{p}, \quad (\text{B7})$$

where W is the total energy of the NN system in its center of mass, and the last expression holds if $p = p' \rightarrow \infty$. We see that it is now possible for $|z_0|$ to be less than unity, so the peak can lie in the region of integration.

To handle this case we need a mapping where the number of Gauss points is strongly distributed near $z = z_0$, where $F(z)$ is sharply peaked. It is convenient to use a mapping which maps the points $y = z_0, 1, -1$ into $z = z_0, 1, -1$ respectively. Such a mapping is

$$z(y) = \begin{cases} z_+(y, b) & z_0 < y < 1 \\ z_-(y, b) & -1 < y < z_0 \end{cases}, \quad (\text{B8})$$

where

$$\begin{aligned} z_+(y, z_0) &= \frac{1}{2}(1 - z_0) \frac{\tanh[\zeta(2y - 1 - z_0)/(1 - z_0)]}{\tanh[\zeta]} \\ &\quad + \frac{1}{2}(z_0 + 1) \\ z_-(y, z_0) &= \frac{1}{2}(1 + z_0) \frac{\tanh[\zeta(2y + 1 - z_0)/(1 + z_0)]}{\tanh[\zeta]} \\ &\quad + \frac{1}{2}(z_0 - 1) \end{aligned} \quad (\text{B9})$$

and ζ is a parameter. This mapping (with $\zeta = 4$) is shown in Fig. 12. Note that each curve distributes the 90 Gauss points so they are densely spaced in the region of $z = z_0$.

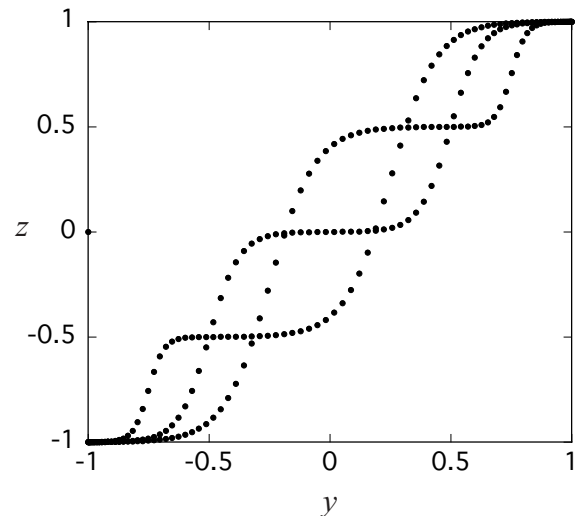


FIG. 12. Plot of the mapping function $z(y)$ defined in Eq. (B8), with $\zeta = 4$. The three curves, each with 90 Gauss points, correspond to $z_0 = -0.5$, $z_0 = 0$, and $z_0 = 0.5$. Each curve flattens out in the region of $z = y = z_0$.

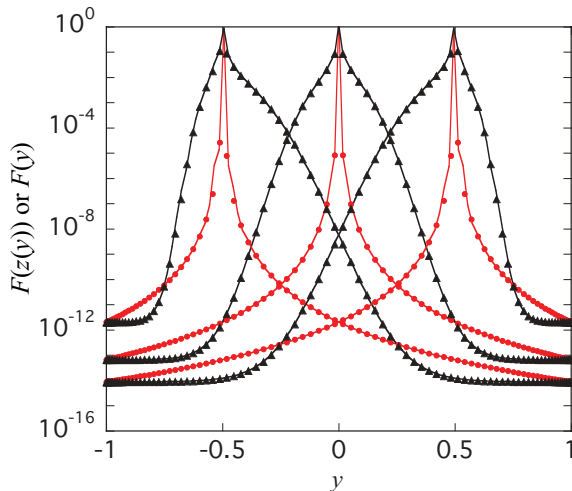


FIG. 13. (Color online) Plot of the meson propagator F (with $m_b = 0.14$, $\Lambda_b = 4$, and $A = 1800$) as a function of both the unmapped variable y and the mapped variable $z(y)$, defined in Eq. (B8), for each of the choices $z_0 = -0.5$, $z_0 = 0$, and $z_0 = 0.5$. The 90 Gauss points distributed between $-1 < y < 1$ are shown on each curve. All the curves peak at $y = z_0$, with the unmapped functions $F(y)$ having narrower peaks than the mapped function $F[z(y)]$.

Both the mapped and unmapped meson propagators are shown in Fig. 13, for each of the cases $z_0 = -0.5, 0$, and 0.5 . Note that the unmapped propagator has a very narrow peak at $y = z_0$ and that this peak is described by only a few Gauss points, even when the total number of points is quite large (90 in this case). The mapped functions, however, have their peaks spread out with many Gauss points distributed around the peak. From this picture, it is easy to see why the mapping allows a more accurate evaluation of the integrals. (In practice, a better result was obtained by dividing the interval into $\ell_1 = [-1, z_0]$ and $\ell_2 = [z_0, 1]$ and doing separate Gaussian integrations over each interval. The points assigned to each interval, n_i , depended on the value of z_0 . If $z_0 < -0.5$, we chose $n_1 = n/4$ and $n_2 = n$, for $-0.5 \leq z_0 < 0$, $n_1 = n/2$ and $n_2 = n - n/4$, for $0 \leq z_0 < 0.5$, $n_1 = n - n/4$ and $n_2 = n/2$, and for $0.5 \leq z_0$, $n_1 = n$ and $n_2 = n/4$, so in all cases we actually used $5n/4$ Gauss points.)

Finally, the angular convergence is illustrated in Tables XV and XVI for a typical matrix element [the isoscalar kernel v_1^{++} defined in Eq. (A12)] at one of the most difficult points: the diagonal elements with $p = p' = p_{\max}$, where $p_{\max} \approx$

30 GeV, the largest momenta used in the numerical solutions of the equations. Table XV shows the convergence of the direct terms (those with $x_0 = 1$) and Table XVI the exchange terms (those with $x_0 = -1$). Note that the direct terms calculated with unmapped integrations (the columns labeled with y in the tables) do not converge and that the convergence of the exchange terms is very marginal. With the mappings (columns labeled z), the direct terms converge even at $n = 20$, but excellent results for the exchange terms requires $n \geq 40$. Because the high-momentum matrix elements only make a small contribution to the solutions of the equation, smaller values of n give very good results for the phase shifts, and these improved angular integrations only affect the accuracy with which the asymptotic deuteron wave functions can be determined.

APPENDIX C: DEUTERON WAVE FUNCTIONS: SOME DETAILS

In this appendix we start by showing in detail how the matrix Eq. (3.26) is derived and then use the equation to study the asymptotic behavior of the wave functions, obtaining the results (3.41). We conclude with a demonstration of how the general normalization condition (3.12) reduces to (3.39).

1. Equations for the partial-wave deuteron wave functions

The equation for the deuteron wave functions, in partial-wave form, can be extracted from the general result (E36) of Ref. I. This equation uses a kernel antisymmetrized under particle exchange (with the superscript $J = 1$ suppressed)

$$\overline{V}_{d,\lambda_1\lambda_2,\lambda'_1\lambda'_2}^{+\rho_2+\rho'_2}(p, k) = \frac{1}{2} \left\{ V_{\text{dir},\lambda_1\lambda_2,\lambda'_1\lambda'_2}^{+\rho_2+\rho'_2}(p, p_0; k) + \delta_{p_0} V_{\text{dir},\lambda_1\lambda_2,\lambda'_1\lambda'_2}^{+\rho_2+\rho'_2}(p, -p_0; k) \right\}. \quad (\text{C1})$$

We will show how the full kernel with well defined parity, Eq. (A1), will emerge automatically, but the phase of the amplitude under particle interchange cannot emerge automatically in the CST (as discussed in Refs. I and II); that phase must be imposed by hand from the start by using the symmetrized kernel (C1). The phase $\delta_{p_0} = \pm 1$ is related but not equal to the phase under particle exchange [because the helicities λ_1, λ_2 are not exchanged in (C1)].

TABLE XV. Convergence of the diagonal element of the *direct* part of the kernel v_1^{++} at the point $p = p' = p_{\max}$. Columns labeled y use no mapping; those labeled $z(y)$ use the mapping function (B5). Here the nucleon form factor has been set to 1. With the correct form factor, all of these matrix elements are reduced by a factor of 7×10^{-15} . (All numbers in this table are multiplied by 10^{-3} .)

n	$J = 0$		$J = 1$		$J = 2$		$J = 3$		$J = 4$		$J = 5$		$J = 6$	
	y	$z(y)$	y	$z(y)$	y	$z(y)$	y	$z(y)$	y	$z(y)$	y	$z(y)$	y	$z(y)$
20	0.14321	-3.3764	-0.38372	-24.288	0.14001	-3.3825	-0.36991	-24.233	0.13277	-3.3967	-0.34585	-24.134	0.12190	-3.4184
30	0.62029	-3.3765	-2.0149	-24.288	0.61385	-3.3827	-1.9816	-24.232	0.59907	-3.3968	-1.9225	-24.133	0.57642	-3.4186
40	1.4365	-3.3765	-5.2403	-24.288	1.4277	-3.3827	-5.1898	-24.232	1.4074	-3.3968	-5.0999	-24.133	1.3762	-3.4186
60	2.5656	-3.3765	-12.496	-24.288	2.5569	-3.3827	-12.435	-24.232	2.5368	-3.3968	-12.326	-24.133	2.5059	-3.4186
90	1.4055	-3.3765	-18.312	-24.288	1.3992	-3.3827	-18.254	-24.232	1.3847	-3.3968	-18.152	-24.133	1.3624	-3.4186

TABLE XVI. Convergence of the diagonal element of the *exchange* part of the kernel v_1^{++} at the point $p = p' = p_{\max}$. Columns labeled y use no mapping; those labeled $z(y)$ use the mapping function (B8). Here the nucleon form factor has been set to 1. With the correct form factor, all of these matrix elements are reduced by a factor of 7×10^{-15} .

n	$J = 0$		$J = 1$		$J = 2$		$J = 3$		$J = 4$		$J = 5$		$J = 6$	
	y	$z(y)$	y	$z(y)$	y	$z(y)$	y	$z(y)$	y	$z(y)$	y	$z(y)$	y	$z(y)$
20	-25.184	-74.786	-5.1601	-14.525	-16.551	-49.079	-2.1999	-6.2065	-2.1638	-6.2802	0.98501	2.7617	8.5080	25.368
30	-37.148	-81.841	-6.0514	-11.903	-24.393	-53.713	-2.5703	-5.0887	-3.1487	-6.8793	1.1690	2.2584	12.580	27.757
40	-43.422	-80.907	-8.1636	-13.401	-28.504	-53.097	-3.4750	-5.7208	-3.6625	-6.7959	1.5681	2.5541	14.716	27.443
60	-54.117	-80.899	-9.9713	-13.594	-35.514	-53.091	-4.2542	-5.8025	-4.5432	-6.7939	1.9029	2.5918	18.356	27.441
90	-67.458	-80.910	-9.1111	-13.607	-44.267	-53.098	-3.8893	-5.8080	-5.6594	-6.7948	1.7357	2.5944	22.885	27.445

Using this kernel, the deuteron bound state equation is

$$\frac{m}{E_p} \Gamma_{\lambda_1 \lambda_2, \lambda_d}^{+\rho_2}(p) = - \int_0^\infty \frac{k^2 dk}{(2\pi)^3} \sum_{\mu_1 \mu_2 \rho} \bar{V}_{d, \lambda_1 \lambda_2, \mu_1 \mu_2}^{+\rho_2, +\rho}(p, k) \times G^\rho(k) \frac{m}{E_k} \Gamma_{\mu_1 \mu_2, \lambda_d}^{+\rho}(k). \quad (\text{C2})$$

The dependence of the amplitudes on the total momentum P has been dropped, and the energy factors of m/E , written explicitly, will be absorbed later when we go to the matrix notation of Eq. (3.26). At this stage the sum is over all helicities μ_1, μ_2 and the intermediate rho spin ρ .

The symmetries of the vertex function are a direct consequence of the symmetries of the kernel. Under parity, the deuteron kernel transforms as

$$\mathcal{P} \bar{V}_{d, \lambda_1 \lambda_2, \mu_1 \mu_2}^{+\rho_2, +\rho}(p, k) \mathcal{P}^{-1} = \rho \rho_2 \bar{V}_{d, -\lambda_1 -\lambda_2, -\mu_1 -\mu_2}^{+\rho_2, +\rho}(p, k). \quad (\text{C3})$$

Hence the equation satisfied by the transformed vertex function,

$$\frac{m}{E_p} \rho_2 \Gamma_{-\lambda_1 -\lambda_2, \lambda_d}^{+\rho_2}(p) = -\rho_2 \int_0^\infty \frac{k^2 dk}{(2\pi)^3} \sum_{\mu_1 \mu_2 \rho} \bar{V}_{d, -\lambda_1 -\lambda_2, \mu_1 \mu_2}^{+\rho_2, +\rho}(p, k) \times G^\rho(k) \frac{m}{E_k} \Gamma_{\mu_1 \mu_2, \lambda_d}^{+\rho}(k) = - \int_0^\infty \frac{k^2 dk}{(2\pi)^3} \sum_{\mu_1 \mu_2 \rho} \bar{V}_{d, \lambda_1 \lambda_2, \mu_1 \mu_2}^{+\rho_2, +\rho}(p, k) \times G^\rho(k) \frac{m}{E_k} \rho \Gamma_{-\mu_1 -\mu_2, \lambda_d}^{+\rho}(k), \quad (\text{C4})$$

is identical to the original equation, showing that the parity relation (3.24) is satisfied. Using this result we may reduce the sum over μ_1, μ_2 from four to two terms. Dropping the redundant $\rho_1 = \rho'_1 = +$ indices, we can write the summation over the two values of μ_1 explicitly and use the parity relations to reduce the equation to a form in which only $\mu_1 = +\frac{1}{2}$ appears:

$$\frac{m}{E_p} \Gamma_{\lambda_1 \lambda_2, \lambda_d}^{\rho_2}(p) = - \int_0^\infty \frac{k^2 dk}{(2\pi)^3} \frac{m}{E_k} \sum_{\mu_2 \rho} G^\rho(k) \{ \bar{V}_{d, \lambda_1 \lambda_2, +\frac{1}{2} \mu_2}^{\rho_2, \rho}(p, k) \times \Gamma_{+\frac{1}{2} \mu_2, \lambda_d}^\rho(k) + \bar{V}_{d, \lambda_1 \lambda_2, -\frac{1}{2} \mu_2}^{\rho_2, \rho}(p, k) \times \Gamma_{-\frac{1}{2} \mu_2, \lambda_d}^\rho(k) \} = - \int_0^\infty \frac{k^2 dk}{(2\pi)^3} \frac{m}{E_k} \sum_{\mu_2 \rho} \{ \bar{V}_{d, \lambda_1 \lambda_2, +\frac{1}{2} \mu_2}^{\rho_2, \rho}(p, k) \times \Gamma_{+\frac{1}{2} \mu_2, \lambda_d}^\rho(k) + \bar{V}_{d, \lambda_1 \lambda_2, -\frac{1}{2} \mu_2}^{\rho_2, \rho}(p, k) \times \Gamma_{-\frac{1}{2} \mu_2, \lambda_d}^\rho(k) \}$$

$$+ \rho \bar{V}_{d, \lambda_1 \lambda_2, -\frac{1}{2}, -\mu_2}^{\rho_2, \rho}(p, k) \times \{ G^\rho(k) \Gamma_{+\frac{1}{2} \mu_2, \lambda_d}^\rho(k) \} = - \int_0^\infty \frac{k^2 dk}{(2\pi)^3} \frac{m}{E_k} \sum_{\mu_2 \rho} \mathbf{V}_{d, \lambda_1 \lambda_2, +\frac{1}{2} \mu_2}^{\rho_2, \rho}(\delta_{p_0}, \rho) \times G^\rho(k) \Gamma_{+\frac{1}{2} \mu_2, \lambda_d}^\rho(k), \quad (\text{C5})$$

where in the last line we have recovered the fully symmetrized kernel of Eq. (A1). The deuteron channel must have the phases $\delta_{p_0} = \rho_2^\lambda$, appropriate to an isospin 0 state, and $\delta_S = \rho_2' = \rho$. Using the notation $\mathbf{V} = \mathbf{V}_{\lambda_2 \mu_2, \delta_{p_0}}^{\rho_2, \rho}$ (and, for the helicities, $\pm \frac{1}{2} \rightarrow \pm$) leads to the identifications

$$\begin{aligned} \mathbf{V}_{+++,+}^{\rho_2+}(p, k) &= v_9^{\rho_2+} & \mathbf{V}_{+++,+}^{\rho_2-}(p, k) &= v_1^{\rho_2-} \\ \mathbf{V}_{+++,+}^{\rho_2+}(p, k) &= v_{11}^{\rho_2+} & \mathbf{V}_{+++,+}^{\rho_2-}(p, k) &= v_3^{\rho_2-} \\ \mathbf{V}_{-+,-,+}^{\rho_2+}(p, k) &= v_{12}^{\rho_2+} & \mathbf{V}_{-+,-,+}^{\rho_2-}(p, k) &= v_8^{\rho_2-} \\ \mathbf{V}_{--,+}^{\rho_2+}(p, k) &= v_{10}^{\rho_2+} & \mathbf{V}_{--,+}^{\rho_2-}(p, k) &= v_6^{\rho_2-} \\ \mathbf{V}_{-+,-,-}^{\rho_2+}(p, k) &= v_{16}^{\rho_2+} & \mathbf{V}_{-+,-,-}^{\rho_2-}(p, k) &= v_4^{\rho_2-} \\ \mathbf{V}_{--,-}^{\rho_2+}(p, k) &= v_{14}^{\rho_2+} & \mathbf{V}_{--,-}^{\rho_2-}(p, k) &= v_2^{\rho_2-}, \end{aligned} \quad (\text{C6})$$

where the kernels are written in the order for coupling to $|\mathbf{\Gamma}_{++}^+\rangle, |\mathbf{\Gamma}_{++}^-\rangle, |\mathbf{\Gamma}_{+-}^+\rangle, |\mathbf{\Gamma}_{+-}^-\rangle$, respectively. Note that they reproduce the matrix (3.28), finishing this discussion.

2. Asymptotic behavior of the wave functions

Starting from Eq. (C5), the wave function (3.22) has the following general behavior at large p

$$\psi_{\lambda_1 \lambda_2, \lambda_d}^\rho(p) = N_d G^\rho(p) \frac{m}{E_p} \Gamma_{\lambda_1 \lambda_2, \lambda_d}^\rho(p) \simeq G^\rho(p) \int_0^\infty dk \mathbf{V}_d(p, k) C(k), \quad (\text{C7})$$

where, in this section, we return to the notation $k \equiv |\mathbf{k}|$ for the magnitude of the three-momentum, and \mathbf{V}_d is the appropriate partial-wave projection of the Feynman OBE amplitudes [shown in detail in Eq. (C6)]. All we need know about the function $C(k)$ is that it provides the convergence for the k integral. The kernels include a factor of m/E_p , which contributes to the large p behavior.

The kernels \mathbf{V}_d are multiplied by nucleon form factors [recall Eq. (3.5)], one of which depends on the intermediate momentum (k) and the other on the final momentum (p). The

final-state form factor, $h(p)$, plays an important role in the asymptotic behavior of the wave function. It depends only on p^2 , which is the square of the mass of the off-shell nucleon

$$\begin{aligned} p^2 &= (M_d - E_p)^2 - \mathbf{p}^2 \\ p^2 &\rightarrow -2M_d p \quad \text{for } p \gg m. \end{aligned} \quad (\text{C8})$$

Hence,

$$h(p^2) \rightarrow \frac{(\Lambda_N^2 - m^2)^4}{16M_d^4 p^4} \quad \text{for } p \gg m, \quad (\text{C9})$$

so the form factor behaves like p^{-4} . This contributes a k^{-4} behavior to the function $C(k)$, which, together with the internal nucleon form factor $h(k)$ that is part of \mathbf{V}_d , assures that the integral over k will converge. Hence we may conclude that the wave functions go like

$$\psi_{\lambda_1 \lambda_2 \lambda_d}^\rho(\mathbf{p}) \rightarrow G^\rho(p) \mathbf{V}_d(p, \bar{k}) C(\bar{k}), \quad (\text{C10})$$

where the mean value theorem has been used to write the integral over k , with \bar{k} a four-vector whose three-vector part has some fixed (unknown) value \bar{k} in the interval $[0, \infty)$.

To finish this discussion we examine the asymptotic behavior of the partial-wave amplitudes for one of the OBE terms in the kernel. If the form factors $h(p)$ and $h(k)$ are removed, this integral has the generic form

$$D_b(p, \bar{k}) = \int_{-1}^1 dz g_b(p) \frac{m}{E_p} \frac{f(\Lambda_b, q)}{m_b^2 + |q^2|} d(z), \quad (\text{C11})$$

where q^2 is the momentum transferred by the boson with mass m_b , form factor $f(\Lambda_b, q)$, and momentum-dependent coupling $g_b(p) \rightarrow p^\ell$ (as $p \rightarrow \infty$, where ℓ is yet to be determined). We have isolated the m/E_p term incorporated into the definition of \mathbf{V}_d , and $d(z)$ includes the z dependence that arises from the generic rotation matrices of the type tabulated in Table XI. In Ref. I the meson form factors have the form

$$f(\Lambda_b, q) = \left[\frac{\Lambda_b^2}{\Lambda_b^2 + |q^2|} \right]^{n_b}, \quad (\text{C12})$$

where $n_b = 4$ for all mesons.

The asymptotic behavior of the integral (C11) depends on whether it is a *direct* or an *exchange* term. The momentum transferred by the mesons in the direct terms

$$\begin{aligned} q^2 &= (E_p - E_k)^2 - (\mathbf{p} - \mathbf{k})^2 \\ &\rightarrow 2m^2 - 2p(E_k - kz) \quad (\text{as } p \rightarrow \infty, \mathbf{k} \text{ fixed}) \end{aligned} \quad (\text{C13})$$

and is never positive, so the integrand is a smooth function of z . The asymptotic behavior, for a fixed $k = k_0$, is therefore

$$D_b^{\text{dir}}(p, \bar{k}) \rightarrow \frac{p^\ell}{p^{n_b+2}}. \quad (\text{C14})$$

As shown below, these terms are vanishingly small compared to the exchange terms.

The exchange terms have a very different structure. Here the momentum transferred by the meson is

$$\begin{aligned} q^2 &= (M_d - E_p - E_k)^2 - (\mathbf{p} + \mathbf{k})^2 \\ &\rightarrow m^2 + (M_d - E_k)^2 - k^2 - 2p(M_d - E_k + kz) \\ &\rightarrow -2p(M_d - E_k + kz) \quad \text{as } p \rightarrow \infty. \end{aligned} \quad (\text{C15})$$

When p is very large, this is zero along a line in the k, z plane given approximately by $M_d - E_k + kz = 0$. Along this line the integrand has a cusp which sits at

$$z_c = -\frac{M_d - E_k}{k} \quad (\text{C16})$$

in the region of integration ($-1 \leq z \leq 1$) for values of k in the interval

$$\frac{M_d^2 - m^2}{2M_d} \equiv k_{\min} < k < \infty. \quad (\text{C17})$$

Contributions from the cusp clearly dominate the integral, because in that region the integrand is constant. Hence, letting $m_b = \Lambda_b$ in order to simplify the determination of the asymptotic p dependence, the exchange integral is approximately

$$\begin{aligned} \frac{D_b^{\text{ex}}(p, \bar{k})}{p^{\ell-1}} &\simeq \int_{-1}^1 dz \left[\frac{\Lambda_b^2}{\Lambda_b^2 + |q^2|} \right]^{n_b+1} d(z) \\ &\simeq \int_{-1}^{z_c} dz \left[\frac{\Lambda_b^2}{\Lambda_b^2 + 2p\bar{k}(z_c - z)} \right]^{n_b+1} d(z) \\ &\quad + \int_{z_c}^1 dz \left[\frac{\Lambda_b^2}{\Lambda_b^2 + 2p\bar{k}(z - z_c)} \right]^{n_b+1} d(z) \\ &\simeq \frac{2\Lambda_b^2}{n_b p \bar{k}} d(z_c), \end{aligned} \quad (\text{C18})$$

where we have integrated by parts, keeping only the leading term (the corrections go like higher powers of $1/p$).

Collecting these results together we see that the reduced partial-wave amplitude, D_b , behaves like

$$D(p, k_0) \rightarrow \begin{cases} p^{\ell-(2+n_b)} & \text{direct} \\ p^{\ell-2} & \text{exchange,} \end{cases} \quad (\text{C19})$$

where n_b is the power of the form factor at the NNb vertex and ℓ the power of the p dependence of the NNb vertex function.

It remains to determine the power ℓ . First, note that the single particle ρ -spin matrix elements given in Table XIV each contribute a factor of $\sqrt{p/(2m)}$ [from the high-momentum limits (A5)]; their product gives one power of p implicitly contained in Eqs. (A45)–(A48).

Next, examine any additional p dependence coming from the two particle ρ -spin matrix elements given in Eqs. (A45)–(A48). There are terms of order p coming from the off-shell energy factors ΔE_i defined in (A24), but the condition (A25) shows that these factors cannot be multiplied together to make a term of order p^2 . Special terms of order p^2 seem to be present in the vector meson exchange terms, but we will show below that these terms cancel. Hence the final result is the product of two terms of order p giving $\ell = 2$.

To show the cancellation of the special p^2 terms contributing to vector-meson exchange, consider the three terms in Eq. (A47) which give rise to them: the product of R_{v1} 's (first term), the p^2 multiplying the D_4 terms (fourth term), and the last term. It is sufficient to look at the exchange-term matrix elements where particle 1 is off-shell in the final state, so that $\Delta E_2 = 0$. At large p use the fact that

$\Delta E_1 \rightarrow (\rho_1 + 1)p$ to obtain

$$\begin{aligned} R_{v1}R_{v1} &\rightarrow p^2 \frac{g_v^2 \kappa_v}{(2m)^2} [\rho_1 \kappa_v - (\rho_1 + 1) \nu_v] \mathbf{1}_1 \mathbf{1}_2 \\ D_4 D_4 &\rightarrow p^2 \left(\frac{g_v \kappa_v}{2m} \right)^2 \mathbf{1}_1 \mathbf{1}_2 \\ R_{v3} D_4 &\rightarrow p^2 \frac{g_v^2 \kappa_v}{(2m)^2} (\kappa_v - \nu_v) (\rho_1 + 1) 2\lambda_1 \boldsymbol{\tau}_1^1 \mathbf{1}_2. \end{aligned} \quad (\text{C20})$$

Adding these together and replacing ρ_1 by ρ_2 [in accordance with the notation for the exchange terms in Eq. (A12)] gives the following compact result

$$g_v(p) \simeq p^2 \frac{g_v^2 \kappa_v}{(2m)^2} (\kappa_v - \nu_v) (\rho_2 + 1) [\mathbf{1}_1 + 2\lambda_1 \boldsymbol{\tau}_1^1] \mathbf{1}_2. \quad (\text{C21})$$

Note that this is zero if $\rho_2 = -1$, insuring immediately that eight of the deuteron kernels (C6) have no special terms of order p^2 . To see that this is also true of the other eight cases set $\rho_2 = 1$ (and recall that $\rho_1' = 1$) and substitute for the matrix elements $\mathbf{1}$ and $\boldsymbol{\tau}^1$ using the results from Table XIV (with $\rho_2 \leftrightarrow \rho_1$) and the high-momentum limits (A5)

$$g_v(p) \simeq 2p^2 \sqrt{\frac{p}{2m}} \frac{g_v^2 \kappa_v}{(2m)^2} (\kappa_v - \nu_v) (1 - 2\lambda_1) (c' - s') \mathbf{1}_2. \quad (\text{C22})$$

This is zero if $\lambda_1 = \frac{1}{2}$, which is true of all of the amplitudes (and, in particular, the remaining 8). We have proved that the special p^2 terms coming from vector meson exchange cancel, and that $\ell = 2$. The conclusion is that the wave functions should go like

$$z^\rho(p) \rightarrow G^\rho(p) h(p^2) D^{\text{ex}}(p, k_0) \quad (\text{C23})$$

which gives the result (3.41) as expected.

To confirm the predictions of Eq. (C6), Fig. 14 shows how the 16 deuteron kernels $V_i(p, k)$ vary with p for a fixed k .

3. Reduction of Eqs. (3.12) to (3.39)

Substituting the expansion (3.16) into Eq. (3.12), and using the orthogonality of the Dirac spinors, compactly written as

$$\bar{u}^\rho(-\mathbf{p}, \lambda) \gamma^0 u^{\rho'}(-\mathbf{p}, \lambda') = \frac{E_p}{m} \delta_{\lambda' \lambda} \delta_{\rho' \rho} \quad (\text{C24})$$

gives

$$\begin{aligned} \delta_{\lambda \lambda'} \delta_{\rho \rho'} &= \int d^3 p \psi_{\lambda_1 \lambda_2, \lambda'_d}^{\rho \dagger}(\mathbf{p}) \psi_{\lambda_1 \lambda_2, \lambda_d}^\rho(\mathbf{p}) - \int \int \frac{d^3 p d^3 p'}{(2\pi)^3} \\ &\times \psi_{\lambda_1 \lambda_2, \lambda_d}^{\rho \dagger}(\mathbf{p}) \Delta \bar{V}^{\rho \rho'}_{\lambda_1 \lambda_2, \lambda'_d} (p, p'; P) \psi_{\lambda'_1 \lambda'_2, \lambda'_d}^{\rho'}(\mathbf{p}'), \end{aligned} \quad (\text{C25})$$

where the $\bar{V}^{\rho \rho'} \equiv \bar{V}^{+\rho, +\rho'}$ includes the factors of $m^2/E_p E_{p'}$ [cf. Eq. (E13) of Ref. I].

Next we express the helicity amplitudes in terms of the wave functions $z_\ell(p)$ using (3.30). To this end recall Eq. (A6) for the two-component helicity spinors and use the convenient

relation $i\sigma_2 \chi_\lambda = -2\lambda \chi_{-\lambda}$ and the matrix elements

$$\begin{aligned} \chi_{-\lambda_2}^\dagger \boldsymbol{\sigma} \cdot \boldsymbol{\xi}_{\lambda_d} i\sigma_2 \chi_{\lambda_1} &= \sqrt{2}^{|\lambda|} d_{\lambda_d, \lambda}^1(\theta) \\ \chi_{-\lambda_2}^\dagger \boldsymbol{\sigma} \cdot \hat{\mathbf{p}} i\sigma_2 \chi_{\lambda_1} &= \chi_{-\lambda_2}^\dagger \chi_{-\lambda_1} = \delta_{\lambda_1, \lambda_2} \\ \hat{\mathbf{p}} \cdot \boldsymbol{\xi}_{\lambda_d} &= d_{\lambda_d, 0}^1(\theta), \end{aligned} \quad (\text{C26})$$

where $\lambda = \lambda_1 - \lambda_2$ and the rotation matrices are

$$\begin{aligned} d_{00}^1(\theta) &= \cos \theta \\ d_{10}^1(\theta) &= d_{0, -1}^1(\theta) = -d_{0, 1}^1(\theta) = -d_{-1, 0}^1(\theta) = -\frac{1}{\sqrt{2}} \sin \theta \end{aligned} \quad (\text{C27})$$

$$d_{11}^1(\theta) = d_{-1, -1}^1(\theta) = \frac{1}{2}(1 + \cos \theta)$$

$$d_{-1, 1}^1(\theta) = d_{1, -1}^1(\theta) = \frac{1}{2}(1 - \cos \theta).$$

The result is

$$\begin{aligned} \psi_{\lambda_1 \lambda_2, \lambda_d}^+(\mathbf{p}) &= \frac{1}{\sqrt{8\pi}} \left\{ \left(u(\mathbf{p}) - \frac{1}{\sqrt{2}} w(\mathbf{p}) \right) \sqrt{2}^{|\lambda|} d_{\lambda_d, \lambda}^1(\theta) \right. \\ &\quad \left. + \delta_{\lambda_1, \lambda_2} \frac{3w(\mathbf{p})}{\sqrt{2}} d_{\lambda_d, 0}^1(\theta) \right\} \\ \psi_{\lambda_1 \lambda_2, \lambda_d}^-(\mathbf{p}) &= -2\lambda_1 \sqrt{\frac{3}{8\pi}} \left\{ \delta_{\lambda_1, \lambda_2} \left(v_s(\mathbf{p}) - \frac{v_t(\mathbf{p})}{\sqrt{2}} \right) d_{\lambda_d, 0}^1(\theta) \right. \\ &\quad \left. + \frac{v_t(\mathbf{p})}{\sqrt{2}} \sqrt{2}^{|\lambda|} d_{\lambda_d, \lambda}^1(\theta) \right\}. \end{aligned} \quad (\text{C28})$$

Explicitly, for each individual case, using $\lambda = \pm \frac{1}{2}$

$$\begin{aligned} \psi_{\lambda \lambda, \lambda_d}^+(\mathbf{p}) &= \frac{1}{\sqrt{8\pi}} (u(\mathbf{p}) + \sqrt{2} w(\mathbf{p})) d_{\lambda_d, 0}^1(\theta) \\ \psi_{\lambda, -\lambda, \lambda_d}^+(\mathbf{p}) &= \frac{1}{\sqrt{4\pi}} \left(u(\mathbf{p}) - \frac{1}{\sqrt{2}} w(\mathbf{p}) \right) d_{\lambda_d, 2\lambda}^1(\theta) \\ \psi_{\lambda \lambda, \lambda_d}^-(\mathbf{p}) &= -2\lambda \sqrt{\frac{3}{8\pi}} v_s(\mathbf{p}) d_{\lambda_d, 0}^1(\theta) \\ \psi_{\lambda, -\lambda, \lambda_d}^-(\mathbf{p}) &= -2\lambda \sqrt{\frac{3}{8\pi}} v_t(\mathbf{p}) d_{\lambda_d, 2\lambda}^1(\theta). \end{aligned} \quad (\text{C29})$$

We now can reduce the first term in the normalization condition (C25). Taking $\lambda_d = \lambda'_d$ and using the normalization condition (3.20) for the d^1 functions gives

$$\begin{aligned} 1 &= 2\pi \int_0^\infty p^2 dp \int_0^\pi \sin \theta d\theta \{ 2|\psi_{\lambda \lambda, \lambda_d}^+(\mathbf{p})|^2 \\ &\quad + 2|\psi_{\lambda, -\lambda, \lambda_d}^+(\mathbf{p})|^2 + 2|\psi_{\lambda \lambda, \lambda_d}^-(\mathbf{p})|^2 + 2|\psi_{\lambda, -\lambda, \lambda_d}^-(\mathbf{p})|^2 \} \\ &= \frac{1}{3} \int_0^\infty p^2 dp \left\{ (u(\mathbf{p}) + \sqrt{2} w(\mathbf{p}))^2 \right. \\ &\quad \left. + 2 \left(u(\mathbf{p}) - \frac{1}{\sqrt{2}} w(\mathbf{p}) \right)^2 + 3v_s^2(\mathbf{p}) + 3v_t^2(\mathbf{p}) \right\} \\ &= \int_0^\infty p^2 dp \{ u^2(\mathbf{p}) + w^2(\mathbf{p}) + v_t^2(\mathbf{p}) + v_s^2(\mathbf{p}) \}, \end{aligned} \quad (\text{C30})$$

in agreement with Eq. (3.39). The derivative terms can be similarly reduced and expressed in terms of the kernels given in Eq. (C6).

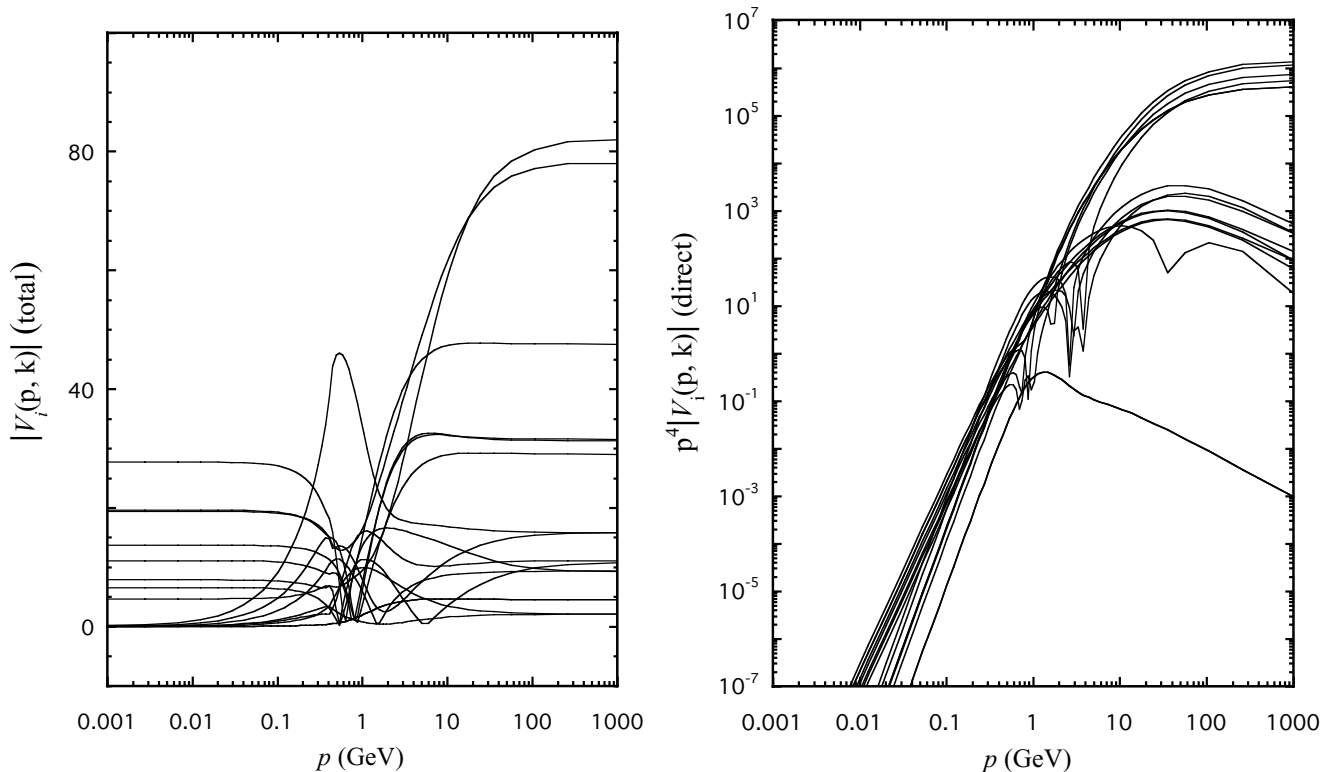


FIG. 14. The p dependence (for k fixed at 816 MeV) of the 16 model WJC-1 deuteron kernels $V_i(p, k)$ defined in Eq. (C6) (but with the nucleon form factors set to unity). The left panel shows clearly that the *total* kernel (direct plus exchange) approaches a constant at large p , as predicted by (C19) with $\ell = 2$. The right panel shows that the asymptotic p dependence of 5 of the *direct* kernels behaves like p^{-4} , again predicted by (C19) with $n_b = 4$ and $\ell = 2$. The other direct kernels decrease more rapidly, presumably because they have an $\ell < 2$.

Before leaving this section we note that the partial-wave deuteron wave functions follow from the normalization condition (3.20) and the definitions (3.21)

$$\begin{aligned}
 \psi_{+++,\lambda_d}^+(\mathbf{p}) &= \sqrt{\frac{3}{8}} \int_{-1}^1 dz |d_{\lambda_d,0}^1(\theta)|^2 (u + \sqrt{2}w) \\
 &= \frac{1}{\sqrt{6}} (u + \sqrt{2}w) \\
 \psi_{+-,\lambda_d}^+(\mathbf{p}) &= \sqrt{\frac{3}{8}} \int_{-1}^1 dz |d_{\lambda_d,1}^1(\theta)|^2 (\sqrt{2}u - w) \\
 &= \frac{1}{\sqrt{6}} (\sqrt{2}u - w) \\
 \psi_{+++,\lambda_d}^-(\mathbf{p}) &= -\frac{3}{\sqrt{8}} \int_{-1}^1 dz |d_{\lambda_d,0}^1(\theta)|^2 v_s = -\frac{1}{\sqrt{2}} v_s \\
 \psi_{+-,\lambda_d}^-(\mathbf{p}) &= -\frac{3}{\sqrt{8}} \int_{-1}^1 dz |d_{\lambda_d,1}^1(\theta)|^2 v_t = -\frac{1}{\sqrt{2}} v_t.
 \end{aligned} \tag{C31}$$

4. Coordinate-space expansion functions

The coordinate-space wave functions discussed in Sec. III G 3 are constructed from the spherical Bessel transforms (3.37), which can be written in terms of the j_0 Bessel

transform

$$\begin{aligned}
 G_\ell^i(r) &= r \sqrt{\frac{2}{\pi}} \int_0^\infty p^2 dp j_\ell(pr) G_\ell^i(p) \\
 &= (-1)^\ell \sqrt{\frac{2}{\pi}} r^{\ell+1} \left(\frac{1}{r} \frac{d}{dr} \right)^\ell \int_0^\infty p^2 dp j_0(pr) \frac{G_\ell^i(p)}{p^\ell} \\
 &= (-1)^\ell \frac{r^{\ell+1}}{M_i^\ell} \left(\frac{1}{r} \frac{d}{dr} \right)^\ell \left[\frac{G_0^i(r)}{r} \right].
 \end{aligned} \tag{C32}$$

Hence, for S and P states it is sufficient to calculate the following Fourier sine transforms:

$$\begin{aligned}
 G_0^i(r) &= \sqrt{\frac{2}{\pi}} \int_0^\infty p dp \sin(pr) G_0^i(p) \\
 &= A_i \left\{ e^{-z_i} - e^{-Z_i} \left[1 + \frac{1}{2} Z_i (1 - R_i^2) \right] \right\} \\
 G_0^n(r) &= \sqrt{\frac{2}{\pi}} \int_0^\infty p dp \sin(pr) G_0^n(p) \\
 &= \frac{2}{3\pi} M_n^2 Z_n^2 K_1(Z_n),
 \end{aligned} \tag{C33}$$

where z_i , Z_i , R_i , and A_i were given in Eq. (3.52) and the modified Bessel functions of the second kind are

$$K_n(z) = \frac{z^n}{(2n-1)!!} \int_1^\infty dt e^{-zt} (t^2 - 1)^{n-\frac{1}{2}}. \tag{C34}$$

The P -state wave functions are obtained by differentiation, as outlined in Eq. (C32). They are

$$G_1^i(r) = -\frac{r}{M_i} \frac{d}{dr} \left[\frac{G_0^i(r)}{r} \right] = A_i \left\{ R_i e^{-z_i} \left[1 + \frac{1}{z_i} \right] - e^{-z_i} \left[1 + \frac{1}{Z_i} + \frac{1}{2} Z_i (1 - R_i^2) \right] \right\}, \quad (\text{C35})$$

$$G_1^n(r) = -\frac{r}{M_n} \frac{d}{dr} \left[\frac{G_0^n(r)}{r} \right] = \frac{2}{3\pi} M_n^2 Z_n^2 K_0(Z_n).$$

Finally, the D -state wave functions are computed by first transforming the functions

$$G_{0d}^i(p) = \sqrt{\frac{2}{\pi}} \frac{m_i^2 M_i^6}{(m_i^2 + p^2)(M_i^2 + p^2)^3} \quad (\text{C36})$$

$$G_{0d}^n(p) = \sqrt{\frac{2}{\pi}} \frac{M_n^7}{(M_n^2 + p^2)^{7/2}}$$

which gives

$$G_{0d}^i(r) = \sqrt{\frac{2}{\pi}} \int_0^\infty p dp \sin(pr) G_{0d}^i(p)$$

$$= B_i \left\{ e^{-z_i} - e^{-Z_i} - \frac{1}{8} Z_i e^{-Z_i} [(5 - 6R_i^2 + R_i^4) + Z_i (1 - R_i^2)^2] \right\}$$

$$G_{0d}^n(r) = \sqrt{\frac{2}{\pi}} \int_0^\infty p dp \sin(pr) G_{0d}^n(p)$$

$$= \frac{2}{15\pi} M_n^2 Z_n^3 K_2(Z_n), \quad (\text{C37})$$

where B_i was given in Eq. (3.52). Differentiating these following (C32) gives

$$G_2^i(r) = \frac{r^2}{M_n^2} \left(\frac{d}{dr} \frac{1}{r} \right)^2 G_{0d}^i(r) = B_i \left\{ R_i^2 e^{-z_i} \left[1 + \frac{3}{z_i} + \frac{3}{z_i^2} \right] - e^{-z_i} \left[1 + \frac{3}{Z_i} + \frac{3}{Z_i^2} + \frac{1}{2} (1 - R_i^2)(1 + Z_i) + \frac{1}{8} Z_i^2 (1 - R_i^2)^2 \right] \right\},$$

$$G_2^n(r) = \frac{r^2}{M_n^2} \left[\frac{d}{dr} \frac{1}{r} \right]^2 G_{0d}^n(r) = \frac{2}{15\pi} M_n^2 Z_n^3 K_0(Z_n). \quad (\text{C38})$$

At large r the functions G_ℓ^i go like

$$\lim_{r \rightarrow \infty} G_0^i(r) = A_i e^{-z_i}$$

$$\lim_{r \rightarrow \infty} G_1^i(r) = A_i R_i e^{-z_i} \left[1 + \frac{1}{z_i} \right] \quad (\text{C39})$$

$$\lim_{r \rightarrow \infty} G_2^i(r) = B_i R_i^2 e^{-z_i} \left[1 + \frac{3}{z_i} + \frac{3}{z_i^2} \right].$$

This is the asymptotic behavior expected for solutions of the Schrödinger equation with orbital angular momentum ℓ . The tail functions fall off like exponentials multiplied by a fractional power of r

$$\lim_{r \rightarrow \infty} G_0^n(r) = \frac{1}{3} \sqrt{\frac{2}{\pi}} M_n^2 Z_n^{3/2} e^{-Z_n}$$

$$\lim_{r \rightarrow \infty} G_1^n(r) = \frac{1}{3} \sqrt{\frac{2}{\pi}} M_n^2 Z_n^{3/2} e^{-Z_n} \quad (\text{C40})$$

$$\lim_{r \rightarrow \infty} G_2^n(r) = \frac{1}{15} \sqrt{\frac{2}{\pi}} M_n^2 Z_n^{5/2} e^{-Z_n}.$$

At small r the functions G_ℓ^i have the expected $r^{\ell+1} \sim Z_i^{\ell+1}$ behavior

$$\lim_{r \rightarrow 0} G_0^i(r) = \frac{1}{2} A_i Z_i (1 - R_i)^2$$

$$\lim_{r \rightarrow 0} G_1^i(r) = \frac{1}{6} A_i Z_i^2 (1 - R_i)^2 (1 + 2R_i) \quad (\text{C41})$$

$$\lim_{r \rightarrow 0} G_2^i(r) = \frac{1}{40} B_i Z_i^3 (1 - R_i)^3 (1 + 3R_i + \frac{8}{3} R_i^2).$$

However, the tail functions contribute some nonanalytic behavior at small r :

$$\lim_{r \rightarrow 0} G_0^n(r) = \frac{2}{3\pi} M_n^2 Z_n.$$

$$\lim_{r \rightarrow 0} G_1^n(r) = -\frac{2}{3\pi} M_n^2 Z_n^2 (\log Z_n - \log 2 + \gamma) \quad (\text{C42})$$

$$\lim_{r \rightarrow 0} G_2^n(r) = -\frac{2}{15\pi} M_n^2 Z_n^3 (\log Z_n - \log 2 + \gamma),$$

where $\gamma = 0.5772$ is Euler's constant.

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