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Hadronic molecules for charmed and bottom baryons near thresholds

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We study hadronic molecules formed by a heavy meson and a nucleon, DN and D^*N ($\bar{B}N$ and \bar{B}^*N) systems. Respecting the heavy quark symmetry and chiral symmetry, we consider the $DN-D^*N$ ($\bar{B}N-\bar{B}^*N$) mixing induced by the one boson exchange potential including the tensor force. We find many bound and resonant states with $J^P = 1/2^{\pm}$, $3/2^{\pm}$, $5/2^{\pm}$, and $7/2^{-}$ in isospin singlet channels, while we find only a few resonant states with $J^P = 1/2^{-}$ in isospin triplet channels. The analysis of DN and D^*N ($\bar{B}N$ and \bar{B}^*N) molecules will be useful to study mass spectra of excited charmed (bottom) baryons with large angular momenta, when their masses are close to the DN and D^*N ($\bar{B}N$ and \bar{B}^*N) thresholds.

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I. INTRODUCTION

Recently, many exotic hadrons observed in experiments have attracted our interest. They are expected to have unexpected structures, because their properties, such as quantum numbers, masses, and decay patterns, cannot be explained by the ordinary picture of hadrons, baryons as qqq and mesons as $q\bar{q}$. As one of the new aspects of these structures, the possibility of multiquarks or hadronic molecules have been extensively discussed. It is considered that, particularly near thresholds, hadronic molecules such as loosely bound or resonant states of meson-meson or meson-baryon emerge as one of the hadronic configurations realized in nature. As typical candidates, the meson-meson molecules for X(3872) and Z_b have been studied in the charm and bottom quark regions [1–8].

The hadronic molecules as bound or resonant states formed by D or D^* mesons (\bar{B} or \bar{B}^* mesons) and a nucleon (N) are also interesting. Since the DN and D^*N ($\bar{B}N$ and \bar{B}^*N) molecules contain ordinary flavor structures of three quarks, they are intimately related to the ordinary heavy baryons, Λ_c , Σ_c , Σ_c^* and their excited states for the charm sector, and Λ_b , Σ_b , Σ_b^* and their excited states for the bottom sector [9]. These excited states have been studied in the quark model extensively [11,12]. However, near the DN and D^*N ($\bar{B}N$ and \bar{B}^*N) thresholds, it is expected that the wave functions of the quark model states have a large overlap with the component of DN and D^*N ($\bar{B}N$ and \bar{B}^*N), and the properties of such baryon states are strongly affected by DN and D^*N ($\bar{B}N$ and \bar{B}^*N) states. There would be even such states that are dominated by the molecular components.

In the DN and D^*N ($\bar{B}N$ and \bar{B}^*N) systems, there are two important symmetries, the heavy quark symmetry and chiral symmetry. The heavy quark symmetry manifests in mass degeneracy of heavy pseudoscalar mesons $P = D, \bar{B}$ and vector mesons $P^* = D^*, \bar{B}^*$ in the heavy quark mass limit [13,14]. Indeed, the mass splitting between a D (\bar{B})

meson and a D^* (\bar{B}^*) meson is small, 140 MeV for D and D^* mesons and 46 MeV for \bar{B} and \bar{B}^* mesons. The small mass splittings in the heavy flavor sectors should be compared with the large mass splittings, ~ 400 MeV for \bar{K} and \bar{K}^* mesons and ~600 MeV for π and ρ mesons, in the light flavor sectors. Because of this, the heavy quark symmetry introduces the mixing of PN and P^*N states, where both P and P^* mesons are considered on the same footing as fundamental degrees of freedom in the dynamics [15]. Then, physical states are given as a superposition of PN and P^*N states. For convenience, in the following, we denote $P^{(*)}$ to stand for P or P^* . Thanks to the mixing of states in $P^{(*)}N$ systems, a long range force between a $P^{(*)}$ meson and a nucleon N is supplied by one pion exchange with the same coupling strengths for $PP^*\pi$ and $P^*P^*\pi$ vertices [16–18].

It is interesting that the pion exchange can exist in $P^{(*)}N$ systems. As a matter of fact, it is known that the pion exchange is important for the binding of atomic nuclei [19]. There the tensor force of the pion exchange that mixes channels with different angular momenta, i.e., L and $L \pm 2$, yields a strong attraction to generate a rich structure of bound as well as resonant states. Because the pion exchange is a long range force, it becomes more significant for states with larger orbital angular momenta. In fact, it was shown that, in the nucleon-nucleon scattering, the phase shifts in the channels with large orbital angular momenta are reproduced almost only by the pion exchange potential [20,21]. The importance of the pion exchange was also discussed in $N\bar{N}$ systems. It was investigated that the properties of bound and resonant states in $N\bar{N}$ systems are dramatically changed if the tensor force is switched off [22,23]. In our previous studies, it was also shown that the pion exchange potential played a significant role for hadronic molecules such as $\bar{D}N$ and BN baryons [16–18], $D\bar{D}$ ($B\bar{B}$) mesons (e.g., Z_b meson) [24], and *DD* (*BB*) mesons [25].

In reality, however, the hadronic molecules of $P^{(*)}N$ do not necessarily correspond to the observed states, because there should be couplings not only to three quark states but also to other meson-baryon states such as $\pi\Lambda_{\rm c},\,\pi\Sigma_{\rm c}$ and $\pi\Sigma_{\rm c}^*$ ($\pi\Lambda_{\rm b}$, $\pi\Sigma_{\rm b}$, and $\pi\Sigma_{\rm b}^*$), and so on. However, we may expect that such couplings are small for DN and $\bar{B}N$ baryons near the thresholds. The reasons are that the wave functions of hadronic molecules are spatially large as compared to the conventional three quark states, and that the transitions, e.g., from DN $(\bar{B}N)$ to $\pi\Sigma_{\rm c}$ $(\pi\Sigma_{\rm b})$, are suppressed by a heavy quark exchange. From those points of view, in the present discussion, we focus on the role of the $D^{(*)}N$ and $\bar{B}^{(*)}N$ sectors and study the bound or resonant states generated by $D^{(*)}N$ and $\bar{B}^{(*)}N$. Recently, there have been several studies of baryon states as $D^{(*)}N(\bar{B}^{(*)}N)$ [26–40] and $\bar{D}^{(*)}\Sigma_c$ $(B^{(*)}\Sigma_b)$ [41–43] molecules. The present study is different from them in that we focus on the states near the thresholds and with a large angular momentum.

This paper is organized as follows. In Sec. II, we briefly describe the interaction between a $D^{(*)}$ or $\bar{B}^{(*)}$ meson and a nucleon. In Sec. III, we solve the coupled-channel Schrödinger equations numerically for bound and resonant states. We discuss the results of the $D^{(*)}N$ and $\bar{B}^{(*)}N$ molecule in Sec. IV, and summarize the present work in the final section.

II. INTERACTIONS

Following our previous studies [16–18], we consider π , ρ , and ω exchange potentials between $P^{(*)}$ and N. The interaction Lagrangians for heavy meson vertices are given by the heavy quark symmetry and chiral symmetry [44–50]:

$$\mathcal{L}_{\pi H H} = i g_{\pi} \operatorname{Tr}[H_b \gamma_{\mu} \gamma_5 A^{\mu}_{ba} \bar{H}_a], \tag{1}$$

$$\mathcal{L}_{vHH} = -i\beta \operatorname{Tr}[H_b v^{\mu}(\rho_{\mu})_{ba} \bar{H}_a] + i\lambda \operatorname{Tr}[H_b \sigma^{\mu\nu} F_{\mu\nu}(\rho)_{ba} \bar{H}_a], \qquad (2)$$

where the subscripts π and v are for the pion and vector $(\rho \text{ and } \omega)$ meson interactions. The heavy meson field H is defined by $H_a = \frac{1+\psi}{2} \big[P_{a\mu}^* \gamma^\mu - P_a \gamma_5 \big]$ with the four-velocity v^μ of a heavy meson, where the subscript a is for light flavors u, d. The conjugate field is given by $\bar{H}_a = \gamma_0 H_a^\dagger \gamma_0$. The axial current of light flavors is written by $A^\mu = \frac{1}{2} (\xi^\dagger \partial^\mu \xi - \xi \partial^\mu \xi^\dagger)$ with $\xi = \exp{(i\hat{\pi}/f_\pi)}$ and the pion decay constant $f_\pi = 132$ MeV. The pion field is defined by

$$\hat{\pi} = \begin{pmatrix} \frac{\pi^0}{\sqrt{2}} & \pi^+ \\ \pi^- & -\frac{\pi^0}{\sqrt{2}} \end{pmatrix}, \tag{3}$$

and the vector meson field by

$$\rho^{\mu} = i \frac{g_V}{\sqrt{2}} \hat{\rho}_{\mu},\tag{4}$$

$$\hat{\rho}_{\mu} = \begin{pmatrix} \frac{\rho^{0}}{\sqrt{2}} & \rho^{+} \\ \rho^{-} & -\frac{\rho^{0}}{\sqrt{2}} \end{pmatrix}_{\mu}, \tag{5}$$

where $g_V \simeq 5.8$ corresponds to the gauge coupling constant from hidden local symmetry [51]. The vector meson field tensor is written by $F_{\mu\nu}(\rho) = \partial_{\mu}\rho_{\nu} - \partial_{\nu}\rho_{\mu} + [\rho_{\mu}, \rho_{\nu}]$. The coupling constants g_{π} , β and λ are summarized in Table I. These coupling constants are essentially the same as those given in Refs. [17,18], except for the signs of g_{π} , β , and λ for vertices of π and ω , which are reversed due to G-parity transformation between $D^{(*)}[\bar{B}^{(*)}]$ and $\bar{D}^{(*)}[B^{(*)}]$. In Eqs. (1) and (2), we consider the static approximation $v^{\mu} = (1, \vec{0})$.

For vertices of π , ρ , and ω mesons to a nucleon, we employ the Bonn model [52] as

$$\mathcal{L}_{\pi NN} = \sqrt{2} i g_{\pi NN} \bar{N} \gamma_5 \hat{\pi} N, \tag{6}$$

$$\mathcal{L}_{\nu NN} = \sqrt{2} g_{\nu NN} \left[\bar{N} \gamma_{\mu} \hat{\rho}^{\mu} N + \frac{\kappa}{2m_N} \bar{N} \sigma_{\mu\nu} \partial^{\nu} \hat{\rho}^{\mu} N \right], \quad (7)$$

where $N = (p, n)^T$ is the nucleon field. The coupling constants for nucleons are summarized in Table I. The pseudoscalar coupling becomes approximately equivalent to the pseudovector coupling with a derivative, if the particles coupled to the pion are on shell, which is approximately the case of hadronic molecules.

When we consider the π , ρ , and ω exchange potentials between the $P^{(*)}$ meson and nucleon, we take into account the internal structure of hadrons by introducing the monopole-type form factors at each vertex;

$$F_{\alpha}(\Lambda, \vec{q}) = \frac{\Lambda^2 - m_{\alpha}^2}{\Lambda^2 + |\vec{a}|^2},\tag{8}$$

with the mass m_{α} , three momentum \vec{q} of the incoming meson $\alpha(=\pi,\rho,\omega)$, and the cutoff parameter Λ . Here, we introduce two cutoff parameters for a heavy meson $[D^{(*)}]$ and $\bar{B}^{(*)}]$ and a nucleon. The cutoff parameter Λ_N for the nucleon is determined to reproduce the properties (the binding energy, the scattering length and effective range) of the deuteron. The cutoff parameter Λ_P for the heavy meson $P^{(*)} = D^{(*)}$, $\bar{B}^{(*)}$ is determined by the ratios of

TABLE I. Masses and coupling constants of mesons $\alpha = \pi$, ρ , ω in Refs. [17,18].

Meson	m_{α}	g_{π}	β	$\lambda \; [\text{GeV}^{-1}]$	$g_{\alpha NN}^2/4\pi$	κ
$\overline{\pi}$	137.27	-0.59		•••	13.6	• • • •
ho	769.9		0.9	0.59	0.84	6.1
ω	781.94	• • •	-0.9	-0.59	20.0	0.0

TABLE II. Cutoff parameters of a nucleon (Λ_N) and heavy mesons $[\Lambda_D \text{ for } D^{(*)} \text{ meson and } \Lambda_B \text{ for } \bar{B}^{(*)} \text{ meson}]$ as employed in Refs. [17,18].

Potential	Λ_N [MeV]	Λ_D [MeV]	Λ_B [MeV]
π	830	1121	1070
$\pi ho \omega$	846	1142	1091

matter radii of the heavy meson and nucleon where the ratio is obtained by a quark model calculation [16–18]. Their numerical values are summarized in Table II.

Here we briefly discuss cutoff dependence. Indeed, numerical results are rather sensitive to the choice of the parameters Λ . Here we have determined them from the size of hadrons, and so a large deviation from the present choice is not appropriate. If we change, however, Λ by, for example, about 10%, the locations of levels change quantitatively, but the structure of spectrum does not change qualitatively.

In the present study, we employ two types of potentials: the π exchange and the $\pi\rho\omega$ exchanges. By comparing the results from them, we show the importance of the one pion exchange potential.

III. NUMERICAL RESULTS

We consider the states with $J^P = 1/2^\pm$, $3/2^\pm$, $5/2^\pm$, and $7/2^\pm$ (total angular momentum J and parity P) with isospin I=0 and 1. Various states with J^P are expanded by coupled channels of ${}^{2S+1}L_J$ (spin S and orbital angular momentum L) as summarized in Table III. The meson exchange potentials among various channels and kinetic terms are summarized in the Appendix. Using the Hamiltonian composed of the kinetic and potential terms, we solve the coupled-channel Schrödinger equations for PN and P^*N channels numerically.

A. Isospin singlet (I = 0)

We present the results for the isosinglet state (I = 0). The energies of bound and resonant states are summarized in Table IV and also presented in Figs. 1 and 2. The results

TABLE III. Various coupled channels for a given quantum number J^P .

J^P	Channels						
1/2 ⁻ 1/2 ⁺ 3/2 ⁻ 3/2 ⁺ 5/2 ⁻ 5/2 ⁺ 7/2 ⁻ 7/2 ⁺	$PN(^2S_{1/2})$ $PN(^2P_{1/2})$ $PN(^2D_{3/2})$ $PN(^2P_{3/2})$ $PN(^2D_{5/2})$ $PN(^2F_{5/2})$ $PN(^2G_{7/2})$ $PN(^2F_{7/2})$	$P^*N(^2S_{1/2})$ $P^*N(^2P_{1/2})$ $P^*N(^4S_{3/2})$ $P^*N(^2P_{3/2})$ $P^*N(^2D_{5/2})$ $P^*N(^4D_{5/2})$ $P^*N(^4D_{7/2})$ $P^*N(^2F_{7/2})$	$P^*N(^4D_{1/2})$ $P^*N(^4P_{1/2})$ $P^*N(^4P_{3/2})$ $P^*N(^4P_{3/2})$ $P^*N(^4D_{5/2})$ $P^*N(^2F_{5/2})$ $P^*N(^2G_{7/2})$ $P^*N(^4F_{7/2})$	$P^*N(^2D_{3/2}) \\ P^*N(^4F_{3/2}) \\ P^*N(^4G_{5/2}) \\ P^*N(^4F_{5/2}) \\ P^*N(^4G_{7/2}) \\ P^*N(^4H_{7/2})$			

are shown for the π and $\pi\rho\omega$ potentials, and the corresponding energy levels are connected by arrows in Figs. 1 and 2. The numerical values are measured from the DN and $\bar{B}N$ thresholds, respectively.

First, we discuss the states with $J^P = 1/2^{\pm}$ and $3/2^{-}$, and then those with $J^P = 3/2^{+}$, $5/2^{\pm}$, and $7/2^{\pm}$. In fact, it will turn out that the results in those two categories have different behaviors.

1.
$$J^P = 1/2^{\pm}$$
 and $3/2^{-}$

For $(I, J^P) = (0, 1/2^-)$, we find bound states in both of $D^{(*)}N$ and $\bar{B}^{(*)}N$. For $D^{(*)}N$, the binding energies are -14.4 MeV for the π potential and -82.5 MeV for the $\pi\rho\omega$ potential. The relative radii are 1.51 and 0.86 fm. respectively. As expected, for a larger binding energy the system becomes smaller. The tensor force from the π exchange which is the main driving force of the DN-D*N mixing plays an important role to form the bound states. In fact, we have verified that neither bound nor resonant states exist when the tensor force from the π exchange is switched off and the $DN-D^*N$ mixing is small. For $J^P = 1/2^-$, the results in the π and $\pi \rho \omega$ potentials are very different, as indicated in Figs. 1 and 2. Since both ρ and ω exchanges are attractive for the $D^{(*)}N$ [$\bar{B}^{(*)}N$] system, the vector meson exchanges contribute to form the deeply bound state of the binding energy around 80 MeV. This contrasts with the previous result for the $\bar{D}^{(*)}N$ and $B^{(*)}N$ systems of truly exotic channels, where the ρ and ω exchanges play a minor role due to the cancellation of these two potentials [17,18].

We estimate the mixing ratio of various channels in the bound states as summarized in Table V. We observe that, for $J^P=1/2^-$ states, the most dominant component is $DN(^2S_{1/2})$ with a fraction 71.8%. The second dominant component is $D^*N(^4D_{1/2})$ with a fraction 20.4%. Therefore, the tensor force which mixes the S wave in $DN(^2S_{1/2})$ and the D wave in $D^*N(^4D_{1/2})$ is important to provide a strong attraction.

For $\bar{B}^{(*)}N$ of $(I,J^P)=(0,1/2^-)$, we also obtain bound states. The binding energies are -57.8 MeV for the π potential and -145.9 MeV for the $\pi\rho\omega$ potential, and the relative radii are 0.99 and 0.76 fm, respectively. Again the results of the two potentials are very different with the same reason for the $D^{(*)}N$ system. As compared to the $D^{(*)}N$ system, the binding energy in the $\bar{B}^{(*)}N$ system is much larger, because heavier particles suppress kinetic energy. For the mixing ratios, we also find a similar tendency as we discussed for the $D^{(*)}N$ system: 56.1% for $\bar{B}N(^2S_{1/2})$, 13.3% for $\bar{B}^*N(^2S_{1/2})$, and 30.6% for $\bar{B}^*N(^4D_{1/2})$.

Let us move to the $(I, J^P) = (0, 1/2^+)$ state. For $D^{(*)}N$, we find one resonance near the DN threshold when the π potential is used. The resonance energy is 1.4 MeV and the half-decay width is 0.2 MeV as shown in Table IV. In the

TABLE IV. Properties of bound and resonant states of $D^{(*)}N$ and $\bar{B}^{(*)}N$ systems. The energies E are either pure real for bound states or complex for resonant states. The complex energies for resonances are written as $E_{\rm re} - i\Gamma/2$ where $E_{\rm re}$ is the resonance energy and $\Gamma/2$ is the half-width. The binding and resonance energies are measured from the lowest threshold (DN and $\bar{B}N$). Root mean square radii are shown only for bound states.

		DN	7	$ar{B}N$	I
$I(J^P)$	Potential	E [MeV]	$\langle r^2 \rangle^{1/2}$ [fm]	E [MeV]	$\langle r^2 \rangle^{1/2}$ [fm]
0(1/2=)	π	-14.4	1.51	-57.8	0.99
$0(1/2^{-})$	$\pi ho\omega$	-82.5	0.86	-145.9	0.76
0(1/2+)	π	1.4 - i0.2		-83.8	0.92
$0(1/2^+)$	$\pi ho\omega$	-81.5	0.85	-185.0	0.75
$0(3/2^{-})$	π	63.5 - i7.9		-38.7	0.99
	$\pi ho\omega$	-13.7	0.89	-127.8	0.76
$0(3/2^+)$	π	23.8 - i118.1		12.9 - i15.5	
	$\pi ho\omega$	26.0 - i44.2		-2.6	1.81
0(5/2-)	π	153.6 - i671.9		63.7 - i177.6	
$0(5/2^{-})$	$\pi ho\omega$	160.0 - i375.4		71.3 - i102.8	
0(5/2+)	π	160.8 - i3.1		46.0 - i1.1	
	$\pi ho\omega$	137.0 - i7.6		20.0 - i0.2	
$0(7/2^{-})$	π	217.7 - i182.4		85.6 - i74.5	
	$\pi ho\omega$	220.8 - i109.1		87.5 - i46.7	
0(7/2+)	π	no		no	
	$\pi ho \omega$	no		no	
$1(1/2^-)$	π	no		no	
	$\pi ho\omega$	147.2 - i105.5	• • •	50.7 - i75.5	• • •

scattering state, we define the resonance energy $E_{\rm re}$ by an inflection point of the phase shift [53], and the decay width Γ is given by $\Gamma = 2/(d\delta/dE)_{E=E_{\rm re}}$. When the $\pi\rho\omega$ potential is used, we find a bound state with a binding energy

-81.5 MeV and with a relative radius 0.85 fm. The mixing ratios of the bound state are 38.8% for $DN(^2P_{1/2})$, 6.0% for $D^*N(^2P_{1/2})$, and 55.2% for $D^*N(^4P_{1/2})$, as shown in Table V. Interestingly, $D^*N(^4P_{1/2})$ is the most dominant

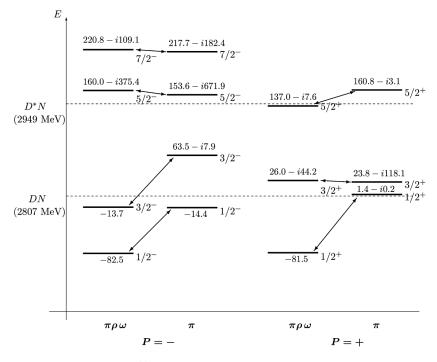


FIG. 1. Energies of bound and resonant states of $D^{(*)}N$ for I=0 with positive parity (P=+) and negative parity (P=-). The energies are measured from the DN threshold. The results for π and $\pi\rho\omega$ potentials are shown, where corresponding states are connected by arrows. The binding energies are given as a real negative value, and the resonance energies $E_{\rm re}$ and decay widths Γ are given as $E_{\rm re} = i\Gamma/2$.

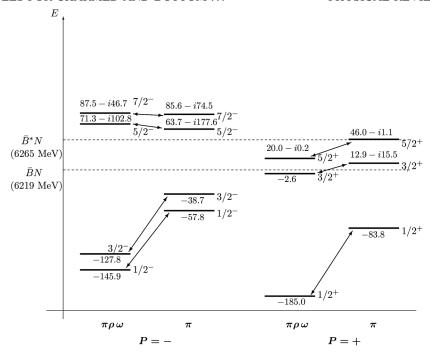


FIG. 2. Energies of bound and resonant states of $\bar{B}^{(*)}N$ for I=0 with positive parity (P=+) and negative parity (P=-). The energies are measured from the $\bar{B}N$ threshold. The same convention is used as Fig. 1.

channel, although the mass of D^*N is heavier than the mass of DN. This is because the attraction of the tensor force is the strongest for the $D^*N(^4P_{1/2})$ channel.

For $\bar{B}^{(*)}N$ of $(I,J^P)=(0,1/2^+)$, we find bound states for both cases when the π and $\pi\rho\omega$ potentials are used. The binding energies are -83.8 MeV for the π potential and -185.0 MeV for the $\pi\rho\omega$ potential, and the relative radii are 0.92 and 0.75 fm, respectively. For the $\pi\rho\omega$ potential, the mixing ratios of the bound state are 28.4% for $\bar{B}N(^2P_{1/2})$, 7.7% for $\bar{B}^*N(^2P_{1/2})$, and 63.9% for $\bar{B}^*N(^4P_{1/2})$. In this case again, $\bar{B}^*N(^4P_{1/2})$ is the most dominant component regardless of its heavy mass.

For $(I, J^P) = (0, 3/2^-)$, we find a resonant DN state for the π potential with the resonance energy 63.5 MeV and the half-decay width 7.9 MeV as shown in Table IV. In contrast, for the $\pi\rho\omega$ potential, we find a bound DN state with the binding energy -13.7 MeV and the relative radius 0.89 fm. For the bound state, the mixing ratios are 19.8% for $DN(^2D_{3/2})$, 62.8% for $D^*N(^4S_{3/2})$, 14.2% for $D^*N(^4D_{3/2})$, and 3.2% for $D^*N(^2D_{3/2})$. Thus, $D^*N(^4S_{3/2})$

is the dominant channel although its mass is heavier than the mass of $DN(^2D_{3/2})$; once again a large attraction due to the tensor force is provided.

For the $\bar{B}^{(*)}N$ state of $(I, J^P) = (0, 3/2^-)$, we obtain bound states for both cases; the π and $\pi\rho\omega$ potentials are used. The binding energies are -38.7 MeV for the π potential and -127.8 MeV for the $\pi\rho\omega$ potential with relative radii 0.99 and 0.76 fm, respectively.

2.
$$J^P = 3/2^+, 5/2^{\pm}, and 7/2^{\pm}$$

For $(I, J^P) = (0, 3/2^+)$, we find resonant states above the threshold. For $D^{(*)}N$, the resonance energies are 23.8 MeV for the π potential and 26.0 MeV for the $\pi\rho\omega$ potential. The half-decay widths are 118.1 and 44.2 MeV, respectively. As compared to the cases of $1/2^\pm$ and $3/2^-$, the results of the π and $\pi\rho\omega$ potentials are not very much different, as indicated in Fig. 1. Since the wave functions are extended due to larger orbital angular momenta of the P wave (L=1) and F wave (L=3) for $J^P=3/2^+$, the long range potential of the π exchange dominates, while

TABLE V. Mixing ratio of each channel in the bound $D^{(*)}N$ and $\bar{B}^{(*)}N$ states for $J^P=1/2^\pm$ and $3/2^\pm$ with I=0 when the $\pi\rho\omega$ potential is employed.

1/2-	$PN(^{2}S_{1/2})$	$P^*N(^2S_{1/2})$	$P^*N(^4D_{1/2})$		1/2+	$PN(^{2}P_{1/2})$	$P^*N(^2P_{1/2})$	$P^*N(^4P_{1/2})$	• • •
DN	$71.8\%^{-2}$	7.8%	20.4%	• • •	DN	38.8%	6.0%	55.2%	• • •
$\bar{B}N$	56.1%	13.3%	30.6%	• • •	$ar{B}N$	28.4%	7.7%	63.9%	• • •
$3/2^{-}$	$PN(^{2}D_{3/2})$	$P^*N(^4S_{3/2})$	$P^*N(^4D_{3/2})$	$P^*N(^2D_{3/2})$	$3/2^{+}$	$PN(^{2}P_{3/2})$	$P^*N(^2P_{3/2})$	$P^*N(^4P_{3/2})$	$P^*N(^4F_{3/2})$
DN	19.8%	62.8%	14.2%	3.2%	DN	• • • •	• • • •	•••	•••
$\bar{B}N$	14.6%	64.7%	16.7%	4.0%	$\bar{B}N$	71.2%	6.7%	7.5%	14.6%

the short range potentials from ρ and ω exchanges are suppressed. For $\bar{B}^{(*)}N$, when the π potential is used, we find a resonant state whose resonance energy is 12.9 MeV and half-decay width is 15.5 MeV, while when the $\pi\rho\omega$ potential is used, we find a loosely bound state of a binding energy -2.6 MeV and a relative radius 1.81 fm. The mixing ratios are 71.2% for $\bar{B}N(^2P_{3/2})$, 6.7% for $\bar{B}^*N(^2P_{3/2})$, 7.5% for $\bar{B}^*N(^4P_{3/2})$, and 14.6% for $\bar{B}^*N(^4F_{3/2})$. The most dominant component is $\bar{B}N(^2P_{3/2})$, and the second dominant one is $\bar{B}^*N(^4F_{3/2})$.

For $(I, J^P) = (0, 5/2^-)$, we obtain resonant states for both cases when the π and $\pi\rho\omega$ potentials are used. For $D^{(*)}N$, the resonance energies are 153.6 MeV for the π potential and 160.0 MeV for the $\pi\rho\omega$ potential, which are above the D^*N threshold. The corresponding half-decay widths are 671.9 and 375.4 MeV, respectively. The difference between the results of the π and $\pi \rho \omega$ potentials is once again small, due to the same reason as before with large angular momenta. For the $\bar{B}^{(*)}N$ state, we also find resonant states above the \bar{B}^*N threshold. The resonance positions are 63.7 MeV for the π potential and 71.3 MeV for the $\pi\rho\omega$ potential. The corresponding half-decay widths are 177.6 and 102.8 MeV, respectively. We observe from Table IV that the widths of the $5/2^-$ states of both DN and $\bar{B}N$ are very large. The reason is that there is only very weak attraction in these channels. Because of the very large widths of order 1 GeV, it is not easy to interpret them physically as particle states.

For $(I, J^P) = (0, 5/2^+)$, we find resonant states with narrow widths for both cases when the π and $\pi \rho \omega$ potentials are used. For $D^{(*)}N$, when the π potential is used, the resonance energy is 160.8 MeV and the half-decay width is $\Gamma/2 = 3.1$ MeV. When the $\pi \rho \omega$ potential is used, the resonance energy is 137.0 MeV and the half-decay width is $\Gamma/2 = 7.6$ MeV. For $\bar{B}^{(*)}N$, we also find resonances

whose energies are 46.0 MeV for the π potential and 20.0 MeV for the $\pi\rho\omega$ potential, with the corresponding half-decay widths 1.1 and 0.2 MeV, respectively. Again, the results for the π and $\pi\rho\omega$ potentials are similar.

For $(I, J^P) = (0, 7/2^-)$, we obtain resonances above the D^*N and \bar{B}^*N thresholds. For $D^{(*)}N$, there exist resonances at 217.7 MeV for the π potential and at 220.8 MeV for the $\pi\rho\omega$ potential, with half-decay widths 182.4 and 109.1 MeV, respectively. For $\bar{B}^{(*)}N$, there also exist resonances whose energies are 85.6 MeV for the π potential and 87.5 MeV for the $\pi\rho\omega$ potential, with half-decay widths 74.5 and 46.7 MeV, respectively.

Finally, for $(I, J^P) = (0, 7/2^+)$, we find no bound nor resonant states.

B. Isospin triplet (I = 1)

Let us move to the states of isospin triplet (I=1). We summarize the results for $D^{(*)}N$ and for $\bar{B}^{(*)}N$ in Table IV and show the energy levels in Fig. 3. As a result, we find resonant states only for $J^P=1/2^-$ when the $\pi\rho\omega$ potential is employed. For $D^{(*)}N$, the resonance energy is 147.2 MeV and the half-decay width is 105.5 MeV. For the $\bar{B}^{(*)}N$ state, we obtain the resonance whose energy is 50.7 MeV and half-width 75.5 MeV. The reason that there are not many states in the isospin triplet channel can be understood as follows; as compared to the isosinglet channel, the attractive force in the isotriplet channel is weak due to small isospin factor, $\vec{\tau}_P \cdot \vec{\tau}_N = -3$ for isosinglet and $\vec{\tau}_P \cdot \vec{\tau}_N = 1$ for isotriplet.

IV. DISCUSSION

A. Contribution from short range interactions

By solving coupled-channel Schrödinger equations for PN and P^*N channels, we find many bound and resonant

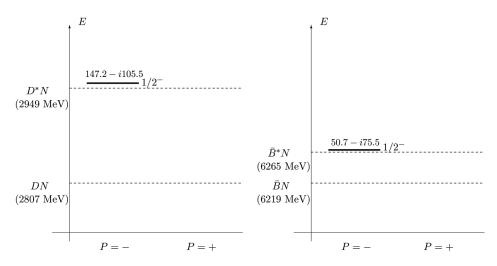


FIG. 3. Energies of bound and resonant states of $D^{(*)}N$ and $\bar{B}^{(*)}N$ for I=1 when the $\pi\rho\omega$ potential is used. The same convention is used as Fig. 1.

states for I=0 and few resonances for I=1. For these states, the tensor force of the π exchange potential plays a significant role to produce them. When we ignore the P^*N channels and solve the Schrödinger equation only with PN channels, we find no bound state nor resonance. Thus, the $PN-P^*N$ mixing and the π exchange potential play an important role to generate a rich structure of molecular states.

The importance of the π exchange potential stands out more in the bottom sector. The small mass difference between \bar{B} and \bar{B}^* mesons helps to yield the strong attraction because it induces the strong $\bar{B}N-\bar{B}^*N$ mixing with the tensor force. Furthermore, the $\bar{B}^{(*)}N$ has a heavier reduced mass and hence the kinetic term is suppressed. For these reasons, the binding energies of $\bar{B}^{(*)}N$ states are larger than these of $D^{(*)}N$ states.

When we compare the results of the π potential and of the $\pi\rho\omega$ potential, they are quite different for $J^P = 1/2^{\pm}$, $3/2^-$, with I=0, where the ρ and ω exchange potentials become important to produce attraction. For $D^{(*)}N$ and $\bar{B}^{(*)}N$ states, both ρ and ω exchange potentials are attractive, and hence they increase the binding energy significantly. On the other hand, for $J^P = 3/2^+$, $5/2^{\pm}$, and $7/2^-$, the results for the $\pi\rho\omega$ potential are similar to those for the π potential. For large J states, the π exchange potential plays a dominant role to generate bound and resonant states, while the ρ and ω exchanges play only a minor one. This is attributed to the fact that these states have large orbital angular momenta. If relevant channels include large orbital angular momenta, the wave functions tend to extend spatially, and the long range force, namely the π exchange potential, becomes important while the short range force is suppressed.

B. Emergent hadronic molecules

In the present study, we discuss the molecular structure formed by the $P^{(*)}N$ bound and resonant states. However, the hadronic molecular picture is not applicable to deeply bound states with small radii. For such states, the constituent hadrons, namely, $D^{(*)}[\bar{B}^{(*)}]$ meson and nucleon, overlap each other, and therefore we have to consider short range effects such as an internal structure of hadrons, channel couplings to conventional three quark states and so on. As a naive criterion for the hadronic molecule. we have shown the relative radii of the bound states as discussed in Ref. [25]. If the size of the bound state is larger than twice of typical hadron size (namely, 1 fm), the state could be well described by a molecular structure. For resonant states, we identify the states as the hadronic molecule. According to the criterion, for the $\pi\rho\omega$ potential, only the bound state for $J^P=3/2^+$ of $\bar{B}^{(*)}N$ constructs a hadronic molecule, where the relative radius is 1.81 fm. Contrary to this, the bound states for $J^P = 1/2^{\pm}$ and $3/2^-$ with I = 0, which have a small

radius and a large binding energy, are not described as simple molecules. We need to consider the short range effects including various channel couplings to do more realistic discussions.

V. SUMMARY

We have investigated heavy baryons as hadronic molecules formed by a heavy meson and a nucleon, $D^{(*)}N$ and $\bar{B}^{(*)}N$. The interaction is given by the meson exchange potential with respecting the heavy quark and chiral symmetries. Many bound states and resonances are found in the isosinglet channel both for $D^{(*)}N$ and $\bar{B}^{(*)}N$ systems. In contrast, there are few resonances in the isotriplet channel due to the small isospin factor. For $J^P = 1/2^{\frac{1}{2}}$ and $3/2^-$, we have found deeply bound states far below the DN or $\bar{B}N$ thresholds with a small radius. In these states, the vector meson exchange potential yields a strong attraction. In order to perform more realistic discussions for such compact states, we need to consider further effects from internal structures of constituent hadrons and channel couplings to quark intrinsic states. On the other hand, for $J^P = 3/2^+$, $5/2^{\pm}$, and $7/2^-$ with large orbital angular momenta, the one pion exchange potential is favored and we have obtained loosely bound states and resonances near the thresholds. There the short range interactions are rather inactive and the long range force of the one pion exchange dominates, where the tensor force plays an important role to generate these states, and they compose the hadronic molecular structure.

It is expected that the hadronic molecular states from a heavy meson and a nucleon are studied in many accelerator facilities, the *D* meson production in the antiproton-nucleon annihilation process in FAIR [54] and the heavy ion collision in RHIC and LHC [55,56]. Furthermore, the exotic baryons are investigated in J-PARC in the coming future.

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APPENDIX: POTENTIALS AND KINETIC TERMS

Interaction potentials are derived by using the Lagrangians in Eqs. (1), (2), (6), and (7) as shown in Refs. [17,18]. The potentials for the coupled-channel systems are given in the matrix form of 3×3 for $J^P = 1/2^{\pm}$ and of 4×4 for the other states.

The π exchange potentials for each J^P are obtained by

$$V_{1/2^{-}}^{\pi} = \begin{pmatrix} 0 & \sqrt{3}V_{C}^{\pi} & -\sqrt{6}V_{T}^{\pi} \\ \sqrt{3}V_{C}^{\pi} & -2V_{C}^{\pi} & -\sqrt{2}V_{T}^{\pi} \\ -\sqrt{6}V_{T}^{\pi} & -\sqrt{2}V_{T}^{\pi} & V_{C}^{\pi} - 2V_{T}^{\pi} \end{pmatrix}, \quad (A1)$$

$$V_{1/2^{+}}^{\pi} = \begin{pmatrix} 0 & \sqrt{3}V_{C}^{\pi} & -\sqrt{6}V_{T}^{\pi} \\ \sqrt{3}V_{C}^{\pi} & -2V_{C}^{\pi} & -\sqrt{2}V_{T}^{\pi} \\ -\sqrt{6}V_{T}^{\pi} & -\sqrt{2}V_{T}^{\pi} & V_{C}^{\pi} - 2V_{T}^{\pi} \end{pmatrix}, \quad (A2)$$

$$V_{3/2^{-}}^{\pi} = \begin{pmatrix} 0 & \sqrt{3}V_{T}^{\pi} & -\sqrt{3}V_{T}^{\pi} & \sqrt{3}V_{C}^{\pi} \\ \sqrt{3}V_{T}^{\pi} & V_{C}^{\pi} & 2V_{T}^{\pi} & V_{T}^{\pi} \\ -\sqrt{3}V_{T}^{\pi} & 2V_{T}^{\pi} & V_{C}^{\pi} & -V_{T}^{\pi} \\ \sqrt{3}V_{C}^{\pi} & V_{T}^{\pi} & -V_{T}^{\pi} & -2V_{C}^{\pi} \end{pmatrix}, \quad (A3)$$

$$V_{3/2^{+}}^{\pi} = \begin{pmatrix} 0 & \sqrt{3}V_{C}^{\pi} & \sqrt{\frac{3}{5}}V_{T}^{\pi} & -3\sqrt{\frac{3}{5}}V_{T}^{\pi} \\ \sqrt{3}V_{C}^{\pi} & -2V_{C}^{\pi} & \frac{1}{\sqrt{5}}V_{T}^{\pi} & -\sqrt{3}V_{T}^{\pi} \\ \sqrt{\frac{3}{5}}V_{T}^{\pi} & \frac{1}{\sqrt{5}}V_{T}^{\pi} & V_{C}^{\pi} + \frac{8}{5}V_{T}^{\pi} & \frac{6}{5}V_{T}^{\pi} \\ -3\sqrt{\frac{3}{5}}V_{T}^{\pi} & -\sqrt{3}V_{T}^{\pi} & \frac{6}{5}V_{T}^{\pi} & V_{C}^{\pi} - \frac{8}{5}V_{T}^{\pi} \end{pmatrix},$$
(A4)

$$V_{5/2^{-}}^{\pi} = \begin{pmatrix} 0 & \sqrt{3}V_{C}^{\pi} & \sqrt{\frac{6}{7}}V_{T}^{\pi} & -\frac{6}{\sqrt{7}}V_{T}^{\pi} \\ \sqrt{3}V_{C}^{\pi} & -2V_{C}^{\pi} & \sqrt{\frac{2}{7}}V_{T}^{\pi} & -2\sqrt{\frac{3}{7}}V_{T}^{\pi} \\ \sqrt{\frac{6}{7}}V_{T}^{\pi} & \sqrt{\frac{2}{7}}V_{T}^{\pi} & V_{C}^{\pi} + \frac{10}{7}V_{T}^{\pi} & \frac{4}{7}\sqrt{6}V_{T}^{\pi} \\ -\frac{6}{\sqrt{7}}V_{T}^{\pi} & -2\sqrt{\frac{3}{7}}V_{T}^{\pi} & \frac{4}{7}\sqrt{6}V_{T}^{\pi} & V_{C}^{\pi} - \frac{10}{7}V_{T}^{\pi} \end{pmatrix},$$
(A5)

$$V_{5/2^{+}}^{\pi} = \begin{pmatrix} 0 & \frac{3}{5}\sqrt{10}V_{T}^{\pi} & \sqrt{3}V_{C}^{\pi} & -2\sqrt{\frac{3}{5}}V_{T}^{\pi} \\ \frac{3}{5}\sqrt{10}V_{T}^{\pi} & V_{C}^{\pi} - \frac{2}{5}V_{T}^{\pi} & \sqrt{\frac{6}{5}}V_{T}^{\pi} & \frac{4}{5}\sqrt{6}V_{T}^{\pi} \\ \sqrt{3}V_{C}^{\pi} & \sqrt{\frac{6}{5}}V_{T}^{\pi} & -2V_{C}^{\pi} & -\frac{2}{\sqrt{5}}V_{T}^{\pi} \\ -2\sqrt{\frac{3}{5}}V_{T}^{\pi} & \frac{4}{5}\sqrt{6}V_{T}^{\pi} & -\frac{2}{\sqrt{5}}V_{T}^{\pi} & V_{C}^{\pi} + \frac{2}{5}V_{T}^{\pi} \end{pmatrix},$$

$$V_{1/2^{+}}^{v} = \begin{pmatrix} V_{C}^{v'} & 2\sqrt{3}V_{C}^{v} & \sqrt{6}V_{C}^{v} \\ 2\sqrt{3}V_{C}^{v} & V_{C}^{v'} - 4V_{C}^{v} & \sqrt{2}V_{T}^{v} \\ \sqrt{6}V_{T}^{v} & \sqrt{2}V_{T}^{v} & V_{C}^{v'} + 2V_{C}^{v} + 2V_{T}^{v} \end{pmatrix},$$

$$(A6)$$

$$(A6)$$

The
$$\pi$$
 exchange potentials for each J^{P} are obtained by
$$V_{1/2^{-}}^{\pi} = \begin{pmatrix} 0 & \sqrt{3}V_{C}^{\pi} & -\sqrt{6}V_{T}^{\pi} \\ \sqrt{3}V_{C}^{\pi} & -2V_{C}^{\pi} & -\sqrt{2}V_{T}^{\pi} \\ -\sqrt{6}V_{T}^{\pi} & -\sqrt{2}V_{T}^{\pi} & V_{C}^{\pi} - 2V_{T}^{\pi} \end{pmatrix}, \quad (A1) \qquad V_{7/2^{-}}^{\pi} = \begin{pmatrix} 0 & 3\sqrt{\frac{3}{7}}V_{T}^{\pi} & \sqrt{3}V_{C}^{\pi} & -\sqrt{\frac{15}{7}}V_{T}^{\pi} \\ 3\sqrt{\frac{3}{7}}V_{T}^{\pi} & V_{C}^{\pi} - \frac{4}{7}V_{T}^{\pi} & \frac{3}{\sqrt{7}}V_{T}^{\pi} & \frac{6}{7}\sqrt{5}V_{T}^{\pi} \\ -\sqrt{\frac{15}{7}}V_{T}^{\pi} & \frac{6}{7}\sqrt{5}V_{T}^{\pi} & -2V_{C}^{\pi} & -\sqrt{\frac{5}{7}}V_{T}^{\pi} \\ -\sqrt{\frac{15}{7}}V_{T}^{\pi} & \frac{6}{7}\sqrt{5}V_{T}^{\pi} & -\sqrt{\frac{5}{7}}V_{T}^{\pi} & V_{C}^{\pi} + \frac{4}{7}V_{T}^{\pi} \end{pmatrix}, \quad (A7)$$

$$V_{1/2^{+}}^{\pi} = \begin{pmatrix} 0 & \sqrt{3}V_{C}^{\pi} & -\sqrt{6}V_{T}^{\pi} \\ \sqrt{3}V_{C}^{\pi} & -2V_{C}^{\pi} & -\sqrt{2}V_{T}^{\pi} \\ -\sqrt{6}V_{T}^{\pi} & -\sqrt{2}V_{T}^{\pi} & V_{C}^{\pi} - 2V_{T}^{\pi} \end{pmatrix}, \quad (A2)$$

$$V_{7/2^{+}}^{\pi} = \begin{pmatrix} 0 & \sqrt{3}V_{C}^{\pi} & V_{T}^{\pi} & -\sqrt{5}V_{T}^{\pi} \\ \sqrt{3}V_{C}^{\pi} & -2V_{C}^{\pi} & \frac{1}{\sqrt{3}}V_{T}^{\pi} & -\sqrt{\frac{5}{3}}V_{T}^{\pi} \\ V_{T}^{\pi} & \frac{1}{\sqrt{3}}V_{T}^{\pi} & V_{C}^{\pi} + \frac{4}{3}V_{T}^{\pi} & \frac{2}{3}\sqrt{5}V_{T}^{\pi} \\ \sqrt{3}V_{T}^{\pi} & V_{C}^{\pi} & -\sqrt{\frac{5}{3}}V_{T}^{\pi} & V_{C}^{\pi} - \frac{4}{3}V_{T}^{\pi} \end{pmatrix}, \quad (A2)$$

$$(A3)$$

$$V_C^{\pi} = \frac{g_{\pi}g_{\pi NN}}{\sqrt{2}m_N f_{\pi}} \frac{1}{3} C_{m_{\pi}} \vec{\tau}_P \cdot \vec{\tau}_N,$$

$$V_T^{\pi} = \frac{g_{\pi}g_{\pi NN}}{\sqrt{2}m_N f_{\pi}} \frac{1}{3} T_{m_{\pi}} \vec{\tau}_P \cdot \vec{\tau}_N.$$
(A9)

The $\vec{\tau}_P$ and $\vec{\tau}_N$ are the isospin matrices for $P^{(*)}$ and N. The functions $C_m = C(r; m)$ and $T_m = T(r; m)$ are given by

$$C(r;m) = \int \frac{d^3q}{(2\pi)^3} \frac{m^2}{\vec{q}^2 + m^2} e^{i\vec{q}\cdot\vec{r}} F(\Lambda_P, \vec{q}) F(\Lambda_N, \vec{q}),$$
(A10)

$$\begin{split} T(r;m)S_{12}(\hat{r}) &= \int \frac{d^3q}{(2\pi)^3} \frac{-\vec{q}^2}{\vec{q}^2 + m^2} S_{12}(\hat{q}) e^{i\vec{q}\cdot\vec{r}} F(\Lambda_P,\vec{q}) F(\Lambda_N,\vec{q}), \end{split} \tag{A11}$$

with $S_{12}(\hat{x}) = 3(\vec{\sigma}_1 \cdot \hat{x})(\vec{\sigma}_2 \cdot \hat{x}) - \vec{\sigma}_1 \cdot \vec{\sigma}_2$, and $F(\Lambda, \vec{q})$ denotes the form factor given in Eq. (8).

The vector meson exchange potentials ($v = \rho, \omega$) are

$$V_{1/2^{-}}^{v} = \begin{pmatrix} V_{C}^{vl} & 2\sqrt{3}V_{C}^{v} & \sqrt{6}V_{T}^{v} \\ 2\sqrt{3}V_{C}^{v} & V_{C}^{vl} - 4V_{C}^{v} & \sqrt{2}V_{T}^{v} \\ \sqrt{6}V_{T}^{v} & \sqrt{2}V_{T}^{v} & V_{C}^{vl} + 2V_{C}^{v} + 2V_{T}^{v} \end{pmatrix},$$
(A12)

$$V_{1/2^{+}}^{v} = \begin{pmatrix} V_{C}^{v\prime} & 2\sqrt{3}V_{C}^{v} & \sqrt{6}V_{T}^{v} \\ 2\sqrt{3}V_{C}^{v} & V_{C}^{v\prime} - 4V_{C}^{v} & \sqrt{2}V_{T}^{v} \\ \sqrt{6}V_{T}^{v} & \sqrt{2}V_{T}^{v} & V_{C}^{v\prime} + 2V_{C}^{v} + 2V_{T}^{v} \end{pmatrix}, \tag{A13}$$

$$V_{3/2^{-}}^{v} = \begin{pmatrix} V_{C}^{v\prime\prime} & -\sqrt{3}V_{T}^{v} & \sqrt{3}V_{T}^{v} & 2\sqrt{3}V_{C}^{v} \\ -\sqrt{3}V_{T}^{v} & V_{C}^{v\prime\prime} + 2V_{C}^{v} & -2V_{T}^{v} & -V_{T}^{v} \\ \sqrt{3}V_{T}^{v} & -2V_{T}^{v} & V_{C}^{v\prime\prime} + 2V_{C}^{v} & V_{T}^{v} \\ 2\sqrt{3}V_{C}^{v} & -V_{T}^{v} & V_{T}^{v} & V_{C}^{v\prime\prime} - 4V_{C}^{v} \end{pmatrix},$$
(A14)

$$V_{3/2^{+}}^{v} = \begin{pmatrix} V_{C}^{v\prime} & 2\sqrt{3}V_{C}^{v} & -\sqrt{\frac{3}{5}}V_{T}^{v} & 3\sqrt{\frac{3}{5}}V_{T}^{v} \\ 2\sqrt{3}V_{C}^{v} & V_{C}^{v\prime} - 4V_{C}^{v} & -\frac{1}{\sqrt{5}}V_{T}^{v} & \frac{3}{\sqrt{5}}V_{T}^{v} \\ -\sqrt{\frac{3}{5}}V_{T}^{v} & -\frac{1}{\sqrt{5}}V_{T}^{v} & V_{C}^{v\prime} + 2V_{C}^{v} - \frac{8}{5}V_{T}^{v} & -\frac{6}{5}V_{T}^{v} \\ 3\sqrt{\frac{3}{5}}V_{T}^{v} & \frac{3}{\sqrt{5}}V_{T}^{v} & -\frac{6}{5}V_{T}^{v} & V_{C}^{v\prime} + 2V_{C}^{v} + \frac{8}{5}V_{T}^{v} \end{pmatrix},$$
(A15)

$$V_{5/2^{-}}^{v} = \begin{pmatrix} V_{C}^{v\prime} & 2\sqrt{3}V_{C}^{v} & -\sqrt{\frac{6}{7}}V_{T}^{v} & \frac{6}{\sqrt{7}}V_{T}^{v} \\ 2\sqrt{3}V_{C}^{v} & V_{C}^{v\prime} - 4V_{C}^{v} & -\sqrt{\frac{2}{7}}V_{T}^{v} & 2\sqrt{\frac{5}{7}}V_{T}^{v} \\ -\sqrt{\frac{6}{7}}V_{T}^{v} & -\sqrt{\frac{2}{7}}V_{T}^{v} & V_{C}^{v\prime} + 2V_{C}^{v} - \frac{10}{7}V_{T}^{v} & -\frac{4}{7}\sqrt{6}V_{T}^{v} \\ \frac{6}{\sqrt{7}}V_{T}^{v} & 2\sqrt{\frac{3}{7}}V_{T}^{v} & -\frac{4}{7}\sqrt{6}V_{T}^{v} & V_{C}^{v\prime} + 2V_{C}^{v} + \frac{10}{7}V_{T}^{v} \end{pmatrix},$$
(A16)

$$V_{5/2^{+}}^{v} = \begin{pmatrix} V_{C}^{v\prime} & -\frac{3}{5}\sqrt{10}V_{T}^{v} & 2\sqrt{3}V_{C}^{v} & 2\sqrt{\frac{3}{5}}V_{T}^{v} \\ -\frac{3}{5}\sqrt{10}V_{T}^{v} & V_{C}^{v\prime} + 2V_{C}^{v} + \frac{2}{5}V_{T}^{v} & -\sqrt{\frac{6}{5}}V_{T}^{v} & -\frac{4}{5}\sqrt{6}V_{T}^{v} \\ 2\sqrt{3}V_{C}^{v} & -\sqrt{\frac{6}{5}}V_{T}^{v} & V_{C}^{v\prime} - 4V_{C}^{v} & \frac{2}{\sqrt{5}}V_{T}^{v} \\ 2\sqrt{\frac{3}{5}}V_{T}^{v} & -\frac{4}{5}\sqrt{6}V_{T}^{v} & \frac{2}{\sqrt{5}}V_{T}^{v} & V_{C}^{v\prime} + 2V_{C}^{v} - \frac{2}{5}V_{T}^{v} \end{pmatrix},$$
(A17)

$$V_{7/2^{-}}^{\nu} = \begin{pmatrix} V_{C}^{\nu\prime} & -3\sqrt{\frac{3}{7}}V_{T}^{\nu} & 2\sqrt{3}V_{C}^{\nu} & \sqrt{\frac{15}{7}}V_{T}^{\nu} \\ -3\sqrt{\frac{3}{7}}V_{T}^{\nu} & V_{C}^{\nu\prime} + 2V_{C}^{\nu} + \frac{4}{7}V_{T}^{\nu} & -\frac{3}{\sqrt{7}}V_{T}^{\nu} & -\frac{6}{7}\sqrt{5}V_{T}^{\nu} \\ 2\sqrt{3}V_{C}^{\nu} & -\frac{3}{\sqrt{7}}V_{T}^{\nu} & V_{C}^{\nu\prime} - 4V_{C}^{\nu} & \sqrt{\frac{5}{7}}V_{T}^{\nu} \\ \sqrt{\frac{15}{7}}V_{T}^{\nu} & -\frac{6}{7}\sqrt{5}V_{T}^{\nu} & \sqrt{\frac{5}{7}}V_{T}^{\nu} & V_{C}^{\nu\prime} + 2V_{C}^{\nu} - \frac{4}{7}V_{T}^{\nu} \end{pmatrix},$$
(A18)

$$V_{7/2^{+}}^{v} = \begin{pmatrix} V_{C}^{v\prime} & 2\sqrt{3}V_{C}^{v} & -V_{T}^{v} & \sqrt{5}V_{T}^{v} \\ 2\sqrt{3}V_{C}^{v} & V_{C}^{v\prime} - 4V_{C}^{v} & -\frac{1}{\sqrt{3}}V_{T}^{v} & \sqrt{\frac{5}{3}}V_{T}^{v} \\ -V_{T}^{v} & -\frac{1}{\sqrt{3}}V_{T}^{v} & V_{C}^{v\prime} + 2V_{C}^{v} - \frac{4}{3}V_{T}^{v} & -\frac{2}{3}\sqrt{5}V_{T}^{v} \\ \sqrt{5}V_{T}^{v} & \sqrt{\frac{5}{3}}V_{T}^{v} & -\frac{2}{3}\sqrt{5}V_{T}^{v} & V_{C}^{v\prime} + 2V_{C}^{v} + \frac{4}{3}V_{T}^{v} \end{pmatrix}, \tag{A19}$$

where $V_C^{v\prime}$, V_C^v , and V_T^v are defined as

$$V_C^{\rho\prime} = \frac{g_V g_{\rho NN} \beta}{\sqrt{2} m_\rho^2} C_{m_\rho} \vec{\tau}_P \cdot \vec{\tau}_N, \tag{A20}$$

$$V_C^{\rho} = \frac{g_V g_{\rho NN} \lambda (1+\kappa)}{\sqrt{2} m_N} \frac{1}{3} C_{m_{\rho}} \vec{\tau}_P \cdot \vec{\tau}_N, \tag{A21}$$

$$V_T^{\rho} = \frac{g_V g_{\rho NN} \lambda (1 + \kappa)}{\sqrt{2} m_N} \frac{1}{3} T_{m_{\rho}} \vec{\tau}_P \cdot \vec{\tau}_N, \tag{A22}$$

$$V_C^{\omega\prime} = \frac{g_V g_{\omega NN} \beta}{\sqrt{2} m_\omega^2} C_{m_\omega},\tag{A23}$$

$$V_C^{\omega} = \frac{g_V g_{\omega NN} \lambda}{\sqrt{2} m_N} \frac{1}{3} C_{m_{\omega}},\tag{A24}$$

$$V_T^{\omega} = \frac{g_V g_{\omega NN} \lambda}{\sqrt{2} m_N} \frac{1}{3} T_{m_{\omega}}.$$
 (A25)

Finally, the kinetic terms are given by

$$K_{1/2^{-}} = \operatorname{diag}\left(-\frac{1}{2\tilde{m}_{P}}\Delta_{0}, -\frac{1}{2\tilde{m}_{P^{*}}}\Delta_{0} + \Delta m_{PP^{*}}, -\frac{1}{2\tilde{m}_{P^{*}}}\Delta_{2} + \Delta m_{PP^{*}}\right),\tag{A26}$$

$$K_{1/2^{+}} = \operatorname{diag}\left(-\frac{1}{2\tilde{m}_{P}}\Delta_{1}, -\frac{1}{2\tilde{m}_{P^{*}}}\Delta_{1} + \Delta m_{PP^{*}}, -\frac{1}{2\tilde{m}_{P^{*}}}\Delta_{1} + \Delta m_{PP^{*}}\right),\tag{A27}$$

$$K_{3/2^{-}} = \operatorname{diag}\left(-\frac{1}{2\tilde{m}_{P}}\Delta_{2}, -\frac{1}{2\tilde{m}_{P^{*}}}\Delta_{0} + \Delta m_{PP^{*}}, -\frac{1}{2\tilde{m}_{P^{*}}}\Delta_{2} + \Delta m_{PP^{*}}, -\frac{1}{2\tilde{m}_{P^{*}}}\Delta_{2} + \Delta m_{PP^{*}}\right), \tag{A28}$$

$$K_{3/2^{+}} = \operatorname{diag}\left(-\frac{1}{2\tilde{m}_{P}}\Delta_{1}, -\frac{1}{2\tilde{m}_{P^{*}}}\Delta_{1} + \Delta m_{PP^{*}}, -\frac{1}{2\tilde{m}_{P^{*}}}\Delta_{1} + \Delta m_{PP^{*}}, -\frac{1}{2\tilde{m}_{P^{*}}}\Delta_{3} + \Delta m_{PP^{*}}\right), \tag{A29}$$

$$K_{5/2^{-}} = \operatorname{diag}\left(-\frac{1}{2\tilde{m}_{P}}\triangle_{2}, -\frac{1}{2\tilde{m}_{P^{*}}}\triangle_{2} + \Delta m_{PP^{*}}, -\frac{1}{2\tilde{m}_{P^{*}}}\triangle_{2} + \Delta m_{PP^{*}}, -\frac{1}{2\tilde{m}_{P^{*}}}\triangle_{4} + \Delta m_{PP^{*}}\right), \tag{A30}$$

$$K_{5/2^+} = \mathrm{diag} \left(-\frac{1}{2\tilde{m}_P} \triangle_3, -\frac{1}{2\tilde{m}_{P^*}} \triangle_1 + \Delta m_{PP^*}, -\frac{1}{2\tilde{m}_{P^*}} \triangle_3 + \Delta m_{PP^*}, -\frac{1}{2\tilde{m}_{P^*}} \triangle_3 + \Delta m_{PP^*} \right), \tag{A31}$$

$$K_{7/2^{-}} = \operatorname{diag}\left(-\frac{1}{2\tilde{m}_{P}}\triangle_{4}, -\frac{1}{2\tilde{m}_{P^{*}}}\triangle_{2} + \Delta m_{PP^{*}}, -\frac{1}{2\tilde{m}_{P^{*}}}\triangle_{4} + \Delta m_{PP^{*}}, -\frac{1}{2\tilde{m}_{P^{*}}}\triangle_{4} + \Delta m_{PP^{*}}\right), \tag{A32}$$

$$K_{7/2^{+}} = \operatorname{diag}\left(-\frac{1}{2\tilde{m}_{P}}\Delta_{3}, -\frac{1}{2\tilde{m}_{P^{*}}}\Delta_{3} + \Delta m_{PP^{*}}, -\frac{1}{2\tilde{m}_{P^{*}}}\Delta_{3} + \Delta m_{PP^{*}}, -\frac{1}{2\tilde{m}_{P^{*}}}\Delta_{5} + \Delta m_{PP^{*}}\right), \tag{A33}$$

where $\triangle_l = \partial^2/\partial r^2 + (2/r)\partial/\partial r - l(l+1)/r^2$, $\tilde{m}_{P^{(*)}} = m_N m_{P^{(*)}}/(m_N + m_{P^{(*)}})$, and $\Delta m_{PP^*} = m_{P^*} - m_P$. The total Hamiltonian is then given by $H_{J^P} = K_{J^P} + V_{J^P}$.

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