Dynamics of hadronic molecule in one-boson exchange approach and possible heavy flavor molecules

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We extend the one pion exchange model at quark level to include the short distance contributions coming from η , σ , ρ and ω exchange. This formalism is applied to discuss the possible molecular states of $D\bar{D}^*/\bar{D}D^*$, $B\bar{B}^*/\bar{B}B^*$, DD^* , BB^* , the pseudoscalar-vector systems with C = B = 1 and C = -B = 1respectively. The " δ function" term contribution and the *S*-*D* mixing effects have been taken into account. We find the conclusions reached after including the heavier mesons exchange are qualitatively the same as those in the one pion exchange model. The previous suggestion that $1^{++} B\bar{B}^*/\bar{B}B^*$ molecule should exist, is confirmed in the one-boson exchange model, whereas DD^* bound state should not exist. The $D\bar{D}^*/\bar{D}D^*$ system can accommodate a 1^{++} molecule close to the threshold, the mixing between the molecule and the conventional charmonium has to be considered to identify this state with X(3872). For the BB^* system, the pseudoscalar-vector systems with C = B = 1 and C = -B = 1, near threshold molecular states may exist. These bound states should be rather narrow, isospin is violated and the I = 0component is dominant. Experimental search channels for these states are suggested.

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I. INTRODUCTION

Since the 1970s it is widely believed that quantum chromodynamics should accommodate a richer spectrum than just $q\bar{q}$ and qqq resonances, many possible nonconventional structures are suggested, e.g. glueballs (gg, ggg, ...), hybrid mesons $(q\bar{q}g)$ and multiquark states $(qq\bar{q}\bar{q}, qqqq\bar{q}, qqqq\bar{q}, qqq\bar{q}\bar{q}\bar{q}\bar{q}\bar{q})$. Unfortunately, so far there is still no uncontroversial evidence for nonconventional states experimentally except the hadronic molecules. The deuteron is a well-known example of hadronic molecule, and the approximate 10⁵ known nuclear levels are all hadronic molecule. In the past few years, many new states have been reported, a striking feature is that some of them are close to the thresholds of certain two hadrons, which inspires the possible interpretation of hadronic molecule.

Hadronic molecule is an old idea, about 30 years ago the possible hadronic molecules consisting of two charm mesons are suggested [1], and $\psi(4040)$ was proposed to be a *P* wave D^*D^* molecule [2]. Since in general a molecule is weakly bound, the separation between the two hadrons in the molecule should be large. We can picture the two hadrons as interacting via a meson exchange potential [3]. At large distance, one pion exchange is dominant. Guided by the binding of deuteron, Tornqvist performed a systematic study of possible deuteronlike two-meson bound states [4,5]. The role of pion exchange in forming hadronic molecules was studied by Ericson and Karl [6]. Recently Close et al. [7] performed a pedagogic analysis of the overall sign, in addition they included the contribution of the " δ function" term which gives a δ function in the effective potential when no regularization is used. In these

original works, only long distance one pion exchange has been considered, and the short distance contributions are neglected. In Ref. [8] Swanson assumed that the short distance dynamics is governed by the one gluon exchange induced constituent quark interchange mechanism, which results in state mixing.

In the model of the nucleon-nucleon interaction, the long range part of the nucleon-nucleon force is quantitatively accounted for by the one π exchange. However, the short and intermediate range interactions are governed by more complex dynamics. Combining the well-established one π exchange with the exchange of heavier bosons (e.g. scalar and vector mesons) to describe the behavior at short distance has been proved to be a very successful approach [9-11]. Physically, the scalar and vector meson exchange describes part of multiple pion exchange effects. For the two π exchange, if they interact and correlate in a P wave state, such a exchange can be modeled by ρ exchange. If the two correlated π pair is in a S-wave state, Durso *et al.* showed that one can approximate them by the exchange of a scalar σ meson [12]. Similarly, the correlated 3π exchange can be approximated by the exchange of one ω meson.

Inspired by the nucleon-nucleon interaction, we shall represent the short distance interactions by the heavier bosons η , σ , ρ , and ω exchange instead of the quark interchange. The effective potential between two hadrons is obtained by summing over the interactions between light quarks or antiquarks as in the original work [4,5,7]. It is well-known that one pion exchange between two light quarks results in two terms: the isospin dependent spinspin interaction and tensor force. After taking into account the heavy bosons exchange, six additional terms appear

including the spin-isospin independent central term, only isospin dependent term, isospin independent spin-spin interaction and tensor force, both isospin dependent and independent spin-orbit interactions. Consequently the situation becomes more complex than the only one pion exchange model. In our model, both the " δ function" term and the S-D mixing effects would be considered, which have been shown to play an important role in the binding [4,5,7]. In this work, we first give a good description of the deuteron in our model, which is an unambiguous hadronic molecule, then apply this formalism to the heavy flavor pseudoscalar-vector (PV) systems. Thus the predictions for the possible heavy flavor PV molecules are base on a solid and reliable foundation. This is a greater advantage over other approaches dealing with the dynamics of hadronic molecule, such as one-boson exchange in the effective field theory [13,14] and residual strong force with pairwise interactions [15,16] etc.

The paper is organized as follows. In Sec. II, the formalism of the one-boson exchange model is presented, the effective potentials from pseudoscalar, scalar and vector meson exchange are given explicitly. In Sec. III, we give the meson parameters involved in our model and the boson-quark couplings which are extracted from the boson-nucleon couplings. The formalism is applied to the deuteron in Sec. IV, the $D\overline{D}^*/\overline{D}D^*$ system and the molecular interpretation of X(3872) are investigated in Sec. V. We further apply the one-boson exchange approach to other heavy flavor PV systems in Sec. VI, and possible molecular states are discussed. Section VII is our conclusions and discussions section. The expressions for the matrix elements of the spin relevant operators are analytically given in the Appendix.

II. THE FORMALISM OF ONE-BOSON EXCHANGE MODEL

The construction of one-boson exchange interaction is constrained by the symmetry principle. To leading order in the boson fields and their derivative, the effective interaction Lagrangian describing the coupling between the constituent quarks and the exchange boson fields is as follows [9-11]

Pseudoscalar:
$$\mathcal{L}_{p} = -g_{pqq}\bar{\psi}(x)i\gamma_{5}\psi(x)\varphi(x)$$

Scalar: $\mathcal{L}_{s} = -g_{sqq}\bar{\psi}(x)\psi(x)\phi(x)$
Vector: $\mathcal{L}_{v} = -g_{vqq}\bar{\psi}(x)\gamma_{\mu}\psi(x)v^{\mu}(x)$
 $-\frac{f_{vqq}}{2m_{q}}\bar{\psi}(x)\sigma_{\mu\nu}\psi(x)\partial^{\mu}v^{\nu}(x).$ (1)

Here m_q is the constituent quark mass, $\psi(x)$ is the constituent quark Dirac spinor field, $\varphi(x)$, $\phi(x)$, and $v^{\mu}(x)$ are the isospin-singlet pseudoscalar, scalar and vector boson fields, respectively. In this work we take $m_q \equiv m_u = m_d \approx 313$ MeV, since we concentrate on constituent up and

down quarks. If the isovector bosons are involved, the couplings enter in the form $\boldsymbol{\tau} \cdot \boldsymbol{\varphi}, \, \boldsymbol{\tau} \cdot \boldsymbol{\phi}$, and $\boldsymbol{\tau} \cdot \boldsymbol{v}^{\mu}$ respectively, where $\boldsymbol{\tau}$ is the well-known Pauli matrices. For the pseudoscalar, another interaction term is allowed $\mathcal{L}'_p = \frac{f_{pqq}}{m_p} \, \bar{\psi}(x) \gamma^{\mu} \gamma_5 \psi(x) \partial_{\mu} \varphi(x)$, where m_p is the exchange pseudoscalar mass, this Lagrangian has been used by Tornqvist [4,5] and Close [7]. By partial integration and using the equation of motion, one can easily show that \mathcal{L}_p and \mathcal{L}'_p are equivalent provided the coupling constants are related by

$$\frac{f_{pqq}}{m_p} = \frac{g_{pqq}}{2m_q}.$$
(2)

From the above effective interactions, the effective potential between two quarks in momentum space can be calculated straightforwardly following the standard procedure. To the leading order in \mathbf{q}^2/m_q^2 , where \mathbf{q} is the momentum transfer, the potentials are

(1) Pseudoscalar boson exchange

$$V_{p}(\mathbf{q}) = -\frac{g_{pqq}^{2}}{4m_{q}^{2}} \frac{(\boldsymbol{\sigma}_{i} \cdot \mathbf{q})(\boldsymbol{\sigma}_{j} \cdot \mathbf{q})}{\mathbf{q}^{2} + \mu_{p}^{2}}$$
$$= -\frac{g_{pqq}^{2}}{12m_{q}^{2}} \left[\frac{\mathbf{q}^{2}}{\mathbf{q}^{2} + \mu_{p}^{2}} \boldsymbol{\sigma}_{i} \cdot \boldsymbol{\sigma}_{j} + \frac{\mathbf{q}^{2}S_{ij}(\hat{\mathbf{q}})}{\mathbf{q}^{2} + \mu_{p}^{2}} \right],$$
(3)

where $S_{ij}(\hat{\mathbf{q}}) = 3(\boldsymbol{\sigma}_i \cdot \hat{\boldsymbol{q}})(\boldsymbol{\sigma}_j \cdot \hat{\boldsymbol{q}}) - \boldsymbol{\sigma}_i \cdot \boldsymbol{\sigma}_j$, we have used $\mu_p^2 = m_p^2 - q_0^2$ instead of m_p^2 to approximately account for the recoil effect [4,5,7].

(2) Scalar boson exchange

$$V_{s}(\mathbf{q}) = -\frac{g_{sqq}^{2}}{\mathbf{q}^{2} + \mu_{s}^{2}} \left(1 + \frac{\mathbf{q}^{2}}{8m_{q}^{2}}\right) - \frac{g_{sqq}^{2}}{2m_{q}^{2}}$$
$$\times \frac{i\mathbf{S}_{ij} \cdot (\mathbf{p} \times \mathbf{q})}{\mathbf{q}^{2} + \mu_{s}^{2}}, \tag{4}$$

where $\mathbf{S}_{ij} \equiv \frac{1}{2}(\boldsymbol{\sigma}_i + \boldsymbol{\sigma}_j)$, $\mu_s^2 = m_s^2 - q_0^2$ with m_s the exchange scalar meson mass, and **p** denotes the total momentum.

(3) Vector boson exchange

$$V_{\nu}(\mathbf{q}) = \frac{g_{\nu qq}^{2}}{\mathbf{q}^{2} + \mu_{\nu}^{2}} - \frac{g_{\nu qq}^{2} + 4g_{\nu qq}f_{\nu qq}}{8m_{q}^{2}} \frac{\mathbf{q}^{2}}{\mathbf{q}^{2} + \mu_{\nu}^{2}} \\ + \frac{(g_{\nu qq} + f_{\nu qq})^{2}}{12m_{q}^{2}} \frac{\mathbf{q}^{2}S_{ij}(\hat{\mathbf{q}}) - 2\mathbf{q}^{2}(\boldsymbol{\sigma}_{i} \cdot \boldsymbol{\sigma}_{j})}{\mathbf{q}^{2} + \mu_{\nu}^{2}} \\ - \frac{3g_{\nu qq}^{2} + 4g_{\nu qq}f_{\nu qq}}{2m_{q}^{2}} \frac{i\mathbf{S}_{ij} \cdot (\mathbf{p} \times \mathbf{q})}{\mathbf{q}^{2} + \mu_{\nu}^{2}}, \quad (5)$$

where $\mu_v^2 = m_v^2 - q_0^2$ approximately reflects the recoil effect with m_v the exchange vector meson mass.

The effective potential in configuration space is obtained by Fourier transforming the momentum space potential. DYNAMICS OF HADRONIC MOLECULE IN ONE-BOSON ...

$$V_i(\mathbf{r}) = \frac{1}{(2\pi)^3} \int d^3 \mathbf{q} e^{i\mathbf{q}\cdot\mathbf{r}} V_i(\mathbf{q}), \tag{6}$$

where i = p, s, and v respectively. However, the resulting potentials are singular, which contains the delta function, so the potentials have to be regularized. Considering the internal structure of the involved hadrons, one usually introduces form factor at each vertex. Here the form factor is taken as

$$F(q) = \frac{\Lambda^2 - m^2}{\Lambda^2 - q^2} = \frac{\Lambda^2 - m^2}{X^2 + \mathbf{q}^2},$$
(7)

where Λ is the so-called regularization parameter, *m*, and *q* are the mass and the four momentum of the exchanged boson, respectively, with $X^2 = \Lambda^2 - q_0^2$. The form factor suppresses the contribution of high momentum, i.e. small

distance. The presence of such a form factor is dictated by the extended structure of the hadrons. The parameter Λ , which governs the range of suppression, can be directly related to the hadron size that is approximately proportional to $1/\Lambda$. However, since the question of hadron size is still very much open, the value of Λ is poorly known phenomenologically, and it is dependent on the model and application. In the nucleon-nucleon interaction, the Λ in the range 0.8–1.5 GeV has been used to fit the data. For the present application to heavy mesons, which have a much smaller size than nucleon, we would expect a larger regularization parameter Λ . We can straightforwardly obtain the effective potentials between two quarks in configuration space. For convenience, the following dimensionless functions are introduced.

$$\begin{split} H_{0}(\Lambda, m_{ex}, \mu, r) &= \frac{1}{\mu r} (e^{-\mu r} - e^{-Xr}) - \frac{\Lambda^{2} - m_{ex}^{2}}{2\mu X} e^{-Xr} \\ H_{1}(\Lambda, m_{ex}, \mu, r) &= -\frac{1}{\mu r} (e^{-\mu r} - e^{-Xr}) + \frac{X(\Lambda^{2} - m_{ex}^{2})}{2\mu^{3}} e^{-Xr} \\ H_{2}(\Lambda, m_{ex}, \mu, r) &= \left(1 + \frac{1}{\mu r}\right) \frac{1}{\mu^{2} r^{2}} e^{-\mu r} - \left(1 + \frac{1}{Xr}\right) \frac{X}{\mu} \frac{1}{\mu^{2} r^{2}} e^{-Xr} - \frac{\Lambda^{2} - m_{ex}^{2}}{2\mu^{2}} \frac{e^{-Xr}}{\mu r} \\ H_{3}(\Lambda, m_{ex}, \mu, r) &= \left(1 + \frac{3}{\mu r} + \frac{3}{\mu^{2} r^{2}}\right) \frac{1}{\mu r} e^{-\mu r} - \left(1 + \frac{3}{Xr} + \frac{3}{X^{2} r^{2}}\right) \frac{X^{2}}{\mu^{2}} \frac{e^{-Xr}}{\mu r} - \frac{\Lambda^{2} - m_{ex}^{2}}{2\mu^{2}} (1 + Xr) \frac{e^{-Xr}}{\mu r} \\ G_{1}(\Lambda, m_{ex}, \tilde{\mu}, r) &= \frac{1}{\tilde{\mu} r} [\cos(\tilde{\mu}r) - e^{-Xr}] + \frac{X(\Lambda^{2} - m_{ex}^{2})}{2\tilde{\mu}^{3}} e^{-Xr} \\ G_{3}(\Lambda, m_{ex}, \tilde{\mu}, r) &= -\left[\cos(\tilde{\mu}r) - \frac{3\sin(\tilde{\mu}r)}{\tilde{\mu}r} - \frac{3\cos(\tilde{\mu}r)}{\tilde{\mu}^{2} r^{2}}\right] \frac{1}{\tilde{\mu} r} - \left(1 + \frac{3}{Xr} + \frac{3}{X^{2} r^{2}}\right) \frac{X^{2}}{\tilde{\mu}^{2}} \frac{e^{-Xr}}{\tilde{\mu} r} - \frac{\Lambda^{2} - m_{ex}^{2}}{2\tilde{\mu}^{2}} (1 + Xr) \frac{e^{-Xr}}{\mu r} \\ \end{cases}$$
(8)

Then the effective potentials between two quarks from one-boson exchange are

(1) Pseudoscalar boson exchange

$$V_{p}(\mathbf{r}) = \begin{cases} \frac{g_{pqq}^{2}}{4\pi} \frac{\mu_{p}^{3}}{12m_{q}^{2}} [-H_{1}(\Lambda, m_{p}, \mu_{p}, r)\boldsymbol{\sigma}_{i} \cdot \boldsymbol{\sigma}_{j} + H_{3}(\Lambda, m_{p}, \mu_{p}, r)S_{ij}(\hat{\mathbf{r}})], & \mu_{p}^{2} > 0\\ \frac{g_{pqq}^{2}}{4\pi} \frac{\tilde{\mu}_{p}^{3}}{12m_{q}^{2}} [-G_{1}(\Lambda, m_{p}, \tilde{\mu}_{p}, r)\boldsymbol{\sigma}_{i} \cdot \boldsymbol{\sigma}_{j} + G_{3}(\Lambda, m_{p}, \tilde{\mu}_{p}, r)S_{ij}(\hat{\mathbf{r}})], & \mu_{p}^{2} = -\tilde{\mu}_{p}^{2} < 0 \end{cases}$$
(9)

with $S_{ij}(\hat{\mathbf{r}}) = 3(\boldsymbol{\sigma}_i \cdot \hat{\mathbf{r}})(\boldsymbol{\sigma}_j \cdot \hat{\mathbf{r}}) - \boldsymbol{\sigma}_i \cdot \boldsymbol{\sigma}_j$. (2) Scalar boson exchange

$$V_{s}(\mathbf{r}) = -\mu_{s} \frac{g_{sqq}^{2}}{4\pi} \bigg[H_{0}(\Lambda, m_{s}, \mu_{s}, r) + \frac{\mu_{s}^{2}}{8m_{q}^{2}} H_{1}(\Lambda, m_{s}, \mu_{s}, r) + \frac{\mu_{s}^{2}}{2m_{q}^{2}} H_{2}(\Lambda, m_{s}, \mu_{s}, r) \mathbf{L} \cdot \mathbf{S}_{ij} \bigg].$$
(10)

Here $\mathbf{L} = \mathbf{r} \times \mathbf{p}$ is the angular momentum operator.

(3) Vector boson exchange

$$V_{v}(\mathbf{r}) = \frac{\mu_{v}}{4\pi} \bigg\{ g_{vqq}^{2} H_{0}(\Lambda, m_{v}, \mu_{v}, r) - \frac{(g_{vqq}^{2} + 4g_{vqq}f_{vqq})\mu_{v}^{2}}{8m_{q}^{2}} H_{1}(\Lambda, m_{v}, \mu_{v}, r) - (g_{vqq} + f_{vqq})^{2} \frac{\mu_{v}^{2}}{12m_{q}^{2}} [H_{3}(\Lambda, m_{v}, \mu_{v}, r)S_{ij}(\hat{\mathbf{r}}) + 2H_{1}(\Lambda, m_{v}, \mu_{v}, r)(\boldsymbol{\sigma_{i}} \cdot \boldsymbol{\sigma_{j}})] - (3g_{vqq}^{2} + 4g_{vqq}f_{vqq}) \frac{\mu_{v}^{2}}{2m_{q}^{2}} H_{2}(\Lambda, m_{v}, \mu_{v}, r)\mathbf{L} \cdot \mathbf{S}_{ij} \bigg\}.$$

$$(11)$$

For I = 1 isovector boson exchange, the above potential should be multiplied by the operator $\tau_i \cdot \tau_j$ in the isospin space. We have included the contribution of the " δ function" term in the above potentials, which gives the delta function when no regularization is used, since this contribution turns out to be important [7]. The effective potential between two hadrons are obtained by summing the interactions between light quarks or antiquarks via one-boson exchange.

III. MESON PARAMETERS AND COUPLING CONSTANTS

As the well-known nuclear-nuclear interaction in the one-boson exchange model, we shall take into account the contributions from pseudoscalar mesons π and η exchange, that from scalar meson σ exchange, and those from vector mesons ρ and ω exchange. The basic input parameters are the boson masses and the effective coupling constants between the exchanged bosons and the constituent quarks. The meson masses with their quantum numbers are taken from the compilation of the Particle Data Group [17]. For the constituent quark-meson coupling constants, one may derive suitable estimates from the phenomenologically known πNN , ηNN , σNN , ρNN and ωNN coupling constants using the Goldberger-Treiman relation. Riska and Brown have demonstrated that the nucleon resonance transition couplings to π , ρ and ω can be derived in the single-quark operator approximation [18], which are in good agreement with the experimental data. Along the same way, we can straightforwardly derive the following relations between the boson-quark couplings and the boson-nucleon couplings,

$$g_{\pi qq} = \frac{3}{5} \frac{m_q}{m_N} g_{\pi NN}, \qquad g_{\eta qq} = \frac{m_q}{m_N} g_{\eta NN}$$

$$g_{\rho qq} = g_{\rho NN}, \qquad f_{\rho qq} = \frac{3}{5} \frac{m_q}{m_N} f_{\rho NN} - \left(1 - \frac{3}{5} \frac{m_q}{m_N}\right) g_{\rho NN}$$

$$g_{\omega qq} = \frac{1}{3} g_{\omega NN}, \qquad f_{\omega qq} = \frac{m_q}{m_N} f_{\omega NN} - \left(\frac{1}{3} - \frac{m_q}{m_N}\right) g_{\omega NN}$$

$$g_{\sigma qq} = \frac{1}{3} g_{\sigma NN}, \qquad (12)$$

where m_N is the nucleon mass. In the present work, the constituent up (down) quark mass $m_{u(d)}$ is taken to be usual

TABLE I. Spin, parity, isospin, *G*-parity, the masses of the exchange bosons, and the meson-nucleon coupling constants in the model.

Boson	$I^G(J^P)$	Mass (MeV)	$g^2/4\pi$	$f^2/4\pi$
π^{\pm}	$1^{-}(0^{-})$	139.57	14.9	
π^0	$1^{-}(0^{-})$	134.98	14.9	
η	$0^+(0^-)$	547.85	3.0	
σ	$0^+(0^+)$	600.0	7.78	
ρ	$1^{+}(1^{-})$	775.49	0.95	35.35
ω	$0^{-}(1^{-})$	782.65	20.0	0.0

value $m_{u(d)} \simeq 313$ MeV, which is about one third of the nucleon mass. The effective boson-nucleon coupling constants are taken from the well-known Bonn model [11], and a typical set of parameters is shown in Table I. The uncertainty of the effective couplings will be taken into account later, all the coupling constants except $g_{\pi NN}$ would be reduced by a factor of 2, since the experimental value for $g_{\pi NN}$ has been determined accurately from pion-nucleon and nucleon-nucleon scatterings. In the following, we shall explore the possible molecular states consisting of a pair heavy flavor pseudoscalar and vector mesons, their masses are taken from Particle Data Group [17]: $m_{D^0} =$ 1864.84 MeV, $m_{D^{\pm}} = 1869.62$ MeV, $m_{D^{*0}} =$ 2006.97 MeV, $m_{D^{*\pm}} = 2010.27$ MeV, $m_{B^0} =$ 5279.53 MeV, $m_{B^{\pm}} = 5279.15$ MeV, and $m_{B^*} =$ 5325.1 MeV.

IV. DEUTERON FROM ONE-BOSON EXCHANGE MODEL

Deuteron is an uncontroversial proton-neutron bound state with J = 1 and I = 0. It has been established that the long distance one pion exchange is the main binding mechanism, and the tensor force plays a crucial role, which results in the ${}^{3}S_{1}$ and ${}^{3}D_{1}$ states mixing. Tornqvist and Close only considered the pion exchange contribution in Refs. [5,7], however, the scalar meson σ exchange and the vector mesons ρ , ω exchange turn out to be important in providing the short distance repulsion and the intermediate range attraction, consequently, we shall take into account the contributions from the heavier boson exchange in the following. The effective potential becomes

$$V^{d}(\mathbf{r}) = V^{d}_{\pi}(\mathbf{r}) + V^{d}_{\eta}(\mathbf{r}) + V^{d}_{\sigma}(\mathbf{r}) + V^{d}_{\rho}(\mathbf{r}) + V^{d}_{\omega}(\mathbf{r})$$

$$\equiv V^{d}_{C}(r) + V^{d}_{S}(r)(\boldsymbol{\sigma}_{1} \cdot \boldsymbol{\sigma}_{2}) + V^{d}_{I}(r)(\boldsymbol{\tau}_{1} \cdot \boldsymbol{\tau}_{2}) + V^{d}_{T}(r)S_{12}(\hat{\mathbf{r}}) + V^{d}_{SI}(r)(\boldsymbol{\sigma}_{1} \cdot \boldsymbol{\sigma}_{2})(\boldsymbol{\tau}_{1} \cdot \boldsymbol{\tau}_{2}) + V^{d}_{TI}(r)S_{12}(\hat{\mathbf{r}})(\boldsymbol{\tau}_{1} \cdot \boldsymbol{\tau}_{2})$$

$$+ V^{d}_{LS}(r)(\mathbf{L} \cdot \mathbf{S}) + V^{d}_{LSI}(r)(\mathbf{L} \cdot \mathbf{S})(\boldsymbol{\tau}_{1} \cdot \boldsymbol{\tau}_{2}), \qquad (13)$$

where $\mathbf{S} = \frac{1}{2}(\boldsymbol{\sigma}_1 + \boldsymbol{\sigma}_2)$ is the total spin, and **L** is the relative angular momentum operator. In the isospin symmetry limit, the components $V_C^d(r)$, $V_S^d(r)$ etc are given by

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$$\begin{split} V_{C}^{d}(r) &= -\frac{g_{\sigma NN}^{2}}{4\pi} m_{\sigma} \bigg[H_{0}(\Lambda, m_{\sigma}, m_{\sigma}, r) + \frac{m_{\sigma}^{2}}{8m_{N}^{2}} H_{1}(\Lambda, m_{\sigma}, m_{\sigma}, r) \bigg] + \frac{g_{\omega NN}^{2}}{4\pi} m_{\omega} H_{0}(\Lambda, m_{\omega}, m_{\omega}, r) \\ &- \frac{g_{\omega NN}^{2} + 4g_{\omega NN} f_{\omega NN}}{4\pi} \frac{m_{\omega}^{3}}{8m_{N}^{2}} H_{1}(\Lambda, m_{\omega}, m_{\omega}, r) \\ V_{S}^{d}(r) &= -\frac{g_{\eta NN}^{2}}{4\pi} \frac{m_{\eta}^{3}}{12m_{N}^{2}} H_{1}(\Lambda, m_{\eta}, m_{\eta}, r) - \frac{(g_{\omega NN} + f_{\omega NN})^{2}}{4\pi} \frac{m_{\omega}^{3}}{6m_{N}^{2}} H_{1}(\Lambda, m_{\omega}, m_{\omega}, r) \\ V_{I}^{d}(r) &= \frac{g_{\rho NN}^{2}}{4\pi} m_{\rho} H_{0}(\Lambda, m_{\rho}, m_{\rho}, r) - \frac{g_{\rho NN}^{2} + 4g_{\rho NN} f_{\rho NN}}{4\pi} \frac{m_{\rho}^{3}}{8m_{N}^{2}} H_{1}(\Lambda, m_{\rho}, m_{\rho}, r) \\ V_{I}^{d}(r) &= \frac{g_{\eta NN}^{2}}{4\pi} \frac{m_{\eta}^{3}}{12m_{N}^{2}} H_{3}(\Lambda, m_{\eta}, m_{\eta}, r) - \frac{(g_{\omega NN} + f_{\omega NN})^{2}}{4\pi} \frac{m_{\omega}^{3}}{12m_{N}^{2}} H_{3}(\Lambda, m_{\omega}, m_{\omega}, r) \\ V_{SI}^{d}(r) &= -\frac{g_{\pi NN}^{2}}{4\pi} \frac{m_{\pi}^{3}}{12m_{N}^{2}} H_{3}(\Lambda, m_{\pi}, m_{\pi}, r) - \frac{(g_{\rho NN} + f_{\rho NN})^{2}}{4\pi} \frac{m_{\rho}^{3}}{6m_{N}^{2}} H_{1}(\Lambda, m_{\rho}, m_{\rho}, r) \\ V_{II}^{d}(r) &= \frac{g_{\pi NN}^{2}}{4\pi} \frac{m_{\pi}^{3}}{12m_{N}^{2}} H_{3}(\Lambda, m_{\pi}, m_{\pi}, r) - \frac{(g_{\rho NN} + f_{\rho NN})^{2}}{4\pi} \frac{m_{\rho}^{3}}{12m_{N}^{2}} H_{3}(\Lambda, m_{\rho}, m_{\rho}, r) \\ V_{LS}^{d}(r) &= -\frac{g_{\pi NN}^{2}}{4\pi} \frac{m_{\pi}^{3}}{12m_{N}^{2}} H_{2}(\Lambda, m_{\sigma}, m_{\sigma}, r) - \frac{3g_{\omega NN}^{2} + 4g_{\omega NN} f_{\omega NN}}{4\pi} \frac{m_{\omega}^{3}}{2m_{N}^{2}} H_{2}(\Lambda, m_{\omega}, m_{\omega}, r) \\ V_{LSI}^{d}(r) &= -\frac{3g_{\rho NN}^{2}}{4\pi} \frac{m_{\eta}^{3}}{2m_{N}^{2}} H_{2}(\Lambda, m_{\sigma}, m_{\sigma}, r) - \frac{3g_{\omega NN}^{2} + 4g_{\omega NN} f_{\omega NN}}{4\pi} \frac{m_{\omega}^{3}}{2m_{N}^{2}} H_{2}(\Lambda, m_{\omega}, m_{\omega}, r) \\ V_{LSI}^{d}(r) &= -\frac{3g_{\rho NN}^{2}}{4\pi} \frac{4g_{\rho NN} f_{\rho NN}}{2m_{N}^{2}} H_{2}(\Lambda, m_{\rho}, m_{\rho}, r). \end{split}$$

In the basis of ${}^{3}S_{1}$ and ${}^{3}D_{1}$ states, the deuteron potential can be written in the matrix form as

$$V^{d} = \left[V_{C}^{d}(r) + V_{S}^{d}(r) - 3V_{I}^{d}(r) - 3V_{SI}^{d}(r)\right] \begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix} + \left[9V_{LSI}^{d}(r) - 3V_{LS}^{d}(r)\right] \begin{pmatrix} 0 & 0 \\ 0 & 1 \end{pmatrix} + \left[V_{T}^{d}(r) - 3V_{TI}^{d}(r)\right] \begin{pmatrix} 0 & \sqrt{8} \\ \sqrt{8} & -2 \end{pmatrix}.$$
(15)

Taking into account the centrifugal barrier from D wave and solving the corresponding two channel Schrödinger equation numerically via the Fortran77 package FESSDE2.2 [19], which can fastly and accurately solve the eigenvalue problem for systems of coupled Schrödinger equations, we find the binding energy $\varepsilon_d \simeq$ 2.25 MeV for the cutoff parameter $\Lambda = 808$ MeV, and the corresponding wave function is presented in Fig. 1. If we reduce half of the effective coupling constants except $g_{\pi NN}$, the binding energy is found to be about 2.28 MeV with $\Lambda = 970$ MeV. From the wave function one can calculate the static properties of deuteron such as the root of mean square radius, the D wave probability, the magnetic moment and the quadrupole moment, which are in agreement with experimental data. We would like to note that the small binding energy of deuteron is a cancellation result of different contributions of opposite signs. The detailed results are listed in Table II, it is obvious the results are sensitive to the regularization parameter Λ , and the same conclusion has been drawn in the one pion exchange model [5,7]. The binding energy variation with respect to Λ is shown in Fig. 2, the dependence is less sensitive than the one pion exchange model. It is obvious that the binding energy variation with Λ is dependent on the coupling constants. For the coupling constants listed in Table I, the binding energy no longer monotonically increases with Λ in contrast with the one pion exchange model. To understand this peculiar behavior, we plot the three components of the deuteron effective potential in Eq. (15) in Fig. 3. We can see that both $V_{11}(\Lambda, r)$ and $V_{22}(\Lambda, r)$ potentials are repulsive, and they increase with



FIG. 1 (color online). The deuteron ${}^{3}S_{1}$ and ${}^{3}D_{1}$ wave function with binding energy $\varepsilon_{d} \simeq 2.25$ MeV and $\Lambda \simeq 808$ MeV.

Λ at short distance. However, at intermediate distance the relation $|V_{12}(\Lambda = 1.2 \text{ GeV}, r)| < |V_{12}(\Lambda = 0.8 \text{ GeV}, r)| < |V_{12}(\Lambda = 0.9 \text{ GeV}, r)| < |V_{12}(\Lambda = 1.6 \text{ GeV}, r)|$ is satisfied, the $V_{12}(\Lambda, r)$ does not monotonically increases with Λ . Therefore the nonmonotonous behavior in Fig. 2(a) mainly comes from the nonmonotonous dependence of

 $V_{12}(\Lambda, r)$ potential on Λ , which is a cancellation result of various contributions. As has been discussed above, the heavy flavor system should admit a larger Λ than the deuteron. Therefore the above values of Λ with which the smaller deuteron binding energy is reproduced, would be assumed to be the lower bound in the following.

TABLE II. The deuteron static properties in the one-boson exchange potential model, where ε_d is the binding energy, $r_{\rm rms}$ is the root of mean square radius, P_S and P_D represent the S-state and D-state probabilities, respectively, μ_d is the magnetic moment, and Q_d denotes the quadrupole moment.

	_				
Λ (MeV)	ε_d (MeV)	r _{rms} (fm)	$P_D:P_S(\%)$	$\mu_d(\mu_N)$	$Q_d \ (\mathrm{fm}^2)$
808	2.25	3.85	5.66:94.34	0.85	0.27
900	5.33	2.77	7.44:92.56	0.84	0.20
1000	4.96	2.87	7.37:92.63	0.84	0.21
		All couplings are reduc	ced by half except $g_{\pi NN}$		
Λ (MeV)	ε_d (MeV)	$r_{\rm rms}$ (fm)	$P_D:P_S$	$\mu_d(\mu_N)$	$Q_d \ (\mathrm{fm}^2)$
970	2.28	3.84	6.52:93.48	0.84	0.28
1100	5.65	2.70	8.92:91.08	0.83	0.20
1200	8.89	2.28	10.26:89.74	0.82	0.16



FIG. 2 (color online). The deuteron binding energy variation with respect to the regularization parameter Λ . Figure 2(a) corresponds to the coupling constants shown in Table I, and Fig. 2(b) for the couplings reduced by half.



FIG. 3 (color online). The three components of the deuteron effective potential in Eq. (13), Fig. 3(a)–3(c) respectively illustrate the $V_{11}(\Lambda, r)$, $V_{12}(\Lambda, r)$ and $V_{22}(\Lambda, r)$ components. The solid line, dashed, dotted and dash-dotted lines correspond to $\Lambda = 0.8$ GeV, 0.9 GeV, 1.2 GeV and 1.6 GeV, respectively.

V. POSSIBLE $D\bar{D}^*/\bar{D}D^*$ HADRONIC MOLECULE AND X(3872)

The narrow charmoniumlike state X(3872) was discovered by the Belle collaboration in the decay $B^+ \rightarrow K^+ +$ X(3872) followed by $X(3872) \rightarrow J/\psi \pi^+ \pi^-$ with a statistical significance of 10.3σ [20]. The existence of X(3872) has been confirmed by CDF [21], D0 [22], and Babar collaboration [23]. the CDF collaboration measured the X(3872) mass to be $(3871.61 \pm 0.16(\text{stat}) \pm$ 0.19(sys)) MeV. Its quantum number is strongly preferred to be 1^{++} [24]. In the one pion exchange model, Tornqvist suggested that X(3872) is a 1⁺⁺ $D\bar{D}^*/\bar{D}D^*$ molecule and isospin is strongly broken [25]. Recently Close et al. reanalyzed X(3872) in the same model, the critical overall sign is corrected and the contribution of the " δ function" term is included [7]. Swanson have taken into account both the long rang pion exchange and short range contribution arising from constituent quark interchange [8]. Recently Zhu *et al.* dynamically studied the binding of X(3872) in the heavy quark effective theory [14]. In this section, we will investigate the 1^{++} $D\bar{D}^*/\bar{D}D^*$ system from the oneboson exchange model at quark level, where the short range interactions are represented by the heavier bosons η, σ, ρ and ω exchange instead of the quark interchange. There is only a sign difference $(-1)^G$ between the quark-quark interaction and quark-antiquark interaction, and the magnitudes are the same, where *G* is the *G*-parity of the exchanged meson. The diagrams contributing to the $D\bar{D}^*$ and $\bar{D}D^*$ interactions are displayed in Fig. 4. Because of the parity conservation, $D\bar{D}^*$ can only scatter into $D^*\bar{D}$ via the pseudoscalar π and η exchange, and $D\bar{D}^*$ scatters into $D\bar{D}^*$ with the scalar σ exchange, whereas the vector mesons ρ and ω exchange contribute to both processes. The effective potential for the 1^{++} $D\bar{D}^*/\bar{D}D^*$ system is

$$V^{X}(\mathbf{r}) = -V^{X}_{\pi}(\mathbf{r}) + V^{X}_{\eta}(\mathbf{r}) + V^{X}_{\sigma}(\mathbf{r}) + V^{X}_{\rho}(\mathbf{r}) - V^{X}_{\omega}(\mathbf{r})$$

$$\equiv V^{X}_{C}(r) + V^{X}_{S}(r)(\boldsymbol{\sigma}_{i} \cdot \boldsymbol{\sigma}_{j}) + V^{X}_{I}(r)(\boldsymbol{\tau}_{i} \cdot \boldsymbol{\tau}_{j})$$

$$+ V^{X}_{T}(r)S_{ij}(\hat{\mathbf{r}}) + V^{X}_{SI}(\boldsymbol{\mu}_{k}, r)(\boldsymbol{\sigma}_{i} \cdot \boldsymbol{\sigma}_{j})(\boldsymbol{\tau}_{i} \cdot \boldsymbol{\tau}_{j})$$

$$+ V^{X}_{TI}(\boldsymbol{\mu}_{k}, r)S_{ij}(\hat{\mathbf{r}})(\boldsymbol{\tau}_{i} \cdot \boldsymbol{\tau}_{j}) + V^{X}_{LS}(r)(\mathbf{L} \cdot \mathbf{S}_{ij})$$

$$+ V^{X}_{LSI}(r)(\mathbf{L} \cdot \mathbf{S}_{ij})(\boldsymbol{\tau}_{i} \cdot \boldsymbol{\tau}_{j})$$
(16)

with *i* and *j* is the index of light quark or antiquark, $\mu_k(k = 1, 2, 3, 4)$ takes four different values due to the mass difference within the *D*, *D*^{*} and π isospin multiplets. The eight components $V_C^X(r)$, $V_S^X(r)$ etc are given by

$$\begin{split} V_{C}^{\chi}(r) &= -\frac{g_{qqq}^{2}}{4\pi} m_{q} \bigg[H_{0}(\Lambda, m_{\sigma}, m_{\sigma}, r) + \frac{m_{\sigma}^{2}}{8m_{q}^{2}} H_{1}(\Lambda, m_{\sigma}, m_{\sigma}, r) \bigg] - \frac{g_{\omega qq}^{2}}{4\pi} m_{\omega} H_{0}(\Lambda, m_{\omega}, m_{\omega}, r) \\ &+ \frac{g_{\omega qq}^{2} + 4g_{\omega qq} f_{\omega qq}}{4\pi} \frac{m_{\omega}^{3}}{8m_{q}^{2}} H_{1}(\Lambda, m_{\omega}, m_{\omega}, r) \\ V_{S}^{\chi}(r) &= -\frac{g_{\eta qq}^{2}}{4\pi} \frac{\mu_{S}^{3}}{12m_{q}^{2}} H_{1}(\Lambda, m_{\eta}, \mu_{S}, r) + \frac{(g_{\omega qq} + f_{\omega qq})^{2}}{4\pi} \frac{\mu_{\eta}^{3}}{6m_{q}^{2}} H_{1}(\Lambda, m_{\omega}, \mu_{\sigma}, r) \\ V_{I}^{\chi}(r) &= \frac{g_{\eta qq}^{2}}{4\pi} \frac{\mu_{S}^{3}}{12m_{q}^{2}} H_{1}(\Lambda, m_{\eta}, \mu_{S}, r) + \frac{(g_{\omega qq} + f_{\omega qq})^{2}}{4\pi} \frac{m_{\eta}^{3}}{8m_{q}^{2}} H_{1}(\Lambda, m_{\omega}, \mu_{\sigma}, r) \\ V_{T}^{\chi}(r) &= \frac{g_{\eta qq}^{2}}{4\pi} \frac{\mu_{S}^{3}}{12m_{q}^{2}} H_{3}(\Lambda, m_{\eta}, \mu_{S}, r) + \frac{(g_{\omega qq} + f_{\omega qq})^{2}}{4\pi} \frac{\mu_{\eta}^{3}}{12m_{q}^{2}} H_{3}(\Lambda, m_{\omega}, \mu_{\tau}, r) \\ V_{SI}^{\chi}(\mu, r) &= \begin{cases} \frac{g_{\eta qq}^{2}}{4\pi} \frac{\mu_{S}^{3}}{12m_{q}^{2}} H_{1}(\Lambda, m_{\pi^{\pm 0}}, \mu, r) - \frac{(g_{\rho qq} + f_{\rho qq})^{2}}{4\pi} \frac{\mu_{\eta}^{3}}{6m_{\eta}^{2}} H_{1}(\Lambda, m_{\rho}, \mu_{6}, r), \quad \mu^{2} > 0 \\ g_{\eta q}^{2} \frac{\mu_{\eta}^{3}}{4\pi} \frac{\mu_{\eta}^{3}}{12m_{\eta}^{2}} G_{3}(\Lambda, m_{\pi^{\pm 0}}, \mu, r) - \frac{(g_{\rho qq} + f_{\rho qq})^{2}}{4\pi} \frac{\mu_{\eta}^{3}}{12m_{\eta}^{2}} H_{3}(\Lambda, m_{\rho}, \mu_{6}, r), \quad \mu^{2} = -\tilde{\mu}^{2} < 0 \end{cases} \\ V_{LS}^{\chi}(r) &= -\frac{g_{\eta qq}^{2}}{4\pi} \frac{m_{\eta}^{3}}{2m_{q}^{2}} H_{2}(\Lambda, m_{\sigma}, m_{\sigma}, r) + \frac{3g_{\omega qq}^{2} + 4g_{\omega qq} f_{\omega qq}}{4\pi} \frac{m_{\eta}^{3}}{2m_{q}^{2}} H_{3}(\Lambda, m_{\rho}, \mu_{6}, r), \quad \mu^{2} = -\tilde{\mu}^{2} < 0 \end{cases} \\ V_{LS}^{\chi}(r) &= -\frac{g_{\eta qq}^{2}}{4\pi} \frac{m_{\eta}^{3}}{2m_{q}^{2}} H_{2}(\Lambda, m_{\sigma}, m_{\sigma}, r) + \frac{3g_{\omega qq}^{2} + 4g_{\omega qq} f_{\omega qq}}{4\pi} \frac{m_{\eta}^{3}}{2m_{q}^{2}}} H_{2}(\Lambda, m_{\omega}, m_{\omega}, r) \\ V_{LSI}^{\chi}(r) &= -\frac{3g_{\rho qq}^{2} + 4g_{\rho qq} f_{\rho qq}}{4\pi} \frac{m_{\eta}^{3}}{2m_{q}^{2}}} H_{2}(\Lambda, m_{\rho}, m_{\rho}, r), \end{cases}$$

FIG. 4. $D\bar{D}^*$ and $\bar{D}D^*$ interaction in one-boson exchange model at quark level, where the thick line represents heavy quark or antiquark, and the thin line denotes light quark or antiquark.

where

$$\mu_{1}^{2} = m_{\pi^{0}}^{2} - (m_{D^{*0}} - m_{D^{0}})^{2},$$

$$\mu_{2}^{2} = m_{\pi^{\pm}}^{2} - (m_{D^{*0}} - m_{D^{\pm}})^{2}$$

$$\mu_{3}^{2} = m_{\pi^{\pm}}^{2} - (m_{D^{*\pm}} - m_{D^{0}})^{2},$$

$$\mu_{4}^{2} = m_{\pi^{0}}^{2} - (m_{D^{*\pm}} - m_{D^{\pm}})^{2}$$

$$\mu_{5}^{2} = m_{\eta}^{2} - (m_{D^{*0}} - m_{D^{0}})^{2},$$

$$\mu_{6}^{2} = m_{\rho}^{2} - (m_{D^{*0}} - m_{D^{0}})^{2},$$

$$\mu_{5}^{2} = m_{\omega}^{2} - (m_{D^{*0}} - m_{D^{0}})^{2}.$$
(18)

These μ^2 parameters approximately represent the recoil effect due to different values of m_D and m_{D^*} as in

Refs. [5,7]. For the η , σ , ρ , and ω exchange processes, the mass difference of m_{D^0} and $m_{D^{\pm}}$ as well as $m_{D^{*0}}$ and $m_{D^{*\pm}}$ are neglected, since they are much smaller comparing with m_{η} , m_{ρ} , and m_{ω} . X(3872) is very close to the $D^0 \bar{D}^{*0}$ threshold, however, it is about 8.3 MeV below the $D^+ D^{*-}$ threshold. Hence, isospin symmetry is drastically broken [25,26]. For the $J^{PC} = 1^{++} D\bar{D}^* / \bar{D}D^*$ system, they can be in *S* wave or *D* wave similar to the deuteron, then the wave function of this system is written as

$$\begin{aligned} |X(3872)\rangle &= \frac{u_1(r)}{r} \frac{1}{\sqrt{2}} |(D^0 \bar{D}^{*0} + \bar{D}^0 D^{*0})_S\rangle \\ &+ \frac{u_2(r)}{r} \frac{1}{\sqrt{2}} |(D^0 \bar{D}^{*0} + \bar{D}^0 D^{*0})_D\rangle \\ &+ \frac{u_3(r)}{r} \frac{1}{\sqrt{2}} |(D^+ D^{*-} + D^- D^{*+})_S\rangle \\ &+ \frac{u_4(r)}{r} \frac{1}{\sqrt{2}} |(D^+ D^{*-} + D^- D^{*+})_D\rangle, \end{aligned}$$
(19)

where the subscript *S* and *D* denote the system in *S* wave and *D* wave, respectively. $u_1(r)$, $u_2(r)$, $u_3(r)$ and $u_4(r)$ are the spatial wave functions. There are four channels coupled with each other as has been shown above, and we might as well choose the basis to be $|1\rangle \equiv \frac{1}{\sqrt{2}} |(D^0 \bar{D}^{*0} + \bar{D}^0 D^{*0})_S\rangle$, $|2\rangle \equiv \frac{1}{\sqrt{2}} |(D^0 \bar{D}^{*0} + \bar{D}^0 D^{*0})_D\rangle$, $|3\rangle \equiv \frac{1}{\sqrt{2}} |(D^+ D^{*-} + D^- D^{*+})_S\rangle$, and $|4\rangle \equiv \frac{1}{\sqrt{2}} |(D^+ D^{*-} + D^- D^{*+})_D\rangle$. Using the analytical formula for the matrix elements presented in the appendix, the effective potential for $1^{++} D \bar{D}^* / \bar{D} D^*$ can be written in the matrix form as

$$V^{X}(r) = \begin{bmatrix} V_{C}^{X}(r) + V_{S}^{X}(r) \end{bmatrix} \begin{pmatrix} 1 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 1 \end{pmatrix} + \begin{bmatrix} V_{I}^{X}(r) + V_{SI}^{X}(\mu_{k}, r) \end{bmatrix} \begin{pmatrix} -1 & 0 & -2 & 0 \\ 0 & -1 & 0 & -2 \\ -2 & 0 & -1 & 0 \\ 0 & -2 & 0 & -1 \end{pmatrix} \\ + V_{T}^{X}(r) \begin{pmatrix} 0 & -\sqrt{2} & 0 & 0 \\ -\sqrt{2} & 1 & 0 & 0 \\ 0 & 0 & 0 & -\sqrt{2} \\ 0 & 0 & -\sqrt{2} & 1 \end{pmatrix} + V_{TI}^{X}(\mu_{k}, r) \begin{pmatrix} 0 & \sqrt{2} & 0 & 2\sqrt{2} \\ \sqrt{2} & -1 & 2\sqrt{2} & -2 \\ 0 & 2\sqrt{2} & 0 & \sqrt{2} \\ 2\sqrt{2} & -2 & \sqrt{2} & -1 \end{pmatrix} \\ + V_{LS}^{X}(r) \begin{pmatrix} 0 & 0 & 0 & 0 \\ 0 & -3/2 & 0 & 0 \\ 0 & 0 & 0 & -3/2 \end{pmatrix} + V_{LSI}^{X}(r) \begin{pmatrix} 0 & 0 & 0 & 0 \\ 0 & 3/2 & 0 & 3 \\ 0 & 0 & 0 & 0 \\ 0 & 3 & 0 & 3/2 \end{pmatrix}.$$

$$(20)$$

In the above equation, the value of μ_k^2 is μ_1^2 for the upperleft 2 × 2 matrix elements, and it is equal to μ_4^2 for the lower-right 2 × 2 matrix elements. There is ambiguity in choosing μ_k^2 value for the processes $D^0 \bar{D}^{*0} \rightarrow D^{*+} D^-$ or $D^+ D^{*-} \rightarrow D^{*0} \bar{D}^0$, accordingly μ_k^2 can take the value μ_2^2 or μ_3^2 for the off-diagonal 2×2 matrix elements, the numerical results for both choices would be given in the following. The different μ_k values is due to the isospin symmetry breaking from the mass difference within the *D*, D^* and π isospin multiplets. Taking into account the

1		1	/	
μ^2	Λ (MeV)	M (MeV)	r _{rms} (fm)	$P_{S}^{00}:P_{D}^{00}:P_{S}^{+-}:P_{D}^{+-}(\%)$
$\overline{\mu_3^2}$	808	3871.6	7.02	90.76:0.56:8.11:0.56
• 5	840	3870.4	2.84	78.23:1.08:19.59:1.11
	850	3869.8	2.45	75.26:1.21:22.29:1.23
	900	3865.9	1.61	65.17:1.89:31.06:1.88
	1000	3849.2	1.08	53.06:4.72:37.65:4.57
μ_2^2	808	3871.7	11.34	94.40:0.38:4.86:0.36
. 2	840	3870.7	3.19	80.74:0.99:17.26:1.01
	850	3870.2	2.68	77.44:1.14:20.28:1.15
	900	3866.4	1.66	66.23:1.85:30.09:1.83
	1000	3849.9	1.09	53.35:4.69:37.43:4.53
	All	couplings except g_{π}	_{NN} are reduced by	half
μ^2	Λ (MeV)	M (MeV)	r _{rms} (fm)	$P_S^{00}: P_D^{00}: P_S^{+-}: P_D^{+-}(\%)$
μ_3^2	970	3869.1	2.13	70.65:1.65:26.02:1.69
• 5	1100	3860.1	1.25	57.24:2.98:36.83:2.95
	1200	3848.2	1.00	51.80:4.40:39.46:4.33
μ_2^2	970	3869.5	2.28	72.56:1.57:24.28:1.60
. 2	1100	3860.8	1.27	57.76:2.94:36.38:2.92
	1200	3849.0	1.01	52.04:4.38:39.29:4.30

TABLE III. The predictions about the mass, the root of mean square radius (rms) and the probabilities of the different components for the $1^{++} D\bar{D}^*/\bar{D}D^*$ molecule.

centrifugal barrier from D wave and solving the four channel coupled Schrödinger equation using the package FESSDE2.2, the numerical results are listed in Table III. It is remarkable that the $1^{++} D\bar{D}^*/\bar{D}D^*$ system could accommodate a molecular state with mass about 3871.6 MeV for $\Lambda = 808$ MeV, it is very close to the central value of X(3872) mass 3871.61 MeV. The corresponding wave function is shown in Fig. 5, it is obvious that the $D^0\bar{D}^{*0} + \bar{D}^0D^{*0}$ component dominates over the $D^+D^{*-} + D^-D^{*+}$ component. Since the spatial wave functions $u_1(r)$ and $u_3(r)$ have the same sign, the same is true for $u_2(r)$ and $u_4(r)$, thus the I = 0 component in this state is predomi-



FIG. 5 (color online). The four components spatial wave functions of the $1^{++} D\bar{D}^*/\bar{D}D^*$ system with $\Lambda = 808$ MeV.

nant, it would be a isospin singlet in the isospin symmetry limit. From the results in Table III, we notice that the predictions about the static properties for the two μ^2 choices are very similar to each other, and the difference is small. The isospin symmetry is strongly broken especially for the states near the threshold. The uncertainties induced by the effective coupling constants are considered, we reduce half of the couplings except $g_{\pi NN}$, and the numerical results are given in Table III as well. For both choices of the coupling constants, the binding energy and other static properties are sensitive to the regularization parameter Λ , and the bound state mass dependence on Λ is displayed in Fig. 6. It is obvious that the bound state mass decreases monotonically with the regularization parameter Λ as in the one pion exchange model. In short summary, the predictions are qualitatively the same as those in the one pion exchange model, even after we have included the contributions from η , σ , ρ , and ω exchange. Since unexpectedly large branch ratio of $X(3872) \rightarrow \psi(2S)\gamma$ recently was reported [27], we have to take into account the mixing between the 1^{++} $D\bar{D}^*/\bar{D}D^*$ molecule and the conventional charmonium state in order to identify this state with X(3872). This is outside the range of the present work.

VI. POSSIBLE MOLECULAR STATES OF OTHER HEAVY FLAVOR PV SYSTEMS

A. $B\bar{B}^*/\bar{B}B^*$ system

For the $1^{++} B\bar{B}^*/\bar{B}B^*$ system, the kinetic energy is greatly reduced due to the heavier mass of *B* meson, and



FIG. 6 (color online). The variation of the $1^{++} D\bar{D}^*/\bar{D}D^*$ bound state mass with respect to Λ . (a) corresponds to the coupling constants shown in Table I, and (b) for the couplings reduced by half.

the interaction potential has features similar to those of the $D\bar{D}^*/\bar{D}D^*$ system except that the former is deeper than the latter. Therefore molecular states should be more easily formed. Following the same procedure as the $D\bar{D}^*/\bar{D}D^*$ case, the numerical results are shown in Table IV, where the μ^2 ambiguity is considered. For the same value of Λ , the $B\bar{B}^*/\bar{B}B^*$ system is really more strongly bound than the $D\bar{D}^*/\bar{D}D^*$ system, its binding energy is a few tens of MeV, and the same was predicted in the one pion exchange

model [5,7] and in the models [14,28]. It is obvious that the predictions about the static properties for the two μ^2 choices are approximately the same. We notice that the isospin symmetry breaking is less stronger than the charm system, this is because the mass difference of B^0 and B^+ is smaller than that of D^0 and D^+ as well as D^{*0} and D^{*+} . It is notable that there may be two molecular states for appropriate values of Λ . The corresponding wave functions for $\Lambda = 1000$ MeV and $\mu^2 = \mu_{bb1}^2 \equiv m_{\pi^{\pm}}^2 - (m_{B^*} - m_{B^0})^2$

TABLE IV. The predictions about the mass, the root of mean square radius (rms) and the probabilities of the different components for the 1⁺⁺ $B\bar{B}^*/\bar{B}B^*$ system with $\mu_{b\bar{b}1}^2 = m_{\pi^{\pm}}^2 - (m_{B^*} - m_{B^0})^2$ and $\mu_{b\bar{b}2}^2 = m_{\pi^{\pm}}^2 - (m_{B^*} - m_{B^+})^2$.

	$1^{++} B\bar{B}^*/\bar{B}B^*$					
μ^2	Λ (MeV)	M (MeV)	r _{rms} (fm)	$P_S^{00}:P_D^{00}:P_S^{+-}:P_D^{+-}(\%)$		
$\mu^2_{h\bar{h}1}$	808	10565.3	0.60	47.70:2.05:48.20:2.05		
001	900	10543.5	0.59	44.11:5.69:44.50:5.70		
	1000	10457.6	0.52	27.52:22.41:27.64:22.44		
		10600.4	1.73	37.10:9.43:44.26:9.22		
$\mu^2_{h\bar{b}2}$	808	10565.3	0.60	47.70:2.05:48.20:2.05		
002	900	10543.5	0.59	44.11:5.69:44.49:5.70		
	1000	10457.6	0.52	27.52:22.41:27.64:22.44		
		10600.4	1.72	37.10:9.43:44.25:9.22		
	All	couplings except g	τ_{NN} are reduced by	y half		
μ^2	Λ (MeV)	M (MeV)	$r_{\rm rms}$ (fm)	$P_S^{00}:P_D^{00}:P_S^{+-}:P_D^{+-}(\%)$		
$\mu^2_{h\bar{h}1}$	970	10544.8	0.55	45.24:4.60:45.55:4.61		
001	1100	10503.9	0.51	40.12:9.78:40.30:9.80		
	1200	10443.1	0.46	33.66:16.27:33.77:16.29		
		10601.9	1.91	38.82:8.03:45.26:7.89		
$\mu_{h\bar{h}2}^2$	970	10544.8	0.55	45.24:4.60:45.55:4.61		
002	1100	10503.9	0.51	40.12:9.78:40.30:9.80		
	1200	10443.1	0.46	33.66:16.27:33.77:16.29		
		10601.9	1.91	38.83:8.04:45.25:7.89		



FIG. 7 (color online). The spatial wave functions of $1^{++} B\bar{B}^*/\bar{B}B^*$ molecule with $\Lambda = 1000$ MeV. There are two bound states with mass 10457.6 MeV and 10600.4 MeV, respectively, Fig. 7(a) is for the first state, and the Fig. 7(b) for the second.

are displayed in Fig. 7, the first state is tightly bound, whereas the second is loosely bound. We notice that the first state is almost an isospin singlet, and the I = 0 component is dominant for the second state. This state can no longer be produced through *B* meson decay because of its large mass, and we have to resort to hadron collider. We can search for this state at Tevatron via $p\bar{p} \rightarrow \pi^+\pi^-\Upsilon(1S)$, and LHC is more promising.

B. DD^* system with C = 2

The interaction potentials arise from the one-boson exchange between two antiquarks instead of a quark and antiquark pair, hence both the π exchange and ω exchange potentials have overall opposite sign relative to the $D\bar{D}^*/\bar{D}D^*$ case. In this case we have four coupled channels $(D^+D^{*0})_S$, $(D^+D^{*0})_D$, $(D^0D^{*+})_S$ and $(D^0D^{*+})_D$. The numerical results are given in Table V. For $\Lambda = 808$ MeV

TABLE V. The predictions about the mass, rms, and the probabilities of the different components for the DD^* system, and P_S^{+0} denotes the probability of *S* wave D^+D^{*0} in the state, and the meaning of P_D^{+0} , P_S^{0+} and P_D^{0+} is similar. Here $\mu_{cc1}^2 = m_{\pi^0}^2 - (m_{D^{*+}} - m_{D^+})^2$ and $\mu_{cc2}^2 = m_{\pi^0}^2 - (m_{D^{*0}} - m_{D^0})^2$.

	DD^* system with $C = 2$					
μ^2	Λ (MeV)	M (MeV)	r _{rms} (fm)	$P_S^{+0}: P_D^{+0}: P_S^{0+}: P_D^{0+}(\%)$		
μ^2_{cc1}	1600	3873.9	3.15	34.61:4.03:56.82:4.54		
	1700	3865.1	1.30	20.29:28.37:22.63:28.71		
	1800	3770.9	0.58	0.19:49.78:0.19:49.84		
		3872.8	2.48	41.47:1.38:55.54:1.63		
μ_{cc2}^2	1600	3873.9	3.16	34.54:4.04:56.87:4.55		
	1700	3865.1	1.30	20.31:28.35:22.65:28.70		
	1800	3771.0	0.58	0.19:49.78:0.19:49.84		
		3872.8	2.49	41.41:1.38:55.58:1.63		
	All	couplings are reduc	ed by half except	$g_{\pi NN}$		
μ^2	Λ (MeV)	M (MeV)	$r_{\rm rms}$ (fm)	$P_{S}^{+0}:P_{D}^{+0}:P_{S}^{0+}:P_{D}^{0+}(\%)$		
μ^2_{cc1}	1900	3873.1	2.53	36.05:5.90:51.55:6.50		
	2000	3870.1	1.82	38.09:8.08:45.32:8.52		
	2100	3865.3	1.44	37.67:10.21:41.57:10.55		
	2200	3858.4	1.19	36.31:12.42:38.60:12.68		
μ^2_{cc2}	1900	3873.1	2.54	36.01:5.91:51.58:6.51		
	2000	3870.1	1.82	38.07:8.08:45.32:8.53		
	2100	3865.4	1.44	37.66:10.22:41.57:10.55		
	2200	3858.4	1.19	36.30:12.42:38.59:12.68		



FIG. 8 (color online). The wave function for the DD^* system assuming $\Lambda = 1600$ MeV and $\mu^2 = \mu_{cc1}^2 \equiv m_{\pi^0}^2 - (m_{D^{*+}} - m_{D^+})^2$, its mass approximately is 3873.9 MeV.

or $\Lambda = 970$ MeV, we find no bound state. A bound state with mass about 3873.9 MeV appears for $\Lambda = 1600$ MeV (about 3873.1 MeV for $\Lambda = 1900$ MeV if the couplings except $g_{\pi NN}$ are reduced half), and the corresponding wave function is shown in Fig. 8. We notice that the wave functions of ${}^{3}S_{1} D^{+}D^{*0}$ and $D^{0}D^{*+}$ have opposite signs, the same is true for the ${}^{3}D_{1} D^{+}D^{*0}$ and $D^{0}D^{*+}$ wave functions, therefore this state would be a isospin singlet in the isospin symmetry limit. We notice that the ${}^{3}D_{1}$ probability are much larger than ${}^{3}S_{1}$ probability for the state with $\Lambda = 1800$ MeV, although there is centrifugal barrier for the *D* wave state. Thus the *S*-*D* mixing effect induced by the tensor force is especially crucial for this state. In short, the bound state of the *DD*^{*} system appears only for the regularization parameter Λ as large as 1600 MeV or 1900 MeV, which is beyond the range of 0.8 to 1.5 GeV favored by the nucleon-nucleon interaction. Moreover, the parameters that allow *X*(3872) to emerge as a $D\bar{D}^*/\bar{D}D^*$ molecule exclude the *DD*^{*} bound state, as can be seen from the results in Sec. V. Consequently we tend to conclude that the *DD*^{*} molecular state may not exist.

C. BB^* system with B = 2

The situation is very similar to the DD^* system except the different mass of D mesons and B mesons, we list the numerical results in Table VI. We find a marginally bound state with mass 10603.9 MeV for $\Lambda = 808$ MeV, which is very close to the BB^* threshold. Its binding energy is much smaller than that of the $1^{++} B\bar{B}^*/\bar{B}B^*$ system, however, the binding energy is less sensitive to Λ than the latter case. Figure 9 displays the wave function of the bound state solution with mass 10602.3 MeV and $\Lambda = 900$ MeV. It is obvious that the B^+B^{*0} and B^0B^{*+} wave functions have the opposite sign, then the I = 0 component is dominant in this state. If the couplings except $g_{\pi NN}$ are reduced by half, a weakly bound state with mass about 10601.5 MeV is found as well assuming $\Lambda = 970$ MeV. These indicates a weakly bound BB^* should exist, This is consistent with the results of Manohar and Wise form heavy quark effective

TABLE VI. The predictions about the mass, rms, and the probabilities of the different components for the *BB*^{*} system with B = 2. P_S^{+0} represents the probability of *S*-wave B^+B^{*0} . Here $\mu_{bb1}^2 = m_{\pi^0}^2 - (m_{B^*} - m_{B^+})^2$ and $\mu_{bb2}^2 = m_{\pi^0}^2 - (m_{B^*} - m_{B^0})^2$.

 BB^* system with B = 2

μ^2	Λ (MeV)	M (MeV)	r _{rms} (fm)	P_S^{+0} : P_D^{+0} : P_S^{0+} : $P_D^{0+}(\%)$
μ^2_{bb1}	808	10603.9	4.09	59.37:6.74:28.49:5.40
	900	10602.3	2.23	43.52:10.84:35.63:10.02
	1000	10598.8	1.61	37.22:14.37:34.48:13.93
	1100	10592.2	1.27	31.33:19.34:30.24:19.09
μ_{bb2}^2	808	10603.9	4.09	59.38:6.74:28.48:5.40
002	900	10602.3	2.23	43.52:10.84:35.63:10.01
	1000	10598.8	1.61	37.22:14.37:34.48:13.93
	1100	10592.2	1.27	31.33:19.34:30.24:19.09
	All	couplings except g	$_{\tau NN}$ are reduced by	y half
μ^2	Λ (MeV)	M (MeV)	r _{rms} (fm)	$P_S^{+0}: P_D^{+0}: P_S^{0+}: P_D^{0+}(\%)$
μ^2_{bb1}	970	10601.5	1.99	39.93:13.08:34.67:12.31
	1000	10600.7	1.82	38.34:14.02:34.30:13.34
	1100	10596.6	1.44	34.37:16.80:32.47:16.36
	1200	10590.3	1.19	31.33:19.33:30.32:19.03
μ_{bh2}^2	970	10601.5	1.99	39.94:13.08:34.67:12.31
002	1000	10600.7	1.82	38.34:14.02:34.30:13.34
	1100	10596.6	1.44	34.37:16.80:32.47:16.36
	1200	10590.3	1.19	31.33:19.33:30.32:19.03



FIG. 9 (color online). The spatial wave function for the *BB*^{*} system assuming $\Lambda = 900$ MeV and $\mu^2 = \mu_{bb1}^2 \equiv m_{\pi^0}^2 - (m_{B^*} - m_{B^+})^2$, its mass is 10602.3 MeV.

theory [29]. For a loosely bound molecule, the leading source of decay is dissociation, to a good approximation the dissociation will proceed via the free space decay of the constituent mesons. The spin-parity forbids its decay into *BB*, therefore the *BB*^{*} molecule is a very narrow state and it mainly decays into *BB* γ .

D. Pseudoscalar-vector system with C = B = 1

This system could have the same quantum as B_c meson or its antiparticle, and it is different from all the systems

discussed above, eight channels instead of four channels are coupled with each other under the one-boson exchange i.e. $(D^+B^{*0})_S$, $(D^+B^{*0})_D$, $(D^0B^{*+})_S$, interaction. $(D^{*+}B^0)_S$, $(D^{*+}B^0)_D$, $(D^{*0}B^+)_S$ and $(D^0 B^{*+})_D,$ $(D^{*0}B^+)_D$. We can investigate the possible bound states along the same line, although it is somewhat lengthy and tedious. There is ambiguity in choosing the μ^2 value as well, for the $DB^* \rightarrow D^*B$ scattering process, we could take $\mu^2 = m_{\text{ex}}^2 - (m_{D^*} - m_D)^2$ or $\mu^2 = m_{\text{ex}}^2 - (m_{B^*} - m_B)^2$, where m_{ex} is the mass of the exchanged boson. Specifically for $D^+B^{*0} \rightarrow D^{*+}B^0$ via π exchange, we can choose $\mu^2 = m_{\pi^0}^2 - (m_{D^{*+}} - m_{D^+})^2$ or $\mu^2 = m_{\pi^0}^2 - m_{\pi^0}^2 (m_{B^{*0}} - m_{B^0})^2$. This ambiguity has been taken into account in our analysis. The numerical results are given in Table VII, For $\Lambda = 808$ MeV, we find no bound state. With the choice $\mu^2 = m_{\text{ex}}^2 - (m_{D^*} - m_D)^2$, a bound state with mass 7189.7 MeV is found for $\Lambda = 850$ MeV, However, this solution disappears if one chooses $\mu^2 =$ $m_{\rm ex}^2 - (m_{B^*} - m_B)^2$. Only when Λ is around 880 MeV, the bound state solutions can be found for both μ^2 choices. The difference of the static properties for the two μ^2 choices is relatively larger than that of the above systems considered, this is because of the larger difference between $m_{D^*} - m_D \simeq 140$ MeV and $m_{B^*} - m_B \simeq 45$ MeV. We notice that the D^0B^{*+} component has the largest probability in the states, since the threshold of D^0B^{*+} is lower than that of D^+B^{*0} , $D^{*+}B^0$ and $D^{*0}B^+$. The wave function of the state with mass about 7185.9 MeV and $\Lambda = 900$ MeV is shown in Fig. 10, it is obvious all the eight components of the spatial wave function have the same sign, conse-

TABLE VII. The predictions for the static properties of the PV system with C = B = 1, where $P_S^{+0}(DB^*)$ denotes the probability of *S*-wave D^+B^{*0} , and $P_S^{+0}(D^*B)$ denotes the probability of *S*-wave $D^{*+}B^0$. Here $\mu_{\tilde{b}c1}^2 = m_{ex}^2 - (m_{D^*} - m_D)^2$ and $\mu_{\tilde{b}c2}^2 = m_{ex}^2 - (m_{B^*} - m_B)^2$ with m_{ex} the exchanged boson mass.

			Th	be pseudoscalar-vector system with $C = B = 1$
μ^2	Λ (MeV)	M (MeV)	r _{rms} (fm)	$P_{S}^{+0}(DB^{*}):P_{D}^{+0}(DB^{*}):P_{S}^{0+}(DB^{*}):P_{D}^{0+}(DB^{*}):P_{S}^{+0}(D^{*}B):P_{D}^{+0}(D^{*}B):P_{S}^{0+}(D^{*}B):P_{D}^{0+}($
$\mu_{\bar{h}c1}^2$	850	7189.7	5.54	8.10:0.02:89.69:0.02:0.87:0.36:0.67:0.26
bui	880	7187.9	2.05	21.88:0.07:72.40:0.08:2.09:0.96:1.73:0.80
	900	7185.9	1.58	27.34:0.11:65.19:0.12:2.54:1.35:2.16:1.19
	1000	7157.2	0.86	36.55:0.01:46.27:0.01:1.39:7.32:1.21:7.26
$\mu_{\bar{h}c^2}^2$	850	no bounded	-	-
• 002	880	7189.5	4.11	10.70:0.02:86.98:0.02:0.79:0.50:0.61:0.39
	900	7188.1	2.18	20.59:0.04:75.09:0.04:1.34:0.98:1.10:0.83
	1000	7161.1	0.88	36.86:0.22:47.18:0.23:0.48:7.33:0.41:7.31
			A	Il couplings are reduced by half except $g_{\pi NN}$
μ^2	Λ (MeV)	M (MeV)	$r_{\rm rms}~({\rm fm})$	$P_{S}^{+0}(DB^{*}):P_{D}^{+0}(DB^{*}):P_{S}^{0+}(DB^{*}):P_{D}^{0+}(DB^{*}):P_{S}^{+0}(D^{*}B):P_{D}^{+0}(D^{*}B):P_{S}^{0+}(D^{*}B):P_{D}^{0+}($
$\mu^2_{\bar{h}c1}$	970	7189.2	3.13	16.76:0.07:78.38:0.08:1.75:0.85:1.45:0.67
bui	1000	7187.6	1.88	25.51:0.13:66.52:0.15:2.78:1.35:2.42:1.15
	1100	7177.1	1.03	34.50:0.42:49.32:0.45:4.94:2.96:4.62:2.81
	1200	7156.3	0.78	34.62:0.72:41.51:0.75:5.85:5.51:5.64:5.42
$\mu_{\bar{h}c2}^2$	970	no bounded	-	
002	1000	7189.6	4.90	11.19:0.03:86.02:0.03:0.90:0.62:0.74:0.49
	1100	7181.9	1.20	34.04:0.20:54.01:0.21:3.26:2.67:3.08:2.53



FIG. 10 (color online). The wave function of the C = B = 1pseudoscalar-vector system with $\Lambda = 900$ MeV and $\mu^2 = m_{ex}^2 - (m_{D^*} - m_D)^2$, the mass of this state is about 7185.9 MeV.

quently this state would be isospin singlet in the isospin symmetry limit. Similar patterns of bound state solutions are predicted if the coupling constants except $g_{\pi NN}$ are reduced by half. This state is difficult to be produced, since both *c* and \bar{b} have to be produced simultaneously. The direct production of this state at hadron collider such as LHC and Tevatron is most promising, and the indirect production via top quark decay is a possible alternative. Once produced, it should be very stable, $DB\pi$ and $DB\gamma$ are the main decay channels.

E. Pseudoscalar-vector system with C = -B = 1

The effective interaction potentials are induced by oneboson exchange between two antiquarks, therefore both the π exchange and ω exchange contributions give opposite sign between the C = B = 1 system and the C = -B = 1system, nevertheless the overall signs of η , σ and ρ exchange potentials remain. We have eight coupled channels as well, $(D^+B^{*-})_S$, $(D^+B^{*-})_D$, $(D^0\bar{B}^{*0})_S$, $(D^0\bar{B}^{*0})_D$, $(D^{*+}B^-)_S$, $(D^{*+}B^-)_D$, $(D^{*0}\bar{B}^0)_S$ and $(D^{*0}\bar{B}^0)_D$ are involved. The numerical results are given in Table VIII. It is remarkable that the μ^2 and Λ dependence of the bound state solutions is similar to the C = B = 1 case. With the same μ^2 and Λ values, the predictions for the static properties of the two systems are not drastically different from each other. Concretely for $\Lambda = 900$ MeV and $\mu^2 = m_{ex}^2 - m_{ex}^2$ $(m_{D^*} - m_D)^2$, we find a bound state with mass 7187.6 MeV for the C = -B = 1 system, and the mass of C = B = 1bound state is 7185.9 MeV, the difference is about 1.7 MeV. The corresponding wave function with $\Lambda = 900$ MeV is plotted in Fig. 11, which can be roughly obtained by reversing the overall sign of the third, fourth, seventh and eighth components of the C = B = 1 system wave function in Fig. 10. To understand the similarity of the predic-

TABLE VIII. The predictions about the static properties of the PV system with C = -B = 1, where $P_S^{+-}(DB^*)$ denotes the probability of *S* wave D^+B^{*-} , and $P_S^{+-}(D^*B)$ denotes the probability of *S* wave $D^{*+}B^{-}$. Here $\mu_{bc1}^2 = m_{ex}^2 - (m_{D^*} - m_D)^2$ and $\mu_{bc2}^2 = m_{ex}^2 - (m_{B^*} - m_B)^2$ with m_{ex} the exchanged boson mass.

	The pseudoscalar-vector system with $C = -B = 1$					
μ^2	Λ (MeV)	M (MeV)	$r_{\rm rms}$ (fm)	$P_{S}^{+-}(DB^{*}):P_{D}^{+-}(DB^{*}):P_{S}^{00}(DB^{*}):P_{D}^{00}(DB^{*}):P_{S}^{+-}(D^{*}B):P_{D}^{+-}(D^{*}B):P_{S}^{00}(D^{*}B):P_{D}^{00}(D^{*}B)(\%)$		
μ_{bc1}^2	880	7189.2	3.11	15.16:0.04:80.15:0.05:2.21:0.45:1.63:0.32		
001	900	7187.6	1.85	23.82:0.08:68.07:0.09:3.83:0.61:3.02:0.48		
	1000	7169.0	0.79	30.56:0.28:46.30:0.30:11.86:0.74:9.27:0.69		
	1050	7148.4	0.51	15.57:0.02:53.99:0.04:22.84:0.10:7.41:0.03		
		7154.3	0.63	53.57:0.36:16.62:0.34:5.65:0.59:22.22:0.66		
μ_{bc2}^2	880	no bounded	-	-		
	900	7189.7	5.17	9.30:0.02:88.09:0.02:1.25:0.27:0.87:0.19		
	1000	7176.2	0.91	30.77:0.20:50.54:0.21:9.62:0.70:7.32:0.65		
	1050	7154.6	0.52	25.22:0.00:45.64:0.01:17.96:0.04:11.14:0.00		
		7163.3	0.70	46.16:0.31:27.72:0.30:7.95:0.64:16.25:0.68		
				All couplings except $g_{\pi NN}$ are reduced half		
μ^2	Λ (MeV)	M (MeV)	$r_{\rm rms}~({\rm fm})$	$P_{S}^{+-}(DB^{*}):P_{D}^{+-}(DB^{*}):P_{S}^{00}(DB^{*}):P_{D}^{00}(DB^{*}):P_{S}^{-}(D^{*}B):P_{D}^{+-}(D^{*}B):P_{S}^{00}(D^{*}B):P_{D}^{00}(D^{*}B)(\%)$		
μ^2_{bc1}	970	7189.4	3.57	15.68:0.06:78.98:0.07:2.29:0.58:1.90:0.44		
	1020	7185.6	1.38	30.23:0.20:56.36:0.21:5.82:1.03:5.25:0.90		
	1100	7173.2	0.82	33.18:0.42:43.08:0.45:10.53:1.24:9.93:1.17		
	1200	7147.4	0.59	30.96:0.68:35.69:0.70:15.05:1.28:14.40:1.25		
μ_{bc2}^2	970	no bounded	-	-		
	1020	7189.2	3.03	18.64:0.07:75.25:0.07:2.60:0.63:2.24:0.51		
	1100	7180.5	1.00	33.61:0.29:47.78:0.30:8.09:1.16:7.68:1.09		
	1200	7158.2	0.64	32.34:0.58:37.41:0.60:13.43:1.28:13.09:1.26		

tions for the C = B = 1 and C = -B = 1 system, we turn to the one π exchange model, the effective potential comprises a spin-spin potential proportional to $(\boldsymbol{\sigma}_i \cdot \boldsymbol{\sigma}_j)(\boldsymbol{\tau}_i \cdot \boldsymbol{\tau}_j)$ and a tensor potential proportional to $S_{ij}(\hat{\mathbf{r}})(\boldsymbol{\tau}_i \cdot \boldsymbol{\tau}_j)$, where the isospin matrix τ_i and the spin matrix σ_i only act on the light quarks. In the basis of the eight channels listed above, these two operators can be written as 8×8 matrices

$$(\boldsymbol{\sigma}_{i} \cdot \boldsymbol{\sigma}_{j})(\boldsymbol{\tau}_{i} \cdot \boldsymbol{\tau}_{j}) \rightarrow \begin{pmatrix} 0 & 0 & 0 & 0 & -1 & 0 & 2 & 0 \\ 0 & 0 & 0 & 0 & 2 & 0 & -1 & 0 \\ 0 & 0 & 0 & 0 & 0 & 2 & 0 & -1 \\ -1 & 0 & 2 & 0 & 0 & 0 & 0 & 0 \\ 2 & 0 & -1 & 0 & 2 & 0 & 0 & 0 & 0 \\ 0 & 2 & 0 & -1 & 0 & 0 & 0 & 0 & 0 \end{pmatrix}$$

$$S_{ij}(\hat{\mathbf{r}})(\boldsymbol{\tau}_{i} \cdot \boldsymbol{\tau}_{j}) \rightarrow \begin{pmatrix} 0 & 0 & 0 & 0 & 0 & \sqrt{2} & 0 & -2\sqrt{2} \\ 0 & 0 & 0 & 0 & \sqrt{2} & -1 & -2\sqrt{2} & 2 \\ 0 & 0 & 0 & 0 & 0 & -2\sqrt{2} & 2 & \sqrt{2} & -1 \\ 0 & \sqrt{2} & 0 & -2\sqrt{2} & 2 & \sqrt{2} & -1 \\ 0 & \sqrt{2} & 0 & -2\sqrt{2} & 0 & 0 & 0 \\ \sqrt{2} & -1 & -2\sqrt{2} & 2 & 0 & 0 & 0 \\ \sqrt{2} & -1 & -2\sqrt{2} & 2 & 0 & 0 & 0 & 0 \\ 0 & -2\sqrt{2} & 0 & \sqrt{2} & 0 & 0 & 0 & 0 \\ -2\sqrt{2} & 2 & \sqrt{2} & -1 & 0 & 0 & 0 & 0 \end{pmatrix},$$

$$(21)$$

For the C = B = 1 pseudoscalar-vector system, the corresponding matrix representations are obtained by replacing 2 and $2\sqrt{2}$ with -2 and $-2\sqrt{2}$ respectively in Eq. (21). It is obvious both operators contribute to only the off-diagonal 4×4 matrix elements. As a result, the eigenvalues of the corresponding Schrödinger equation for the C = B = 1 and C = -B = 1 cases are exactly the same, if the small



FIG. 11 (color online). The wave function of the C = -B = 1pseudoscalar-vector system with $\Lambda = 900$ MeV and $\mu^2 = m_{ex}^2 - (m_{D^*} - m_D)^2$, its mass is about 7187.6 MeV.

mass difference within the isospin multiplets is neglected, and the eingen-wave function of one system can be obtained from another by reversing the overall sign of the third, fourth, seventh and eighth components. Therefore the heavy bosons η , σ , ρ and ω exchange contributes to effective potential, and the pion exchange contribution is still dominant. In short summary, even after including shorter distance contributions from η , σ , ρ and ω exchange, the results obtained are qualitatively the same as those in the one π exchange model. The same conclusion has been reached for all the system consider above.

VII. CONCLUSIONS AND DISCUSSIONS

Motivated by the nucleon-nucleon interaction, we have represented the short range interaction by heavier mesons η , σ , ρ and ω exchange. The effective potentials between two hadrons are obtained by summing the interactions between light quarks or antiquarks via one-boson exchange. The potential becomes more complicated than that in the one pion exchange model, and there are six additional terms which are proportional to 1, $\tau_i \cdot \tau_j$, $\sigma_i \cdot \sigma_j$, $S_{ij}(\hat{\mathbf{r}})$, $\mathbf{L} \cdot \mathbf{S}_{ij}$ and $(\mathbf{L} \cdot \mathbf{S}_{ij})(\tau_i \cdot \tau_j)$ respectively.

We first apply the one-boson exchange formalism to the deuteron, then generalize to $D\bar{D}^*/\bar{D}D^*$, $B\bar{B}^*/\bar{B}B^*$, DD^* , BB^* , PV systems with C = B = 1 and C = -B = 1. S-D mixing effects has been taken into account, and the un-

certainties from the regularization parameter Λ and effective coupling constants are considered. We find the conclusions reached are qualitatively the same as those in the one pion exchange model. This implies that the long range π exchange effects dominate the physics of a weakly bound hadronic molecule, and we can safely use one pion exchange model to qualitatively discuss the binding of molecule candidates. Since the predictions for the binding energy and other static properties are sensitive to the regularization parameter Λ and the effective couplings, we are not able to predict the binding energies very precisely. If the potential is so strong that binding energy is large enough, we would be quite confident that such bound state must exist. However, the exact binding energy will depend on the details of the regularization and the effective couplings involved. Our results indicate that the $1^{++} B\bar{B}^*/\bar{B}B^*$ molecule should exist, whereas DD^* bound state does not exist. For $\Lambda = 808$ MeV (970 MeV), the binding energy, D wave probability and other static properties of deuteron are produced, meanwhile near threshold $1^{++} D\bar{D}^*/\bar{D}D^*$ molecule is predicted. To identify this state with X(3872), the mixing between this $D\bar{D}^*/\bar{D}D^*$ molecule and the conventional charmonium state should be further considered to be consistent with the recent experimental data on $X(3872) \rightarrow \psi(2S)\gamma$ [27]. For the BB^{*} system, the PV systems with C = B = 1 and C = -B = 1, near threshold molecular states may exist. Similar to the $1^{++} D\bar{D}^*/\bar{D}D^*$ molecule, these states should be rather stable, isospin is drastically broken, and the I = 0 component is dominant. Direct production of the above doubly heavy states at Tevatron and LHC is the most promising way. We can search for the 1⁺⁺ $B\bar{B}^*/\bar{B}B^*$ molecule via $p\bar{p} \rightarrow$ $\pi^+\pi^-\Upsilon(1S)$ at Tevatron. The BB^* bound state mainly decays into $BB\gamma$ if it really exists. The dominant decay channels of the heavy flavor PV bound state with C = B =1 are $DB\pi$ and $DB\gamma$, and the possible heavy flavor PV bound state with C = -B = 1 mainly decays into $D\bar{B}\pi$ and $D\bar{B}\gamma$.

In our model, the involved parameters include the effective quark-boson couplings, the masses of the exchanged bosons and the hadrons inside the molecule. Therefore this model is quite general, it can be widely used to dynamically study the possible molecular candidates. We will further apply the one-boson exchange model to baryonantibaryon system etc, and compare the predictions with the recent experimental observations [30].

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APPENDIX: THE MATRIX ELEMENTS OF THE SPIN RELEVANT OPERATORS

For initial state consisting of two mesons A and B, with relative angular momentum L, total spin S and total angular momentum J, its wave function is written as

$$|(AB)LS, JM_{J}\rangle = \sum_{M_{L},M_{S}} \langle LM_{L}; SM_{S} | JM_{J}\rangle | LM_{L}\rangle | SM_{S}\rangle$$
$$= \sum_{S_{13},S_{24}} \hat{S}_{A} \hat{S}_{B} \hat{S}_{13} \hat{S}_{24} \begin{cases} 1/2 & 1/2 & S_{A} \\ 1/2 & 1/2 & S_{B} \\ S_{13} & S_{24} & S \end{cases}$$
$$\times | L(S_{13}S_{24})S, JM_{J}\rangle, \qquad (A1)$$

where $\hat{S} = \sqrt{2S + 1}$. For the convenience of calculating the matrix elements of the spin-orbit operator $\mathbf{L} \cdot \mathbf{S}_{24}$, we can recouple the state as

$$|(AB)LS, JM_{J}\rangle = \sum_{S_{13}, S_{24}, J_{LS}} (-1)^{L+S+J} \hat{S}_{A} \hat{S}_{B} \hat{S}_{13} \hat{S}_{24} \hat{S} \hat{J}_{LS} \times \left\{ \begin{array}{ccc} L & S_{24} & J_{LS} \\ S_{13} & J & S \end{array} \right\} \left\{ \begin{array}{ccc} 1/2 & 1/2 & S_{A} \\ 1/2 & 1/2 & S_{B} \\ S_{13} & S_{24} & S \end{array} \right\} \times |(LS_{24})J_{LS}S_{13}, JM_{J}\rangle.$$
(A2)

In the same way, we can recouple the final state $|(A'B')L'S', J'M'_{j}\rangle$ via the Wigner 6-*j* and 9-*j* coefficients. In the following, we shall present the matrix elements of four light quark operators involved in the work, which is helpful to calculating the matrix representation of the effective interactions.

(1) The unit operator 1

Using Eq. (A1), it is obvious that

$$\langle (A'B')L'S', J'M'_{J}|\mathbf{1}|(AB)LS, JM_{J} \rangle = \delta_{LL'}\delta_{SS'}\delta_{JJ'}\delta_{M_{J}M'_{J}} \sum_{S_{13},S_{24}} \hat{S}_{A}\hat{S}'_{A}\hat{S}_{B}\hat{S}'_{B}\hat{S}^{2}_{13}\hat{S}^{2}_{24} \begin{cases} 1/2 & 1/2 & S_{A} \\ 1/2 & 1/2 & S_{B} \\ S_{13} & S_{24} & S \end{cases} \begin{cases} 1/2 & 1/2 & S_{A} \\ 1/2 & 1/2 & S_{B} \\ S_{13} & S_{24} & S \end{cases} \\ = \delta_{LL'}\delta_{SS'}\delta_{S_{A}S'_{A}}\delta_{S_{B}S'_{B}}\delta_{JJ'}\delta_{M_{J}M'_{J}}.$$

$$(A3)$$

(2) The spin-spin operator $\boldsymbol{\sigma}_2 \cdot \boldsymbol{\sigma}_4$

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$$\langle (A'B')L'S', J'M'_{J} | \boldsymbol{\sigma}_{2} \cdot \boldsymbol{\sigma}_{4} | (AB)LS, JM_{J} \rangle = \delta_{LL'} \delta_{SS'} \delta_{JJ'} \delta_{M_{J}M'_{J}} \sum_{S_{13}, S_{24}} \hat{S}_{A} \hat{S}'_{A} \hat{S}_{B} \hat{S}'_{B} \hat{S}^{2}_{13} \hat{S}^{2}_{24} [2S_{24}(S_{24}+1)-3] \\ \times \begin{cases} 1/2 & 1/2 & S_{A} & 1/2 & 1/2 & S'_{A} \\ 1/2 & 1/2 & S_{B} \} \{1/2 & 1/2 & S'_{B} \\ S_{13} & S_{24} & S & S_{13} & S_{24} & S \end{cases},$$
(A4)

where the spin operators σ_2 and σ_4 only act on the light quarks and antiquarks.

(3) The spin-orbit operator $\mathbf{L} \cdot \mathbf{S}_{24}$

$$\langle (A'B')L'S', J'M'_{J}|\mathbf{L} \cdot \mathbf{S}_{24}|(AB)LS, JM_{J} \rangle = \delta_{LL'} \delta_{JJ'} \delta_{M_{J}M'_{J}} \sum_{S_{13}, S_{24}, J_{LS}} (-1)^{S+S'+2L+2J} \hat{S}_{A} \hat{S}'_{A} \hat{S}_{B} \hat{S}'_{B} \hat{S}'_{B} \hat{S}'_{B} \hat{S}_{13} \hat{S}_{24}^{2} \hat{J}_{LS}^{2} \frac{1}{2}$$

$$\times \left[J_{LS}(J_{LS}+1) - L(L+1) - S_{24}(S_{24}+1) \right]$$

$$\times \left\{ \begin{array}{ccc} L & S_{24} & J_{LS} \\ S_{13} & J & S \end{array} \right\} \left[\begin{array}{ccc} L & S_{24} & J_{LS} \\ S_{13} & J & S' \end{array} \right] \left\{ \begin{array}{ccc} 1/2 & 1/2 & S_{A} & 1/2 & 1/2 & S'_{A} \\ 1/2 & 1/2 & S_{B} \right\} \{1/2 & 1/2 & S'_{B} \\ S_{13} & S_{24} & S & S_{13} & S_{24} & S \end{array} \right\},$$

$$(A5)$$

where $\mathbf{S}_{24} = \frac{1}{2}(\boldsymbol{\sigma}_2 + \boldsymbol{\sigma}_4)$, **L** is the relative spatial angular momentum. The matrix elements of $\mathbf{L} \cdot \mathbf{S}_{24}$ can be calculated by the Wigner-Echart theorem [31], and the same result has been obtained.

(4) The tensor operator S₂₄(**r**) ≡ 3(**σ**₂ · **r**)(**σ**₄ · **r**) - **σ**₂ · **σ**₄
 It can be checked that the tensor operator S₂₄(*r*) is proportional to the scalar product of two rank-2 tensor operators Y_{2m} and S⁽²⁾_m with m = 0, ±1, ±2, where Y_{2m} is the spherical harmonic function of degree 2, and the five components of S⁽²⁾_m are

$$S_{2}^{(2)} = \frac{1}{2}S_{2+}S_{4+}, \qquad S_{1}^{(2)} = -\frac{1}{2}(S_{20}S_{4+} + S_{2+}S_{40}), \qquad S_{0}^{(2)} = -\frac{\sqrt{6}}{12}(S_{2-}S_{4+} - 4S_{20}S_{40} + S_{2+}S_{4-})$$
$$S_{-1}^{(2)} = \frac{1}{2}(S_{2-}S_{40} + S_{20}S_{4-}), \qquad S_{-2}^{(2)} = \frac{1}{2}S_{2-}S_{4-}.$$
(A6)

Here $S_{2+} = \frac{1}{2}(\sigma_{2x} + i\sigma_{2y})$, $S_{20} = \frac{1}{2}\sigma_{20}$ and $S_{2-} = \frac{1}{2}(\sigma_{2x} - i\sigma_{2y})$. The same convention applies to $S_{4,\pm}$ and S_{40} , the spin operators σ_2 and σ_4 only act on the light quark and antiquarks. Using the Wigner-Echart theorem, the matrix element of this tensor operator can be obtained, although it is somewhat lengthy.

$$\langle (A'B')L'S', J'M'_{J}|S_{24}(\hat{\mathbf{r}})|(AB)LS, JM_{J}\rangle = \delta_{JJ'}\delta_{M_{J}M'_{J}}\frac{2}{3}\sqrt{30}\sum_{S_{13},S_{24}}\delta_{S_{24},1}(-1)^{J+L+L'+2S'+S_{13}+S_{24}}\hat{S}_{A}\hat{S}'_{A}\hat{S}_{B}\hat{S}'_{B}\hat{S}'_{L}\hat{L}'\hat{S}^{2}_{13}\hat{S}^{4}_{24} \\ \times \left\{ \begin{matrix} L' & S' & J \\ S & L & 2 \end{matrix} \right\} \left[\begin{matrix} S_{24} & S' & S_{13} \\ S & S_{24} & 2 \end{matrix} \right] \left(\begin{matrix} L' & 2 & L \\ 0 & 0 & 0 \end{matrix} \right) \\ \times \left\{ \begin{matrix} 1/2 & 1/2 & S_{A} \\ 1/2 & 1/2 & S_{B} \\ S_{13} & S_{24} & S \end{matrix} \right\} \left\{ \begin{matrix} 1/2 & 1/2 & S'_{A} \\ 1/2 & 1/2 & S'_{B} \\ S_{13} & S_{24} & S \end{matrix} \right\} \left\{ \begin{matrix} 1/2 & 1/2 & S'_{A} \\ 1/2 & 1/2 & S'_{B} \\ S_{13} & S_{24} & S' \end{matrix} \right\}.$$
 (A7)

The above expression is apparently different from the results in Ref. [7], However, the numerical results of all the matrix elements are the same.

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