

# Supplemental Material

Fang Qin (覃昉)<sup>1,2</sup> and Pengfei Zhang<sup>3,4,\*</sup>

<sup>1</sup>*Shenzhen Institute for Quantum Science and Engineering and Department of Physics,  
Southern University of Science and Technology (SUSTech), Shenzhen 518055, China*

<sup>2</sup>*CAS Key Laboratory of Quantum Information, University of Science and Technology of China,  
Chinese Academy of Sciences, Hefei, Anhui 230026, China*

<sup>3</sup>*Walter Burke Institute for Theoretical Physics, California Institute of Technology, Pasadena, California 91125, USA*

<sup>4</sup>*Institute for Quantum Information and Matter,  
California Institute of Technology, Pasadena, California 91125, USA*

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In this Supplemental Material, we provide the detailed calculations for the results in the main text.

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## SI. DIMER PROPAGATOR MATRIX

We consider finite total momentum  $Q$  for each two-body pairing state. As shown in Fig. S1, the inverse of the dimer propagator matrix is given by the Dyson equation

$$\begin{aligned}
 D^{-1}(E_0, Q) &= \begin{pmatrix} (ig_S)^{-1} & 0 \\ 0 & (ig_P)^{-1} \end{pmatrix} - \begin{pmatrix} \Pi_{SS}(E_0, Q) & \Pi_{SP}(E_0, Q) \\ \Pi_{PS}(E_0, Q) & \Pi_{PP}(E_0, Q) \end{pmatrix} \\
 &= \begin{pmatrix} (ig_S)^{-1} - \Pi_{SS}(E_0, Q) & -\Pi_{SP}(E_0, Q) \\ -\Pi_{PS}(E_0, Q) & (ig_P)^{-1} - \Pi_{PP}(E_0, Q) \end{pmatrix}, \tag{S1}
 \end{aligned}$$

where the polarization bubble is given by (the derivations are given in the Appendix)

$$\Pi_{\alpha\beta}(E_0, Q) = - \int \frac{dp dp_0}{(2\pi)^2} \frac{p^{l_\alpha + l_\beta}}{2} \text{Tr} \left[ G^T(p_0, Q/2 + p) \sigma_\alpha G(E_0 - p_0, Q/2 - p) \sigma_\beta^\dagger \right], \tag{S2}$$

$E_0 = Q^2/(4m) + k^2/m$  is the total energy,  $\alpha, \beta \in \{S, P\}$ ,  $l_S = 0$ ,  $l_P = 1$ , and Tr denotes the trace over the spin degrees of freedom.

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\* pengfeizhang.physics@gmail.com

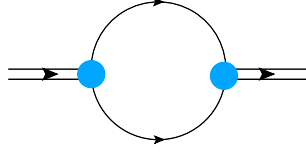


FIG. S1. Feynman diagrams for the matrix elements of the dimer-atom interaction operator. The single line denotes the bare atom propagator matrix  $G$ , the double lines denote the matrix elements of the dimer propagator matrix  $D_{\alpha\beta}$  with  $\alpha, \beta \in \{S, P\}$ , and the blue dot represents the interaction vertex:  $-i\sigma_\alpha$  or  $-i\sigma_\beta$ .

For convenience, one can also write the dimer propagator matrix as

$$D(E_0, Q) = \frac{1}{\det[D^{-1}(E_0, Q)]} \begin{pmatrix} (ig_P)^{-1} - \Pi_{PP}(E_0, Q) & \Pi_{SP}(E_0, Q) \\ \Pi_{PS}(E_0, Q) & (ig_S)^{-1} - \Pi_{SS}(E_0, Q) \end{pmatrix} = \begin{pmatrix} D_{SS}(E_0, Q) & D_{SP}(E_0, Q) \\ D_{PS}(E_0, Q) & D_{PP}(E_0, Q) \end{pmatrix}, \quad (\text{S3})$$

where  $\det[D^{-1}(E_0, Q)]$  is the determinant of  $D^{-1}(E_0, Q)$ .

In the absence of the Raman coupling, the  $s$ -wave polarization bubble is given by

$$\Pi_{SS}(E_0, Q) = - \int \frac{dpdp_0}{(2\pi)^2} \frac{i}{p_0 - (Q/2 + p)^2/(2m) + i0^+} \frac{i}{E_0 - p_0 - (Q/2 - p)^2/(2m) + i0^+} = -\frac{m}{2k}, \quad (\text{S4})$$

where we use  $E_0 = Q^2/(4m) + k^2/m$ .

The 1D scattering amplitude can be written as [1, 2]  $f_{1D}(k) = -1/(1 + i \cot \delta_k) \simeq -1/(1 + ik a_s)$ , where  $\delta_k$  is the scattering phase shift. With the  $s$ -wave  $T$  matrix  $T_s(k) = ik f_{1D}(k)/m_r$  and  $D_{SS}(k) = iT_s(k)$ , one can get

$$a_s = -\frac{2}{mg_S}, \quad (\text{S5})$$

where  $a_s$  is the effective 1D  $s$ -wave scattering length.

In the absence of the Raman coupling, the  $p$ -wave polarization bubble is given by

$$\begin{aligned} \Pi_{PP}(E_0, Q) &= -\frac{1}{2} \int \frac{dpdp_0}{(2\pi)^2} \frac{ip^2}{p_0 - (Q/2 + p)^2/(2m) + i0^+} \frac{i}{E_0 - p_0 - (Q/2 - p)^2/(2m) + i0^+} \\ &= -\frac{1}{2} \int \frac{dp}{2\pi} \frac{ip^2}{E_0 - Q^2/(4m) - p^2/m + i0^+} = \frac{im}{2} \int_{-\Lambda}^{\Lambda} \frac{dp}{2\pi} \frac{(p^2 - k^2) + k^2}{[p - (k + i0^+)][p + (k + i0^+)]} \\ &= \frac{im}{2} \left( \frac{\Lambda}{\pi} + i\frac{k}{2} \right), \end{aligned} \quad (\text{S6})$$

where  $\Lambda$  is the ultraviolet momentum cutoff and we use  $E_0 = Q^2/(4m) + k^2/m$ .

Therefore, the  $p$ -wave  $T$  matrix is given by

$$\frac{T_{pp}(E_0, Q)}{qq'} = \frac{D_{PP}(E_0, Q)}{i} = \frac{1}{\frac{1}{g_P} - i\Pi_{pp}(E_0, Q)} = \frac{1}{\frac{1}{g_P} + \frac{m\Lambda}{2\pi} + i\frac{mk}{4}}. \quad (\text{S7})$$

Comparing Eq. (S7) with [3]

$$\frac{T_{pp}(E_0, Q)}{qq'} = \frac{1}{\frac{m}{4a_p} + i\frac{mk}{4}}, \quad (\text{S8})$$

one can get

$$\frac{m}{4a_p} = \frac{1}{g_P} + \frac{m\Lambda}{2\pi}, \quad (\text{S9})$$

where  $a_p$  is the effective 1D  $p$ -wave scattering length.



FIG. S2. Feynman diagrams for the matrix elements of the dimer local operator  $\varphi_\alpha^\dagger(R)\varphi_\beta(R)$  and its derivatives  $\varphi_\alpha^\dagger(R) [i\partial_t + \partial_R^2/(4m)]^u (-i\partial_R)^v \varphi_\beta(R)$  with  $u, v = 0, 1, 2, 3, \dots$ . The open dot represents the operators.

In the presence of the Raman fields, the bubble with zero total momentum ( $Q = 0$ ) can be calculated as

$$\begin{aligned} \Pi_{SS}(E_0, 0) &= \frac{-m^3\Omega^2}{2(mE_0k_0^2 - k_0^4 + m^2\Omega^2)\sqrt{mE_0 + i0^+ - k_0^2}} - \frac{mk_0^2}{4} \left[ \frac{\sqrt{mE_0 + k_0^2 + 2\sqrt{mE_0k_0^2 + m^2\Omega^2}}}{m^2\Omega^2 + k_0^2(mE_0 + \sqrt{mE_0k_0^2 + m^2\Omega^2})} \right. \\ &\quad \left. + \frac{\sqrt{mE_0 + i0^+ + k_0^2 - 2\sqrt{mE_0k_0^2 + m^2\Omega^2}}}{m^2\Omega^2 + k_0^2(mE_0 - \sqrt{mE_0k_0^2 + m^2\Omega^2})} \right] \approx \frac{-m}{2\sqrt{mE_0}}, \end{aligned} \quad (\text{S10})$$

$$\begin{aligned} \Pi_{PP}(E_0, 0) &= \frac{im}{2} \left( \frac{\Lambda}{\pi} + i \frac{\sqrt{mE_0 + i0^+ - k_0^2}}{2} \right) - \frac{m^3\Omega^2}{16} \left( \frac{\sqrt{mE_0 + k_0^2 + 2\sqrt{mE_0k_0^2 + m^2\Omega^2}}}{mE_0k_0^2 + m^2\Omega^2 + k_0^2\sqrt{mE_0k_0^2 + m^2\Omega^2}} \right. \\ &\quad \left. + \frac{\sqrt{mE_0 + i0^+ + k_0^2 - 2\sqrt{mE_0k_0^2 + m^2\Omega^2}}}{mE_0k_0^2 + m^2\Omega^2 - k_0^2\sqrt{mE_0k_0^2 + m^2\Omega^2}} \right) \approx \frac{im}{2} \left( \frac{\Lambda}{\pi} + i \frac{\sqrt{mE_0 + i0^+ - k_0^2}}{2} \right), \end{aligned} \quad (\text{S11})$$

$$\begin{aligned} \Pi_{SP}(E_0, 0) &= \Pi_{PS}(E_0, 0) \\ &= \frac{m^2k_0\Omega\sqrt{mE_0 + i0^+ - k_0^2}}{4(mE_0k_0^2 - k_0^4 + m^2\Omega^2)} - \frac{m^2k_0\Omega}{8} \left( \frac{\sqrt{mE_0 + k_0^2 + 2\sqrt{mE_0k_0^2 + m^2\Omega^2}}}{mE_0k_0^2 + m^2\Omega^2 + k_0^2\sqrt{mE_0k_0^2 + m^2\Omega^2}} \right. \\ &\quad \left. + \frac{\sqrt{mE_0 + i0^+ + k_0^2 - 2\sqrt{mE_0k_0^2 + m^2\Omega^2}}}{mE_0k_0^2 + m^2\Omega^2 - k_0^2\sqrt{mE_0k_0^2 + m^2\Omega^2}} \right) \approx \frac{\sqrt{m}k_0\Omega}{8E_0^{-3/2}}, \end{aligned} \quad (\text{S12})$$

where  $k_0$  and  $\Omega$  are treated perturbatively up to the  $k_0^2$  and  $\Omega$  order. Note that Eq. (S9) can be used to cancel the divergence of the  $p$ -wave bubble (S11), and the  $s$ - and  $p$ -wave mixing bubble (S12) needs both finite  $k_0$  and  $\Omega$ , as expected from having a non-trivial SOC.

## SII. OTHER UNIVERSAL RELATIONS

### A. Adiabatic relations

The traditional  $s$ - and  $p$ -wave adiabatic relations are given by [2, 3]

$$\frac{C_{SS}}{2m} \equiv \frac{\partial E}{\partial a_s} = - \int dR \left\langle \frac{\partial \mathcal{L}(R)}{\partial a_s} \right\rangle = \frac{m}{2} \int dR \left\langle \varphi_S^\dagger(R)\varphi_S(R) \right\rangle, \quad (\text{S13})$$

$$\frac{C_{PP}}{4m} \equiv - \frac{\partial E}{\partial a_p^{-1}} = \int dR \left\langle \frac{\partial \mathcal{L}(R)}{\partial a_p^{-1}} \right\rangle = \frac{m}{4} \int dR \left\langle \varphi_P^\dagger(R)\varphi_P(R) \right\rangle, \quad (\text{S14})$$

where  $E$  is the total energy of the many-body system,  $\mathcal{L}(R)$  is the density of the Lagrangian (??),  $C_{SS}$  ( $C_{PP}$ ) is the 1D  $s(p)$ -wave contact.

When SOC is present, there are two new parameters  $k_0$  and  $\Omega$ . One can define two new contacts  $C_\lambda$  and  $C_\Omega$  as

$$C_\lambda \equiv - \int dR \left\langle \frac{\partial \mathcal{L}(R)}{\partial k_0} \right\rangle, \quad (\text{S15})$$

$$C_\Omega \equiv - \int dR \left\langle \frac{\partial \mathcal{L}(R)}{\partial \Omega} \right\rangle. \quad (\text{S16})$$

Here,  $C_\lambda$  and  $C_\Omega$  refer to only single-atom operators which give nonzero matrix elements in the single-atom sector. The momentum distribution under single-particle states is just a delta function, so that  $C_\lambda$  and  $C_\Omega$  will not contribute to the large-momentum tail, which is different from  $C_{SS}$  and  $C_{PP}$  [4]. However, both  $k_0$  and  $\Omega$  have nonzero energy scale, so that they would appear in the pressure relation and virial theorem.

### B. Pressure relation

For a uniform gas, the pressure relation can be derived following the expression of the Helmholtz free energy density  $\mathcal{F} = F/L$  which can be expressed in terms of [2, 5, 6]

$$\mathcal{F}(T, n_\uparrow, n_\downarrow, a_s, a_p, k_0, \Omega) = \frac{k_F^3}{2m} f\left(\frac{2mT}{k_F^2}, \frac{n_\uparrow}{k_F}, \frac{n_\downarrow}{k_F}, a_s k_F, a_p k_F, \frac{k_0}{k_F}, \frac{2m\Omega}{k_F^2}\right), \quad (\text{S17})$$

where  $L$  is the length along the  $x$  direction,  $f$  is a dimensionless function,  $T$  is the temperature,  $n = n_\uparrow + n_\downarrow = k_F/\pi$  is the Fermi particle number density, and  $k_F$  is the Fermi wave vector.

Equation (S17) implies the scaling behavior of the Helmholtz free energy density as follows:

$$\tilde{\lambda}^3 \mathcal{F}(T, n_\uparrow, n_\downarrow, a_s, a_p, k_0, \Omega) = \mathcal{F}\left(\tilde{\lambda}^2 T, \tilde{\lambda} n_\uparrow, \tilde{\lambda} n_\downarrow, \tilde{\lambda}^{-1} a_s, \tilde{\lambda}^{-1} a_p, \tilde{\lambda} k_0, \tilde{\lambda}^2 \Omega\right), \quad (\text{S18})$$

where  $\tilde{\lambda}$  is a dimensionless and arbitrary parameter.

Taking the derivative of Eq. (S18) with respect to  $\tilde{\lambda}$  at  $\tilde{\lambda} = 1$ , we have

$$3\mathcal{F} = \left(2T \frac{\partial}{\partial T} + n_\uparrow \frac{\partial}{\partial n_\uparrow} + n_\downarrow \frac{\partial}{\partial n_\downarrow} - a_s \frac{\partial}{\partial a_s} - a_p \frac{\partial}{\partial a_p} + k_0 \frac{\partial}{\partial k_0} + 2\Omega \frac{\partial}{\partial \Omega}\right) \mathcal{F}. \quad (\text{S19})$$

Replacing the free energy density  $\mathcal{F}$  in the left side of Eq. (S19) by  $n_\uparrow \mu_\uparrow + n_\downarrow \mu_\downarrow - \mathcal{P}$  and substituting  $S = -\partial F/\partial T$  and  $\mu_\sigma = \partial F/\partial n_\sigma$  into Eq. (S19), one gets

$$3(n_\uparrow \mu_\uparrow + n_\downarrow \mu_\downarrow - \mathcal{P}) = -2TS + n_\uparrow \mu_\uparrow + n_\downarrow \mu_\downarrow - a_s \frac{\partial \mathcal{F}}{\partial a_s} - a_p \frac{\partial \mathcal{F}}{\partial a_p} + k_0 \frac{\partial \mathcal{F}}{\partial k_0} + 2\Omega \frac{\partial \mathcal{F}}{\partial \Omega}, \quad (\text{S20})$$

where  $\mathcal{P}$  is the pressure density,  $S$  is the entropy, and  $\mu_\sigma$  is the chemical potential with spin  $\sigma$ .

Using the adiabatic relations (S13), (S14), (S15) and (S16), we can get the pressure relation as

$$\mathcal{P} = 2\mathcal{E} + \frac{a_s C_{SS}}{2mL} + \frac{C_{PP}}{4ma_p L} - \frac{k_0 C_\lambda}{L} - \frac{2\Omega C_\Omega}{L}, \quad (\text{S21})$$

where  $\mathcal{E} = E/L$  is the energy density and we use  $E = F + TS$  and

$$-a_p \frac{\partial E}{\partial a_p} = -a_p \frac{\partial E}{\partial a_p^{-1}} \frac{\partial a_p^{-1}}{\partial a_p} = -\frac{C_{PP}}{4ma_p}. \quad (\text{S22})$$

### C. Virial theorem

For an atomic gas in a harmonic potential  $V_T = m\omega^2 x^2/2$  with the trapping frequency  $\omega$ , the free energy can be expressed in terms of [2, 5, 6]

$$F(T, \omega, a_s, a_p, k_0, \Omega, N_\uparrow, N_\downarrow) = \omega \tilde{f}(T/\omega, \omega/\omega, a_s/a_{\text{ho}}, a_p/a_{\text{ho}}, k_0 a_{\text{ho}}, \Omega/\omega, N_\uparrow, N_\downarrow), \quad (\text{S23})$$

where  $N = N_\uparrow + N_\downarrow$  is the particle number,  $a_{\text{ho}} = \sqrt{2/(m\omega)}$  is the harmonic oscillator length and the dimensionless function  $\tilde{f}$  is dependent on the dimensionless ratios  $T/\omega$ ,  $a_s/a_{\text{ho}}$ ,  $a_p/a_{\text{ho}}$ ,  $k_0 a_{\text{ho}}$ ,  $\Omega/\omega$ , and particle numbers  $N_\uparrow$  and  $N_\downarrow$ .

With Eq. (S23), we can get the scaling law

$$\tilde{\lambda} F(T, \omega, a_s, a_p, k_0, \Omega, N_\uparrow, N_\downarrow) = F(\tilde{\lambda} T, \tilde{\lambda} \omega, \tilde{\lambda}^{-1/2} a_s, \tilde{\lambda}^{-1/2} a_p, \tilde{\lambda}^{1/2} k_0, \tilde{\lambda} \Omega, N_\uparrow, N_\downarrow), \quad (\text{S24})$$

where  $\tilde{\lambda}$  is a dimensionless and arbitrary parameter.

The derivative of Eq. (S24) with respect to  $\tilde{\lambda}$  at  $\tilde{\lambda} = 1$  gives

$$F = \left( T \frac{\partial}{\partial T} + \omega \frac{\partial}{\partial \omega} - \frac{1}{2} a_s \frac{\partial}{\partial a_s} - \frac{1}{2} a_p \frac{\partial}{\partial a_p} + \frac{1}{2} k_0 \frac{\partial}{\partial k_0} + \Omega \frac{\partial}{\partial \Omega} \right) F. \quad (\text{S25})$$

Substituting  $E = F + TS$  and  $S = -\partial F / \partial T$  into Eq. (S25), one gets

$$E = \left( \omega \frac{\partial}{\partial \omega} - \frac{1}{2} a_s \frac{\partial}{\partial a_s} - \frac{1}{2} a_p \frac{\partial}{\partial a_p} + \frac{1}{2} k_0 \frac{\partial}{\partial k_0} + \Omega \frac{\partial}{\partial \Omega} \right) E, \quad (\text{S26})$$

which, together with the Hellmann-Feynman theorem and the adiabatic relations (S13), (S14), (S15) and (S16), gives

$$E = 2\langle V_T \rangle - \frac{a_s C_{SS}}{4m} - \frac{C_{PP}}{8ma_p} + \frac{k_0 C_\lambda}{2} + \Omega C_\Omega. \quad (\text{S27})$$

### III. CONTACTS IN TWO-BODY BOUND STATES

The inverse of the two-body  $T$  matrix can be written as

$$T^{-1}(E_0, Q) \sim iD^{-1}(E_0, Q) = \begin{pmatrix} g_S^{-1} - i\Pi_{SS}(E_0, Q) & \delta_{SP} - i\Pi_{SP}(E_0, Q) \\ \delta_{PS} - i\Pi_{PS}(E_0, Q) & g_P^{-1} - i\Pi_{PP}(E_0, Q) \end{pmatrix}, \quad (\text{S28})$$

where  $\delta_{SP} = \delta_{PS}$  is a phenomenological parameter which is used to describe the bare coupling between  $s$ - and  $p$ -wave interactions on the two-body level. It corresponds to add a term  $\sum_Q \left( \delta_{SP} \varphi_{Q,S}^\dagger \varphi_{Q,P} + \delta_{PS} \varphi_{Q,P}^\dagger \varphi_{Q,S} \right)$  in the Lagrangian. We introduce this additional coupling to derive the contact using the adiabatic relation.

The binding energy  $E_b$  can be calculated by the pole of the  $T$ -matrix as  $\det(T^{-1}(E_b, 0)) = 0$ , where  $E_b = -\kappa^2/m$  with momentum  $k = i\kappa$  [3]. We consider the case where  $k_0$  and  $\Omega$  can be treated perturbatively. To the  $k_0^2$  and  $\Omega$  order, we find

$$\begin{pmatrix} g_S^{-1} - i\Pi_{SS}(E_b, 0) & \delta_{SP} - i\Pi_{SP}(E_b, 0) \\ \delta_{PS} - i\Pi_{PS}(E_b, 0) & g_P^{-1} - i\Pi_{PP}(E_b, 0) \end{pmatrix} = \begin{pmatrix} -\frac{ma_s}{2} + \frac{m}{2\sqrt{-mE_b}} & \delta_{SP} + \frac{\sqrt{mk_0\Omega}}{8(-E_b)^{3/2}} \\ \delta_{SP} + \frac{\sqrt{mk_0\Omega}}{8(-E_b)^{3/2}} & \frac{m - a_p m \sqrt{-mE_b + k_0^2}}{4a_p} \end{pmatrix}, \quad (\text{S29})$$

We assume  $a_s > 0$  and  $a_p > 0$ . Without the coupling  $\Omega$  and  $\delta_{sp}$ , the  $s$ - and  $p$ -wave dimers decouple. The binding energies are

$$E_b^{(s)} = -1/(ma_s^2), \quad E_b^{(p)} = k_0^2/m - 1/(ma_p^2). \quad (\text{S30})$$

When  $E_b^{(s)}$  and  $E_b^{(p)}$  are almost degenerate, we expect strong mixing between the  $s$ -wave and the  $p$ -wave dimer. As a result, we expect a small off-diagonal term is important only when

$$1/(a_s^0)^2 = 1/(a_p^0)^2 - k_0^2. \quad (\text{S31})$$

we thus approximate the off-diagonal terms as  $\delta_{SP} + \frac{k_0\Omega m^2 (a_s^0)^3}{8}$ . We further expand the diagonal terms near their own pole as

$$T^{-1}(E_b, 0) \sim \begin{pmatrix} \frac{m^2 a_s^3}{4} \left( E_b + \frac{1}{ma_s^2} \right) & \delta_{SP} + \frac{k_0\Omega m^2 (a_s^0)^3}{8} \\ \delta_{SP} + \frac{k_0\Omega m^2 (a_s^0)^3}{8} & \frac{m^2 a_p}{8} \left( E_b + \frac{1}{ma_p^2} - \frac{k_0^2}{m} \right) \end{pmatrix}. \quad (\text{S32})$$

Therefore, the pole of the  $T$ -matrix as  $\det(T^{-1}(E_b, 0)) = 0$  gives:

$$E_b^{(\pm)} = -\frac{1}{2a_p^2 a_s^2 m^2} \left\{ [a_s^2 + a_p^2 (1 - a_s^2 k_0^2)] m \pm \sqrt{a_s^4 m^2 + a_p^4 m^2 (1 + a_s^2 k_0^2)^2 - 2a_p^2 a_s^2 m^2 (1 + a_s^2 k_0^2) + 2a_p^3 a_s [8\delta_{SP} + (a_s^0)^3 k_0 \Omega m^2]^2} \right\}. \quad (\text{S33})$$

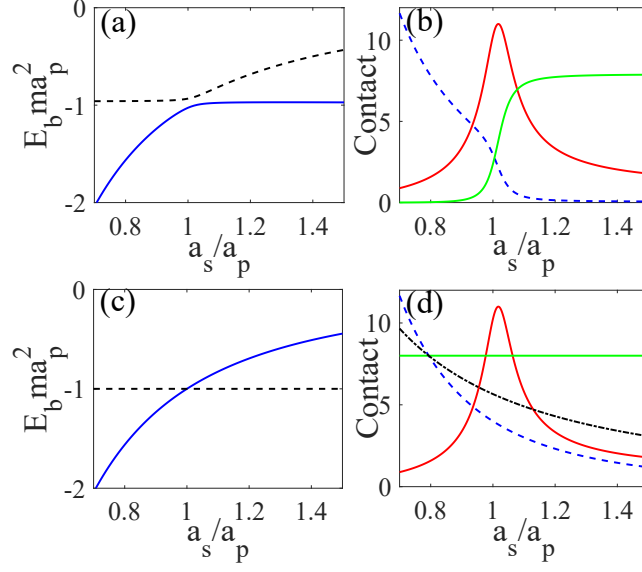


FIG. S3. (a) Dimensionless two-body banding energy versus  $a_s/a_p$  with SOC. The black dashed curve denotes  $E_b^{(+)}ma_p^2$  and the blue solid curve denotes  $E_b^{(-)}ma_p^2$ . (b) Dimensionless two-body contacts versus  $a_s/a_p$  with SOC. (c) Dimensionless two-body banding energy versus  $a_s/a_p$  without SOC. The black dashed curve denotes  $E_b^{(p)}ma_p^2$  and the blue solid curve denotes  $E_b^{(s)}ma_p^2$ . (d) Dimensionless two-body contacts versus  $a_s/a_p$  without SOC. As a comparison, we also plot the  $C_{SP}$  with finite SOC (the same curve as (b)). The red solid curve denotes  $C_{SP}a_p^2$ , the blue dashed curve denotes  $C_{SS}a_p^3$ , the green solid curve denotes  $C_{PP}a_p$ , and the black DotDashed curve denotes  $\sqrt{C_{SS}C_{PP}a_p^2}$ . Here, we choose the SOC parameters as  $k_0a_p = 0.2$  and  $m\Omega a_p^2 = 0.3$ .

The explicit formula for all contacts are given in the supplementary material. A plot for  $E_b^{(\pm)}$  and contacts for  $E_b^{(-)}$  are shown in Fig. S3 (a) and (c). Away from the degenerate point,  $E_b^{(\pm)}$  approaches  $E_b^{(s)}$  or  $E_b^{(p)}$ . Comparing Fig. S3 (a) with (c), it is found that the SOC parameters can open a gap between the two banding energies  $E_b^{(+)}$  and  $E_b^{(-)}$ .

Further, with the adiabatic relations (S13), (S14), we can get the two-body contacts as

$$C_{SS} \equiv 2m \frac{\partial E_b^{(-)}}{\partial a_s} = \frac{2}{a_s^3} + \frac{-2a_s^2 + 2a_p^2(1 + a_s^2 k_0^2) + 3a_p a_s (a_s^0)^6 k_0^2 m^2 \Omega^2}{a_s^2 \sqrt{a_s^2 \{ [a_s^2 - a_p^2(1 + a_s^2 k_0^2)]^2 + 2a_p^3 a_s (a_s^0)^6 k_0^2 m^2 \Omega^2 \}}}, \quad (\text{S34})$$

$$C_{PP} \equiv -4m \frac{\partial E_b^{(-)}}{\partial a_p^{-1}} = \frac{4}{a_p} + \frac{2\{a_s[2a_s^2 - 2a_p^2(1 + a_s^2 k_0^2)] + a_p^3 (a_s^0)^6 k_0^2 m^2 \Omega^2\}}{a_p a_s \sqrt{[a_s^2 - a_p^2(1 + a_s^2 k_0^2)]^2 + 2a_p^3 a_s (a_s^0)^6 k_0^2 m^2 \Omega^2}}, \quad (\text{S35})$$

$$C_{SP} = C_{PS} \equiv -2m^2 \frac{\partial E_b^{(-)}}{\partial \delta_{SP}} \Big|_{\delta_{SP}=0} = \frac{16a_p (a_s^0)^3 k_0 m \Omega}{a_s \sqrt{[a_s^2 - a_p^2(1 + a_s^2 k_0^2)]^2 + 2a_p^3 a_s (a_s^0)^6 k_0^2 m^2 \Omega^2}}, \quad (\text{S36})$$

where we use  $\delta_{SP} = 0$ ,  $a_s^0 = a_s$ , and  $E_b^{(-)}$  to calculate the contacts as shown in Fig. S3(b). Consequently, for the diagonal components of the contact matrix, we have  $C_{SS} \approx 0$  for  $a_s/a_p \gg 1$  and  $C_{PP} \approx 0$  for  $a_s/a_p \ll 1$ . Near the degenerate point  $a_s/a_p \sim 1$ , we see a peak for  $C_{SP}$ , indicating a large mixing between  $s$ - and  $p$ -wave dimers as expected.

Without SOC, one can have

$$C_{SS} \equiv 2m \frac{\partial E_b^{(s)}}{\partial a_s} = \frac{4}{a_s^3}, \quad (\text{S37})$$

$$C_{PP} \equiv -4m \frac{\partial E_b^{(p)}}{\partial a_p^{-1}} = \frac{8}{a_p}, \quad (\text{S38})$$

$$C_{SP} = C_{PS} = 0. \quad (\text{S39})$$

Moreover, we also calculate the amplitude of the hybridized new contacts compared to the  $s$ - and  $p$ -wave ones without SOC as shown in Fig. S3 (d) to give the possibility of the measurement.

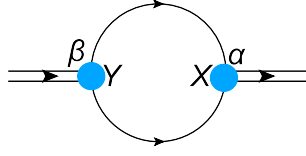


FIG. S4. Feynman diagram for the polarization bubble  $\Pi_s$ . Here,  $X = (x_1, t_1)$  and  $Y = (x_2, t_2)$ .

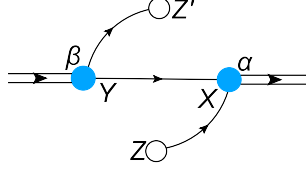


FIG. S5. Feynman diagram for the matrix of the operator  $\Psi^{\dagger T}(Z)\Psi^T(Z')$ . Here,  $X = (x_1, t_1)$ ,  $Y = (x_2, t_2)$ ,  $Z = (x_3, t_3)$ , and  $Z' = (x_4, t_4)$ .

#### SIV. DERIVATION OF SOME FORMULAS

##### A. Derivations of Eq. (S2)

The polarization bubble can be calculated by the diagram in Fig. S4 as

$$\begin{aligned}
\Pi_{\alpha\beta}(E_0, Q) &= \int dXdY \text{Tr} \left\langle \mathcal{T} \left[ \frac{1}{2} \Psi^T(X) \sigma_\alpha \Psi(X) \right] \left[ \frac{1}{2} \Psi^\dagger(Y) \sigma_\beta^\dagger \Psi^{\dagger T}(Y) \right] \right\rangle \\
&= \frac{1}{4} \int dXdY \text{Tr} \left[ \langle \mathcal{T} \overbrace{\Psi^T(X) \sigma_\alpha \Psi(X) \Psi^\dagger(Y) \sigma_\beta^\dagger \Psi^{\dagger T}(Y)} + \langle \mathcal{T} \overbrace{\Psi^T(X) \sigma_\alpha \Psi(X) \Psi^\dagger(Y) \sigma_\beta^\dagger \Psi^{\dagger T}(Y)} \rangle \right] \\
&= \frac{2}{4} \int dXdY \text{Tr} \langle \mathcal{T} \overbrace{\Psi^T(X) \sigma_\alpha \Psi(X) \Psi^\dagger(Y) \sigma_\beta^\dagger \Psi^{\dagger T}(Y)} \rangle \text{ (two equivalent contractions),} \\
&= \frac{1}{2} \int dXdY \text{Tr} \left\langle \mathcal{T} [\Psi^T(X)]_{1a} (\sigma_\alpha)_{ab} [\Psi(X)]_{b1} [\Psi^\dagger(Y)]_{1m} (\sigma_\beta^\dagger)_{mn} [\Psi^{\dagger T}(Y)]_{n1} \right\rangle \\
&= \frac{1}{2} \int dXdY \text{Tr} \left\langle \mathcal{T} [\Psi^T(X)]_{1a} [\Psi^{\dagger T}(Y)]_{n1} (\sigma_\alpha)_{ab} [\Psi(X)]_{b1} [\Psi^\dagger(Y)]_{1m} (\sigma_\beta^\dagger)_{mn} \right\rangle \\
&= \frac{1}{2} \int dXdY \text{Tr} \langle \mathcal{T} [\Psi(X)]_{a1} [\Psi^\dagger(Y)]_{1n} (\sigma_\alpha)_{ab} [G(X-Y)]_{bm} (\sigma_\beta^\dagger)_{mn} \rangle \\
&= \frac{1}{2} \int dXdY \text{Tr} [G(X-Y)]_{an} (\sigma_\alpha)_{ab} [G(X-Y)]_{bm} (\sigma_\beta^\dagger)_{mn} \\
&= \frac{1}{2} \int dXdY \text{Tr} [G^T(X-Y)]_{na} (\sigma_\alpha)_{ab} [G(X-Y)]_{bm} (\sigma_\beta^\dagger)_{mn} \\
&= \int dXdY \frac{1}{2} \text{Tr} \left[ G^T(X-Y) \sigma_\alpha G(X-Y) \sigma_\beta^\dagger \right], \tag{S40}
\end{aligned}$$

where  $\int dXdY = \frac{1}{2} \int d(X+Y)d(X-Y)$ ,  $\Psi = (\psi_\uparrow, \psi_\downarrow)^T$ ,  $X = (x_1, t_1)$ ,  $Y = (x_2, t_2)$ , and we use the definition  $G(X-Y) \equiv \langle \mathcal{T} \Psi(X) \Psi^\dagger(Y) \rangle$ .

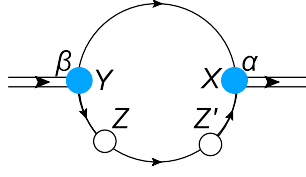


FIG. S6. (Color online) Feynman diagram for the matrix element of  $\mathcal{O}_{\sigma_3}(R+x, t)\mathcal{O}_{\sigma'_3}^\dagger(R, 0)$ .

### B. Derivations of Eq. (14)

As shown in Fig. S5, the expectation of the operator  $\Psi^{\dagger T}(Z)\Psi^T(Z')$  is calculated as follows:

$$\begin{aligned}
\langle O_{\alpha_o} | \Psi^{\dagger T}(Z)\Psi^T(Z') | I_{\alpha_i} \rangle &= \left( \begin{array}{l} \langle O_{\alpha_o} | \psi_\uparrow^\dagger(Z)\psi_\uparrow(Z') | I_{\alpha_i} \rangle \\ \langle O_{\alpha_o} | \psi_\downarrow^\dagger(Z)\psi_\downarrow(Z') | I_{\alpha_i} \rangle \end{array} \right) \\
&= (-i)^2 D_{\alpha_o\alpha}(E_0, Q) D_{\beta\alpha_i}(E_0, Q) \int dXdY \left\langle \mathcal{T} \left[ \frac{1}{2} \Psi^T(X) \sigma_\alpha \Psi(X) \right] \Psi^{\dagger T}(Z) \Psi^T(Z') \left[ \frac{1}{2} \Psi^\dagger(Y) \sigma_\beta^\dagger \Psi^\dagger(Y) \right] \right\rangle \\
&= -\frac{D_{\alpha_o\alpha}(E_0, Q) D_{\beta\alpha_i}(E_0, Q)}{4} \int dXdY \\
&\quad \left[ \langle \mathcal{T} \overbrace{\Psi^T(X) \sigma_\alpha \Psi(X)} \overbrace{\Psi^{\dagger T}(Z) \Psi^T(Z')} \overbrace{\Psi^\dagger(Y) \sigma_\beta^\dagger \Psi^\dagger(Y)} \rangle + \langle \mathcal{T} \overbrace{\Psi^T(X) \sigma_\alpha \Psi(X)} \overbrace{\Psi^{\dagger T}(Z) \Psi^T(Z')} \overbrace{\Psi^\dagger(Y) \sigma_\beta^\dagger \Psi^\dagger(Y)} \rangle \right. \\
&\quad \left. + \langle \mathcal{T} \overbrace{\Psi^T(X) \sigma_\alpha \Psi(X)} \overbrace{\Psi^{\dagger T}(Z) \Psi^T(Z')} \overbrace{\Psi^\dagger(Y) \sigma_\beta^\dagger \Psi^\dagger(Y)} \rangle + \langle \mathcal{T} \overbrace{\Psi^T(X) \sigma_\alpha \Psi(X)} \overbrace{\Psi^{\dagger T}(Z) \Psi^T(Z')} \overbrace{\Psi^\dagger(Y) \sigma_\beta^\dagger \Psi^\dagger(Y)} \rangle \right] \\
&\sim \int dXdY \langle \mathcal{T} \overbrace{\Psi^T(X) \sigma_\alpha \Psi(X)} \overbrace{\Psi^{\dagger T}(Z) \Psi^T(Z')} \overbrace{\Psi^\dagger(Y) \sigma_\beta^\dagger \Psi^\dagger(Y)} \rangle \text{ (four equivalent contractions)}, \\
&= \int dXdY \langle \mathcal{T} [\Psi^T(X)]_{1n} (\sigma_\alpha)_{nm} [\Psi(X)]_{m1} [\Psi^{\dagger T}(Z)]_{j1} [\Psi^T(Z')]_{1j} [\Psi^\dagger(Y)]_{1a} (\sigma_\beta^\dagger)_{ab} [\Psi^{\dagger T}(Y)]_{b1} \rangle \\
&= \int dXdY \langle \mathcal{T} [\Psi(X)]_{n1} [\Psi^\dagger(Z)]_{1j} (\sigma_\alpha)_{nm} [\Psi(X)]_{m1} [\Psi^\dagger(Y)]_{1a} (\sigma_\beta^\dagger)_{ab} [\Psi(Z')]_{j1} [\Psi^\dagger(Y)]_{1b} \rangle \\
&= \int dXdY [G(X-Z)]_{nj} (\sigma_\alpha)_{nm} [G(X-Y)]_{ma} (\sigma_\beta^\dagger)_{ab} [G(Z'-Y)]_{jb} \\
&= \int dXdY [G^T(X-Z)]_{jn} (\sigma_\alpha)_{nm} [G(X-Y)]_{ma} (\sigma_\beta^\dagger)_{ab} [G^T(Z'-Y)]_{bj} \\
&= \int dXdY G^T(X-Z) \sigma_\alpha G(X-Y) \sigma_\beta^\dagger G^T(Z'-Y), \tag{S41}
\end{aligned}$$

where  $\int dXdY = \frac{1}{2} \int d(X+Y)d(X-Y)$ , we label the field operator  $\Psi = (\psi_\uparrow, \psi_\downarrow)^T$ ,  $X = (x_1, t_1)$ ,  $Y = (x_2, t_2)$ ,  $Z = (x_3, t_3)$ ,  $Z' = (x_4, t_4)$ , and we use the definition  $G(X-Y) \equiv \langle \mathcal{T} \Psi(X) \Psi^\dagger(Y) \rangle$ .

Therefore, we have

$$\begin{aligned}
&\left( \begin{array}{l} \langle O_{\alpha_o} | \psi_\uparrow^\dagger(Z)\psi_\uparrow(Z') | I_{\alpha_i} \rangle \\ \langle O_{\alpha_o} | \psi_\downarrow^\dagger(Z)\psi_\downarrow(Z') | I_{\alpha_i} \rangle \end{array} \right)^T = [\langle O_s | \Psi^{\dagger T}(Z)\Psi^T(Z') | I_{\alpha_i} \rangle]^T \\
&= -D_{\alpha_o\alpha}(E_0, Q) D_{\beta\alpha_i}(E_0, Q) \int dXdY G(Z'-Y) \sigma_\beta G^T(X-Y) \sigma_\alpha^\dagger G(X-Z), \tag{S42}
\end{aligned}$$

where we use  $(ABCDE)^T = E^T D^T C^T B^T A^T$ .



### C. Derivations of Eq. (18)

Equation (18) in the main text can be calculated by the diagram in Fig. S6 as

$$\begin{aligned}
& \langle O_{\alpha_o} | \mathcal{T} \mathcal{O}_{\sigma_3}(R+x, t) \mathcal{O}_{\sigma'_3}^\dagger(R, 0) | I_{\alpha_i} \rangle \\
& \sim (-i)^2 D_{\alpha_o \alpha}(E_0, Q) D_{\beta \alpha_i}(E_0, Q) \int dX dY \left\langle \mathcal{T} \left[ \frac{1}{2} \Psi^T(X) \sigma_\alpha \Psi(X) \right] [\Psi^\dagger(Z) \Psi(Z)] [\Psi^\dagger(Z') \Psi(Z')] \left[ \frac{1}{2} \Psi^\dagger(Y) \sigma_\beta^\dagger \Psi^{\dagger T}(Y) \right] \right\rangle \\
& \sim \frac{1}{4} \int dX dY \left[ \langle \mathcal{T} \overbrace{\Psi^T(X) \sigma_\alpha \Psi(X) \Psi^\dagger(Z) \Psi(Z) \Psi^\dagger(Z') \Psi(Z') \Psi^\dagger(Y) \sigma_\beta^\dagger \Psi^{\dagger T}(Y)} \rangle \right. \\
& \quad + \langle \mathcal{T} \overbrace{\Psi^T(X) \sigma_\alpha \Psi(X) \Psi^\dagger(Z) \Psi(Z) \Psi^\dagger(Z') \Psi(Z') \Psi^\dagger(Y) \sigma_\beta^\dagger \Psi^{\dagger T}(Y)} \rangle \\
& \quad + \langle \mathcal{T} \overbrace{\Psi^T(X) \sigma_\alpha \Psi(X) \Psi^\dagger(Z) \Psi(Z) \Psi^\dagger(Z') \Psi(Z') \Psi^\dagger(Y) \sigma_\beta^\dagger \Psi^{\dagger T}(Y)} \rangle \\
& \quad \left. + \langle \mathcal{T} \overbrace{\Psi^T(X) \sigma_\alpha \Psi(X) \Psi^\dagger(Z) \Psi(Z) \Psi^\dagger(Z') \Psi(Z') \Psi^\dagger(Y) \sigma_\beta^\dagger \Psi^{\dagger T}(Y)} \rangle \right] \\
& = \int dX dY \langle \mathcal{T} \overbrace{\Psi^T(X) \sigma_\alpha \Psi(X) \Psi^\dagger(Z) \Psi(Z) \Psi^\dagger(Z') \Psi(Z') \Psi^\dagger(Y) \sigma_\beta^\dagger \Psi^{\dagger T}(Y)} \rangle \text{ (four equivalent contractions),} \\
& = \int dX dY \langle \mathcal{T} [\Psi^T(X)]_{1a} (\sigma_\alpha)_{ab} [\Psi(X)]_{b1} [\Psi^\dagger(Z)]_{1c} [\Psi(Z)]_{c1} [\Psi^\dagger(Z')]_{1d} [\Psi(Z')]_{d1} [\Psi^\dagger(Y)]_{1m} (\sigma_\beta^\dagger)_{mn} [\Psi^{\dagger T}(Y)]_{n1} \rangle \\
& = \int dX dY \langle \mathcal{T} [\Psi(X)]_{a1} (\sigma_\alpha)_{ab} [G(X-Z)]_{bc} [G(Z-Z')]_{cd} [G(Z'-Y)]_{dm} (\sigma_\beta^\dagger)_{mn} [\Psi^\dagger(Y)]_{1n} \rangle \\
& = \int dX dY \langle \mathcal{T} [\Psi(X)]_{a1} [\Psi^\dagger(Y)]_{1n} \rangle (\sigma_\alpha)_{ab} [G(X-Z)]_{bc} [G(Z-Z')]_{cd} [G(Z'-Y)]_{dm} (\sigma_\beta^\dagger)_{mn} \\
& = \int dX dY [G(X-Y)]_{an} (\sigma_\alpha)_{ab} [G(X-Z)]_{bc} [G(Z-Z')]_{cd} [G(Z'-Y)]_{dm} (\sigma_\beta^\dagger)_{mn} \\
& = \int dX dY [G^T(X-Y)]_{na} (\sigma_\alpha)_{ab} [G(X-Z)]_{bc} [G(Z-Z')]_{cd} [G(Z'-Y)]_{dm} (\sigma_\beta^\dagger)_{mn} \\
& = \int dX dY [G^T(X-Y) \sigma_\alpha G(X-Z) G(Z-Z') G(Z'-Y) \sigma_\beta^\dagger] \\
& \sim G_0(Z-Z') \int dX dY [G^T(X-Y) \sigma_\alpha G(X-Z) G(Z'-Y) \sigma_\beta^\dagger] \\
& = G_0(Z-Z') \int dX dY [G(Z'-Y) \sigma_\beta^\dagger G^T(X-Y) \sigma_\alpha G(X-Z)]. \tag{S43}
\end{aligned}$$

### D. $T$ matrix for $s$ wave without SOC

As shown in Fig. S7(a), the two-body  $T$  matrix for  $s$  wave is given by [7]

$$-iT_s(E_0, Q) = \frac{-ig_S}{1 - (-ig_S)\Pi_s(E_0, Q)}. \tag{S44}$$

As shown in Fig. S7(c), the dimer propagator matrix for  $s$  wave is given by

$$D_{SS}(E_0, Q) = \frac{ig_S}{1 - (ig_S)\Pi_{SS}(E_0, Q)}, \tag{S45}$$

With  $D_{SS}(E_0, Q) = iT_s(E_0, Q)$ , one can get  $\Pi_{SS}(E_0, Q) = -\Pi_s(E_0, Q)$ .

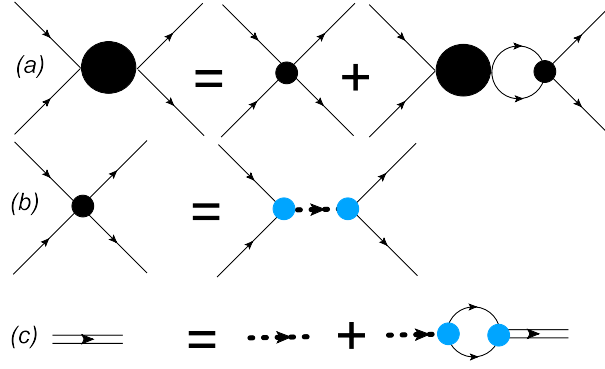


FIG. S7. (a) Diagrams for the  $T$  matrix of  $s$ -wave without SOC. (b) Diagrams for the interaction vertex of  $s$ -wave. (c) Diagrams for the dimer propagator of  $s$ -wave. The black disk represents the  $s$ -wave  $T$  matrix:  $-iT_s$ , the single line denotes the bare atom propagator  $G$ , the black dot represents the interaction vertex:  $-ig_S$ , the black dotted line denotes the bare dimer propagator  $ig_S$ , the double lines denote the dimer propagator  $D$ , and the blue dot represents the interaction vertex:  $-i$ .

### E. $T$ matrix for $p$ wave without SOC

The two-body  $T$  matrix for  $p$  wave is given by [7]

$$\frac{-iT_p(E_0, Q)}{qq'} = \frac{-ig_P}{1 - (-ig_P)\Pi_p(E_0, Q)}. \quad (\text{S46})$$

The dimer propagator matrix for  $p$  wave is given by

$$D_{PP}(E_0, Q) = \frac{ig_P}{1 - (ig_P)\Pi_{PP}(E_0, Q)}, \quad (\text{S47})$$

With  $D_{PP}(E_0, Q) = iT_p(E_0, Q)/(qq')$ , one can get  $\Pi_{PP}(E_0, Q) = -\Pi_p(E_0, Q)$ .

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