Electron-K-phonon interaction in twisted bilayer graphene

Chao-Xing Liu^{1,2,*} Yulin Chen,³ Ali Yazdani,² and B. Andrei Bernevig^{2,4,5,*}

¹Department of Physics, The Pennsylvania State University, University Park, Pennsylvania 16802, USA

²Department of Physics, Princeton University, Princeton, New Jersey 08544, USA

³Department of Physics, University of Oxford, Oxford OX1 3PU, United Kingdom

⁴Donostia International Physics Center, P. Manuel de Lardizabal 4, 20018 Donostia-San Sebastian, Spain ⁵IKERBASQUE, Basque Foundation for Science, Plaza Euskadi 5, 48009, Bilbao, Spain

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We develop an analytic theory to describe the interaction between electrons and *K* phonons and study its influence on superconductivity in the *bare bands* of twisted bilayer graphene (TBG). We find that, due to symmetry and the two-center approximation, only one optical *K* phonon (~160 meV) of graphene is responsible for the intervalley electron-phonon interaction. This phonon has recently been found in angular-resolved photoemission spectroscopy to be responsible for replicas of the TBG flat bands. By projecting the interaction to the TBG flat bands, we perform the full symmetry analysis of the phonon-mediated attractive interaction and pairing channels in the Chern basis, and show that several channels are guaranteed to have gapless order parameters. From the linearized gap equations, we find that the highest T_c pairing induced by this phonon is a singlet gapped *s*-wave inter-Chern-band order parameter, followed closely by a gapless nematic *d*-wave intra-Chern-band order parameter. We justify these results analytically, using the topological heavy-fermion mapping of TBG which has allowed us to obtain an analytic form of a phonon-mediated attractive interaction and to analytically solve the linearized and T = 0 gap equations. For the intra-Chern-band Coulomb interaction can be screened sufficiently enough—around the chiral flat-band limit. While the flat-band Coulomb interaction can be screened sufficiently enough—around the Van Hove singularities—to allow for electron-phonon based superconductivity, it is unlikely that this effect can be maintained in the lower density of states excitation bands around the correlated insulator states.

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I. INTRODUCTION

Superconductivity in twisted bilayer graphene (TBG) appears within its phase diagram around the correlated insulator states [1–19]. Among the mechanisms suggested for superconductivity are phonons, spin fluctuations, skyrmions, and others [20–45]. Based on a recent experiment that suggests a strong coupling between the graphene K phonon and the flat bands in TBG [46], we perform a comprehensive analysis of the electron-K-phonon (e-K-ph) interaction and the resulting phonon-mediated superconductivity on the bare flat bands of TBG. We develop an exhaustive numerical, analytical, and symmetry based description of the e-K-ph interaction in TBG and the symmetry classifications of the order parameter, and find the competing singlet gapped inter-Chern-band channel and nematic gapless intra-Chern-band channel. Armed with the heavy-fermion description of TBG [47–55], the form factors of the K-phonon induced attractive interaction can be analytically computed and matched well to full numerical calculations. An analysis of the Coulomb screening shows that, due to the high density of states (DOS) of flat bands, the Coulomb interaction might be strongly renormalized down near the Van Hove singularities. However, it remains unclear if the Hartree-Fock bands of the correlated insulator, with the lower DOS, can provide a similar result.

II. MODEL HAMILTONIAN FOR ELECTRON-PHONON INTERACTION IN TBG

We consider the deformation potential type of theory, described by a tight-binding (TB) model for the electron Hamiltonian with the hopping parameters depending on the atom positions $\tilde{\mathbf{R}}_{\alpha}^{l} = \mathbf{R}^{l} + \tau_{\alpha}^{l} + \mathbf{u}^{l}(\mathbf{R}_{\alpha}^{l})$ with a displacement field $\mathbf{u}^{l}(\mathbf{R}_{\alpha}^{l} = \mathbf{R}^{l} + \tau_{\alpha}^{l})$, where \mathbf{R}^{l} and τ_{α}^{l} label the lattice vector and the sublattice atom position ($\alpha = A, B$) at the layer *l*, respectively. By treating **u** as a perturbation, we expand the intralayer Hamiltonian up to the linear order in u [Supplemental Material (SM) Sec. II [56]]. We only keep u-independent terms for the interlayer Hamiltonian for TBG, thus focusing on the intralayer electron-phonon (e-ph) interaction in this paper. As only the Dirac bands appear around the Fermi energy close to $\pm \mathbf{K}_D = \pm \frac{4\pi}{3a_0}(1,0)$ in the Brillouin zone (BZ) with the lattice constant a_0 in graphene, we also expand the Hamiltonian around $\eta \mathbf{K}_D$ ($\eta = \pm$ labeling two valleys) and focus on Dirac electrons around two valleys. Our full Hamiltonian consists of three parts,

$$H = H_{\rm el} + H_{\rm ph} + H_{\rm eph}.$$
 (1)

Here, $H_{\rm el}$ describes the Dirac electrons located at valley $\eta = \pm$ momentum $\eta \mathbf{K}_D$ that are coupled through interlayer tunnelings and is given by the Bistritzer-MacDonald (BM) model

^{*}These authors contributed equally to this work.



FIG. 1. (a) Phonon dispersion of graphene. The irreps for phonon modes at Γ and \mathbf{K}_D are labeled. Inset: BZ of graphene. (b) MBZ of TBG. (c) and (d) show the momentum dependence of the normalized gap function $|\Delta_{\mathbf{k}}|$ for the inter-Chern-band A_1 singlet (or the A_2 triplet) channel and intra-Chern-band pairing 2D E_2 singlet channel, respectively. The inset in (c) shows the $|\Delta_{\mathbf{k}}|$ along the dashed line $k_y = 0$ for both the inter-Chern-band (red) and intra-Chern-band (blue) channels. The momenta Γ_M , \mathbf{K}_M , \mathbf{M}_M are labeled in MBZ in (b) and (d).

[57] (SM Sec. V [56]),

$$\hat{H}_{\rm el} = \sum_{\eta s} \sum_{\mathbf{k} \in \rm MBZ} \sum_{\alpha \alpha'} \sum_{\mathbf{Q}, \mathbf{Q}'} h_{\mathbf{Q}\alpha, \mathbf{Q}'\alpha'}^{(\eta)}(\mathbf{k}) c_{\mathbf{k}, \mathbf{Q}, \alpha, \eta, s}^{\dagger} c_{\mathbf{k}, \mathbf{Q}', \alpha', \eta, s}, \quad (2)$$

where $c_{\mathbf{k},\mathbf{Q},\alpha,\eta,s}$ is the fermion annihilation operator, **k** is a momentum in the moiré Brillouin zone (MBZ) [Fig. 1(b)], α is the sublattice index, and s is spin. The vector \mathbf{Q} belongs to the lattice set $Q_{l\eta} = \{ l\eta \mathbf{q}_2 + n_1 \mathbf{b}_{M1} + n_2 \mathbf{b}_{M2} | n_{1,2} \in \mathbb{Z} \},\$ where *l* is the layer index, $\mathbf{q}_2 = k_{\theta}(\frac{\sqrt{3}}{2}, \frac{1}{2}), \mathbf{b}_{M1} = k_{\theta}(\frac{\sqrt{3}}{2}, \frac{3}{2}),$ $\mathbf{b}_{M2} = k_{\theta}(-\frac{\sqrt{3}}{2}, \frac{3}{2})$, and $k_{\theta} = 2|\mathbf{K}_D|\sin\frac{\theta}{2}$ with θ the twist angle. $h_{\mathbf{O}\alpha,\mathbf{O}'\alpha'}^{(\eta)}(\mathbf{k})$ is given in SM Sec. V A [56]. \hat{H}_{el} exhibits C_{6v} and time-reversal symmetries, generated by valley-switching $\pi/6$ rotation along the z axis (\hat{C}_{6z}), time reversal (\hat{T}), and π rotation along the y axis (C_{2y}), and valley-preserving π rotation along the x axis (\hat{C}_{2x}), and the composite antiunitary $C_{2z}T$. In addition, \hat{H}_{el} has a unitary particle-hole (\hat{P}) symmetry, as well as a chiral symmetry \hat{C} in the limit with vanishing AA region hopping $(w_0 = 0)$ [19]. A full discussion of symmetry of the BM model [58,59] is found in SM Sec.V B [56] (see also Refs. [60–63] therein).

 $H_{\rm ph}$ describes the intralayer in-plane phonon modes. Outof-plane phonon modes are decoupled from Dirac electrons for the intralayer e-ph interaction. The dynamical matrix for a single-layer graphene is derived in SM Sec. III [56] based both on symmetry considerations and the microscopic model, up to the next-nearest-neighbor interaction. The resulting inplane phonon dispersion in Fig. 1(a) reproduces that in the literature [64–68] (SM Secs. III and IV B [56]). The phonon modes at Γ and ηK_D can induce intravalley and intervalley e-ph interactions, respectively. In this paper we focus on the $\eta \mathbf{K}_D$ phonons. At $\eta \mathbf{K}_D$, we have one A_1 (~160 meV), one A_2 (~140 meV), and one two-dimensional (2D) E mode (~150 meV) of the C_{3v} group. Based on the deformation potential theory, we derive the e-ph interaction H_{eph} by expanding the TB Hamiltonian treating both the momentum and phonon displacement field \mathbf{u} as perturbations. For the e-ph interaction, we only keep the dominant zeroth order in momentum for the $\eta \mathbf{K}_D$ phonons. We find, due to both symmetry and the two-center approximation (SM Sec. II E [56]), that only the A_1 phonons at \mathbf{K}_D can scatter an electron from \mathbf{K}_D to $-\mathbf{K}_D$ [68]. The corresponding Hamiltonian reads

$$H_{\text{intervall}}^{\text{op},A_1} \approx \frac{\gamma_3}{\sqrt{2N_G M \omega_{A1}}} \sum_{\tilde{\mathbf{k}},\tilde{\mathbf{k}}',\eta,\alpha\beta} (b_{-\eta \mathbf{K}_D+\tilde{\mathbf{k}}-\tilde{\mathbf{k}}',A_1} + b_{\eta \mathbf{K}_D-\tilde{\mathbf{k}}+\tilde{\mathbf{k}}',A_1}^{\dagger}) c_{\tilde{\mathbf{k}}+\eta \mathbf{K}_D,\alpha}^{\dagger} (\sigma_x)_{\alpha\beta} c_{\tilde{\mathbf{k}}'-\eta \mathbf{K}_D,\beta}, \quad (3)$$

where \mathbf{k} is the electron momentum away from $\eta \mathbf{K}_D$, N_G is the number of atomic unit cells, M is the atomic mass, ω_{A1} is the A_1 phonon frequency, and b and c are phonon and electron annihilation operators. The material-dependent parameter γ_3 can be derived from the hopping potential as $\gamma_3 = 2i \sum_{\mathbf{G}} e^{i(\tau_A - \tau_B) \cdot \mathbf{G}} (\mathbf{G} + \mathbf{K}_D)_y t(\mathbf{G} + \mathbf{K}_D, 0) \approx 17 \text{ eV/ Å}$, where \mathbf{G} is the reciprocal lattice vector and $t(\mathbf{q})$ is the Fourier transform of the π -bond hopping function between two carbon p_z orbitals in graphene [Eq. (6) in SM Sec. I [56]]. Our next step is to rewrite the electron momentum \tilde{k} into the MBZ by $\tilde{\mathbf{k}} = \mathbf{k} - \mathbf{Q}_{l\eta}$ with $\mathbf{k} \in \text{MBZ}$, so that $c_{\mathbf{k},\mathbf{Q}_{l\eta},\alpha,\eta,s} = c_{\eta \mathbf{K}_D' + \mathbf{\hat{k}},\alpha,l,s}$ and $\sum_{\mathbf{\hat{k}}} \rightarrow \sum_{\mathbf{k} \in \text{MBZ}} \sum_{\mathbf{Q}_{l\eta}}$, where we have added the spin index *s* and layer index *l*. Finally, we project the e-ph interaction $H_{\text{intervall}}^{\text{op},A_1}$ into the flat bands of the BM Hamiltonian as

$$H_{\text{intervall}}^{\text{op},A_1} \approx \frac{1}{\sqrt{N_G}} \sum G_{\mathbf{k},\mathbf{k}',\mathbf{Q}_{-l\eta}}^{\eta n n'l} \gamma_{\mathbf{k},n,\eta,s}^{\dagger} \gamma_{\mathbf{k}',n',-\eta,s} \times (b_{-\eta\mathbf{K}_D+\mathbf{k}-\mathbf{k}'-\mathbf{Q}_{-l\eta},l,A_1} + b_{\eta\mathbf{K}_D-\mathbf{k}+\mathbf{k}'+\mathbf{Q}_{-l\eta},l,A_1}^{\dagger}),$$
(4)

where the summation includes \mathbf{k} , \mathbf{k}' , n, n', η , s, l, $\mathbf{Q}_{-l\eta}$, $\gamma_{\mathbf{k},n,\eta,s}^{\dagger} = \sum_{\mathbf{Q}\alpha} u_{\mathbf{Q}\alpha;n\eta}(\mathbf{k}) c_{\mathbf{k},\mathbf{Q},\eta,\alpha s}^{\dagger}$ with $u_{\mathbf{k},\mathbf{Q}_{l\eta},\alpha,\eta}^{n}$ the eigenstates of $h_{\mathbf{O}\alpha,\mathbf{O}\alpha'}^{(\eta)}(\mathbf{k})$. The matrix element

$$G_{\mathbf{k},\mathbf{k}',\mathbf{Q}_{-i\eta}}^{\eta nn'l} = \frac{\gamma_3}{\sqrt{2M\omega_{A_1}}} \sum_{\mathbf{Q}'_{i\eta},\alpha\beta} \times u_{\mathbf{k},\mathbf{Q}'_{i\eta},\alpha,\eta}^{n\star} \sigma_{\alpha\beta}^{x} u_{\mathbf{k}',\mathbf{Q}'_{i\eta}-\mathbf{Q}_{-i\eta},\beta,-\eta}^{n\star}$$
(5)

characterizes the e-ph interaction strength for TBG and can be evaluated numerically (and later analytically), as shown in SM Sec. VI F [56]. We focus on two flat bands (per valley per spin) of TBG, labeled by $n = \pm$. Instead of the eigenstate basis, we work on the so-called "Chern-band" basis, defined by

$$u_{\mathbf{k},\mathbf{Q},\alpha,\eta}^{e_{Y}} = \frac{1}{\sqrt{2}} (u_{\mathbf{k},\mathbf{Q},\alpha,\eta}^{n=+} + ie_{Y} u_{\mathbf{k},\mathbf{Q},\alpha,\eta}^{n=-}), \tag{6}$$

with $e_Y = \pm 1$. $u_{\mathbf{k},\mathbf{Q},\alpha,\eta}^{e_Y}$ carries the Chern number ± 1 . On the Chern-band basis, the expressions for the e-ph interaction can be obtained by replacing the *n*, *n'* indices in Eqs. (4) and

(5) with e_Y, e_Y' indices and $u_{\mathbf{k},\mathbf{Q},\alpha,\eta}^n$ in Eq. (5) with $u_{\mathbf{k},\mathbf{Q},\alpha,\eta}^{e_Y}$. Discrete symmetries can constrain the form of the function $G_{\mathbf{k},\mathbf{k}',\mathbf{Q}-l\eta}^{\eta e_Y e_Y' l}$, as discussed in SM Sec.VI D [56]. In particular, in the chiral limit $w_0 = 0$ one can show that $G_{\mathbf{k},\mathbf{k}',\mathbf{Q}-l\eta}^{\eta e_Y e_Y' l} = \delta_{e_Y,e_Y'} G_{\mathbf{k},\mathbf{k}',\mathbf{Q}-l\eta}^{\eta e_Y e_Y l}$ has a diagonal form on the Chern-band basis, and this approximation will be adopted below for the discussion of possible superconducting channels.

III. PHONON-MEDIATED ELECTRON-ELECTRON INTERACTION AND SYMMETRY CLASSIFICATION OF SUPERCONDUCTING PAIRING CHANNELS

We next apply the Schrieffer-Wolff transformation [69] to integrate out the phonon modes and obtain the phononmediated electron-electron (el-el) interaction [25,32]. We focus on the Cooper pair channel of the attractive interaction, which takes the form

$$H_{ee} = -\frac{1}{N_M} \sum_{\mathbf{k}, \mathbf{k}', s, s_1, e_Y, e_Y'} V_{\mathbf{k}, \mathbf{k}'}^{\eta, e_Y, e_Y'}$$
$$\times \gamma_{\mathbf{k} e_Y \eta s}^{\dagger} \gamma_{-\mathbf{k} e_Y', -\eta s_1}^{\dagger} \gamma_{-\mathbf{k}' e_Y', \eta s_1} \gamma_{\mathbf{k}' e_Y, -\eta s}, \qquad (7)$$

where

$$V_{\mathbf{k},\mathbf{k}'}^{\eta,e_{Y},e_{Y}'} = \frac{1}{N_{0}\omega_{A_{1}}} \sum_{\mathbf{G}_{M},l} G_{\mathbf{k},\mathbf{k}',-l\eta\mathbf{q}_{2}+\mathbf{G}_{M}}^{\eta,e_{Y},l} G_{-\mathbf{k},-\mathbf{k}',l\eta\mathbf{q}_{2}-\mathbf{G}_{M}}^{-\eta,e_{Y}',l},$$

with G_M the moiré reciprocal lattice vectors, N_M the number of moiré unit cells, and N_0 the number of atomic unit cells in one moiré unit cell $(N_G = N_0 \times N_M)$. Discrete symmetries constrain the form of the interaction parameter $V_{\mathbf{k},\mathbf{k}'}^{\eta,e_Y,e_Y'}$. The ones leaving the momentum $(\mathbf{k}, \mathbf{k}')$ unchanged are as follows: (1) $\hat{C}_{2z}\hat{P}$: $V_{\mathbf{k},\mathbf{k}'}^{\eta,e_Y,e'_Y} = V_{\mathbf{k},\mathbf{k}'}^{-\eta,e_Y,e'_Y}$; (2) $\hat{C}_{2z}\hat{T}$: $V_{\mathbf{k},\mathbf{k}'}^{\eta,e_Y,e'_Y} =$ $V_{\mathbf{k},\mathbf{k}'}^{\eta,-e_{\gamma},-e_{\gamma}'\star}$; and (3) the combination of index reshuffling and \hat{P} symmetry: $V_{\mathbf{k},\mathbf{k}'}^{\eta,e_Ye_Y} = V_{\mathbf{k},\mathbf{k}'}^{-\eta,e_Y'e_Y}$. These three symmetry operations reduce the number of the independent components of the V function for a fixed $(\mathbf{k}, \mathbf{k}')$ from eight complex parameters to one real $(V_{\mathbf{k},\mathbf{k}'}^{+,+-})$ and one complex parameter $(V_{\mathbf{k},\mathbf{k}'}^{+,++})$. Other discrete symmetries, including \hat{P} , reshuffling, Hermiticity, \hat{C}_{3z} , and \hat{C}_{2z} , relate the V function at different (**k**, **k**'). In particular, \hat{C}_{3z} guarantees $V_{\mathbf{K}_{M},0}^{\eta,e_{Y},e_{Y}} = 0$ for the intra-Chernband channels. The projected Coulomb interaction into the flat bands of the BM model possesses a large $U(4) \times U(4)$ spinvalley continuous symmetry [18,58,70]. The el-el interaction (7) breaks this symmetry down to the $U(2)_{e_Y=+} \times U(2)_{e_Y=-}$ in the chiral limit and further to a total spin SU(2) together with a valley charge $U(1) \otimes U(1)$ (SM Sec. VI E [56]).

At the mean-field level, the attractive interaction (7) is decomposed into the fermion bilinear form $H_{\Delta} = \hat{\Delta} + \hat{\Delta}^{\dagger}$ with

$$\hat{\Delta} = \sum \gamma_{\mathbf{k}, e_{Y_1}, \eta, s_1}^{\dagger} \Delta_{\mathbf{k}; e_{Y_1} s_1, e_{Y_2} s_2}^{\eta} \gamma_{-\mathbf{k}, e_{Y_2}, -\eta, s_2}^{\dagger}, \qquad (8)$$

where the summation above includes the indices **k**, e_{Y_1} , e_{Y_2} , s_1 , s_2 , η , and the gap function

$$\Delta^{\eta}_{\mathbf{k};e_{Y_{1}}s_{1},e_{Y_{2}}s_{2}} = -\frac{1}{N_{M}} \sum_{\mathbf{k}'} V^{\eta e_{Y_{1}}e_{Y_{2}}}_{\mathbf{k}\mathbf{k}'} \langle \gamma_{-\mathbf{k}'e_{Y_{2}}\eta s_{2}} \gamma_{\mathbf{k}'e_{Y_{1}}-\eta s_{1}} \rangle.$$
(9)

Since the interaction V function does not involve spin, we can decompose $\Delta^{\eta}_{\mathbf{k};e_{Y_1}s_1,e_{Y_2}s_2} = \sum_{S,M} \Delta^{\eta,SM}_{\mathbf{k};e_{Y_1}e_{Y_2}} S^{SM}_{s_1s_2}$, where S = 0 for the spin singlet and S = 1 ($M = -S, \ldots, S$) for the spin triplet (SM Sec. VI G 1 [56]).

The gap function can be classified according to the discrete symmetries. The C_{6v} group includes four 1D irreducible representations (irreps), e.g., $A_{1,2}$ and $B_{1,2}$, and two 2D irreps, $E_{1,2}$. The 1D irreps $A_{1,2}$ and $B_{1,2}$ channels differ by their \hat{C}_{2z} eigenvalues, $\lambda_{C_{2z}} = +1$ for $A_{1,2}$ and $\lambda_{C_{2z}} = -1$ for $B_{1,2}$. Combining \hat{C}_{2z} and reshuffling symmetries leads to $\Delta^{\eta}_{\mathbf{k};e_{Y_1},e_{Y_2}} = \lambda_{C_{2z}} \Delta^{\eta}_{\mathbf{k};e_{Y_2},e_{Y_1}} \text{ for the spin singlet and } \Delta^{\eta}_{\mathbf{k};e_{Y_1},e_{Y_2}} =$ $-\lambda_{C_{2z}}\Delta^{\eta}_{\mathbf{k};e_{Y_2},e_{Y_1}}$ for the spin triplet. Thus, for intra-Chern-band pairing $(e_{Y_1} = e_{Y_2})$, the $A_{1,2}$ channel must be a spin singlet while the $B_{1,2}$ channel must be a spin triplet. Furthermore, the rotation C_{3z} ensures the existence of nodes at \mathbf{K}_M for the gap function of any 1D irrep intra-Chern-band channel $(\Delta_{\mathbf{K}_{M};e_{Y},e_{Y}}^{\eta}=0)$, while the inter-Chern-band channel does not have such a constraint. The 2D irreps E_1 and E_2 have different \hat{C}_{2z} eigenvalues, $\lambda_{C_{2z}} = +1$ for E_2 and $\lambda_{C_{2z}} = -1$ for E_1 , similarly to the 1D irrep case. Consequently, the E_2 channel must be a spin singlet while the E_1 channel must be a spin triplet for intra-Chern-band pairings. \hat{C}_{3z} guarantees nodes at Γ_M for both intra- and inter-Chern-band channels, and it requires additional nodes at \mathbf{K}_M for the inter-Chern-band channels for both 2D $E_{1,2}$ pairings. Besides discrete symmetries, the continuous $U(2)_{e_y=1} \times U(2)_{e_y=-1}$ spin symmetry in the chiral limit guarantees the singlet and triplet pairings of the inter-Chern-band channel to be degenerate in the chiral flat-band limit. The full symmetry analysis of the gap functions can be found in SM Sec. VI G [56] (see also Refs. [71,72] therein).

IV. GAP EQUATIONS AND SELF-CONSISTENT SOLUTION OF PAIRING CHANNELS

The linearized gap equation (LGE) for the attractive interaction (7) can be derived by evaluating $\langle \gamma_{-\mathbf{k}'e_{Y_2}\eta s_2} \gamma_{\mathbf{k}'e_{Y_1}-\eta s_1} \rangle$ in Eq. (9) and expanding it to linear order of the gap function. In the chiral flat-band limit, e.g., the bandwidth is much smaller than the critical temperature T_c , the LGE is derived as

$$2k_B T \Delta_{\mathbf{k};e_{Y_1}e_{Y_2}}^{\eta,SM} = \frac{1}{N_M} \sum_{\mathbf{k}'} V_{\mathbf{k},\mathbf{k}'}^{\eta e_{Y_1}e_{Y_2}} \Delta_{\mathbf{k}';e_{Y_1}e_{Y_2}}^{-\eta,SM}.$$
 (10)

This is an eigenequation problem for the matrix $V_{\mathbf{k},\mathbf{k}'}^{\eta e_{Y_1}e_{Y_2}}$: The T_c is determined by the largest eigenvalue and the symmetry of the gap function is determined by that of its eigenvector. As mentioned, the only two independent components of the V function (complex $V_{\mathbf{k},\mathbf{k}'}^{+++}$ and real $V_{\mathbf{k},\mathbf{k}'}^{++-}$) lead to two independent LGEs for the intra- and inter-Chern-band channels, respectively. The form of the LGE suggests that all the gap functions are doubly degenerate at T_c in the flatband limit. They belong either to two degenerate 1D irreps or one 2D irrep. We first numerically solve these two LGEs from Eq. (10), and find the forms of the gap functions with the largest eigenvalues, as shown in Fig. 1. Our numerical calculations show $k_B T_c \sim 0.21$ meV for the inter-Chern-band channel, slightly larger than $k_B T_c \sim 0.16$ meV for the intra-Chern-band channel. For the inter-Chern-band channels, the gap function is almost a constant in Fig. 1(a), featuring a fully gapped s-wave pairing with even \hat{C}_{2z} parity (A_1 or A_2 irrep). In the chiral flat-band limit, spin singlet and triplet pairings are degenerate, as required by the continuous $U(2) \times U(2)$ spin symmetry (SM Sec. VI E [56]). Including kinetic energy splits this degeneracy and makes the spin singlet A_1 irrep channel to have the highest T_c . For the intra-Chern-band channel, one can see nodes appearing at the Γ_M in Fig. 1(b). As our previous symmetry analysis shows that the gap function should have nodes at \mathbf{K}_M for the 1D irrep $(A_{1,2}, B_{1,2})$ and Γ_M for the 2D irrep $(E_{1,2})$, numerical results should correspond to a 2D irrep. Numerically analyzing the symmetry property of the gap function suggests that the intra-Chern-band channel belongs to the 2D E_2 irrep with a spin singlet. Full numerical results are discussed in SM Secs. VI H 2 and VI H 3 [56].

Our results for the intra-Chern-band channels reveal a *d*-wave character of the gap. Using the heavy-fermion formalism of TBG [47], we analytically obtain $V_{\mathbf{k},\mathbf{k}'}^{\eta,e_Y,e_Y}$,

$$V_{\mathbf{k},\mathbf{k}'}^{\eta,e_{Y},e_{Y}} = U_{e_{Y},\mathbf{k}}^{*}U_{e_{Y},\mathbf{k}'}, \quad U_{e_{Y},\mathbf{k}} = \frac{\sqrt{V_{0}}}{k^{2}+b^{2}}k_{e_{Y}}^{2}, \quad (11)$$

with $k_{e_Y} = k_x + ie_Y k_y$ ($e_Y = \pm$). This interaction allows us to solve the LGE analytically to obtain the T_c ,

$$k_B T_c = \frac{\tilde{V}_0}{2}, \quad \tilde{V}_0 = \frac{1}{N_M} \sum_{\mathbf{k}} V_0 \frac{k^4}{(k^2 + b^2)^2},$$
 (12)

where V_0 and b are material-dependent parameters. The corresponding self-consistent gap function takes the d-wave form

$$\begin{pmatrix} \Delta_{\mathbf{k},e_{Y}e_{Y}}^{+,00} \\ \Delta_{\mathbf{k},e_{Y}e_{Y}}^{-,00} \end{pmatrix} = \Delta_{e_{Y}} \frac{k_{-e_{Y}}^{2}}{k^{2} + b^{2}} \begin{pmatrix} 1 \\ 1 \end{pmatrix},$$
(13)

with $e_Y = \pm$ and Δ_{e_Y} a parameter to be determined. Time reversal, if it exists, requires $(\Delta_{\mathbf{k};--}^{-,00}, \Delta_{\mathbf{k};--}^{+,00}) = (\Delta_{\mathbf{k};++}^{+,00}, \Delta_{\mathbf{k};++}^{-,00})^*$. The *d*-wave nature of the gap function suggests the possibility of the nodal superconductivity. However, one should note that the single-particle Hamiltonian is *not* diagonal in the Chern-band basis. The Bogoliubov–de Gennes (BdG) spectrum must be checked with kinetic energy added. The BdG Hamiltonian for the intra-Chern-band pairing is block diagonal and one block $\mathcal{H}_{BdG}^{+,+}$ on the basis $(\gamma_{\mathbf{k},e_Y=\pm,+,s=\uparrow}^{\dagger}, \gamma_{-\mathbf{k},e_Y=\pm,-,s=\downarrow})$ reads

$$H_{\text{BdG}}^{+,+}(\mathbf{k}) = \begin{pmatrix} h_{+}(\mathbf{k}) & \Delta_{\mathbf{k}}^{+} \\ (\Delta_{\mathbf{k}}^{+})^{\dagger} & -h_{-}^{*}(-\mathbf{k}) \end{pmatrix},$$
(14)

with $h_{\eta}(\mathbf{k}) = (d_{0,\eta}(\mathbf{k}) - \mu)\zeta^{0} + d_{x,\eta}(\mathbf{k})\zeta^{x}$ and $\Delta_{\mathbf{k}}^{+} = \text{Diag}[\Delta_{\mathbf{k},++}^{+,0}, \Delta_{\mathbf{k},--}^{+,0}]$. Here, $d_{0,\eta}(\mathbf{k}) = [\epsilon_{+,\eta}(\mathbf{k}) + \epsilon_{-,\eta}(\mathbf{k})]/2$ and $d_{x,\eta}(\mathbf{k}) = [\epsilon_{+,\eta}(\mathbf{k}) - \epsilon_{-,\eta}(\mathbf{k})]/2$, where $\epsilon_{\pm,\eta}(\mathbf{k})$ are the eigenenergies for the two low-energy flat bands (per valley per spin) of the BM model $\hat{H}_{el}(2)$. The corresponding energy spectrum can possess nodes when the pairing amplitudes of two Chern-band channels are equal, $|\Delta_{e_{Y}=+}| = |\Delta_{e_{Y}=-}| = \Delta_{0}$, which corresponds to the Euler pairing discussed in Refs. [32,73]. Point nodes appear at the location defined by two conditions, (1) $\cos[(\Phi_{\mathbf{k},-} - \Phi_{\mathbf{k},+})/2] = 0$, where $\Phi_{\mathbf{k},e_{Y}} = \varphi_{e_{Y}} - 2e_{Y}\theta_{\mathbf{k}}$ with $\Delta_{e_{Y}} = \Delta_{0}e^{i\varphi_{e_{Y}}}$ and $k_{e_{Y}} = ke^{ie_{Y}\theta_{\mathbf{k}}}$, and (2) $d_{x,\mathbf{k}}^{2} = (d_{0,\mathbf{k}} - \mu)^{2} + \Delta_{0,\mathbf{k}}^{2}$ with $\Delta_{0,\mathbf{k}} = \Delta_{0}\frac{k^{2}}{k^{2}+k^{2}}$, as discussed in SM Sec. VI H 5 [56]. The first condition determines the momentum angle for the nodes while the second gives the momentum amplitude, thus together fixing



FIG. 2. (a) The superconductor order parameter amplitudes $|\Delta_{\pm}|$ (red circles and blue crosses) and the ground state energy (black dots) as a function of μ . The superconducting phase has nodes in the shadowed regime. (b) and (c) show the BdG spectrum with and without nodes at $\mu = 0.04$ and 0.14 meV, respectively. The single-particle bandwidth is set around 0.3 meV.

the location of point nodes in the 2D momentum space. We next solve the self-consistent gap equation at zero temperature for the interaction form (11). With the gap function ansatz $\Delta_{\mathbf{k};e_Y} = \Delta_{e_Y} \frac{k_{e_Y}^2}{k^2+b^2}$, we find a self-consistent gap equation

$$\Delta_{e_{Y}} = \frac{V_{0}}{N_{M}} \sum_{\mathbf{k}', e_{Y_{1}}} \frac{k'_{e_{Y}}^{2}}{k'^{2} + b^{2}} u_{-\mathbf{k}', e_{Y}e_{Y_{1}}} w_{-\mathbf{k}', e_{Y}e_{Y_{1}}}^{*}, \qquad (15)$$

where $\psi_{\mathbf{k},e_{Y_1}} = (u_{\mathbf{k},\pm,e_{Y_1}}, w_{\mathbf{k},\pm e_{Y_1}}) (e_{Y_1} = \pm)$ are the eigenwave functions with the positive eigenenergies of the BdG Hamiltonian $H^{+,+}_{BdG}(\mathbf{k})$ (14). Figure 2(a) shows the chemical potential dependence of the gap functions and the condensation energy. The Euler pairing $|\Delta_+| = |\Delta_-|$ is always energetically favored for a non-flat-bandwidth ~0.3 meV, quite different from chiral *d*-wave pairing in doped graphene [26,74,75]. For the chemical potential μ below 0.1 meV, a nodal superconductor phase with four point nodes [Fig. 2(b)] located at the positions determined by two conditions discussed above [32]. With increasing μ , four nodes move towards Γ_M and eventually a gapped superconductor phase [Fig. 2(c)] appears for $\mu > 0.1$ meV.

The energy scale of the Coulomb interaction in TBG is \sim 24 meV [58], much larger than the estimated energy scale of e-K-ph mediated attractive interaction \sim 0.3 meV [76]. Near the Van Hove singularities of flat bands, the screening can significantly reduce the Coulomb interaction to a similar order as the e-K-ph mediated interaction due to the large DOS (SM Sec. VI H 7 [56]), thus making superconductivity from this mechanism possible. If, however, the DOS is that of

the Hartree-Fock bands of correlated insulators, the screening might not be enough to reduce the Coulomb interaction. Hence, superconductivity from this K-phonon flat bare band mechanism could appear only when the correlated insulator states are suppressed and the Coulomb interaction is strongly screened [77], which is consistent with the TBG experiments with different Coulomb screenings [78].

V. CONCLUSION

In conclusion, we develop a theory for the projected e-K-ph interaction of the flat bands and the resulting superconductor pairing channels in TBG. We find the inter-Chern-band *s*-wave singlet pairing and the intra-Chern-band *d*-wave nematic singlet pairing have the highest T_c , and the T_c of the inter-Chern-band channel is slightly higher than the intra-Chen-band channel. The intra-Chen-band channel can have nodes in a large parameter regime. From the estimate of the screened Coulomb interaction, we argue that this mechanism requires the correlated insulators to be suppressed.

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