Tunneling and fluctuating electron-hole Cooper pairs in double bilayer graphene

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A strong low-temperature enhancement of the tunneling conductance between graphene bilayers has been reported recently and interpreted as a signature of equilibrium electron-hole pairing, first predicted in bilayers more than 40 years ago but previously unobserved. Here we provide a detailed theory of conductance enhanced by fluctuating electron-hole Cooper pairs, which are a precursor to equilibrium pairing, that accounts for specific details of the multiband double graphene bilayer system which supports several different pairing channels. Above the equilibrium condensation temperature, pairs have finite temporal coherence and do not support dissipationless tunneling. Instead they strongly boost the tunneling conductivity via a fluctuational internal Josephson effect. Our theory makes predictions for the dependence of the zero bias peak in the differential tunneling conductance on temperature and electron-hole density imbalance that capture important aspects of the experimental observations. In our interpretation of the observations, cleaner samples with longer disorder scattering times would condense at temperatures T_c up to ~50 K, compared to the record $T_c \sim 1.5$ K achieved to date in experiment.

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I. INTRODUCTION

The possibility of Cooper pairing in a system with spatially separated electrons and holes in semiconductor quantum wells was first anticipated more than 40 years ago [1,2]. According to theory, strong Coulomb interactions allow pairing at elevated temperatures, which would provide a physical realization of dipolar superfluidity that is potentially relevant for applications. The paired state is fragile however, and can be suppressed by disorder [3,4] or by Fermi-line mismatches due to the differences between electron and hole anisotropies [5,6] that are always present in conventional semiconductors [7]. In fact equilibrium pairing has until recently been observed only in the presence of strong magnetic fields that quench the kinetic energies of electrons and holes and drive the system to the regime of strong correlations [8,9].

Recent progress in fabricating single-atomic-layer twodimensional materials has renewed interest in electron-hole pairing [11–25]. Graphene-based two-dimensional electron systems not only have high mobility and almost perfect electron-hole symmetry, but they also make it possible to fabricate closely spaced, and therefore strongly interacting, independently gated and contacted double layer structures. Very recently low-temperature enhancement of the tunneling conductance between graphene bilayers has been observed at matched concentrations of electrons and holes [10]. A typical conductance trace is presented in Fig. 1, where we see a tunneling conductance that appears to diverge at $T_0 \approx 1.5$ K, signaling equilibrium pair condensation. This observation provides the first clear experimental signature of equilibrium electron-hole pair condensation in the absence of a magnetic field [26,27].

Enhanced tunneling conductance has been observed previously in semiconductor bilayers in the strong field quantum Hall regime [28] and has been interpreted as an internal Josephson effect [8]. The differential conductance does not diverge, however, and instead has a sharp peak at zero bias. The property that the conductance peak width is smaller than temperature, and smaller than the single-electron scattering rate (i.e., the Landau level width), nevertheless points to a collective origin of the peak. Bilayers in the quantum Hall regime are predicted to support *dissipationless* Josephson-like tunneling currents in the presence of long-range electron-hole coherence [29,30]. The development of a quantitative theory of enhanced tunneling in quantum Hall systems [31–34] that fully explains the peak width has been challenged by the importance of inhomogeneity and disorder and by strong

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FIG. 1. (a) Temperature dependence of the tunneling conductance $G_{\rm T}$ between graphene bilayers at zero voltage bias (V = 0). The red curve corresponds to the calculation that incorporates the effect of fluctuating Cooper pairs [Eq. (27)] and accurately fits the experimental data presented by blue dots [Fig. 3(b) in Ref. [10]]. The purple dashed curve corresponds to the model of noninteracting electrons and holes [Eq. (26)]. The fitting details and parameters are presented in Sec. V. Fluctuating Cooper pairs above the critical temperature $T_0 = 1.5$ K strongly enhance $G_{\rm T}$ and are responsible for its critical behavior $G_{\rm T} \sim (T - T_0)^{-2}$. The latter is derived and discussed in Sec. IV E and reasonably matches with the experimental data, as is clearly seen in the inset (b).

interactions in the presence of dispersionless Landau levels. Phase fluctuations that are inevitably present due to the two-dimensional Berezinskii-Kosterlitz-Thouless nature of the phase transition [35–38] also play a role. The theory of enhanced tunneling is simpler at a zero magnetic field, at least in the weak coupling regime where the electron-hole pairing energy is small compared to the Fermi energy, allowing experiments to be explained more fully, as we demonstrate below.

The enhancement of the tunneling conductance in the double bilayer graphene system has been observed over a wide density range 4×10^{10} – 10^{12} cm⁻², where electronic correlations vary from moderate to weak [39]. The Bardeen-Cooper-Schrieffer (BCS) theory of electron-hole pairing has greater validity at weaker pairing. True internal Josephson behavior in this case occurs only below a critical temperature T_0 and is preceded by enhancement of the tunneling conductance that diverges as T_0 is approached as illustrated in Fig. 1. This critical behavior has been predicted by one of the authors [40] and has been interpreted as a *fluctuational* internal Josephson effect. It originates from partly coherent fluctuating electron-hole Cooper pairs [40-42] that are a precursor of equilibrium pairing and reminiscent of Aslamazov-Larkin and related effects in superconductors [43–45]. Above T_0 fluctuating Cooper pairs have a finite coherence time [19,46-48] and cannot support a dissipationless tunneling current. While the recent observations do qualitatively agree with earlier theory, the double bilayer graphene system has some important differences compared to the double parabolic electron gas models considered previously. These are related to the system's wellknown 2π momentum space Berry phases. We show here that



FIG. 2. (a) Schematic of a double bilayer graphene system which specifies our sublayer labeling. External gates induce an excess of electrons (holes) in the top (bottom) bilayer and open a gap in the spectrum of isolated bilayers. (b) Electronic structure of the system in the case of the matched concentration of electrons and holes that most strongly favors electron-hole Cooper pairing.

accounting properly for these differences provides a better account of the low-temperature tunneling anomalies.

In the present work we have developed a theory of the fluctuational internal Josephson effect in a system with closely spaced graphene bilayers. As we show below, the presence of valley and sublattice degrees of freedom provides three competing electron-hole channels for both intravalley and intervalley Cooper pairs. We show that three channels are nearly independent and have have different condensation temperatures and different sublattice structure. The experimental enhancement of the tunneling conductance by fluctuating Cooper pairs can be explained only by the presence of competing channels that dominate in different temperature ranges. In the vicinity of T_0 , the tunneling conductance at zero bias is predicted to have a critical divergence $G_{\rm T} \sim$ $(T - T_0)^{-2}$, which matches the experimental data well, as we see in Fig. 1. The calculated dependence of the tunneling conductivity on interlayer voltage bias and carrier-density imbalance also match the experimental data [10] reasonably. We conclude that the observed enhancement of the tunneling conductance in double bilayer graphene is well explained by our fluctuational internal Josephson effect theory.

The paper is organized as follows. In Sec. II we introduce a model that describes the low-energy physics of two closely spaced graphene bilayers. Section III is devoted to a description of fluctuating Cooper pairs above the critical temperature T_0 . In Sec. IV we use these results as a starting point for a theory of the tunneling conductance. In Sec. V we compare our calculations with the recent experimental data. Finally in Sec. VI we discuss limitations of our theory and aspects of the experimental data that are still not well understood and present our conclusions.

II. MODEL

A. Weak and strong coupling regimes

The system of interest contains two graphene bilayers separated by an insulator as sketched in Fig. 2(a). An external electric field perpendicular to the bilayers induces an excess of electrons in the top bilayer (t) and their deficit in the bottom bilayer (b). Inevitably, it also results in gaps 2|u| in isolated bilayer graphene spectra. The latter are in Fig. 2(b) in the

case of matched concentrations for electrons and holes, the most favorable regime for the Cooper pairing. Double bilayer physics is very rich and has been considered in a number of recent papers [19–21,23,39] that aim to provide realistic predictions of the critical temperature T_0 based on the microscopic model. Here we follow a different route and consider a minimal phenomenological model that accounts for disorder and describes the instability of the system towards electronhole pairing in the weak coupling regime and can explain the observed enhancement of the tunneling conductance between graphene bilayers [10].

The physics of electron-hole Copper pairing depends on three dimensionless parameter: Wigner-Seitz interaction strength parameter $r_{\rm s} = me^2/\kappa \hbar^2 k_{\rm F}$ that scales the ratio of interactions and kinetic energy in an individual bilayer [49], a parameter $k_{\rm F}d$ that scales the interlayer Coulomb interactions respect to the intralayer one, and a parameter $|u|/\epsilon_{\rm F}$ that determines the low-energy spectrum of individual bilayers. Here $k_{\rm F}$ and $\epsilon_{\rm F}$ are Fermi wave vector and energy for electrons and holes, m is their effective constant, and κ and d are a dielectric constant for the spacer between bilayers and its thickness, respectively. If $r_s \ll 1$, $k_F d \gg 1$, and $|u|/\epsilon_F \ll$ 1 the system is in the weak coupling regime with pairing correlations only in the vicinity of Fermi level for electrons and holes. This regime can be described by the BCS theory for electron-hole pairing. The nature of the strong coupling regime $(r_s \gg 1 \text{ and } k_F d \sim 1)$ depends on the ratio $|u|/\epsilon_F$. If $|u| \gg \epsilon_{\rm F}$, the state is a Bose-Einstein condensate (BEC) of indirect excitons that represent a bound state of electron and hole. It has been argued that the most favorable conditions for observation of electron-hole condensation are reached near the midpoint of the BEC-BCS crossover [19-21.39.50]. In the opposite case the paired state is a multiband BCS-like paired state [15–18,51] where pairing correlations also span to remote bands (valence band in the layer with excess of electrons and conduction band in the layer with excess of holes).

Enhanced conductance has been seen over a range of electron and hole densities that covers $4 \times 10^{10}-10^{12}$ cm⁻² and corresponds to $r_s = 3.1-0.72$ and $k_Fd = 0.07-0.35$ [52]. This range suggests to the transition from moderate to weak coupling regime occurs with increasing of electron and hole concentrations. In the experimental setup [10] the Fermi energy ϵ_F of charge carriers and the gap 2|u| in the spectrum of individual bilayers are not controlled independently, but with the same gate (that produces the electric field perpendicular to bilayers). While the density can be tuned within a wide range, the ratio $|u|/\epsilon_F \approx d_{BG}/d \approx 0.15$ is approximately fixed. Here d_{BG} is the thickness of individual bilayers. This favors the multiband BCS-like state at strong coupling and extends the region of the applicability for the phenomenological weak coupling BCS theory that will be employed below.

B. Phenomenological model in the weak coupling regime

The spectrum sketched in Fig. 2(b) has spin and valley degeneracy. The spin degrees of freedom simply added a factor of 2 in the tunneling conductance when the condensed state preserves spin invariance and do not need to be treated explicitly. The low energy states in bilayer graphene are

concentrated around two inequivalent valleys K (v = 1) and K' (v = -1) situated at the corners of the first Brillouin zone and are described by the two-band Hamiltonian [53]

$$H_{0} = \sum_{\mathbf{p}} [\hat{\psi}_{t\mathbf{p}}^{+}(\hat{h}_{t\mathbf{p}} - \mu_{t})\hat{\psi}_{t\mathbf{p}} + \hat{\psi}_{b\mathbf{p}}^{+}(\hat{h}_{b\mathbf{p}} + \mu_{b})\hat{\psi}_{b\mathbf{p}}], \quad (1)$$

where $\hat{\psi}_{\mathbf{p}} = {\psi_{1\mathbf{p}}, \psi_{2\mathbf{p}}}$ is a spinor of annihilation operators for electrons in both layers [top (t) and bottom (b)] with sublattice indexes $\sigma = 1, 2$, that are numerated according to the sketch in Fig. 2(a). We assume that there is a deficit of electrons in the bottom layer, but it is instructive not to perform the transformation to field operators of holes. $\mu_t = \epsilon_F + h$ and $\mu_b = \epsilon_F - h$ characterize the electric potentials in the top and bottom bilayers. Here ϵ_F is the average of the electron and hole Fermi energies at neutrality, while $2h \ll \epsilon_F$ is their difference. The matrices \hat{h}_{tp} and \hat{h}_{bp} are

$$\hat{h}_{t\mathbf{p}} = \begin{pmatrix} u & \frac{p_{\tilde{b}_t}^2}{2m} \\ \frac{p_{\tilde{v}_t}^2}{2m} & -u \end{pmatrix}, \qquad \hat{h}_{b\mathbf{p}} = \begin{pmatrix} -u & \frac{p_{\tilde{v}_b}^2}{2m} \\ \frac{p_{\tilde{v}_b}^2}{2m} & u \end{pmatrix}. \tag{2}$$

Here $v = \pm 1$ and $\bar{v} = \mp 1$ are valley indexes, *m* is the electron mass, and $p_v = p_x + ivp_y$. The electric field perpendicular to bilayers opens a gap 2|u| separating conduction $\epsilon_{cp} = \epsilon_p$ and valence $\epsilon_{vp} = -\epsilon_p$ bands in each bilayer with $\epsilon_p = \sqrt{u^2 + (p^2/2m)^2}$. In the weak coupling regime the pairing correlations appear in the vicinity of Fermi lines for electrons and holes, and the presence of remote bands (valence band in the layer with excess of electrons and conduction band in the layer with excess of holes) can be neglected. In this regime only the conduction band of the top bilayer and the valence band of the bottom bilayer are relevant, and the corresponding spinor wave functions are

$$|\mathbf{tcp}\rangle = \begin{pmatrix} c_{\mathbf{p}} \ e^{-iv_{b}\phi_{\mathbf{p}}} \\ s_{\mathbf{p}} \ e^{iv_{b}\phi_{\mathbf{p}}} \end{pmatrix}, \quad |\mathbf{bvp}\rangle = \begin{pmatrix} c_{\mathbf{p}}e^{iv_{b}\phi_{\mathbf{p}}} \\ -s_{\mathbf{p}}e^{-iv_{b}\phi_{\mathbf{p}}} \end{pmatrix}.$$
(3)

Here $\phi_{\mathbf{p}}$ is the polar angle; $c_{\mathbf{p}} = \cos(\vartheta_{\mathbf{p}}/2)$ and $s_{\mathbf{p}} = \sin(\vartheta_{\mathbf{p}}/2)$ with $\cos(\vartheta_{\mathbf{p}}) = u/\epsilon_{\mathbf{p}}$. The spinors have valleydependent chirality $\pm v_{t(b)}\phi_{\mathbf{p}}$ that defines a sublattice structure of fluctuating electron-hole Cooper pairs as will be shown below. We introduce disorder with the help of phenomenological scattering rates $\gamma_{t(b)}$. It is important in this theory to observe that because the electron and hole components of the Cooper pair are spatially separated and have opposite charges, both short- and long-range Coulomb disorder lead to pair-breaking [3,4].

In experiment [10] the relative angle θ between graphene bilayers can be adjusted. Since valleys *K* and *K'* reside at the corners of the first Brillouin zone, the valley momenta in the two layers do not match in the presence of a twist. For momentum-conserving tunneling, the current is maximized when the layers are aligned ($\theta = 0$) or twisted by $\theta = n\pi/3$. For even *n*, valley K (K') in one layer is aligned with valley K (K') in the other layer, whereas for *n* odd valley K (K') in one layer is aligned with valley K' (K) in the other layer. When states are labeled by their momenta relative to the Brillouin-zone corners, the tunneling Hamiltonian for θ close to $n\pi/3$ is

$$H_{t} = T^{+} + T = \sum_{\mathbf{p}} [\hat{\psi}_{bp}^{+} \hat{t}^{+} \hat{\psi}_{t\mathbf{p}+\mathbf{Q}} + \hat{\psi}_{t\mathbf{p}+\mathbf{Q}}^{+} \hat{t} \hat{\psi}_{bp}].$$
(4)

Here $\mathbf{Q} = \mathbf{Q}_{\theta} + \mathbf{Q}_{\mathbf{B}}$ is the momentum splitting between valleys in the different layers. The twist contribution at small relative angle $\theta \ll 1$ can be approximated as $\mathbf{Q}_{\theta} = -[\mathbf{q}_{\mathbf{K}} \times \mathbf{e}_{z}]\theta$ where $\mathbf{q}_{\mathbf{K}}$ is the momentum for the Brillouin-zone corner in bilayer graphene. It has opposite directions for valleys K and K', and its magnitude is $q_{\mathbf{K}} = 4\pi\hbar/3a_{0}$ with a_{0} the corresponding Bravais lattice period. The contribution induced by an in-plane magnetic field $\mathbf{B}_{||}$ is the same for two valleys and is equal to $Q_{\mathbf{B}} = ed[\mathbf{B}_{||} \times \mathbf{e}_{z}]/\hbar c$. Because each bilayer is represented by a two-band model, the matrix \hat{t} has four matrix elements, which we treat in a phenomenological way below, with the expectation that since t_{22} corresponds to the tunneling between adjacent sublayers whereas t_{11} to the tunneling between remote sublayers, $|t_{11}| \ll |t_{12}| \approx |t_{21}| \ll |t_{22}|$.

III. FLUCTUATING COOPER PAIRS

A. The Cooper instability

Due to Coulomb interactions between electrons and holes, the double bilayer system is unstable towards Cooper pairing. Here we omit repulsive interactions within each graphene bilayer since its main effect in the considered weak coupling regime is a simple renormalization of the quasiparticle spectra. The interbilayer attraction is

$$H_{\rm int} = \sum_{\mathbf{pp'q}} \sum_{\sigma_t \sigma_b} U_{\mathbf{q}} \psi^+_{\mathbf{t}, \mathbf{p+q}, \sigma_t} \psi^+_{\mathbf{b}, \mathbf{p'-q}, \sigma_b} \psi_{\mathbf{b}\mathbf{p'}\sigma_b} \psi_{\mathbf{t}\mathbf{p}\sigma_t}, \qquad (5)$$

where U_q is the screened Coulomb potential estimated in our previous work [14–16]. Below we employ a multipole decomposition of the interaction and set the momenta magnitudes

to the Fermi momentum so that U_q reduces to a constant U_l for each orbital angular momentum channel *l*. The set of U_l parameters are also treated as phenomenological parameters with the expectation that the *s*-wave moment $U_s \equiv U_0$ is largest. Since valleys are well separated in momentum space we neglect intervalley scattering for electrons and holes.

The instability of a double layer system towards electronhole Cooper pairing is signaled by divergence of the electronhole channel scattering vertex $\Gamma_{\sigma_b'\sigma_b}^{\sigma_t'\sigma_t}(\omega, \mathbf{p}', \mathbf{p}, \mathbf{q})$ at zero frequency ω . Note that in our approximation scattering conserves valley indices for electrons (v_t) and holes (v_b) . The Γ vertex satisfies the Bethe-Saltpeter equation presented in Fig. 3(a) and is algebraic within the multipole approximation. It is instructive to combine σ_t and σ_b into a single index $|\sigma_t \sigma_b\rangle$ that varies between 1 and 4 as $|1\rangle = |11\rangle$, $|2\rangle = |12\rangle$, $|3\rangle =$ $|21\rangle$, and $|4\rangle = |22\rangle$. With this definition the Bethe-Saltpeter equation can be written (see Appendix A for details) in a compact matrix form:

$$\hat{\Gamma}_{l'l} = U_l \,\delta_{l'l} + \sum_{l''} U_{l'} \,\hat{M}_{l'-l''} \,\Pi \,\hat{\Gamma}_{l''l}, \tag{6}$$

where momentum and frequency dependence are suppressed, *l* (*l'*) is the orbital momentum for the relative motion of two particles before (after) scattering, and $\hat{\Gamma}_{l'l}$ is the corresponding scattering matrix. We have separated a factor of $\Pi(\omega, \mathbf{q})$ which also appears as the single-step pair propagator in the Cooper ladder sum of a bilayer system without sublattice degrees of freedom:

$$\Pi(\omega, \mathbf{q}) = N_{\rm F} \left(\ln\left[\frac{\epsilon_{\rm c}}{2\pi T}\right] - \frac{1}{2} \sum_{\zeta=\pm} \left\langle \Psi\left\{\frac{1}{2} + \frac{i[\omega + \zeta(h + v_{\rm F}q\cos\phi_{\rm p})] + \gamma}{4\pi T}\right\} \right\rangle_{\phi_{\rm p}} \right),\tag{7}$$

where ϵ_c is an energy cutoff that is required for momentum-independent interactions. The average $\langle \cdots \rangle_{\phi_p}$ is calculated respect to a polar angle ϕ_p . $\Psi(x)$ is the logarithmic derivative of the Γ function (or the digamma function), while $\zeta = \pm 1$ is the summation index. $\gamma = \gamma_t + \gamma_b$ is the pair-breaking rate, which is the sum of the scattering rates for electrons γ_t and holes γ_b . The expression for $\Pi(\omega, \mathbf{q})$ in Eq. (7) is well known [43,44] from previous work on systems without layer degrees of freedom. The chiral nature of the bilayer graphene charge carriers is captured by the nontrivial matrix form factor \hat{M}_l in the Bethe-Salpheter Eq. (6). The matrix form factor \hat{M}_l is defined as the multipole moment of the two-particle matrix element

$$\hat{M}_{\sigma_{b}^{\prime},\sigma_{b}}^{\sigma_{t}^{\prime},\sigma_{t}}(\phi_{\mathbf{p}}) = \left\langle \sigma_{t}^{\prime} | \mathrm{tc}\mathbf{p} + \frac{\mathbf{q}}{2} \right\rangle \left\langle \mathrm{tc}\mathbf{p} + \frac{\mathbf{q}}{2} | \sigma_{t} \right\rangle \left\langle \sigma_{b} | \mathrm{bv}\mathbf{p} - \frac{\mathbf{q}}{2} \right\rangle \left\langle \mathrm{bv}\mathbf{p} - \frac{\mathbf{q}}{2} | \sigma_{b}^{\prime} \right\rangle.$$

$$\tag{8}$$

Corrections to the form factor \hat{M}_l due to finite Cooper pair momentum q, $|\Delta M| = q^2 |u|/4p_F^2 \epsilon_F$, are negligible in the weak coupling regime since $q \ll p_F$. As a result, the matrix (8) can be approximated as

$$\hat{M}(\phi_{\mathbf{p}}) = \begin{pmatrix} c^4 & -c^3 s e^{-2iv_b\phi_{\mathbf{p}}} & c^3 s e^{-2iv_t\phi_{\mathbf{p}}} & -c^2 s^2 e^{-2i(v_t+v_b)\phi_{\mathbf{p}}} \\ -c^3 s e^{2iv_b\phi_{\mathbf{p}}} & c^2 s^2 & -c^2 s^2 e^{2i(v_b-v_t)\phi_{\mathbf{p}}} & cs^3 e^{-2iv_t\phi_{\mathbf{p}}} \\ c^3 s e^{2iv_t\phi_{\mathbf{p}}} & -c^2 s^2 e^{2i(v_t-v_b)\phi_{\mathbf{p}}} & c^2 s^2 & -cs^3 e^{-2iv_b\phi_{\mathbf{p}}} \\ -c^2 s^2 e^{2i(v_t+v_b)\phi_{\mathbf{p}}} & cs^3 e^{2iv_t\phi_{\mathbf{p}}} & -cs^3 e^{2iv_b\phi_{\mathbf{p}}} & s^4 \end{pmatrix}.$$
(9)

Here the coefficient c and s correspond to the coherence factors c_p and s_p in (3) evaluated at the average Fermi energy ϵ_F for electrons and holes and are given by

$$c^{2} = \frac{1}{2} \left(1 + \frac{u}{\epsilon_{\rm F}} \right), \qquad s^{2} = \frac{1}{2} \left(1 - \frac{u}{\epsilon_{\rm F}} \right). \tag{10}$$

The matrix form factor (9) shapes the sublattice structure of fluctuating electron-hole Cooper pairs in double bilayer graphene. Note that it couples scattering channels $\hat{\Gamma}_{l'l}$ with different orbital momenta. Importantly, \hat{M}_l has only even harmonics $l = 0, \pm 2, \pm 4$, which forbid scattering between states with even and odd orbital momenta. For isotropic Coulomb



FIG. 3. (a) Bethe-Saltpeter equation in the electron-hole channel. Divergence of the many-body vertex $\Gamma(\omega = 0, \mathbf{Q}_0)$ signals the double bilayer electron-hole pairing instability with a momentum \mathbf{Q}_0 . (b) The noninteracting tunneling response function $\chi_0(\omega, \mathbf{q})$ introduced in Eq. (18). (c) The tunneling response function $\chi(\omega, \mathbf{q})$ with the renormalized vertex t' that captures the effect of fluctuating Cooper pairs. (d) Renormalization of t' that illustrates Eq. (21).

interactions, the *s*-wave moment U_s is expected to be largest. It is instructive to start by neglecting all other moments. In that case only $\hat{\Gamma}_{00}$ is nonzero. The latter depends on the *s*-wave moment of the form factor \hat{M}_0 , which has a different form for intravalley and intervalley Cooper pairs. We discuss these two cases separately below.

B. Intravalley Cooper pairs

For intravalley $(v_t = 1 \text{ and } v_b = 1)$ electron-hole Cooper pairs the *s*-wave moment of the form factor $\hat{M}_0 = \langle \hat{M}(\phi_p) \rangle_{\phi_p}$ is given by

$$\hat{M}_0 = \begin{pmatrix} c^4 & 0 & 0 & 0\\ 0 & c^2 s^2 & -c^2 s^2 & 0\\ 0 & -c^2 s^2 & c^2 s^2 & 0\\ 0 & 0 & 0 & s^4 \end{pmatrix}.$$
 (11)

The scattering problem decouples into the three channels identified in Ref. [21], and the corresponding scattering vertex is given by

$$\frac{\hat{\Gamma}_{00}}{U_{\rm s}} = \begin{pmatrix} \frac{1}{L_{11}} & 0 & 0 & 0\\ 0 & \frac{1-c^2s^2U_{\rm s}\Pi}{L_{12-21}} & \frac{-c^2s^2U_{\rm s}\Pi}{L_{12-21}} & 0\\ 0 & \frac{-c^2s^2U_{\rm s}\Pi}{L_{12-21}} & \frac{1-c^2s^2U_{\rm s}\Pi}{L_{12-21}} & 0\\ 0 & 0 & 0 & \frac{1}{L_{22}} \end{pmatrix}, \qquad (12)$$

TABLE I. Coupling constants λ_{α} for intravalley Cooper pairs.

Channel, α	s-wave, λ^{s}	<i>d</i> -wave, λ^d
11	$c^4 N_{\rm F} U_{\rm s}$	$2c^2c^2N_{\rm F}U_{\rm d}$
22	$s^4 N_{\rm F} U_{\rm s}$	$2c^2c^2N_{\rm F}U_{\rm d}$
12-21	$2s^2c^2N_{ m F}U_{ m s}$	$(c^4 + s^4)N_{\rm F}U_{\rm d}$



FIG. 4. Dependence of the s-wave coupling constants λ_{α}^{s} for different pairing channels on the displacement field parameter *u*. For intravalley (a) Cooper pairs a competition between channels is possible, but in the considered regime $|u| \ll \epsilon_{\rm F}$ the mixed one 12-21 dominates. For intervalley Cooper pairs (b) the hierarchy between channels does not depend on ratio between |u| and $\epsilon_{\rm F}$, and the mixed channel 11-22 is the dominant one.

where $L_{\alpha} = 1 - \lambda_{\alpha}^{s} \Pi / N_{F}$ is a dimensionless inverse Cooper propagator for channel α and λ_{α}^{s} is the corresponding coupling constant specified in Table I. L_{α} vanishes at the critical temperature T_{α} for the electron-hole pairing instability in channel α .

In the absence of disorder and electron-hole density imbalances, the critical temperatures are given by $\bar{T}_{\alpha} = 2e^{C} \Lambda \exp[-1/\lambda_{\alpha}^{s}]/\pi$, where C = 0.577 is the Euler constant. Although the coupling constant values λ_{α} can be fine-tuned by the displacement field, as is presented in Fig. 4(a), their hierarchy is universal for the case $|u| \ll \epsilon_{\rm F}$. The coupling constant $\lambda_{12-21}^{\rm s} \approx 1/2$ is almost twice as large as the constants $\lambda_{11(22)}^{\rm s} \approx 1/4$, ensuring domination of the mixed channel 12-21. Physically the presence of two sub lattice combinations ($|12\rangle$ and $|21\rangle$) doubles the number of states that take part in the Cooper pairing.

C. Intervalley Cooper pairs

For intervalley $(v_t = 1 \text{ and } v_b = -1)$ electron-hole Cooper pairs the *s*-wave moment of the form factor $\hat{M}_0 = \langle \hat{M}(\phi_{\mathbf{p}}) \rangle_{\phi_{\mathbf{p}}}$ is given by

$$\hat{M}_0 = \begin{pmatrix} c^4 & 0 & 0 & -c^2 s^2 \\ 0 & c^2 s^2 & 0 & 0 \\ 0 & 0 & c^2 s^2 & 0 \\ -c^2 s^2 & 0 & 0 & s^4 \end{pmatrix}.$$
 (13)

TABLE II. Coupling constants λ_{α} for intervalley Cooper pairs.

Channel, α	<i>s</i> -wave, λ^s	<i>d</i> -wave, λ^{d}
12	$c^2 s^2 N_{\rm F} U_{\rm s}$	$(c^4 + s^4)N_{\rm F}U_{\rm d}$
21	$c^2 s^2 N_{\rm F} U_{\rm s}$	$(c^4 + s^4)N_{\rm F}U_{\rm d}$
11-22	$(c^4 + s^4)N_{\rm F}U_{\rm s}$	$2c^2s^2N_{\rm F}U_{\rm d}$

This case also decouples into three independent channels, with scattering vertex

$$\frac{\hat{\Gamma}_{00}}{U_{\rm s}} = \begin{pmatrix} \frac{1-s^4 U_{\rm s}\Pi}{L_{11-22}} & 0 & 0 & \frac{-c^2 s^2 U_{\rm s}\Pi}{L_{11-22}} \\ 0 & \frac{1}{L_{12}} & 0 & 0 \\ 0 & 0 & \frac{1}{L_{21}} & 0 \\ \frac{-c^2 s^2 U_{\rm s}\Pi}{L_{11-22}} & 0 & 0 & \frac{1-c^4 U_{\rm s}\Pi}{L_{11-22}} \end{pmatrix}.$$
 (14)

Interestingly, the sublattice structure of the Cooper pairs and the corresponding coupling constants are different in intravalley and intervalley cases. The latter are presented in Table II, and their dependence on displacement field parameter u is shown in Fig. 4(b). In that case the hierarchy between coupling constants is universal and does not depend on the ratio between |u| and $\epsilon_{\rm F}$. The mixed channel 11-22 has the highest critical temperature. The separation of scattering problem into three channels is not an artifact of the *s*-wave truncation, but is maintained when higher multipole momenta of interactions U_1 are taken into account. When the *d*-wave interaction $U_d \equiv U_{\pm 2}$ is nonzero (*s*- and *p*-wave momenta are decoupled, and the latter is irrelevant) the scattering matrix $\Gamma_{l'l}$ is nonzero for l = -2, 0, 2. We show below that the only effect of the *d*wave momentum U_d on the tunneling conductance between graphene bilayers is a renormalization of coupling constants $\lambda_{\alpha} = \lambda_{\alpha}^s + \lambda_{\alpha}^d$. The *d*-wave coupling constants λ_{α}^d are summarized in Tables I and II.

D. Disorder and density imbalances

Disorder and electron-hole density imbalances both reduce the critical temperature \overline{T}_0 (index 0 corresponds to the channel α that has the largest critical temperature of electron-hole pairing). The latter can stabilize the Fulde-Ferrell-Larkin-Ovchinnikov (FFLO) pairing state [5,6,54–57] with finite Cooper pair momentum \mathbf{Q}_0 . (References [56,57] also outlines a stabilization of the Sarma phases [58] within the BEC-BCS crossover. These phases are unstable in the weak-coupling regime considered here.) The critical temperature T_0 and the instability momentum \mathbf{Q}_0 for a channel with the highest critical temperature satisfy the equation $L_0(0, \mathbf{Q}_0) = 0$, which can be recast as

$$\ln\left[\frac{T_0}{\bar{T}_0}\right] + \frac{1}{2} \sum_{\zeta=\pm} \left\langle \Psi\left[\frac{1}{2} + \frac{i\zeta(h+v_{\rm F}Q_0\cos\phi_{\rm p}) + \gamma}{4\pi T}\right] - \Psi\left(\frac{1}{2}\right) \right\rangle_{\phi_{\rm p}} = 0.$$
⁽¹⁵⁾

The dependence of the critical temperature T_0 on the pairbreaking rate γ and the electron-hole imbalance expressed as a difference in Fermi energies, h, are illustrated in Fig. 5(a). The electron-hole pair instability is suppressed when the pair-breaking rate exceeds a critical value $\gamma \approx 1.78 \bar{T}_0$. If the rate does not exceed $\gamma \approx T_0$ the Fulde-Ferrell-Larkin-Ovchinnikov (FFLO) state with finite Cooper pair momentum \mathbf{Q}_0 is stabilized by finite imbalance. The dependence of the instability momentum \mathbf{Q}_0 is presented in Fig. 5(b). Its magnitude can be approximated by $v_F Q_0 \approx h$, corresponding to the difference between the Fermi momenta for electrons and holes.

E. Coherence time and length of fluctuating Cooper pairs

In the absence of electron-hole imbalance, the Cooper propagator L_0^{-1} of the dominating channel at small frequencies and momenta simplifies to

$$L_0^{-1}(\omega, \mathbf{q}) = \frac{1}{\lambda_0 N_{\rm F}} \frac{1}{i\omega\tau - \epsilon_{\mathbf{q}}}, \qquad \epsilon_{\mathbf{q}} = \epsilon + \frac{\xi^2 \mathbf{q}^2}{\hbar^2}, \quad (16)$$

where $\epsilon = \ln [T/T_0]$ is the energy scale that is required to create a uniform fluctuating Cooper pair. It vanishes at the critical temperature T_0 and is linear, $\epsilon \approx \Delta T/T_0$, in its vicinity $\Delta T \approx T_0$. Here $\Delta T = T - T_0$. τ and ξ are characteristic time

and spatial scales for Cooper pairs, which are given by

$$\tau = \frac{\hbar \Psi' \left(\frac{1}{2} + \frac{\gamma}{4\pi T}\right)}{4\pi T}, \quad \xi = \frac{\hbar v_F |\Psi'' \left(\frac{1}{2} + \frac{\gamma}{4\pi T}\right)|^{\frac{1}{2}}}{8\pi T}.$$
 (17)

They are connected to the coherence time and length of Cooper pairs by $\tau^* = \tau/2\epsilon$ and $\xi^* = \xi/\sqrt{\epsilon}$, which diverge at the critical temperature for electron-hole condensation T_0 . The Cooper propagator (16) has its only pole on the imaginary frequency axis at $\omega_{\mathbf{q}} = -i/2\tau_{\mathbf{q}}^*$ with $\tau_{\mathbf{q}}^* = \tau/2\epsilon_{\mathbf{q}}$, reflecting the dissipative nature of Cooper pairs dynamics. Due to the presence of a finite temporal coherence time τ^* , fluctuating Cooper pairs do not provide a dissipationless Josephson current but do strongly enhance the tunneling conductance at zero voltage bias.

IV. TUNNELING CONDUCTIVITY

A. Linear response theory for the tunneling conductance

When interlayer tunneling amplitudes are treated in leading order of perturbation theory, the interbilayer tunneling conductance at finite voltage bias V is [59–61]

$$G_{\rm T}(V) = \frac{8Ae^2}{\hbar} \frac{{\rm Im}[\chi(eV, \mathbf{Q})]}{eV}.$$
 (18)

Equation (18) accounts for the fourfold degeneracy due to the presence of valley and spin degrees of freedom, and A is the sample area. Here $\chi(\Omega)$ is the retarded correlation function



FIG. 5. Dependence of the critical temperature T_0 (a) and the instability momentum of Cooper pairs Q_0 on the pair-breaking rate γ and electron-hole imbalance, parameterized by the difference between electron and hole Fermi energies 2h. When disorder is weak $\gamma \lesssim T_0$ the imbalance can stabilize the Fulde-Ferrell-Larkin-Ovchinnikov (FFLO) state with the finite Cooper pair momentum.

corresponding to the imaginary-time ordered correlation function constructed from tunneling operators:

$$\chi(\tau, \mathbf{q}) = -\langle T_{\mathrm{M}}T(\tau, \mathbf{q})T^{+}(0, \mathbf{q})\rangle.$$
(19)

 $\chi(\omega, \mathbf{q})$ can be constructed from $\chi(\tau, \mathbf{q})$ by the usual Fourier transform and analytical continuation steps. Without Coulomb interactions between electrons and holes the response function $\chi(\omega, \mathbf{q})$ corresponds to the single electron-hole loop diagram depicted in Fig. 3(b) and is given by

$$\chi_0(\omega, \mathbf{q}) = \hat{t}^+ \hat{M}_0 \,\Pi \,\hat{t},\tag{20}$$

where $\hat{t} = \{t_{11}, t_{12}, t_{21}, t_{22}\}$ is the vector of tunneling matrix elements in the compact representation, and $\Pi \equiv \Pi(\omega, \mathbf{q})$ is a single step pair propagator in the Cooper ladder defined in Eq. (7). In the presence of Coulomb interactions, the singleparticle Green functions and the tunneling vertex need to be renormalized. The renormalization of Green functions results in a dip of the density of states at the Fermi level [42], which does not produce any singularities in the tunneling conductivity and thus is unimportant and can be neglected. The renormalized tunneling vertex $\hat{t} \rightarrow \hat{t}' \equiv \hat{t}'(\omega, \mathbf{q})$ diverges at the critical temperature T_0 and is responsible for the drastic enhancement of the tunneling vertex is presented in Fig. 3(d), and the corresponding equation for \hat{t}' can be written as

$$\hat{t}'_{l} = t \,\delta_{l0} + \sum_{l'} U_l \,\hat{M}_{l-l'} \,\Pi \,\hat{t}'_{l'}.$$
(21)

The matrix form factor \hat{M} couples even orbital channels, and we neglect all multipole moments except for *s* and *d*. (The *p*-wave multipole moment is decoupled from the *s*-wave one and is irrelevant.) Since the form factors \hat{M} are different for intravalley and intervalley Cooper pairs, we again consider these two cases separately.

B. Intravalley tunneling

Without Coulomb interactions between electrons and holes the response function $\chi(\omega, \mathbf{q})$ is a sum of three noninterfering terms that correspond to three channels α introduced in Sec. III and identified in Ref. [21]:

$$\chi^{0} = (c^{4}|t_{11}|^{2} + s^{4}|t_{22}|^{2} + |t_{12} - t_{21}|^{2}c^{2}s^{2}) \Pi.$$
(22)

The three channels are not coupled by Coulomb interactions and the response function $\chi(\omega, \mathbf{q})$ is given by

$$\chi = \left(\frac{c^4 |t_{11}|^2}{L_{11}} + \frac{s^4 |t_{22}|^2}{L_{22}} + \frac{|t_{12} - t_{21}|^2 c^2 s^2}{L_{12-21}}\right) \Pi, \quad (23)$$

where L_{α} is the inverse Cooper propagator for each channel and $\lambda_{\alpha} = \lambda_{\alpha}^{s} + \lambda_{\alpha}^{d}$ is the corresponding coupling constant. Importantly the only role of the *d*-wave interaction moment is the renormalization the coupling constant λ_{α} .

C. Intervalley tunneling

When opposite valleys are aligned by a twist angle between bilayers close to $\theta = \pi/3$, the response function $\chi(\omega, \mathbf{q})$ for noninteracting electrons and holes is

$$\chi_0 = \left(c^2 s^2 |t_{12}|^2 + c^2 s^2 |t_{21}|^2 + |c^2 t_{11} - s^2 t_{22}|^2\right) \Pi.$$
(24)

It again is a sum of three noninterfering terms that correspond to three channels α . Coulomb interactions do not couple the channels, and the response function $\chi(\omega, \mathbf{q})$ is given by

$$\chi = \left(\frac{c^2 s^2 |t_{12}|^2}{L_{12}} + \frac{c^2 s^2 |t_{21}|^2}{L_{21}} + \frac{|c^2 t_{11} - s^2 t_{22}|^2}{L_{11-22}}\right) \Pi.$$
 (25)

Each channel acquires its own Cooper propagator L_{α} with the coupling constant $\lambda_{\alpha} = \lambda_{\alpha}^{s} + \lambda_{\alpha}^{d}$.

D. Tunneling conductance

Due to the remarkable conservation of momentum for electron tunneling, the conductance $G_{\rm T}$ is observable only if graphene bilayers are aligned with twist angle $\theta = 0$ or $\theta = \pi/3$. In the former case the tunneling is intravalley, while in the latter case it is intervalley. Therefore we refer to the cases $\theta = 0$ and $\theta = \pi/3$ as to intravalley and intervalley alignments, respectively. As we explain below, there are important differences in the enhancement of tunneling conductance by fluctuating the Cooper pair in these two cases.

Within the model of noninteracting electrons and holes the tunneling conductance is dominated by tunneling between adjacent sublayers t_{22} and can be approximated (for both for both intra- and intervalley alignments) as

$$G_{\rm T}^0 = \frac{8Ae^2}{\hbar} \frac{s^4 |t_{22}|^2 \text{Im}[\Pi(eV, \mathbf{Q})]}{eV}.$$
 (26)

In the presence of interactions, the enhancement of tunneling conductance by fluctuating Cooper pairs works very differently for intra- and intervalley alignments. The reason is a drastic difference in the sublattice structure for the pairing correlations. For intravalley fluctuating Cooper pairs, the dominating channel 12-21 does not involve pairing correlations at adjacent sublayers. As result, the corresponding contribution of the channel 12-21 is weaker than that of 22, since the latter involves tunneling between adjacent sublayers, except in the vicinity of the critical temperature T_0 . The tunneling conductance is governed by the competition of two channels and can be approximated as

$$G_{\rm T} = \frac{8Ae^2}{\hbar} \left(\frac{s^4 |t_{22}|^2}{|L_{22}|^2} + \frac{|t_{12} - t_{21}|^2 c^2 s^2}{|L_{12} - 21|^2} \right) \frac{{\rm Im}[\Pi(eV, \mathbf{Q})]}{eV}.$$
 (27)

As we explain in the next section, the competition between channels is essential for the explanation of the experimental data [10], which in our interpretation strongly depend on mutual orientation of graphene bilayers.

For the intervalley alignment, the dominant channel for fluctuating Cooper pairs involves pairing correlations at adjacent sublayers. As a result, the tunneling conductance can be well approximated by a single term corresponding to the channel 11-22 and is given by

$$G_{\rm T} = \frac{8Ae^2}{\hbar} \frac{s^4 |t_{22}|^2}{|L_{1-22}|^2} \frac{{\rm Im}[\Pi(eV, \mathbf{Q})]}{eV}.$$
 (28)

The intervalley alignment case has not been studied experimentally yet, but according to our theory it is most favorable for observations of the fluctuational internal Josephson effect.

E. Critical behavior of tunneling conductance

At matched concentrations of electrons and holes the voltage dependence of tunneling conductance G_T in the vicinity of the critical temperature T_0 acquires a Lorentzian shape which is governed by the factor

$$F(eV, \mathbf{Q}) = \frac{\mathrm{Im}[\Pi]}{eV|L_0|^2} = \frac{\tau}{(eV\tau)^2 + \epsilon_{\mathbf{Q}}^2},$$
(29)

where $\epsilon_0 = \epsilon + \xi^2 \mathbf{Q}^2 / \hbar^2$ and $\epsilon = \ln(T/T_0)$ can be interpreted as an energy of fluctuating Cooper pairs. In the absence of a valley splitting $\mathbf{Q} = 0$, the amplitude of the peak has a long high-temperature tail $F(0, 0) = \tau / \ln^2[T/T_0]$. It diverges near the critical temperature T_0 as a function of $\Delta T = T - T_0$ in the critical manner as $F(0, 0) \approx \tau T_0^2 / \Delta T^2$ with an index 2. Its width at half maximum $eV_{\rm HM} = 1/\tau^* = 2\Delta T/T_0\tau$ is equal to the inverse coherence time τ^* of fluctuating Cooper pairs and vanishes linearly at T_0 . In the presence of a valley splitting that can be induced by in-plane magnetic field or relative twist, fluctuating Cooper pairs with finite momentum Q are probed in tunneling experiments. Temperature dependence of the peak width is modified as $eV_{HM} = [1 + (\xi^* \mathbf{Q})^2]/\tau^*$. It has a universal form as a function of coherence time $\tau_* = 2\tau/\epsilon$ and length $\xi^* = \xi / \sqrt{\epsilon}$ for fluctuating Cooper pairs, that allow one to extract them from the experimental data in the presence of an in-plane magnetic field.

For the case of the intervalley alignment ($\theta = \pi/3$), the temperature dependence of the tunneling conductance (28) is well approximated by Eq. (29) in a wide temperature range. For the case of intravalley alignment ($\theta = 0$) channels 12-21 and 22 compete with each other. As a result, tunneling conductance Eq. (27) cannot be approximated by a simple analytical expression over as wide of a temperature range. It is

well approximated by the critical behavior given by Eq. (29) only in the narrow temperature range where the contribution of the channel 21-21 dominates.

V. COMPARISON WITH EXPERIMENT

In Ref. [10] the twist angle between graphene layers can in principle be tuned to access the intravalley ($\theta = 0$) and intervalley ($\theta = \pi/3$) tunneling cases, although experimental results for tunneling conductance $G_{\rm T}$ have so far been reported only for one alignment. A preliminary analysis of the data suggests that a divergent zero bias peak appears on the top of a background with a weaker temperature dependence. This behavior can be explained by the competition between channels 12-21 and 22. This is not surprising since top and bottom graphene bilayers originate from the same flake and are aligned with the twist angle $\theta = 0$. Calculations for the intravalley alignment are presented below and compared with the experimental data, and ones for the intervalley one are presented in Appendix B.

To fit the experimental data, the tunneling conductance $G_{\rm T}$ has been calculated with the help of Eq. (27). The model has a large number of fitting parameters. To adjust their values and compare our results with the experimental data we use the following strategy, which involves six steps.

(1) The experimental data have already been carefully analyzed [10] within a model of noninteracting charge carriers. In the wide range of concentrations the noninteracting model explains the tunneling data very well except in the case of opposite polarity charge carriers with nearly equal electron and hole densities. Based on the fits to experimental data away from matched electron and hole densities we can confidently assign values for the adjacent layer tunneling amplitude, $|t_{22}| = 30 \ \mu eV$ and the disorder-broadening energies $\gamma_{t(b)} = 4$ meV. Note that the measured disorder broadening parameter $\gamma_{t(b)}$ is much larger than T_0 , where T_0 is the temperature at which the tunneling conductance appears to diverge experimentally. It is immediately clear therefore, even before performing a detailed analysis, that the condensation temperature must be strongly suppressed by disorder.

(2) The colossal enhancement of the tunneling conductance has been observed in the wide density range $4 \times 10^{10} - 10^{12}$ cm⁻². The strength of the interactions in the system decreases with doping, and we have chosen the elevated doping level $n = 7.4 \times 10^{11}$ cm⁻² because of the detailed experimental data available in this case. This doping level corresponds to the average concentration of electrons and holes $n = (n_e + n_h)/2$, while the electron-hole imbalance $\Delta n \ll n$ can be present.

(3) The transport experiments with similar double bilayer graphene samples and with similar gating geometry have already been performed [62]. The doping level of charge carriers $\epsilon_{\rm F}$ and the gap 2|u| in the electronic spectrum of bilayer graphene are independent and are controlled by the same gate. At the considered average density $n = 7.4 \times 10^{11}$ cm⁻², the effective mass parameter of bilayer graphene can be approximated as $m \approx 0.04 m_0$, while the gap is equal to $2|u| \approx 6.6$ meV and is much smaller than the corresponding Fermi energy $\epsilon_{\rm F} \approx 20$ meV. It follows that $|u|/\epsilon_{\rm F} \approx 0.16$, which

implies sublayer polarization within the bilayers $c^2 \approx 0.58$ and $s^2 \approx 0.42$ to be modest.

(4) The bare critical temperature without disorder $\bar{T}_0 \approx$ 50 K can be recalculated from the actual critical temperature at the considered doping level $T_0 = 1.5$ K and the Cooper pair-breaking rate $\gamma = \gamma_t + \gamma_b = 8$ meV, which has been chosen above with help of Eq. (15). The system is in the regime of strong pair breaking $T_0 \ll \bar{T}_0 \sim \gamma$, and the value $\gamma/T_0 \approx 1.74$ is very close to the critical value 1.78. Experimentally, singular behavior of the tunneling conductance is observed only in the cleanest samples.

(5) The bare critical temperature T_0 of Cooper pairing in the weak-coupling regime is given by $\overline{T}_0 = 2e^{C} \Lambda \exp[-1/\lambda_0]/\pi$, where C = 0.577 is the Euler constant. Here $\lambda_0 = \lambda_0^{\rm s} + \lambda_0^{\rm d}$ corresponds to the mixed channel 12-21 that dominates in the considered regime $|u| \ll \epsilon_{\rm F}$. Approximating the high energy cutoff as $\epsilon_{\rm c} \approx 2\epsilon_{\rm F}$ and employing the values of \overline{T}_0 and $\epsilon_{\rm F}$ chosen above we get $\lambda_0 \approx 0.44$. The applicability condition for the weak-coupling approach $\lambda_0 \ll 1$ is not well satisfied, but this value for λ_0 corresponds to moderate coupling regime $\lambda_0 \lesssim 1$, which justifies the approximations used in our theory.

(6) The relative contribution to the tunneling conductance $G_{\rm T}$ of channels 12-21 and 22 can be characterized by a dimensionless parameter $r = |t_{12} - t_{21}|^2 / |t_{22}|^2$ that we treat in the phenomenological way. Along with the sample area A, r and A are only free parameters of the model that have not yet been assigned. We adjust them by fitting the measured temperature dependence for tunneling conductance at zero voltage bias V = 0 (and also at matched concentrations of electrons and holes and zero in-plane magnetic field), which is presented in Fig. 1, with the theory that incorporates the effect of fluctuating electron-hole Cooper pairs [Eq. (27)]. The area A can be be obtained by matching the high-temperature behavior of $G_{\rm T}$ since the contribution of 12-21 in this case is negligible small, resulting in $A \approx 397 \ \mu m^2$. The value of $r \approx 7.6 \times 10^{-5}$ is obtained by fitting the singular behavior of the tunneling conductance in the vicinity of the critical temperature. The corresponding theoretical curve is also presented in Fig. 1 and matches the experimental data reasonably well over a wide temperature range.

In Fig. 1 we also present calculations within the model of noninteracting electrons and holes [Eq. (26)]. This model severely underestimates the tunneling conductance G_T in the case of matched electron and hole concentrations, where interactions are crucial to explain enhanced tunneling conductance at low temperature and singular behavior in the vicinity of the critical temperature T_0 for Cooper pair condensation.

We will keep all parameters chosen above in further calculations and investigate an impact of finite voltage bias between bilayers V, electron-hole imbalance Δn , and in-plane magnetic field B at tunneling conductance. We will also present only results of the theory that incorporates the effect of fluctuating Cooper pairs.

A comparison between theory and experiment for the voltage dependence of the tunneling conductance between graphene bilayers $G_{\rm T}$ is presented in Fig. 6. The zero bias peaks emerge with a decreasing temperature on a top of a smooth background that corresponds to the channel 22. The width of the background $eV_{\rm HM}$ is governed by the single-particle energy scales $2\pi T$ and γ and is approximately equal



FIG. 6. Dependence of the tunneling conductance $G_{\rm T}$ on bias voltage V at different temperatures T. The concentrations of electrons and holes do match. The solid curves correspond to calculations, while the dotted lines are experimental data [Fig. S2(b) in Ref. [10]]. For clarity the curves at adjacent temperatures are offset by $\Delta G_{\rm T} = 12$ mS. The theory that incorporates the effect of fluctuating Cooper pairs [Eq. (27)] reasonably fits temperature dependence of the peak height but overestimates its width and does not capture its asymmetry. Possible origins of the asymmetry are discussed in Sec. VI.

to the largest of them. The width of the zero bias peak in the vicinity of T_0 is much smaller than the single-particle disorder scale, which demonstrates its collective origin. While the temperature dependence of peak height is well fit by the theory, the width dependence is captured only qualitatively and is overestimated by a factor of 2. The experimental data also exhibit a voltage asymmetry that becomes more prominent at low temperatures. Within our phenomenological model an asymmetrical voltage dependence of the tunneling conductance can be obtained if the scattering rates of electrons and holes $\gamma_{t(b)}$ that define the Cooper pair-scattering rate as $\gamma = \gamma_t + \gamma_b$ are energy dependent. The simple linear dependence $\gamma = \gamma + \gamma' \omega$ with a phenomenological parameter γ' does not capture the observed asymmetry, however. A quantitative understanding of the asymmetry requires a microscopic understanding of disorder mechanisms that is outside the scope of the present work.

The comparison between theory and experiment for the dependence of the tunneling conductance $G_{\rm T}$ on the electronhole imbalance Δn and the voltage bias V are summarized in Fig. 7, and a zero-bias-voltage cut of the comparison made in this color plot at zero voltage bias is presented in Fig. 8. The theoretical curves again agree reasonably well with the data. A density imbalance splits the Fermi lines of the electrons and holes and disfavors their Cooper pairing. As seen in Fig. 5, the FFLO state with a finite Cooper pair momentum cannot be stabilized in the strong pair-breaking regime realized in the experiment. The critical temperature T_0 is maximal for Cooper pairs with zero Cooper pair momentum and decreases monotonically in the presence of imbalance and vanishes if the latter exceeds the critical temperature. As a result, the dependence of tunneling conductance on imbalance Δn and the in-plane magnetic filed B that is presented in Fig. 9 is smooth and featureless. Figure 9 presents theoretical curves



FIG. 7. Dependence of the tunneling conductance $G_{\rm T}$ on voltage bias V and electron-hole density imbalance Δn . Their average density $n = (n_{\rm e} + n_{\rm h})/2 = 7.4 \times 10^{11} \,{\rm cm}^{-2}$ is fixed, and the temperature is 3.5 K. The top subplot (a) corresponds to experiment [Fig. 5(c) in Ref. [10]], and the bottom one (b) to theory. The cut of this plot at zero voltage bias V = 0 is presented as Fig. 8.

since the experimental data for magnetic field and densitybalance dependence at this temperature are not yet available. The theory qualitatively explains the decrease of peak height with a magnetic field studied experimentally at lower temperatures $T_0 \approx 1.5$ K but considerably overestimates its effect. In this case the system is in the paired state whose behavior lies outside the range of validity of the present theory. Other less fundamental limitations might also explain this discrepancy as we discuss in more detail in the next section. We conclude that our theory of the fluctuational internal Josephson effect, combined with specific features of the multiple-channel structure of pairing in double bilayer graphene provides a reasonable overall description of experiment.

VI. DISCUSSION

Our theory of the internal fluctuational Josephson effect does not account for interactions between fluctuating Cooper pairs. The Gaussian nature of the theory we employ is more clearly seen within an alternate derivation of the tunneling conductance. The latter employs the auxiliary field approach and is presented in Appendix C. Interactions between fluctuating pairs can be safely omitted in the wide range of temperatures $\Delta T \sim T_0$ and are important only in the critical region $\Delta T_{Gi} \lesssim Gi T_0$ where fluctuations are large and strongly interfere with each other. Here $Gi = T_0/E_F$ is the Ginzburg number calculated in Appendix D. The double



FIG. 8. The dependence of tunneling conductance $G_{\rm T}$ at zero voltage bias V = 0 on the electron-hole density imbalance Δn . Their average density $n = (n_{\rm e} + n_{\rm h})/2 = 7.4 \times 10^{11} \text{ cm}^{-2}$ is fixed, and the temperature is 3.5 K. The solid curve is theory, and the dotted curve is experimental data extracted from the V = 0 line from Fig. 5(c) in Ref. [10]. The theory captures the curve profile reasonably well but does not capture the asymmetry. Possible origins of the asymmetry are discussed in Sec. VI.

bilayer graphene system studied experimentally [10] is close to the weak coupling regime, and the critical region $\Delta T_{\text{Gi}} = T_0^2/E_{\text{F}} \approx 10 \text{ mK}$ is much smaller than the temperature range $\Delta T \approx 4 \text{ K}$ where the tunneling conductance is strongly enhanced. The picture of noninteracting fluctuating Cooper pairs is therefore well justified to address the basic phenomena identified in experiment.

Fluctuating Cooper pairs in conventional superconductors alter the thermodynamics of the normal state only in the critical region ΔT_{Gi} . The conductivity and the magnetic susceptibility [43] not only aresingular at the critical temperature, however, but have long high-temperature tails [45]. The high-temperature tail for the diamagnetic susceptibility $\chi \sim$ $\chi_{\rm L}/\ln^2[T/T_0]$, where $\chi_{\rm L}$ is the Landau diamagnetic susceptibility in the normal state, was predicted [63–65] theoretically and observed in experiments [45]. The reason is the paired state is superconducting and mediates the perfect diamagnetism that makes even a small number of fluctuating Cooper pairs important. Similarly the equilibrium-paired state of spatially separated electrons and holes provides a fluctuational internal Josephson effect that colossally enhances the interlayer tunneling. That is why even a small density of fluctuating Cooper pairs can make a strong impact on the tunneling conductance above the critical temperature T_0 which also has the high-temperature tail proportional to $1/\ln^2[T/T_0]$, which is clearly seen in Fig. 1.

In the vicinity of the critical temperature T_0 zero bias peak shape is universal and governed by the factor $F(eV, \mathbf{Q})$ given by Eq. (29). It can be rewritten as $F(eV, \mathbf{Q}) =$ $\text{Im}[L_0^{-1}(eV, \mathbf{Q})]/eV$. Thus an imaginary part of the Cooper propagator $L_0^{-1}(eV, \mathbf{Q})$, which can be interpreted as Cooper pair susceptibility [66], is directly probed in tunneling experiments [67]. It should be noted that Cooper pair susceptibility of a superconductor can also be probed in tunneling Josephson junction in which one side is near its critical temperature while the other is well below its critical temperature [68,69]. The



FIG. 9. Theoretical dependence of tunneling conductance $G_{\rm T}$ on in-plane magnetic *B* and electron-hole density imbalance Δn . Their average density $n = (n_{\rm e} + n_{\rm h})/2 = 7.4 \times 10^{11} {\rm cm}^{-2}$ is fixed and the temperature is 3.5 K. The dependence of $G_{\rm T}$ is smooth and featureless, which demonstrates for the set of parameters used for fitting (the regime of strong pair breaking $\gamma \gtrsim T_0$) the FFLO state with finite Cooper pair momentum is not stabilized by the density imbalance Δn .

junction does not support dissipationless Josephson tunneling current, but the tunneling current at finite voltage is strongly enhanced by fluctuating Cooper pairs that grow in the vicinity of T_0 . The latter has been observed experimentally [70,71].

Only the enhancement of intravalley tunneling that corresponds to the alignment ($\theta = 0$) has been reported so far. In this case the instability happens in one channel, while the main contribution to the tunneling conductance comes form the different one. As a result the divergent contribution to the conductance due to fluctuating Cooper pairs appears on the top of background with weak temperature dependence that dominates at higher temperatures. We predict an enhancement of intervalley tunneling ($\theta = \pi/3$) to be much more profound because Cooper pairs in the dominating pairing channel involve electrons and holes settled at adjacent sublayers. We present calculations for this case in Appendix B.

The Fulde-Ferrell-Larkin-Ovchinnikov (FFLO) state with finite Cooper pair momentum was predicted in conventional superconductors more than 60 years ago [54,55]. It requires the splitting of Fermi surfaces/lines for pairing electrons with opposite spins. The splitting can be induced by a magnetic field provided that its paramagnetic effect is larger than its diamagnetic one (Chandrasekhar-Clogston limit [72,73]). This condition is rarely satisfied even in layered conventional superconductors subjected to an in-plane magnetic field. There are few observations in heavy-fermion and organic superconductors where FFLO state signatures have been claimed but are still debated. (See Ref. [74] and Refs. [75,76] for reviews of progress in solid state and cold atom systems.) So far the FFLO state has not been unambiguously identified. In double bilayer graphene the densities of electrons and holes can be controlled separately in a way that opens the FFLO state up for experimental study in a condensed matter system [5,6]. The FFLO can be unambiguously identified if it appears from the dependence of the zero-bias peak on imbalance and an inplane magnetic field, since the latter makes it possible to probe Cooper pairs with finite momentum. In the vicinity of the instability to the uniform paired state, the tunneling conductance monotonically decreases with the in-plane magnetic field (as presented in Fig. 9). In the vicinity of an instability to the FFLO state the tunneling conductance achieves a maximum at finite field-induced momentum shift $\mathbf{Q}_{\rm B} = \mathbf{Q}_0$ where \mathbf{Q}_0 is the corresponding momentum of Cooper pairs. We discuss how to distinguish these states in more detail in Appendix B, where calculations for the intervalley tunneling are presented.

The sensitivity of Cooper pairing to a disorder opens the possibility of a granular electron-hole state in the presence of its strong long-range variations. In this state the pairing happens in disconnected or weakly coupled regions with a minimal amount of disorder and does not support the spatial coherence. It makes the transport properties of the system including Coulomb drag effect to be different from ones in the uniform paired state. A tunneling conductance in the granular state is still colossally enhanced since the latter requires temporal coherence of Cooper pairs but not the spatial one.

The interpretation of experiment provided by our theory suggested that the pairing critical temperature would be substantial if samples with weaker disorder could be fabricated. This finding is perhaps a bit surprising since the experiments are for the most part conducted in the weak to moderate coupling regime ($r_s = 0.72-3.3$) where some researchers have argued that critical temperatures should be strongly suppressed by screening, especially accounting for spin and valley degeneracy [17,77] (see also arguments that this approach considerably underestimates the critical temperature [15,16,19,20]). Our theory also suggests that high pairing temperatures should be achievable in double singlelayer graphene systems, since there is nothing in its structure that puts single layers at a disadvantage relative to bilayer. Moreover, for the linear spectrum in monolaver graphene the Wigner-Seitz radius is defined in a different way, $r_s = 2.19/\kappa$, and could achieve even larger value $r_s = 1.1$ if hBN is used as a spacer between graphene sheets. Here κ is the corresponding dielectric constant. Future experimental work which seeks to weaken pair breaking by disorder and which explores double single-layer graphene systems as well is therefore important.

In summary, the theory of the fluctuational internal Josephson effect developed here explains the anomalies in the tunneling conductance between graphene bilayers observed experimentally at equal electron and hole densities, including their dependence on temperature, bias voltage bias, and electron-hole imbalance. Some aspects of the observations are nevertheless not understood. First, the observed dependence of the tunneling conductance on bias voltage has an asymmetry between positive and negative bias that becomes more prominent with decreasing temperature. At first glance the asymmetry is unexpected and surprising since the electronic spectrum of two graphene bilayers with matched concentrations of electrons and holes is symmetric, as is clearly seen in Fig. 2. The symmetry can be broken by Coulomb impurities if most of them are of the same charge. For example, positive charges (ionized donors) provide repulsive scattering for holes and attractive scattering for electrons. Our model takes the scattering rates for electrons and holes $\gamma_{t(b)}$ to be momentum and energy independent and ignores these common complications. An asymmetry can be introduced in a phenomenological way by making the assumption that the Cooper pair-breaking time is energy dependent $\gamma(\omega)$. Approximating it by a linear function does predict an asymmetry of the tunneling conductance that grows with decreasing temperature, but the shape of the experimental curves is not captured by this simple ansatz. To clarify whether or not the observed asymmetry can be explained by the presence of charge impurities, a more microscopic description of their scattering characteristics is needed, and this is outside of the scope of the present work. Second, it is not clear whether or not our theory can capture the dependence of tunneling conductance on magnetic field since more experimental data are needed. Comparison with data obtained at $T \approx 1.5$ K suggests that the theory considerably overestimates the effect of a magnetic field. Nevertheless, at such temperatures the system is in the paired state or in the critical regime, which is outside of the applicability range of the theory of Gaussian fluctuations. This discrepancy can be due to other reasons. Bilayer graphene along with other two-dimensional systems have long-range density variations, and if the corresponding length is smaller than $\hbar/Q_{\rm B}$, the effect of the magnetic cannot be reduced just to the relative shift of dispersion for electrons from different layers. To better understand capabilities of the theory more experimental data are needed.

To conclude, we have developed a theory of the fluctuational internal Josephson effect in the system of two closely spaced graphene bilayers. The presence of valley and sublattice degrees of freedom provides three competing electronhole channels for both intravalley and intervalley Cooper pairs. We show that three channels are nearly independent and have different critical temperatures of the condensation and sublattice structures. The observed enhancement of the tunneling conductance can be explained only by the presence of competing channels that dominate in different temperature ranges. The theory reasonably captures the dependence of the conductance on temperature, voltage bias between bilayers, and electron-hole imbalance. We also argue that the enhancement is much stronger for intervalley tunneling than for the intravalley one that has been reported recently. We also discuss how to distinguish the uniform state and the FFLO state with finite Cooper pair momentum that can be stabilized in the system by an electron-hole imbalance.

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APPENDIX A: BETHE-SALPETHER EQUATION

Here we present a detailed derivation of Eq. (6). The Bethe-Saltpeter equation that is illustrated in Fig. 3 can be written as

$$\Gamma^{\sigma_i'\sigma_t}_{\sigma_b'\sigma_b}(ip_n, \mathbf{p}', \mathbf{p}) = U_{\mathbf{p}'-\mathbf{p}}\delta^{\sigma_i'\sigma_t}_{\sigma_b'\sigma_b} + T\sum_{i\omega_n\mathbf{p}''} U_{\mathbf{p}'-\mathbf{p}''}G^{\mathsf{t}}_{\sigma_i'\sigma_i''}(ip_n+i\omega_n, \mathbf{p}''_+)G^{\mathsf{b}}_{\sigma_b''\sigma_b'}(i\omega_n, \mathbf{p}''_-)\Gamma^{\sigma_i''\sigma_t}_{\sigma_b''\sigma_b}(ip_n, \mathbf{p}'', \mathbf{p}).$$
(A1)

Here $\mathbf{p}_{\pm} = \mathbf{p} \pm \mathbf{q}/2$; $\omega_n = (2n+1)\pi/T$ and $p_n = 2n\pi/T$ are fermionic and bosonic Matsubara frequencies, respectively. The electron Green function $\hat{G}^{t(b)}(\omega_n, \mathbf{p})$ in the sublattice space can be presented as

$$G_{\sigma'\sigma}^{t}(i\bar{\omega}_{n}^{t},\mathbf{p}) = \frac{\langle \sigma'|\mathbf{tcp}\rangle\langle\mathbf{tcp}|\sigma\rangle}{i\omega_{n} - \epsilon_{\mathbf{cp}} + i\gamma_{t}\mathrm{sgn}[\omega_{n}]}, \quad G_{\sigma'\sigma}^{b}(i\omega_{n},\mathbf{p}) = \frac{\langle \sigma'|\mathbf{bvp}\rangle\langle\mathbf{bvp}|\sigma\rangle}{i\omega_{n} - \epsilon_{\mathbf{vp}} + i\gamma_{b}\mathrm{sgn}[\omega_{n}]}, \tag{A2}$$

where $\gamma_{t(b)}$ are scattering rates for electrons (holes). We have also neglected the presence of the valence band in the layer with excess of electrons and the presence of the conduction band in the layer with the excess of holes. The product of Green functions that appears in (A1) can be written with as

$$G_{\sigma_i'\sigma_i''}^{\mathsf{t}}(ip_n+i\omega_n,\mathbf{p}_+)G_{\sigma_b''\sigma_b'}^{\mathsf{b}}(i\omega_n,\mathbf{p}_-)=M_{\sigma_b'\sigma_b''}^{\sigma_i'\sigma_t''}(\mathbf{p}'')C(ip_n,\mathbf{q},i\omega_n,\mathbf{p}),$$

where the matrix form factor

$$M^{\sigma'_{t}\sigma''_{t}}_{\sigma'_{b}\sigma''_{b}}(\mathbf{p}) = \langle \sigma'_{t} | tc \mathbf{p}_{+} \rangle \langle tc \mathbf{p}_{+} | \sigma''_{t} \rangle \langle \sigma''_{b} | tv \mathbf{p}_{-} \rangle \langle bv \mathbf{p}_{-} | \sigma'_{b} \rangle$$

reflects the chiral nature of charge carriers in bilayer graphene, while $C(ip_n, \mathbf{q}, i\omega_n, \mathbf{p})$ contains information only about the energy spectrum for electrons and holes and is given by

$$C(ip_n, \mathbf{q}, i\omega_n, \mathbf{p}) = \frac{1}{(i\omega_n + ip_n - \epsilon_{c\mathbf{p}_+} + i\gamma_t \text{sgn}[\omega_n + p_n])(i\omega_n - \epsilon_{v\mathbf{p}_-} + i\gamma_b \text{sgn}[\omega_n])}.$$
(A3)

In the weak coupling regime that we consider in the paper, pairing correlations do appear in the vicinity of Fermi lines for electrons and holes. As a result the vertex $\hat{\Gamma}(ip_n, \phi_{\mathbf{p}'}, \phi_{\mathbf{p}})$, the Fourier transform of interactions $U(\phi_{\mathbf{p}'} - \phi_{\mathbf{p}})$, and the form factor $\hat{M}(\phi_{\mathbf{p}})$ can be safely approximated by their values at the Fermi level ($|\mathbf{p}| = p_F$ and $|\mathbf{p}'| = p_F$) and depend only on the corresponding polar angles ($\phi_{\mathbf{p}}$ and $\phi_{\mathbf{p}'}$). After decomposition over multipole momenta $\hat{\Gamma}_{l',l}$, U_l , and \hat{M}_l Eq. (A1) becomes algebraic and can be presented in a compact form

$$\hat{\Gamma}_{l'l} = U_{l'} \delta_{l'l} \hat{1} + \sum_{l_1 l_2} U_{l'} M_{l'-l_1-l_2} \Pi_{l_2} (ip_n, \mathbf{q}) \hat{\Gamma}_{l_2, l}, \qquad (A4)$$

$$\Pi_l(ip_n, \mathbf{q}) = T \sum_{i\omega_n, \mathbf{p}} e^{-il\phi_{\mathbf{p}}} C(ip_n, \mathbf{q}, i\omega_n, \mathbf{p})$$

Its *s*-wave component $\Pi_0(ip_n, \mathbf{q}) \equiv \Pi(ip_n, \mathbf{q})$ has the ultraviolet logarithmic divergence that is usually present in the weak coupling pairing theories and coincides with the singlestep pair propagator in the Cooper ladder sum of a bilayer system without sublattice degrees of freedom. Its detailed derivation can be found in textbooks [43,44], and its explicit expression is presented in the main part as Eq. (7). Therefor all information about the chiral nature of the bilayer graphene charge carriers is hidden in the nontrivial matrix form factor \hat{M}_l . At finite *l* the value $\Pi_l(ip_n, \mathbf{q})$ is nonzero only at finite Cooper pair momentum and is much smaller than $\Pi_0(ip_n, \mathbf{q})$ and can be neglected. As a result, Eq. (A4) reduces to Eq. (6).

APPENDIX B: INTERVALLEY ALIGNMENT AND IDENTIFICATION OF THE FFLO STATE

In the main text calculations for intravalley tunneling ($\theta =$ 0) are discussed, while ones for the intervalley alignment ($\theta =$ $\pi/3$) are presented here. The tunneling conductance is dominated by the channel 11-22 because of fluctuating Cooper pairs in the dominating channel involve correlations at adjacent sublayers. The tunneling conductance can be approximated as Eq. (28). We use the same set of parameters that have been used above to fit the experimental data except for pairbreaking rate γ . For the latter we use $\gamma = 2, 4$, and 8 meV. The first two values correspond to cleaner samples compared to ones that have been studied experimentally [10]. The temperature dependence of tunneling conductance is presented in Fig. 10. Its enhancement of the conductance is considerably stronger than that for the intervalley tunneling. The pair-breaking rate γ determines the critical temperature of pair condensation but weakly influences the temperature dependence of tunneling conductance above it.



FIG. 10. The temperature dependence of the intervalley tunneling conductance $G_{\rm T}$ between graphene bilayers at zero voltage bias (V = 0). Three curves correspond to pair-breaking rates $\gamma = 2$, 4, and 8 meV. The pair-breaking rate induced by scattering at impurities reduces the critical temperature T_0 of electron-hole condensation but weakly affects the critical behavior above T_0 .



FIG. 11. The dependence of tunneling conductance $G_{\rm T}$ on inplane magnetic *B* and the electron-hole density imbalance Δn . Three subplots correspond to T = 45 K (a), 25 K (b), and 5 K (c). The dashed line corresponds to the phase boundary of equilibrium electron-hole paired state. In (b) and (c) the conductance achieves a maximum at a finite value of magnetic field *B*, which demonstrates that fluctuating Cooper pairs with finite momentum $\mathbf{Q} \approx \mathbf{Q}_{\rm B}$ are the most intensive and the system is in the vicinity of the instability to the FFLO state. The presence of a kink in the phase boundary in (c) clearly demonstrates that the system is unstable towards the equilibrium FFLO state.

For the pair-breaking rate $\gamma = 2$ meV the critical temperature is $T_0 \approx 42$ K. According to the phase diagram (Fig. 5) the ratio $\gamma/T_0 \approx 0.52$ is small enough to stabilize the Fulde-Ferrell-Larkin-Ovchinnikov (FFLO) state by the electron-hole imbalance. The in-plane magnetic field results in a relative shift of electronic dispersions between layers and makes it possible to probe fluctuating Cooper pairs with finite momentum $Q_{\rm B} = e dB_{||}/\hbar c$. The dependence of tunneling conductance at zero voltage bias on the electron-hole imbalance and magnetic field at temperature T = 55 K is presented in Fig. 11(a). The tunneling conductance has a monotonic dependence demonstrating that the system is far from the instability to the FFLO state, and Cooper pair fluctuations with zero momentum $\mathbf{Q} = 0$ are the most intensive. The dependence at T = 25 K is shown in Fig. 11(b) with a dashed line that denotes a phase boundary of a paired state. Since the critical imbalance required to suppress the pairing instability decreases with the magnetic field the paired state is the uniform BCS one. Nevertheless, the tunneling conductance achieves a maximum at the finite value of the magnetic field, which shows that the system is close to the FFLO state and fluctuations with finite Cooper momentum are the most intensive. The dependence at T = 5 K is presented in Fig. 11(c). The dashed line that denotes the phase boundary is nonmonotonous and has a kink at the magnetic field finite value. It clearly demonstrates that the equilibrium FFLO state is stabilized by the electronhole imbalance.

APPENDIX C: BOSONIC PICTURE OF THE FLUCTUATIONAL INTERNAL JOSEPHSON EFFECT

Here we present a bosonic picture of the fluctuational internal Josephson effect that can be employed with a help of the field integral formalism. We consider the contact interactions U between electrons and holes that correspond to the truncation of all multipole momenta of interactions except the s-wave one. In the main text we have demonstrated that they are unimportant. The correlation function of tunneling operators $\chi(\omega, \mathbf{q})$ that defines the tunneling conductance $G_{\rm T}$ according to Eq. (18) can be extracted from the corresponding imaginary time correlation function $\chi(\tau, \mathbf{r})$ given by

$$\chi(\tau, \mathbf{r}) = \frac{1}{Z} \frac{\delta^2 Z}{\delta \Lambda_{\tau, \mathbf{r}} \delta \Lambda_{0, 0}^+},$$
(C1)

where $Z[\Lambda_{\tau,\mathbf{r}}^+, \Lambda_{\tau,\mathbf{r}}]$ is the statistical sum with an auxiliary bosonic field $\Lambda_{\tau,\mathbf{r}}$ introduced to the action *S* of the system as

$$S = \int_{0}^{\beta} d\tau \int \mathbf{dr} \Big[\hat{\psi}_{t}^{+} (\partial_{\tau} + \hat{h}_{t\mathbf{r}} - \mu_{t}) \hat{\psi}_{t} + \hat{\psi}_{b}^{+} (\partial_{\tau} + \hat{h}_{b\mathbf{r}} + \mu_{b}) \hat{\psi}_{b} + \hat{\psi}_{b}^{+} \Lambda^{+} \hat{t}^{+} \hat{\psi}_{t} + \hat{\psi}_{t}^{+} \hat{t} \Lambda \hat{\psi}_{b} + U \psi_{ts_{t}}^{+} \psi_{bs_{b}}^{+} \psi_{ts_{t}} \Big],$$

where $\hat{\psi}_t \equiv \hat{\psi}_{t\tau \mathbf{r}}$ and $\hat{\psi}_b \equiv \hat{\psi}_{b\tau \mathbf{r}}$ are spinor fermionic fields for electrons from top (t) and bottom (b) layers, respectively, with labeling described in Sec. II. If they are integrated out and the corresponding action is expanded in the lowest order in tunneling matrix elements \hat{t} , the Fourier transform χ_q appears in the action as $S = -\sum_q \chi_q |\Lambda_q|^2$. Here $q = \{q_n, \mathbf{q}\}$ with bosonic Matsubara frequency $q_n = 2\pi T n$. For noninteracting electrons and holes, calculations are straightforward and result in

$$\chi_q^0 = \hat{t}^+ \hat{M}_0 \Pi_q \, \hat{t}. \tag{C2}$$

After analytical continuation we get Eq. (20). In the case of interacting electrons and holes it is instructive to start with the Hubbard-Stratonovich transformation. It eliminates interactions but introduces the bosonic field $\hat{\Delta}_{\tau \mathbf{r}}$ corresponding to electron-hole Cooper pairs. The action *S* is modified as

$$S = \int_{0}^{\beta} d\tau \int \mathbf{dr} \bigg[\hat{\psi}_{t}^{+} (\partial_{\tau} + \hat{h}_{tr} - \mu_{t}) \hat{\psi}_{t} + \hat{\psi}_{b}^{+} (\partial_{\tau} + \hat{h}_{br} + \mu_{b}) \hat{\psi}_{b} + \hat{\psi}_{b}^{+} \hat{\Delta}'^{+} \hat{\psi}_{t} + \hat{\psi}_{t}^{+} \hat{\Delta}' \hat{\psi}_{b} + \frac{1}{U} \text{tr} (\hat{\Delta}^{+} \hat{\Delta}) \bigg], \qquad (C3)$$

where $\hat{\Delta}'_{\tau \mathbf{r}} = \hat{\Delta}_{\tau \mathbf{r}} + \hat{t} \Lambda_{\tau \mathbf{r}}$. Above the critical temperature T_0 of the electron-hole pairing the saddle point of the action is trivial $\langle \hat{\Delta} \rangle = 0$, and the field $\hat{\Delta}$ corresponds to Cooper pair fluctuations [86]. In the wide temperature range $\Delta T_{Gi} \lesssim \Delta T \sim T_0$ outside the critical regime $\Delta T \lesssim \Delta T_{Gi}$ fluctuations can be approximated by the noninteracting Gaussian theory. Here $\Delta T_{Gi} = Gi T_0$ with Ginzburg number $Gi = T_0/E_F$ calculated in Appendix D. Integrating out fermions and expanding the action up to the second order in the bosonic field $\hat{\Delta}'$ results in

$$S = \sum_{q} \left[\frac{\hat{\Delta}_{q}^{+} \hat{\Delta}_{q}}{U} - \hat{\Delta}_{q}^{\prime +} \hat{M}_{0} \Pi_{q} \hat{\Delta}_{q}^{\prime} \right] = \sum_{q} \left[\hat{\Delta}_{q}^{+} \hat{\Gamma}_{q}^{-1} \hat{\Delta}_{q} - \hat{\Delta}_{q}^{+} \hat{M}_{0} \Pi_{q} t \Lambda_{q} - \Lambda_{q}^{+} \hat{t}^{+} \hat{M}_{0} \Pi_{q} \hat{\Delta}_{q} - \Lambda_{q}^{+} \chi_{q}^{0} \Lambda_{q} \right].$$
(C4)

Here all matrices are in the compact representation, and $\hat{\Gamma}_q$ is the scattering vertex calculated in Sec. III. Some of its components vanish at the critical temperature T_0 of the electron-hole Cooper pairing. The last term corresponds to the response χ_q^0 function for noninteracting electrons and holes that is given by Eq. (C2). The action represents the bosonic picture of the Josephson effect and is valid outside the weak coupling regime. The action (C4) is quadratic in the bosonic field Δ_q , and after its integration we get the tunneling response function χ_q given by

$$\chi_q = \chi_q^0 + \hat{t}^+ \hat{M}_0 \Pi_q \hat{\Gamma}_q^{-1} \hat{M}_0 \Pi_q \hat{t}.$$
 (C5)

With the help of Eqs. (12) and (11) we get the tunneling response function for the intravalley tunneling (23). In the

same way with the help of Eqs. (14) and (13) we get the response function for the intervalley one (25).

APPENDIX D: GINZBURG CRITERION

The developed theory of the fluctuational internal Josephson effect implies that Cooper pair fluctuations are Gaussian and do not interact with each other. The interactions can be safely omitted in the wide range of temperatures $\Delta T \sim$ T_0 except the critical region $\Delta T \leq \Delta T_{\text{Gi}}$ where fluctuations are overgrown are strongly interfere with each other. The range ΔT_{Gi} can be estimated from the Ginzburg criterion [87], which compares the contribution of Gaussian fluctuations to the heat capacity $C_{\text{FL}} = T_0/4\pi\xi^2\Delta T$ with predictions of the mean-field theory below the critical temperature

 $Gi = T_0/E_{\rm F}$. It does not depend explicitly on the pair-breaking rate γ but only on the critical temperature T_0 . It should be noted that the Ginzburg criterion can be derived microscopically in a more strict way by explicit analysis of the role of grows with decrease of temperatures and dominates in interactions between fluctuating Cooper pairs [88]. the temperature range $\Delta T_{Gi} = GiT_0$ with Ginzburg number

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spectrum becomes nonparabolic and has the Mexican-hat shape. This enhances the role of interactions and is very important in the strong coupling regime [51].

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