# Robustness of Majorana fermions in proximity-induced superconductors

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In two-dimensional chiral *p*-wave superconductors, the zero-energy Majorana fermion excitations trapped at vortex cores are protected from the thermal effects by the minigap,  $\Delta^2/\epsilon_F$  ( $\Delta$ : bulk gap,  $\epsilon_F$ : Fermi energy), which is the excitation gap to the higher energy bound states in the vortex cores. Robustness to thermal effects is guaranteed only when  $T \ll \Delta^2/\epsilon_F \sim 0.1$  mK, which is a very severe experimental constraint. Here we show that when *s*-wave superconductivity is proximity-induced on the surface of a topological insulator or a spin-orbit-coupled semiconductor, as has been recently suggested, the minigaps of the resultant non-Abelian states can be orders of magnitude larger than in a chiral *p*-wave superconductor. Specifically, for interfaces with sufficient barrier transparencies, the minigap can be as high as  $\sim \Delta \gg \Delta^2/\epsilon_F$ , where  $\Delta$  is the bulk gap of the *s*-wave superconductor.

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#### I. INTRODUCTION

Topological quantum computation (TQC) requires the existence of a two-dimensional (2D) topologically ordered state whose lowest energy excitations follow non-Abelian statistics.<sup>1</sup> If the appropriate many-body ground-state wave function-e.g., Pfaffian states in fractional quantum systems<sup>1,2</sup> and chiral Hall *p*-wave  $(p_x + ip_y)$ superconductor/superfluid<sup>3,4</sup>—is a linear combination of states from a degenerate subspace, then a pairwise exchange of the particle coordinates can unitarily rotate the wave function in the degenerate subspace. This exact non-Abelian statistical property can be used to perform quantum gate operations, which are, in principle, fault tolerant.1,5,6 More importantly, these non-Abelian particles, the Majorana fermions, being half-fermions, are new particles in nature distinct from ordinary Dirac fermions, which are of obvious intrinsic fundamental interest.<sup>7</sup>

In practice, a key requirement for TQC is that the degenerate ground-state subspace must be separated from the other excited states by a nonzero energy gap so that thermal effects cannot hybridize the topological quasiparticle states with the other higher energy, nontopological, states in the system.<sup>1</sup> In 2D  $p_x + ip_y$  superconductors (SC), such as possibly SrRuO<sub>4</sub>, where the zero-energy Majorana fermion excitations trapped in the vortex cores are the topological quasiparticle states, this gap is given by the so-called minigap,  $\sim \delta_0 \sim \Delta^2 / \epsilon_F$ , where  $\Delta$  is the bulk superconducting gap and  $\epsilon_F$  is the Fermi energy.<sup>8</sup> Since  $\delta_0 < 0.1$  mK is a very small energy scale for typical *p*-wave superconductors, the requirement  $T \ll \delta_0$  constitutes the real bottle neck for TOC, even if the best possible 2D  $p_x + ip_y$  superconductor-based platform were realized in the laboratory. Here we show that, in a class of newly proposed TQC platforms, involving Majorana fermions in multilayer structures where s-wave superconductivity is proximity induced on a host topological insulator (TI) (Refs. 9 and 10) or a spin-orbit-coupled semiconductor,<sup>11</sup> the minigap can be enhanced by several orders of magnitude. Given that a strong proximity effect in such superconductorsemiconductor structures has already been experimentally demonstrated,<sup>12,13</sup> it is realistic to decrease *T* to satisfy  $T \ll \delta_0$ , since  $\delta_0$  can be made as high as 1 K (i.e., a fraction of order unity of  $\Delta$ ), which is the bulk gap in the *s*-wave superconductor.

To derive these results, we explicitly analyze the microscopic model of the proximity effect between a TI surface and an s-wave superconductor by applying the conventional tunneling formalism.<sup>14</sup> We find that, in addition to the superconducting gap  $\Delta$ , the interface transparency (denoted by  $\lambda$ below) given by the interlayer tunneling amplitude controls the strength of the proximity effect on the TI surface. Our main result is that for barriers with transparency satisfying  $\epsilon_F \gg \lambda \gtrsim U, \Delta$ , where U is the Fermi level on the TI surface, the excitation gap above the non-Abelian quasiparticle states can be  $\sim \Delta \gg \Delta^2 / \epsilon_F$ . We show this by applying our central result, Eq. (13), on the excitation gaps in the two most important structures on a TI surface, a line junction [Eq. (16)] and a vortex [Eq. (19)], which contain Majorana modes. Note that, as discussed earlier,<sup>9</sup> the Majorana modes in a line junction and a vortex are the two most essential elements of a putative TQC architecture on the TI surface. The dramatic increase in the excitation gap on the TI surface greatly enhances the robustness of the topological quasiparticles to thermal decoherence effects, which may bring non-Abelian statistics and TQC to the realm of realistic, achievable, temperature regimes in the laboratory.

The paper is organized as follows. In Sec. II, we describe the microscopic model we consider for the proximity effect at a TI-SC interface. In Sec. III, we derive the proximityinduced effective pair potential with renormalized parameters. In Secs. IV and V we described the line-junction geometry that may be used to manipulate superconducting quasiparticles such as Majorana fermions at a TI-SC interface. In Sec. VI, we show that the minigap of a vortex formed at a TI-SC interface can be orders of magnitude larger than in a vortex in an intrinsic (not proximity-induced) superconductor such as SrRuO<sub>4</sub>.

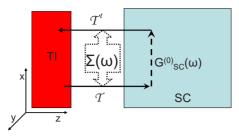


FIG. 1. (Color online) Proximity induced pairing on the TI surface. The (red) region on the left is the TI, and the (blue) region on the right is an *s*-wave superconductor. "Integrating out" the superconducting degrees of freedom produces the self-energy  $\Sigma$  on the TI surface, where  $\Sigma$  is given by the tunneling Hamiltonian  $\mathcal{T}$  and the Green's function  $G_{\rm SC}^{(0)}$  of the superconductor (see text for details).

## **II. MICROSCOPIC MODEL FOR PROXIMITY EFFECT**

An interesting property of superconductors is that they can induce superconductivity in a normal metal in contact with the superconductor.<sup>15</sup> This is referred to as the superconducting proximity effect. As we will show in the rest of the text, the superconducting proximity effect allows a greater degree of control of the superconducting quasiparticle spectrum, than is possible by simply modifying the superconductor, where the superconductivity is intrinsically derived from the quasiparticle spectrum itself. The superconducting properties of a normal-superconductor (NS) interface can be described by the self-consistent Bogoliubov-de Gennes (BdG) equations at the interface.<sup>15,16</sup>

The BdG equations at an NS interface can be written in terms of a Nambu spinor wave function,  $\Psi(\mathbf{r}) = [u_{\uparrow}(\mathbf{r}), u_{\downarrow}(\mathbf{r}), v_{\downarrow}(\mathbf{r}), -v_{\uparrow}(\mathbf{r})]^T$  which is finite on both the superconductor and normal side of the interface. In order to write these equations more compactly it is convenient to introduce the Nambu matrices  $\tau_{x,y,z}$ , which are identical to the Pauli spin matrices  $\sigma_{x,y,z}$ , except that they operate on the (u,v) part of the spinor  $\Psi(\mathbf{r})$ . Thus the spinor  $\Psi(\mathbf{r})$  exists in the tensor product space  $\sigma_{\alpha} \otimes \tau_{\beta}$ . The BdG equations for the quasiparticle wave functions can be written in terms of the  $\sigma \otimes \tau$  matrices in the form  $H_{\text{BdG}}\Psi(\mathbf{r}) = E\Psi(\mathbf{r})$ , where  $H_{\text{BdG}}$  is a  $4 \times 4$  BdG Hamiltonian.

The normal-superconductor interface can be considered to be a planar geometry (Fig. 1) with the coordinates r=(r,z), where r and z are the in-plane [r=(x,y)] and out-of-plane coordinates for the interface (the NS interface is at z=0). In this paper, we will restrict ourselves to the case where the SC in our system is a conventional s-wave superconductor such as Al, and the normal part of our system is the metallic surface state of a TI such as Bi<sub>2</sub>Te<sub>3</sub>.<sup>17-19</sup> The BdG Hamiltonian describing the s-wave superconductor is of the form

$$H_{\rm SC} = \left(-\frac{\boldsymbol{\nabla}_r^2 + \partial_z^2}{2m^*} - \boldsymbol{\varepsilon}_F\right) \boldsymbol{\tau}_z + \boldsymbol{\Delta}_s(r) \boldsymbol{\tau}_x,\tag{1}$$

where the  $m^*$  is the effective mass of the superconductor,  $\varepsilon_F$  is the Fermi energy of the superconductor, and  $\Delta_s(r)$  is the pairing potential in the superconductor. The pairing potential  $\Delta_s(r)$  inside the *s*-wave superconductor is generated by a pairing interaction V(r) that can be taken to be pointlike.<sup>15</sup> In

the self-consistent BdG equations, the pairing potential satisfies the constraint

$$\Delta_s(\boldsymbol{r}) = V(\boldsymbol{r}) \sum_n u_{n,\uparrow}^*(\boldsymbol{r}) v_{n,\downarrow}(\boldsymbol{r}) [1 - 2f(E_n)], \qquad (2)$$

where  $(u_n, v_n)$  are components of the four-spinor eigenvectors  $\Psi_n(\mathbf{r})$  which satisfy the BdG equations  $H_{\rm SC}\Psi_n(\mathbf{r}) = E_n\Psi_n(\mathbf{r})$ . Here  $f(E_n)$  is the Fermi function.

The normal part of the TI-SC system, which in our case consists of the metallic surface band<sup>17–19</sup> of a TI (in anticipation of the proximity-induced superconductivity) can be described by a BdG Hamiltonian

$$H_{\rm TI} = [\iota v \,\boldsymbol{\sigma} \cdot \boldsymbol{\nabla}_r - U] \otimes \tau_z, \tag{3}$$

where  $U = \varepsilon_F - \int dz |\phi(z)|^2 V_{gate}(z)$  is the Fermi level of the TI surface where  $V_{gate}(z)$  is the gate potential and  $\phi(z)$  is the *z*-dependent electron wave function (with momenta close to the Dirac point) of the TI surface states. Here *v* is the effective electron velocity on the TI surface. Note that apart from the factor  $\tau_z$ , this is the Dirac-Hamiltonian describing the surface state of the topological insulator in the normal state. The factor  $\tau_z$  is present to account for the independent electron part (represented by *u* in the spinor  $\Psi$ ) and hole part (represented by *v* in the spinor  $\Psi$ ).

Since there is no tunneling so far between the TI and SC, the TI may be described as being normal with no superconductivity in the TI. The introduction of a tunneling term Twhich transfers electrons between the SC and TI leads to a finite value for the order parameter  $\langle \psi_{\sigma}(r)\psi_{\sigma'}(r')\rangle$  on the TI surface.<sup>15</sup> However it is crucial to note that despite the existence of an order parameter in the normal part of the system (TI), the microscopic pairing potential  $\Delta_{\text{TI}}(r)=0$ . This follows from the self-consistency condition in the TI

$$\Delta_{\mathrm{TI}}(\mathbf{r}) = V_{\mathrm{TI}}(\mathbf{r}) \langle \psi_{\uparrow}(\mathbf{r})\psi_{\downarrow}(\mathbf{r}) \rangle = 0 \tag{4}$$

since the attractive pairing interaction in the TI vanishes  $[V_{\text{TI}}(\mathbf{r})=0]$ . This is consistent with the de Gennes boundary conditions at the interface which requires  $\Delta(\mathbf{r})/N(\mathbf{r})V(\mathbf{r})$  to be continuous across the TI-SC interface.<sup>15</sup> Here  $\Delta(\mathbf{r})$ ,  $V(\mathbf{r})$ , and  $N(\mathbf{r})$  stand for the microscopic pairing potential, pairing interaction and density of states at the Fermi level on both sides of the interface. Note that since  $\Delta_{\text{TI}}$  and  $V_{\text{TI}}$  are both zero on the TI side of the interface, the ratio can be finite which allows  $\frac{\Delta_{\text{TI}}}{V_{\text{TI}}N_{\text{TI}}} = \frac{\Delta_s}{V_s N_s}$ . However, even though there is no microscopic pair potential on the TI side, the superconducting pair potential. In the next few paragraphs, we show explicitly how the superconductor can be integrated out to yield an effective Hamiltonian on the TI surface which has an effective pairing potential.

## III. EFFECTIVE PAIRING POTENTIAL ON THE TI SURFACE

The complete BdG Hamiltonian describing the TI-SC interface including the tunneling term<sup>14</sup> (Fig. 1) is defined by the Hamiltonian:  $H_{\text{total}} = H_{\text{TI}} + H_{\text{SC}} + \mathcal{T} + \mathcal{T}^{\dagger}$ . Here,  $H_{\text{TI}}$  and  $H_{\text{SC}}$ are the Hamiltonians describing the TI surface and the *s*-wave superconductor, respectively.  $\mathcal{T}$  describes the tunneling from the TI surface to the superconductor and  $\mathcal{T}^{\dagger}$  describes the tunneling in the opposite direction. The excitation spectrum of the interface can be determined from the BdG equation

$$(H_{\text{total}} - E)\Psi(\mathbf{r}) = 0, \qquad (5)$$

where  $\Psi(\mathbf{r})$  is the Nambu spinor, and  $H_{\text{total}}$  is written as a  $4 \times 4$  matrix in the Nambu basis.

The tunneling Hamiltonian  $H_t = T + T^{\dagger}$  coupling the 2D TI surface states with the superconductor can be explicitly written in the Nambu space as

$$\mathcal{T}(r;r'z') = \tau_z \int d^2 \mathbf{k} dk_z \chi(z';\mathbf{k}k_z) \langle \chi(\mathbf{k}) | \mathcal{T} | \phi \rangle e^{i\mathbf{k} \cdot (r-r')}.$$
 (6)

Here the momenta are measured relative to the Dirac cone momentum M and the tunneling matrix element in the integrand can be approximately written as<sup>20</sup>

$$\langle \chi(\boldsymbol{k}) | \mathcal{T} | \boldsymbol{\phi} \rangle = \frac{i}{m} [ \boldsymbol{\phi}(z) \partial_z \chi(z; \boldsymbol{k}, k_z) - \chi(z; \boldsymbol{k}, k_z) \partial_z \boldsymbol{\phi}(z) ] |_{z=0},$$

where,  $\chi(z; \mathbf{k}, k_z)$  is the single-particle eigenfunction in the superconductor and  $\phi(z)$  is as defined before.

In order to solve the BdG equation at the TI-SC interface, we decompose the wave function as  $\Psi = \psi_{TI} + \psi_{SC}$ . Decomposing the BdG equation [Eq. (5)] we obtain

$$(H_{\rm TI} - E)\psi_{\rm TI} + \mathcal{T}^{\dagger}\psi_{\rm SC} = 0, \qquad (7)$$

$$(H_{\rm SC} - E)\psi_{\rm SC} = -\mathcal{T}\psi_{\rm TI}.$$
(8)

Substituting the wave function  $\psi_{SC}$  from Eq. (8) in Eq. (7) we get the effective BdG equation on the TI surface

$$[H_{\rm TI} + \Sigma(rr';\omega) - \omega]\psi_{\rm TI} = 0.$$
(9)

Here the self-energy  $\Sigma$  on the TI surface (Fig. 1) is given by

$$\Sigma(\boldsymbol{r}\boldsymbol{r}',\omega) = -\int d\boldsymbol{r}_1 d\boldsymbol{r}_2 \mathcal{T}^{\dagger}(\boldsymbol{r},\boldsymbol{r}_1) G_{\rm SC}^{(0)}(\boldsymbol{r}_1,\boldsymbol{r}_2;\omega) \mathcal{T}(\boldsymbol{r}_2,\boldsymbol{r}'), \tag{10}$$

where  $G_{SC}^{(0)}(\mathbf{r}_1, \mathbf{r}_2; \omega) = (H_{SC} - \omega)^{-1}$  is the Green function matrix in the superconductor.

The effective superconducting pairing potential induced at the surface of a TI appears in the form of the anomalous part of the self-energy,  $\Sigma$ , which can be written as

$$\Sigma_A(\boldsymbol{r}\boldsymbol{r}',\omega) = -\int d\boldsymbol{r}_1 d\boldsymbol{r}_2 \mathcal{T}^{\dagger}(\boldsymbol{r},\boldsymbol{r}_1) F_{\rm SC}^{(0)}(\boldsymbol{r}_1,\boldsymbol{r}_2;\omega) \mathcal{T}(\boldsymbol{r}_2,\boldsymbol{r}'),$$
(11)

where  $F_{SC}^{(0)}(\mathbf{r}_1, \mathbf{r}_2; \omega) = \int dt e^{i\omega t} \langle T\psi(\mathbf{r}_1, 0)\psi(\mathbf{r}_2, t) \rangle$  is the anomalous part of the Green function in the superconductor that represents the superconducting order parameter. Since the anomalous part of the self-energy is proportional to the anomalous part of the Green function, the complex phase of  $\Sigma_A$  must also equal the complex phase of the superconductor.

The Green's function for the superconductor can be written in terms of the normal state eigenbasis of the superconductor as

$$G_{\rm SC}^{(0)}(\boldsymbol{rr'};\omega) = \sum_{n} \chi_n(\boldsymbol{r})\chi_n(\boldsymbol{r'})[(\boldsymbol{\epsilon}_n - \boldsymbol{\varepsilon}_F)\boldsymbol{\tau}_z + \Delta\boldsymbol{\tau}_x - \omega]^{-1},$$
(12)

where the normal state eigenfunctions  $\chi_n$  are taken to be real and spin independent. Using Eqs. (6) and (12) in Eq. (10) and then Fourier transforming to the momentum space the self-energy on the TI surface takes the form

$$\Sigma(rr';\omega) = -\sum_{n} \chi_{n}(\boldsymbol{r}_{1})\chi_{n}(\boldsymbol{r}_{2})\mathcal{T}^{\dagger}(r,\boldsymbol{r}_{1})\mathcal{T}(\boldsymbol{r}_{2},r')\frac{\omega\tau_{0} + (\boldsymbol{\epsilon}_{n} - \boldsymbol{\varepsilon}_{F})\tau_{z} + \Delta\tau_{x}}{(\boldsymbol{\epsilon}_{n} - \boldsymbol{\varepsilon}_{F})^{2} + \Delta^{2} - \omega^{2}}.$$
(13)

This equation is more conveniently written in terms of the tunneling density of states  $\Lambda$  on the TI as

$$\Sigma(rr';\omega) = -\int d\epsilon \frac{\omega\tau_0 + \epsilon\tau_z + \Delta\tau_x}{\epsilon^2 + \Delta^2 - \omega^2} \Lambda(rr';\epsilon), \quad (14)$$

where the tunneling density of states  $\Lambda$  from the TI to the SC relative to the Fermi level is given by

$$\Lambda(rr';\epsilon) = \sum_{n} \delta(\epsilon - \epsilon_n + \epsilon) \chi_n(\mathbf{r}_1) \chi_n(\mathbf{r}_2) \mathcal{T}^{\dagger}(r,\mathbf{r}_1) \mathcal{T}(\mathbf{r}_2,r').$$
(15)

Note that the matrix  $\Lambda$  is the operator that can be used to calculate the time a state  $\phi(x)$  on the TI would take to diffuse

into the SC layer when in the normal state through the expression  $\tau^{-1} = \int dr dr' \phi(r) \phi^*(r') \Lambda(rr'; \epsilon)$ .

The scale of the spatial and energy dependence of the matrix  $\Lambda(rr'; \epsilon)$  can be understood by observing that it is dependent only on the wave functions  $\chi_n$  of the superconductor in the normal state at the energy  $\epsilon$ . Defining the Fourier transform of  $\Lambda(rr'; \epsilon)$  as  $\Lambda(r,k;\epsilon) = \int dr' \Lambda(r+r'/2, r - r'/2;\epsilon)e^{-ikr'}$ , similar to the quasiclassical approximation, we obtain a function in position r and wave-vector k space. Given that the original definition of  $\Lambda$  involved only the normal state band structure of the superconductor,  $\Lambda(rk;\epsilon)$  varies on a scale  $k \sim k_F$  and  $\epsilon \sim \epsilon_F$ . Since we are interested in slow variations  $\xi^{-1} \ll k_F$  and energies  $U, \Delta \ll \epsilon_F$ , we can ignore the k and  $\epsilon$  dependence of  $\Lambda$  and take it to be a function of only r. Thus we will assume  $\Lambda(rk;\epsilon) \approx \lambda(r) = \Lambda(rk_F;\epsilon_F)$ .

Fourier transforming back to real space leads to the real space relation  $\Lambda(rr'; \epsilon) \sim \lambda(r) \delta(r-r')$  which is local in r, r' and independent of  $\epsilon$ . Within this approximation we find

$$\Sigma(r,\omega) \approx \lambda(r) \frac{(-\omega\tau_0 + \Delta\tau_x)}{\sqrt{|\Delta|^2 - \omega^2}}.$$
(16)

Using  $H_{\text{TI}}$  from Eq. (2) and the local (*k*-independent) selfenergy from Eq. (16), we can now straightforwardly rewrite Eq. (9) as an effective BdG equation for the TI surface

$$[\tilde{v}(\omega)\iota\boldsymbol{\sigma}\cdot\boldsymbol{\nabla}\tau_{z}-\tilde{U}(\omega)\tau_{z}+\tilde{\Delta}(\omega)\tau_{x}-\omega]\psi_{\mathrm{TI}}=0, \quad (17)$$

 $\widetilde{U}(\omega) = Z(\omega)U,$  $\widetilde{\Delta}(\omega)$ where  $\widetilde{v}(\omega) = Z(\omega)v,$ and  $=\lambda\Delta/(\sqrt{\Delta^2-\omega^2}+\lambda).$ Here, the factor  $Z(\omega)$  $=\sqrt{\Delta^2-\omega^2}/(\sqrt{\Delta^2-\omega^2}+\lambda)$ .  $\tilde{v}(\omega)$ ,  $\tilde{U}(\omega)$ , and  $\tilde{\Delta}(\omega)$  are the renormalized velocity, Fermi level, and superconducting gap on the TI surface, respectively. Below we will be interested only in the low-energy states with energies  $E \ll \Delta$  (typically  $0.1\Delta < E < 0.5\Delta$ ). In this case, we can approximate the frequency-dependent parameters in Eq. (17) with their values at  $\omega = 0$ 

$$\widetilde{v}(\omega) \approx \frac{v}{1 + \frac{\lambda}{\Delta}} = \widetilde{v}, \quad \widetilde{U}(\omega) \approx \frac{U}{1 + \frac{\lambda}{\Delta}} = \widetilde{U},$$
$$\widetilde{\Delta}(\omega) \approx \frac{\lambda}{1 + \frac{\lambda}{\Delta}} = \widetilde{\Delta}.$$
(18)

The renormalization of the parameters described in Eq. (18)gives the central results of this paper which can be understood as arising from the virtual propagation of the electron in the superconductor. This is consistent with the estimate of the fraction of time spent in the TI, which is  $t_{TI}/(t_{TI}+t_{SC})$  $\sim (1 + \lambda/\Delta)^{-1} = Z(0)$ , since we can estimate life-time of an electron in the TI as  $t_{\text{TI}} \sim \lambda^{-1}$  and the virtual excitation of an electron in the SC as  $t_{\text{SC}} \sim \Delta^{-1}$ . Thus, in addition to inducing a pairing potential  $\tilde{\Delta}(\omega)$  on the TI surface, the proximity effect also renormalizes the velocity on the surface of the TI to  $v \rightarrow \tilde{v}(\omega) = Z(\omega)v$  and the background potential to U  $\rightarrow U(\omega) = Z(\omega)U$ . One can notice that the TI electron density *n* is proportional to  $(\tilde{U}/\tilde{v})^2$  and remains constant as expected. In the weak tunneling regime  $\lambda \ll \varepsilon_F$ , our tunneling matrix approach to the proximity effect is consistent with previous calculations<sup>21</sup> for the superconductor-semiconductor system using the Blonder-Tinkham-Klapwijk model.<sup>16</sup> Since the parameter  $\lambda$  is determined by purely electronic energy scales,  $\lambda$ can be larger than  $\Delta$  and the retardation effects discussed above lead to substantial renormalizations of the original parameters.

The expression for the proximity-induced self-energy, Eq. (18), was derived for a single superconducting island in contact with a TI surface. The proximity-induced pairing was then localized to the region where  $\lambda(r)$  was nonzero. In the case of different superconducting islands covering different parts of the TI surface, one must add the self-energy contribution from each of the separate islands. The pair-potential  $\Delta$ 

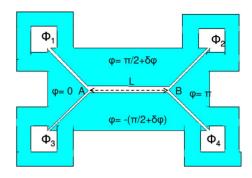


FIG. 2. (Color online) A trijunction-pair geometry of superconducting islands deposited on the TI surface (top view) to confine and manipulate Majorana fermions. The superconducting islands are connected strong superconducting loops enclosing fluxes  $\Phi_{n,m}$ . The values of the fluxes  $\Phi_{1,2}=\Phi_{3,4}=\frac{\pi}{2}+\delta\phi$ ,  $\Phi_{2,3}=\frac{\pi}{2}-\delta\phi$ , and  $\Phi_{4,1}=-3\frac{\pi}{2}-\delta\phi$  which satisfy  $\Sigma_n\Phi_n=0$  control the phases of the superconducting island. For the given superconducting phase configuration and  $\delta\phi=\frac{\pi}{6}$  the structure contains a vortex with a trapped Majorana state only on trijunction A. By changing  $\delta\phi$  to  $-\frac{\pi}{6}$ , the discrete vortex together with the Majorana state is transferred to the trijunction B. The Majorana state formed on the one-dimensional line junction (of length *L*) connecting A and B in the intermediate stage with  $\delta\phi=0$ .

is then taken to be a constant for each island but varies between different islands. In particular, a region of the TI which is in contact with a superconducting island [through a nonzero  $\lambda(r)$ ] inherits a superconducting pairing potential with a phase which has the same value as the island it is in contact with. Furthermore, the effective pair potential, vanishes [since  $\lambda(x)=0$ ] if the TI surface is not locally in contact with the superconductor (for example, the TI regions in between two superconducting islands has vanishing pair potential). Below we will apply the formulas contained in Eq. (18) to estimate the excitation gaps above the Majorana topological excitations that have been discussed for the TI surface.<sup>9</sup>

#### IV. MAJORANA SYSTEM ON TI SURFACE

The proximity effect discussed in the previous sections allows one to induce superconductivity on a semiconducting substrate such as a TI surface state. This leads to the possibility of manipulating the position and energies of superconducting quasiparticles on such a surface state. One of the most interesting scenarios that this kind of manipulation of superconducting quasiparticles might lead to is the possibility of creating and braiding Majorana fermions at such TI/SC interfaces.<sup>9</sup>

The trijunction geometry shown in Fig. 2 provides a way to create and braid a Majorana fermions on the TI surface: two trijunctions (A and B in Fig. 2) of superconducting layers with distinct local phases separated by a line junction of length *L*. The superconducting islands can be considered to have distinct phases for barrier transparencies much less than unity corresponding to  $\lambda \ll \epsilon_F$ . The line-junction separating the superconducting islands consists of a region of width *W* which is not in direct contact with either of the superconducting islands. Therefore the effective pairing potential vanishes in this region and, in principle, the line junction can support multiple low-energy transverse modes with energy much less than  $\Delta$ . This can be avoided by taking the junction width to satisfy  $W \sim \tilde{v}/\tilde{\Delta} = v/\lambda$  such that the confinement energy  $\tilde{v}/W$  creates a sufficiently large energy spacing between the various transverse modes. For the transparencies assumed in the paper  $\Delta \ll \lambda \ll \epsilon_F$ , the width W, which is required to be smaller than the effective coherence length is much smaller than the coherence length of the superconductor in the clean limit but much larger than the Fermi wavelength of the superconductor. Ideally, one should adjust the barrier transparency to set  $\lambda$  to the lowest value such that the line-junctions width W is within the resolution of the fabrication technology.

Such a geometry can be used to trap zero-energy Majorana fermions on the TI surface at discrete vortices formed by the phase configurations of the superconducting islands either at A or B. By varying the phase differences between the islands it is also possible to move the discrete vortex from A to B and thus transport the bound Majorana state. The phase differences between neighboring superconducting islands is controlled by connecting the islands by superconducting loops with fluxes threaded through them.

For barrier transparencies between the TI and the SC, such that  $\lambda \gg \Delta$ , but still much less than unity (which implies  $\lambda \ll \epsilon_F$ ), the Josephson current between the islands mediated by the single channel TI surface state splits up into a small current density in the bulklike superconducting loops. This adds only small gradients in phase to the gradient resulting from the vector potential of the magnetic flux. Thus the difference in the phases on the islands are controlled by the applied fluxes according to the equation

$$\phi_i - \phi_k = \Phi_{ik}.\tag{19}$$

Therefore, for the application of the fluxes shown in Fig. 2 leads to the desired phase configuration on the islands.

The trijunction setup discussed in the previous paragraphs operates on the principle that if one considers a single linejunction separating two superconductors (such as the one joining A and B) with a phase difference of  $\pi$ , there are exactly one pair of zero energy and zero momentum Majorana states trapped between the islands. At finite but small momenta, one can use lowest order  $k \cdot p$  perturbation theory to check that these modes disperse into a pair of linearly dispersing nonchiral Majorana modes with equal and opposite velocities. Off-setting the phase difference from  $\pi$  by  $\delta\phi$ leads to a mass gap in the linear Dirac spectrum of this pair of Majorana modes. Thus the effective Hamiltonian describing such a line-junction is given by

$$H_{junc} = -\imath \frac{\tilde{\upsilon} \tilde{\Delta}^2}{(\tilde{U}^2 + \tilde{\Delta}^2)} \sigma_z \nabla + m(\delta \phi) \sigma_x, \qquad (20)$$

where  $\sigma$  here is an effective pseudospin degree of freedom that keeps track of the mode index.<sup>9</sup> The term proportional to *m* is an effective phase-difference-dependent mass term. Following the Jackiw-Rebbi index theorem<sup>22</sup> it is clear that a localized zero-energy solution results at any point where the sign of the mass  $m\delta\phi$  changes sign. Such changes in signs can be engineered at the ends A and B of the line junction joining A and B by appropriately tuning the fluxes. As it turns out, the appropriate mass changes are generated at the ends A and B whenever there are discrete vortices in the phases at A and B.<sup>9</sup>

With the configuration of the phases on the superconducting islands as shown in Fig. 2, and for  $\delta \phi = \frac{\pi}{6}$ , the total phase change around the trijunction A is  $2\pi$ . Trijunction A then acts as a discrete vortex and there is a localized zero-energy Majorana state confined to A. In this configuration, there is no vortex or zero-energy mode at B. It can be easily checked from Fig. 2 that the roles of A and B are reversed if  $\delta \phi = -\frac{\pi}{6}$ : now B contains a vortex and a localized Majorana mode while A is topologically trivial. In both cases, the spectrum of the line junction connecting A and B is gapped with the excitation gap controlled by  $\delta \phi$ . To avoid hybridization of the localized states at A and B, the length *L* must exceed the size of the localized states themselves

$$L > \xi \sim \tilde{\nu} / \tilde{\Delta}, \tag{21}$$

where  $\xi$  is the decay length of the Majorana states on the TI surface.

It is now clear that the Majorana states trapped at the discrete vortices can be braided by tuning the phase  $\delta\phi$  through zero. For  $\delta\phi=0$  the phase change across the line junction is  $\pi$ , and for the arrangement of the phases as shown in Fig. 2, there is a single zero-energy *extended* Majorana mode on the line junction. When  $\delta\phi$  is tuned from  $\frac{\pi}{6}$  to 0 to  $-\frac{\pi}{6}$ , the Majorana mode shifts from A to the line junction and finally to B. For  $\delta\phi=0$ , the other low-energy delocalized modes on the line junction follow a dispersion given by<sup>9</sup>

$$\omega(q) \approx \pm q \tilde{\upsilon} \tilde{\Delta}^2 / (\tilde{U}^2 + \tilde{\Delta}^2).$$
(22)

Below we will consider two types of excitation gaps which control the thermal robustness of the above Majorana system. First, in the line junction of length L the gap  $E_g$  $\approx \frac{\tilde{v}}{L} \tilde{\Delta}^2 / (\tilde{U}^2 + \tilde{\Delta}^2)$  that follows from Eq. (22) protects the *de*localized zero-energy Majorana mode from thermal decoherence.  $E_g$  controls the thermal robustness of the Majorana fermions while they are braided in TQC. We show below by explicit analytic arguments that it is possible to make  $E_g$  $\sim \Delta$  by appropriately designing the TI-SC interface. The thermal robustness of the (stationary) topological qubits themselves, on the other hand, is determined by the energy gap ( $\Delta E$ ) above the zero-energy *localized* Majorana states within the discrete vortex cores. We will show by rigorous numerical calculations that even this scale  $\Delta E \sim \Delta$ , making the entire TQC architecture surprisingly robust to thermal decoherence effects.

#### **V. EXCITATION GAP IN LINE JUNCTION**

For a line junction of length L, the gap  $E_g$  is given by [see Eq. (22)]  $E_g \approx \frac{\tilde{v}}{L} \tilde{\Delta}^2 / (\tilde{U}^2 + \tilde{\Delta}^2)$ . Now, for barriers with transparency such that  $\lambda \gg U, \Delta$ , we get  $\tilde{\Delta} \gg \tilde{U}$  [Eq. (18)] and the

factor multiplying  $\tilde{v}/L$  in  $E_g$  reduces to unity. Even if  $U \sim \lambda$ , which should be possible experimentally, this factor is still of order unity. To maximize  $E_g$ , we need to take the minimum allowed value of the length of the line junction,  $L_m \sim \tilde{v}/\tilde{\Delta}$  [Eq. (21)]. Therefore, the maximum  $E_g$  attainable on the TI surface is given by

$$E_g \approx \frac{\tilde{v}}{L_m} = \tilde{\Delta} = \frac{\lambda}{1 + \frac{\lambda}{\Lambda}},$$
 (23)

which, in the case of transparency  $\lambda \gg \Delta$ , reduces to  $\Delta$  itself.

### VI. EXCITATION GAP IN VORTEX

To determine the excitation gap  $\Delta E$  within a vortex core numerically, we consider the BdG Hamiltonian on the surface of a TI sphere with a vortex and an antivortex at the poles<sup>23</sup>

$$H = [\tilde{v}\hat{R} \cdot (\sigma \times p) - \tilde{U}]\tau_z + \tilde{\Delta}(\mathbf{r})\tau_x, \qquad (24)$$

which can be written in angular coordinates as

$$H = \left[ -\frac{\tilde{v}}{R} L \cdot \sigma - \tilde{U} \right] \tau_z + \tilde{\Delta}(\theta) \{ \cos \phi \tau_x + \sin \phi \tau_y \}.$$
(25)

 $\widetilde{\Delta}(\theta)$ Here R is the radius of the sphere,  $=\Delta \tanh\{R \sin \theta / \xi_v\}$  and  $\xi_v$  is the size of the vortex core. In the above Hamiltonian, we have approximated discrete vortices by regular ones with continuously varying phases. The resultant azimuthal symmetry allows us to decouple the equations into sets indexed by m with a basis of spinor spherical harmonics of the form  $(Y_{l,m+1}, Y_{l,m+2}, Y_{l,m}, Y_{l,m+1})^T(\theta, \phi)$ . We expect the minigap of such a continuous vortex to be qualitatively similar to the discrete vortex in Fig. 2. We find that the m=-1 channel contains a pair of decaying and oscillating solutions which are spatially localized at the two poles. The corresponding eigenenergies exponentially decay to zero with the radius of the sphere, indicating that, in the limit when the vortices are far-separated, the eigenenergies are exactly zero. On the other hand, the spectrum of the other m channels qualitatively resembles the m=-1 channel with the important difference that the eigenenergy of the first pair of excited states does not vanish as the radius of the sphere increases. This eigenenergy gives us the excitation gap in the vortex core.

Assuming the vortex core size to be equal to  $\xi_v = \xi = \tilde{v}/\Delta$ = $v/\lambda$ , the numerical results for the minigap (Fig. 3) can be fit by the analytic form

$$\Delta E \approx 0.83 \tilde{\Delta}^2 / \sqrt{\tilde{\Delta}^2 + \tilde{U}^2}.$$
 (26)

The above result is interesting since it approaches the value  $\Delta E \sim \tilde{\Delta}$  (avoiding the approximation  $\Delta E \ll \Delta$  in Eq. (18) yields  $\Delta E \approx 0.6\Delta$  instead of  $0.8\Delta$ ) for  $\tilde{U} \sim \tilde{\Delta}$ . Such a large minigap is made possible for proximity-induced superconductors because of the ability to tune the chemical potential on the TI surface independent of the chemical potential of

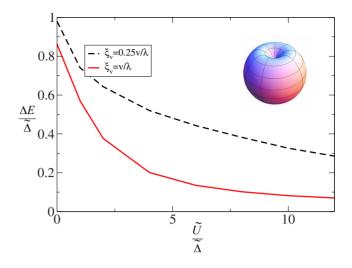


FIG. 3. (Color online) Numerical results for the vortex minigap  $\Delta E$  plotted against the renormalized Fermi level  $\tilde{U}$  on the TI surface ( $\Delta E$ ,  $\tilde{U}$  scaled by  $\tilde{\Delta}$ ). The solid (red) line gives the mini-gap when the vortex core size  $\xi_v = \xi = \tilde{v}/\tilde{\Delta} = v/\lambda$ , as is appropriate in a regular vortex with a continuously varying phase. The dashed (black) line shows that the excitation gap is even larger when the vortex core size is smaller, as is expected in a discrete vortex (see Fig. 1). The inset shows the TI sphere with a vortex and an antivortex (with reduced superconducting amplitudes at the vortex cores) situated at the north and the south poles.

the superconductor. This is unlike the case of more conventional *s*-wave and chiral *p*-wave superconductors where the chemical potential  $\tilde{U} \gg \tilde{\Delta}$  and the minigap obeys the scaling

$$\Delta E \sim 0.83 \frac{\tilde{\Delta}^2}{\tilde{U}} \tag{27}$$

similar to the classic result of Caroli, de Gennes, and Matricon.<sup>8</sup>

In fact this is expected since the scaling of Caroli, de Gennes, and Matricon can be established using fairly generic semiclassical arguments by considering the core of the vortex to be a normal region of radius  $\xi$  with no superconductivity which can host quasiparticle states close to the Fermi level  $\epsilon_{F}$ . Considering a quasiparticle state at zero angular momentum that is confined in the vortex core, the state at the next allowed angular momentum would have a relative energy  $v_F(|k|-k_F) = v_F(\sqrt{k_F^2 + k_t^2} - k_F)$ . Here  $k_t \sim \xi^{-1}$  is the transverse momentum from the angular momentum. For  $k_t \ll k_F$ , this leads to the energy splitting  $v_F k_t^2 / k_F \sim v_F / k_F \xi^2$ . For a vortex, the quasiparticles are confined within a decay length of  $\xi = v_F / \Delta$ . The leads to the relation  $\delta_0 \sim \Delta^2 / \epsilon_F$ . Since this argument applies to any system with weak superconductivity, it also holds for the TI/SC system as verified by our numerics.

Thus the staggering difference between the minigap estimate in a chiral *p*-wave superconductor  $\delta_0 = 10^{-5}$  meV  $\sim 0.1$  mK and the estimate of 1 K for the proximity-induced system arises entirely from our ability to control the Fermi energy  $\epsilon_F$  in the system. Lowering  $\epsilon_F$  lowers  $k_F$  which leads

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to a lower density of quasiparticle states at the Fermi level which in turn reduces the number of midgap states in the vortex core.

Of course, our analysis of the minigap in proximity induced superconductors only accounts for possible quasiparticle states in the vortex core that were localized at the TI surface. Therefore the geometry must ensure that there are no quasiparticles in the superconductor itself. This can be accomplished by using a discrete vortex as in the trijunction geometry. We believe that the estimates for the continuous vortex carry over to the discrete vortex case.

So far the vortex core size has been taken to be on the order of the coherence length in the TI,  $\xi_v = \xi = \tilde{v} / \tilde{\Delta} = v / \lambda$ . If the vortex core size is taken to be smaller, as is expected for a discrete vortex, the numerical calculations lead to an even larger  $\Delta E$  (Fig. 3). As is clear from Eq. (27), for  $\tilde{U} \leq \tilde{\Delta}$ , the excitation gap in a vortex can be of order  $\tilde{\Delta}$ , which is  $\sim \Delta$  for chemical potential on the surface of the TI tuned such that  $U \leq \lambda$  [see Eq. (18)] which can be of the scale of 100 meV (i.e., of the scale set by  $\epsilon_F$ ). The constraint  $U \leq \Delta$  $\sim 0.5$  meV within the Dirac point, that is obtained without the central result Eq. (18), is difficult to achieve in experiments because of impurity-induced density fluctuations as in graphene.<sup>24</sup> The ability to tune the chemical potential U on the TI surface independent of the Fermi energy of the superconductor  $\epsilon_F$  leads to a significant enhancement over the case of a chiral p-wave superconductor where the chemical potential  $U \sim \epsilon_F \gg \Delta$  such that the minigap from Eq. (27) scales as  $\Delta E \sim \Delta^2 / \epsilon_F$  consistent with previous estimates.<sup>8</sup>

## VII. CONCLUSION

In conclusion, we have shown that Majorana fermion excitations in proximity-induced s-wave superconducting systems are much more robust to thermal decoherence effects than in regular chiral *p*-wave superconductors. In the latter system, the excitation gap protecting the Majorana modes, the so-called excitation gap, is given by  $\sim \Delta^2 / \epsilon_F$ , which is a prohibitively low-energy scale  $\sim 0.1$  mK. On the other hand, for proximity-induced s-wave superconducting systems,<sup>9,11</sup> which have generated a lot of recent interest, 25-28 and in the case of sufficient transparency of the barriers, the minigap can be made as high as  $\sim \Delta \sim 1$  K. The possible orders of magnitude enhancement of the minigap in these systems helps bring the observation of non-Abelian statistics to the realm of realistic, accessible, temperature regimes in experiments. Thus, the proposal of Fu-Kane<sup>9</sup> and that of Sau et al.<sup>11</sup> with appropriate control of the proximity effect and feature sizes appear at this state to provide the most robust platforms for the observation of Majorana fermions and the implementation of TQC.

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