

Nodal arc of disordered Dirac fermions and non-Hermitian band theory

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We show that Dirac fermion systems in two dimensions generally exhibit a disorder-induced nodal arc replacing the nodal point and tilted Dirac cone, provided that the two components of the Dirac fermion correspond to two distinct orbitals unrelated by symmetry. This result is explicitly demonstrated using renormalization-group analysis in a disordered Dirac model that we introduce, where the disorder potential acts differently on the two orbitals. As we show by numerical simulations and self-consistent Born approximation calculation, this drives the system into a new strongly disordered phase.

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Disordered Dirac fermions in two dimensions have been studied for decades in a variety of contexts [1–3], including the integer quantum Hall transition [4] and disordered unconventional superconductivity [5–9]. The interest in this field has been reinforced since the discovery of graphene and topological insulators [10–21]. In general, massless Dirac fermions in solids arise from various types of internal degrees of freedom, such as electron spin in topological insulator surface states, the two sublattices in graphene, or particle/hole excitations in cuprates. In these three cases, the two components of Dirac fermions, i.e., the two degenerate states at the Dirac point, belong to a symmetry doublet associated with time-reversal, spatial inversion, and particle-hole symmetry, respectively.

A second kind of massless Dirac fermions exists in solids, whose two components correspond to distinct degrees of freedom unrelated by any symmetry. For example, Dirac fermions on the (001) surface of topological crystalline insulators SnTe and $\text{Pb}_{1-x}\text{Sn}_x\text{Se}$ are comprised of the cation Sn/Pb orbital and the anion Te/Se orbital [22]. In heavy fermion semimetals, Dirac fermions can emerge from the hybridization of f and d bands [23,24], which have very different masses.

In the presence of disorder, the two distinct orbitals, from which the Dirac fermion is formed, are generally expected to have different scattering rates. Then, the self-energy of the Dirac fermion in the disorder-averaged single-particle Green's function acquires an *orbital-dependent* imaginary part. The existence of two scattering rates—a generic property of Dirac fermion of the second kind—unnoticed in previous studies, has an important consequence that is only recognized very recently. As shown by one of the authors [23], in such case the imaginary part of self-energy not only broadens the energy spectrum, but also alters the energy-momentum dispersion. It transforms the original Dirac point into a “nodal arc”—a line of band degeneracy without fine-tuning. The two ends of this nodal arc are exceptional points, where the inverse of the Green's function becomes nondiagonalizable. This bulk nodal arc connecting a pair of exceptional points is topologically robust and unique to non-Hermitian band theory recently developed for finite-lifetime quasiparticles [25]. It is also shown that finite-lifetime effects lead to a flat band

or a nodal line segment in type-II Weyl semimetals in three dimensions [26].

In this Rapid Communication, we report the finding of a nodal arc and a new universality class in disordered two-dimensional (2D) Dirac fermions of the second kind, comprised of two distinct orbitals that do not belong to any symmetry doublet. Our model includes a random potential that acts on the two orbitals differently. This type of random potential, not considered in the standard treatment of disordered Dirac fermions, appears naturally in real materials (see below). Our renormalization-group (RG) analysis shows the disorder is marginally relevant, driving the system into a strongly disordered phase. The disorder reduces the quasiparticle weights of the two orbitals at low energy by different amounts. As a result, it generates a tilt of the Dirac cone, even when it is initially absent. In the disordered phase, the two orbitals acquire different lifetimes, which inevitably leads to a bulk nodal arc replacing the Dirac point in the clean limit. The nodal arcs and the tilt of Dirac cones are also observed directly in our numerical simulations.

We consider a 2D Dirac Hamiltonian

$$H_0(\mathbf{r}) = \psi^\dagger(\mathbf{r})[-i(v_x\sigma_z - w\sigma_0)\partial_x - iv_y\sigma_x\partial_y]\psi(\mathbf{r}), \quad (1)$$

and the disorder of the form

$$H_{\text{dis}}(\mathbf{r}) = V(\mathbf{r})\psi^\dagger(\mathbf{r})\eta\psi(\mathbf{r}), \quad (2)$$

where $\psi = (\psi_1, \psi_2)^T$ is a two-component fermion field, σ_i ($i = x, y, z$) are the Pauli matrices, η is a 2×2 Hermitian matrix depending on the type of disorder to be specified below, and $V(\mathbf{r})$ is a random function. The velocity parameter w describes the tilt of the Dirac cone along the x direction in the absence of disorder, where the velocity of the steep and gentle sides of the cone are $v_x \pm w (> 0)$ depending on the sign of w . The velocity along the y direction is given by v_y . We restrict the tilt to satisfy $w^2 < v_x^2$, so that the Fermi surface remains closed. Such massless and tilted Dirac fermions described by H_0 appear in a number of materials, including (001) surface states of SnTe [27,28] and organic conductor α -(BEDT-TTF) $_2$ I $_3$ [BEDT-TTF=bis(ethylenedithio)tetrathiafulvalene] [29–31].

We assume that the random function $V(\mathbf{r})$ is spatially uncorrelated and obeys a Gaussian distribution $P[V] \propto \exp[-\int d^2r V^2(\mathbf{r})/(2\Delta)]$. The random function is characterized by $\langle V(\mathbf{r})V(\mathbf{r}') \rangle = \Delta\delta(\mathbf{r} - \mathbf{r}')$, where Δ (> 0) characterizes the strength of disorder.

The 2×2 matrix η specifies the type of disorder potential. Previous studies of disordered Dirac systems considered the cases where the orbitals forming Dirac fermions are related by a symmetry. In real materials where the two orbitals are unrelated by a symmetry, however, disorder strengths for the two orbitals are allowed to be different; for example, η may have a form $\eta = A\sigma_0 + B\sigma_z$ with two independent constants A and B . Such a form is excluded when the two orbitals are related by a symmetry since the symmetry is restored after disorder averaging.

For simplicity, we henceforth consider the limit $\eta = \eta_{11} \equiv (\sigma_0 + \sigma_z)/2$, where only the $\sigma_z = +1$ orbital is disordered. This highly asymmetric case is relevant for the surface state of topological crystalline insulator $\text{Pb}_{1-x}\text{Sn}_x\text{Te}$, where the two components $\sigma_z = \pm 1$ of the surface Dirac fermion correspond to Sn/Pb and Te orbitals respectively. Since the Pb sites are substituted with Sn, the disorder potential on the $\sigma_z = +1$ orbital is naturally much stronger.

In the following, we will show that the disorder characterized by η_{11} ($A/B = 1$) is marginally relevant in the RG sense. We also confirmed that the disorder is marginally relevant and our conclusion remains valid for a wide range of the ratio A/B . The analysis for general cases will be reported elsewhere [32].

We study disorder-averaged electron spectral function and density of states, which can be computed from the disorder-averaged Green's function: $\bar{G}(\omega) = \langle G(\omega) \rangle \equiv [\omega - H_0 - \Sigma(\omega)]^{-1}$, where $G = (\omega - H_0 - H_{\text{dis}})^{-1}$ is the one-particle Green's function before disorder average, and Σ is the self-energy—a non-Hermitian 2×2 matrix. Note that the translational invariance is statistically recovered after the disorder average. The poles of \bar{G} on the complex plane determine the quasiparticle energy spectrum and lifetime in the presence of disorder.

We use the replica method to evaluate disorder averages. With the replica method, we take the disorder average and obtain the Euclidean action

$$S = \sum_a \int d\tau d^2r \psi_a^\dagger [\partial_\tau - i(v_x\sigma_z - w\sigma_0)\partial_x - iv_y\sigma_x\partial_y] \psi_a - \sum_{ab} \frac{\Delta}{2} \int d\tau d\tau' d^2r (\psi_a^\dagger \eta \psi_a)(\tau) (\psi_b^\dagger \eta \psi_b)(\tau'), \quad (3)$$

where $a, b (= 1, \dots, n)$ are replica indices. The quartic term, effectively working as an interaction between replicas, is generated by the disorder average.

Now we examine whether the parameters are relevant or irrelevant, by a perturbative RG calculation to one-loop order (Fig. 1). The calculations give the scale dependence of parameters, arising from resummations of logarithmic divergences.

The two-leg diagram [Fig. 1(a)] gives the renormalization of quasiparticle weight $Z = (1 - \text{Re } \partial\Sigma/\partial\omega)^{-1}$. We have two different quasiparticle weights Z_{11} and Z_{22} for the two orbitals $\sigma_z = \pm 1$, respectively. Here, the self-energy $\Sigma(\omega)$ depends only on frequency ω , but not on momentum. Therefore, the renormalization of the velocity parameters v_x , v_y , and w owes

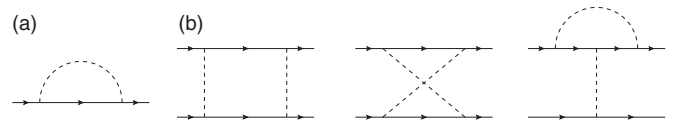


FIG. 1. One-loop diagrams that contribute to the RG equations. Solid and dashed lines correspond to the fermion field and disorder, respectively. The two-point diagram (a) contributes to the field renormalization, and the four-point diagrams (b) give corrections to the disorder strength.

solely to the renormalization of the quasiparticle weights. The four-leg diagrams [Fig. 1(b)] bear disorder strength renormalization. After evaluating the one-loop diagrams, we obtain the following set of RG equations [33]:

$$\begin{aligned} \frac{dv_x}{dl} &= -\alpha v_x, & \frac{dv_y}{dl} &= -\alpha \frac{v_x}{v_x - w} v_y, & \frac{dw}{dl} &= \alpha v_x, \\ \frac{d\Delta}{dl} &= 4\alpha \frac{w}{v_x - w} \Delta, \\ \frac{dZ_{11}}{dl} &= -2\alpha \frac{v_x}{v_x - w} \Delta, & \frac{dZ_{22}}{dl} &= 0, \end{aligned} \quad (4)$$

where $\alpha = \Delta/[4\pi\sqrt{(v_x^2 - w^2)v_y^2}]$ is the dimensionless quantity, $l = \ln(\Lambda/\epsilon)$, Λ is the UV energy cutoff, and ϵ is the energy scale of interest.

Even when there is initially no tilt of the Dirac cone, the tilt $w(> 0)$ is generated by the disorder. This is seen from the decreasing quasiparticle weight of the $\sigma_z = +1$ orbital as the energy scale goes down. For $k_y = 0$, this orbital corresponds to the energy branch of $(v_x - w)k_x$. In low energies, the decreasing Z_{11} , accompanied by the change of $\text{Re } \Sigma(\omega)$, reduces $(v_x - w)$. In contrast, since the $\sigma_z = -1$ orbital is free from disorder, Z_{22} remains constant and so does $(v_x + w)$. The tilt $w > 0$ increases the density of states of the $\sigma_z = +1$ orbital, which helps the disorder strength Δ grow as the energy approaches the Dirac point. Thus, Δ is marginally relevant, driving the system into a disordered phase. In the course of increasing Δ , the tilt keeps growing whereas the two velocity parameters v_x and v_y decrease. Those flows stop as $w \rightarrow v_x$, where the gentle slope of the Dirac cone becomes flat and the increasing density of states further drives the flow to the strongly disordered phase.

To study the properties of the disordered phase, the self-energy Σ is evaluated by the self-consistent Born approximation. Within this approximation, the self-consistent equation becomes

$$\Sigma(\omega) = \Delta \int' \frac{d^2k}{(2\pi)^2} \eta \bar{G}(\mathbf{k}, \omega) \eta. \quad (5)$$

\int' indicates an integration with the cutoff Λ . In the present model with $\eta = \eta_{11}$, the self-energy $\Sigma(\omega)$ is nonzero only for the $\sigma_z = +1$ orbital; $\Sigma(\omega) = \Sigma_{11}(\omega)\eta_{11}$. The self-energy is calculated for the two regions $|\epsilon| \ll |\Sigma|$ and $|\epsilon| \gg |\Sigma|$, which are separated at the energy scale

$$\Gamma_0 = 2\Lambda \sqrt{\frac{v_x - w}{v_x + w}} \exp\left(-\frac{v_x - w}{v_x + w} \frac{1}{\alpha}\right). \quad (6)$$

Γ_0 corresponds to the energy scale where the one-loop RG analysis breaks down. Close to Γ_0 , the velocity $(v_x - w)$ approaches zero, which destroys perturbative expansion with respect to α .

The retarded self-energy $\Sigma_{11}^R(\omega)$ is obtained from Eq. (5) in a series of ω for $|\omega| \ll \Gamma_0$, and in a series of α by iterations for $|\omega| \gg \Gamma_0$:

$$\Sigma_{11}^R(\omega) = \begin{cases} -\frac{2v_x(v_x - w)}{(v_x + w)^2} \frac{\omega}{\alpha} - i\Gamma_0, & |\omega| \ll \Gamma_0, \\ -\alpha \frac{v_x}{v_x - w} \left[\omega \ln \left(\frac{(v_x^2 - w^2)\Lambda^2}{v_x^2 \omega^2} \right) + i\pi|\omega| \right], & |\omega| \gg \Gamma_0. \end{cases} \quad (7)$$

Importantly, it shows that the Green's function has a finite imaginary part even at $\omega = 0$, appearing only in the $\sigma_z = +1$ orbital. The energy spectrum of the quasiparticles is obtained as zeros of \tilde{G} , i.e., $\det[E - H_0(\mathbf{k}) - \Sigma(E)] = 0$. The quasiparticle energy dispersion becomes now complex valued because of the non-Hermitian component of Σ . Furthermore, the parameters of the Hamiltonian are renormalized by the reduced quasiparticle weight Z_{11} , giving the eigenvalues for $|E| \ll \Gamma_0$ as

$$E_{\pm} = \frac{(\tilde{v}_2 - v_1)k_x - i\tilde{\Gamma}_0}{2} \pm \sqrt{\left[\frac{(\tilde{v}_2 + v_1)k_x - i\tilde{\Gamma}_0}{2} \right]^2 + v_y^2 k_y^2} \quad (8)$$

with $v_1 = v_x + w$, $\tilde{v}_2 = Z_{11}(v_x - w)$, and $\tilde{\Gamma}_0 = Z_{11}\Gamma_0$.

In the clean limit, two linearly dispersing energy bands touch at a Dirac point. With disorder, the quasiparticle energy

obtains an imaginary component, and the real parts of the two energy branches coalesce along the line $|k_y| \leq \tilde{\Gamma}_0/(2v_y)$ ($k_x = 0$). Here, we can observe the formation of a nodal arc in a disordered Dirac model, which lies along the y direction and terminates at exceptional points $(0, \pm\tilde{\Gamma}_0/(2v_y))$.

A nodal arc is absent when the two orbitals are related by a symmetry and disorder is characterized by $\eta = \sigma_0$ or σ_i . In such cases, the energy dispersion is simply smeared by the same finite lifetime for the two orbitals because of the symmetry. The self-consistent equation (5) also confirms that the self-energy is proportional to σ_0 .

It is important to point out that the formation of the nodal arc by disorder is observed even at zero temperature $T = 0$, because plane-wave (or Bloch wave) states as quasiparticles are not eigenstates due to the lack of translational symmetry by disorder. Therefore, unlike the arc due to interactions, a nodal arc formed by disorder is independent from thermal effects, which contribute largely to thermal broadening of energy dispersions.

Next, we perform numerical simulations on a square lattice with the periodic boundary conditions to confirm the observations above. Because of the fermion doubling problem, the lattice model necessarily has a pair of Dirac cones in the Brillouin zone. The tight-binding Hamiltonian that we use is

$$\begin{aligned} \hat{H}_0 &= -t \sum_{(ij)} c_i^\dagger \sigma_z c_j + t_y \sum_i (i c_{i+y}^\dagger \sigma_x c_i + \text{H.c.}) \\ &\quad + u \sum_i c_i^\dagger \sigma_z c_i, \\ \hat{H}_{\text{dis}} &= \sum_i V_i c_i^\dagger \eta_{11} c_i. \end{aligned} \quad (9)$$

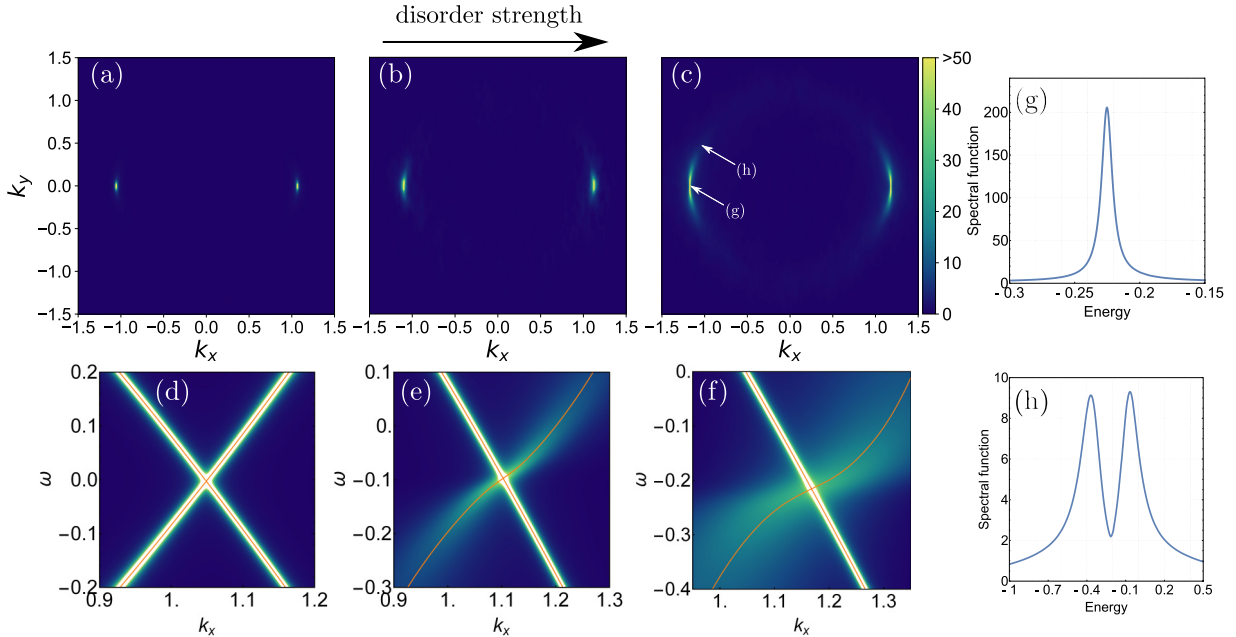


FIG. 2. (a)–(c) Spectral function $A(\mathbf{k}, \omega)$ slices across the relevant fragment of the Brillouin zone for ω corresponding to the open contour Fermi surface, calculated for $t = 1$, $t_y = 0.25$, $u = 3.0$ and disorder strengths $V_0 = 0.5, 3.0, 4.5$, respectively. As the disorder strength increases, the nodal arc becomes longer. (d)–(f) Spectral function $A(\mathbf{k}, \omega)$ slices for $k_y = 0$ and the same parameters as respective panels above. As the disorder increases, the Dirac cone tilts. (g), (h) Energy dependence of the spectral function for a single point in \mathbf{k} space for a point in the middle of the arc and a point on the faint contour away from the nodal arc, chosen as indicated by white arrows on panel (c).

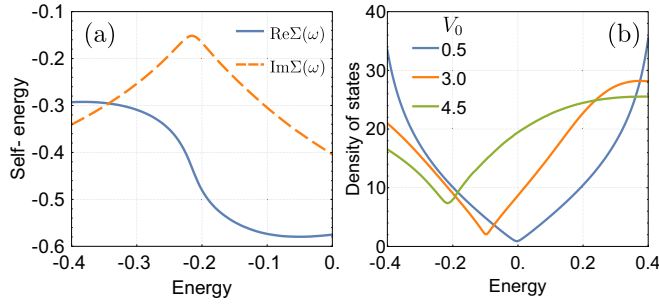


FIG. 3. (a) The self-energy $\Sigma(\omega)$ calculated using self-consistent Born approximation with parameters corresponding to Fig. 2(c). (b) Total integrated density of states $\rho(\omega)$. The position of the minimum, which shifts to lower energies, corresponds to the energy of the nodal arc.

For $0 < u < 4t$, \hat{H}_0 has two Dirac points at $k_x a = \pm \arccos[(u - 2t)/(2t)]$, $k_y a = 0$, where a is a lattice constant. We assume that there is no tilt of the Dirac cones in the clean limit. By a series expansion near the $k_x a = \arccos[(u - 2t)/(2t)] > 0$ point, we can establish the connection to the Dirac model from Eq. (1) by $v_x = \text{sgn}(u - 2t)\sqrt{(4t - u)u}$, $v_y = -2t_y$, $w = 0$. For the purpose of numerical simulations, we choose a random potential V_i obeying a uniform distribution over an interval $[-V_0/2, V_0/2]$, and its spatial correlation function is $\langle V(\mathbf{r}_i)V(\mathbf{r}_j) \rangle = \frac{V_0^2}{12}\delta(\mathbf{r}_i - \mathbf{r}_j)$. In all the simulations the lattice consists of 800×800 sites. We choose $t = 1$, $t_y = 0.25$, $u = 3$, $a = 1$ and change the disorder strength to observe the evolution of nodal arcs. We use MUMPS [34] and KWANT [35] packages for the numerical calculations.

For a particular disorder distribution, we can compute the retarded Green's function $G^R(\mathbf{r}, \mathbf{r}', \omega) = (\omega + i\delta - H_0 - H_{\text{dis}})^{-1}(\mathbf{r}, \mathbf{r}', \omega)$, where δ is a small positive quantity. The disorder average is taken by computing G^R for 2000 independent disorder realizations. (A result for a single disorder realization is presented in the Supplemental Material (SM) [33].) We can then reinstate translational symmetry in the averaged quantity $\bar{G}^R(\mathbf{r} - \mathbf{r}', \omega) = \langle G^R(\mathbf{r}, \mathbf{r}', \omega) \rangle$, which allows the Fourier transformation $\bar{G}^R(\mathbf{k}, \omega)$. The spectral function is $A(\mathbf{k}, \omega) = -\frac{1}{\pi} \text{Im Tr } \bar{G}^R(\mathbf{k}, \omega)$.

Figures 2(a)–2(c) show the spectral function $A(\mathbf{k}, \omega)$ at the energy where the two bands touch, with the disorder strength $V_0 = 0.5, 3.0, 4.5$, respectively. We can observe the nodal arcs extending in the y direction, with stronger disorder yielding longer nodal arcs. The curvature of the nodal arcs reflects the energy dispersion of the tight-binding model \hat{H}_0 , which corresponds to a higher-order effect with respect to k in the linearized model, Eq. (1).

The disorder-averaged Green's function can be alternatively obtained from the self-consistent equation (5) in successive iterations of numerical integrations. We replace the parameter Δ with $V_0^2/12$ and the integration is now over the Brillouin zone. Using the self-energy obtained by the self-consistent calculation, we compute spectral

functions, which are in excellent agreement with those in Figs. 2(a)–2(c) (see SM [33]). This supports the conclusions that the self-energy is largely momentum independent and the important effects are due to its energy dependence instead. We also obtain the slices of $A(\mathbf{k}, \omega)$ in the $k_y = 0$ plane [Figs. 2(d)–2(f)], with solid lines indicating the position of the poles of $\bar{G}^R(\mathbf{k}, \omega)$. We confirm that the tilt of Dirac cones is generated by the disorder, even when it is initially absent. Since only the orbital with a gentle slope of the energy dispersion is affected by the disorder, only its component of the spectral function is smeared by $\text{Im } \Sigma$, and its peaks are smaller due to the decreasing quasiparticle weight Z .

Furthermore, we check that the contour we observe in Fig. 2(c) is indeed a nodal arc by determining the spectral function as a function of energy for particular \mathbf{k} values on and off the arc as indicated by the white arrows. In Figs. 2(g) and 2(h), we observe a single peak in the spectral function for the point on the contour and two peaks on the faint structure outside of the arc.

The self-energy $\Sigma(\omega)$ obtained from the self-consistent Born approximation is shown in Fig. 3(a). Note that $\text{Re } \Sigma_{11}(\omega)$ shifts the band crossing point to $\omega - \text{Re } \Sigma_{11}(\omega)/2 = 0$, since $\Sigma(\omega) = \Sigma_{11}(\omega)(\sigma_0 + \sigma_z)/2$. Crucially, close to the band touching point, $\text{Re } \Sigma_{11}$ is linearly dependent on energy and the slope increases with disorder strength causing the tilting of the Dirac cone, which is consistent with the results of the RG analysis. Moreover, at the band touching point, $\text{Im } \Sigma$ has a dip, but it remains finite, supporting the existence of the nodal arcs. This result agrees with the solution of the self-consistent equation for the linearized model. The shift of the band touching point is also evident in the density of states $\rho(\omega) = \int_{\text{BZ}} d^2k A(\mathbf{k}, \omega)$, which is depicted in Fig. 3(b). The minimum of $\rho(\omega)$ moves to lower energy values with larger disorder strength. The minima match the solutions of $\omega - \text{Re } \Sigma_{11}(\omega)/2 = 0$ (see Fig. S1 in the SM [33]).

In summary, we studied a disordered 2D Dirac system with two orbitals not belonging to a symmetry doublet, and reported a nodal arc and a new universality class in disordered 2D Dirac fermions. By the RG analysis we have shown that asymmetric disorder for the two orbitals is marginally relevant, and in the strongly disordered phase, the inequality of the quasiparticle weights and lifetimes of the two orbitals yields a tilt of the Dirac cone and a bulk nodal arc, replacing the Dirac point. The nodal arc is formed by the non-Hermitian effect of the self-energy generated by the disorder, which is present even at $T = 0$. This conclusion is supported by numerical calculations using a tight-binding model and self-consistent Born approximation, which show the appearance of a nodal arc in the spectral function.

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M.P. and H.I. contributed equally to this work.

[1] E. Fradkin, *Phys. Rev. B* **33**, 3257 (1986); **33**, 3263 (1986).

[2] S. Guruswamy, A. LeClair, and A. W. W. Ludwig, *Nucl. Phys. B* **583**, 475 (2000).

- [3] For a review, see, e.g., F. Evers and A. D. Mirlin, *Rev. Mod. Phys.* **80**, 1355 (2008).
- [4] A. W. W. Ludwig, M. P. A. Fisher, R. Shankar, and G. Grinstein, *Phys. Rev. B* **50**, 7526 (1994).
- [5] A. A. Nersisyan, A. M. Tsvelik, and F. Wenger, *Nucl. Phys. B* **438**, 561 (1995).
- [6] T. Senthil and M. P. A. Fisher, *Phys. Rev. B* **61**, 9690 (2000).
- [7] M. Bocquet, D. Serban, and M. R. Zirnbauer, *Nucl. Phys. B* **578**, 628 (2000).
- [8] A. Altland, B. D. Simons, and M. R. Zirnbauer, *Phys. Rep.* **359**, 283 (2002).
- [9] M. S. Foster, Hong-Yi Xie, and Yang-Zhi Chou, *Phys. Rev. B* **89**, 155140 (2014).
- [10] E. McCann, K. Kechedzhi, V. I. Fal'ko, H. Suzuura, T. Ando, and B. L. Altshuler, *Phys. Rev. Lett.* **97**, 146805 (2006).
- [11] M. Koshino and T. Ando, *Phys. Rev. B* **73**, 245403 (2006).
- [12] P. M. Ostrovsky, I. V. Gornyi, and A. D. Mirlin, *Phys. Rev. B* **74**, 235443 (2006).
- [13] I. L. Aleiner and K. B. Efetov, *Phys. Rev. Lett.* **97**, 236801 (2006).
- [14] A. Altland, *Phys. Rev. Lett.* **97**, 236802 (2006).
- [15] P. M. Ostrovsky, I. V. Gornyi, and A. D. Mirlin, *Phys. Rev. Lett.* **98**, 256801 (2007).
- [16] J. H. Bardarson, J. Tworzydło, P. W. Brouwer, and C. W. J. Beenakker, *Phys. Rev. Lett.* **99**, 106801 (2007).
- [17] K. Nomura, M. Koshino, and S. Ryu, *Phys. Rev. Lett.* **99**, 146806 (2007).
- [18] S. Ryu, C. Mudry, H. Obuse, and A. Furusaki, *Phys. Rev. Lett.* **99**, 116601 (2007).
- [19] M. Noro, M. Koshino, and T. Ando, *J. Phys. Soc. Jpn.* **79**, 094713 (2010).
- [20] R. S. K. Mong, J. H. Bardarson, and J. E. Moore, *Phys. Rev. Lett.* **108**, 076804 (2012).
- [21] L. Fu and C. L. Kane, *Phys. Rev. Lett.* **109**, 246605 (2012).
- [22] I. Zeljkovic, Y. Okada, C.-Y. Huang, R. Sankar, D. Walkup, W. Zhou, M. Serbyn, F. Chou, W.-F. Tsai, H. Lin, A. Bansil, L. Fu, M. Z. Hasan, and V. Madhavan, *Nat. Phys.* **10**, 572 (2014).
- [23] V. Kozii and L. Fu, [arXiv:1708.05841](https://arxiv.org/abs/1708.05841).
- [24] Y. Xu, C. Yue, H. Weng, and X. Dai, *Phys. Rev. X* **7**, 011027 (2017).
- [25] H. Shen, B. Zhen, and L. Fu, *Phys. Rev. Lett.* **120**, 146402 (2018).
- [26] A. A. Zyuzin and A. Yu. Zyuzin, *Phys. Rev. B* **97**, 041203(R) (2018).
- [27] Y. Tanaka, Z. Ren, T. Sato, K. Nakayama, S. Souma, T. Takahashi, K. Segawa, and Y. Ando, *Nat. Phys.* **8**, 800 (2012).
- [28] I. Sodemann and Liang Fu, *Phys. Rev. Lett.* **115**, 216806 (2015).
- [29] A. Kobayashi, S. Katayama, Y. Suzumura, and H. Fukuyama, *J. Phys. Soc. Jpn.* **76**, 034711 (2007); A. Kobayashi, Y. Suzumura, and H. Fukuyama, *ibid.* **77**, 064718 (2008).
- [30] M. O. Goerbig, J.-N. Fuchs, G. Montambaux, and F. Piéchon, *Phys. Rev. B* **78**, 045415 (2008).
- [31] H. Isobe and N. Nagaosa, *J. Phys. Soc. Jpn.* **81**, 113704 (2012); *Phys. Rev. Lett.* **116**, 116803 (2016).
- [32] H. Isobe, M. Papaj, and L. Fu (unpublished).
- [33] See Supplemental Material at <http://link.aps.org/supplemental/10.1103/PhysRevB.99.201107> for the derivation of RG equations and the details of the numerical simulations.
- [34] P. R. Amestoy, I. S. Duff, and J. S. Koster, and J. Y. L'Excellent, *SIAM. J. Matrix Anal. Appl.* **23**, 15 (2001).
- [35] C. W. Groth, M. Wimmer, A. R. Akhmerov, and X. Waintal, *New J. Phys.* **16**, 063065 (2014).

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