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PT-Symmetric Real Dirac Fermions and Semimetals

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Recently, Weyl fermions have attracted increasing interest in condensed matter physics due to their rich phenomenology originated from their nontrivial monopole charges. Here, we present a theory of real Dirac points that can be understood as real monopoles in momentum space, serving as a real generalization of Weyl fermions with the reality being endowed by the PT symmetry. The real counterparts of topological features of Weyl semimetals, such as Nielsen-Ninomiya no-go theorem, 2D subtopological insulators, and Fermi arcs, are studied in the PT symmetric Dirac semimetals and the underlying reality-dependent topological structures are discussed. In particular, we construct a minimal model of the real Dirac semimetals based on recently proposed cold atom experiments and quantum materials about PT symmetric Dirac nodal line semimetals.

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Introduction.—The discovery of topological insulators [1,2] has galvanized the condensed matter community into research of topological phenomena in a vast variety of quantum matters. One class of materials that have attracted particular attention in recent years are the gapless topological systems, such as Weyl and various Dirac semimetals, Dirac nodal line semimetals, and nodal topological superconductors. In this new topological paradigm of gapless solid state physics, Weyl fermions are of the most fundamental status [3–5], in the sense that a Weyl point can be interpreted as a unit monopole of the $U(N)$ Berry bundle of the band structure in momentum space [3], which, unlike other topological gapless modes, does not rely on any symmetry. The nontrivial topological charges of Weyl points as momentum space monopoles put strong constraints on the global band structure of a Weyl semimetal, leading to two primary consequences. First, a Weyl semimetal conforms the Nielsen-Ninomiya (NN) no-go theorem that Weyl points generically appear in pairs of opposite unit $U(N)$ monopole charges. This is due to the fact that the total monopole charge has to vanish due to the orientability and closeness of the first Brillouin zone (BZ) [6,7], and a multiply charged monopole may unstably split into unit ones under perturbations [5,8,9]. Second, the Chern number or Thouless-Kohmoto-Nightingale-den Nijs (TKNN) invariant of a two-dimensional (2D) subsystem [10,11] has to jump by a unit when moving across a Weyl point, and consequently, the gapless chiral states originated from the Chern numbers of these 2D subsystems form Fermi arcs on the surface of a Weyl semimetal connecting the projections of Weyl points in the bulk [12], which are experimentally observable [13–17].

In this Letter, we present a real generalization of the Weyl semimetal through the combined PT symmetry with $(PT)^2 = 1$, where $P$ indicates the inversion symmetry and $T$ the time reversal. Note that PT has to be broken for a Weyl semimetal unless gauge potentials are present [18]. Analogous to the innumerable interesting phenomena embedded in the classification of topological insulators and superconductors [19,20], the real counterpart of the Weyl point, which we call a real Dirac point, is of particular interest in the recently established classification of PT and CP (charge conjugate C) symmetric topological gapless systems [21]. We show that the PT symmetry endows a reality condition on the band structure inducing a real Berry bundle over the BZ, and the real Dirac point is actually a unit monopole for the $O(N)$ Berry bundle in contrast to the $U(N)$ bundle associated to a Weyl point. Historically, the real monopole first appeared in an entirely different context about $SO(3)$ gauge field theory in Polyakov’s classic work [22]. While the real monopole nature of the real Dirac point guarantees that all features of Weyl semimetals find their real counterparts in the real Dirac semimetal including the no-go theorem and gapless chiral Fermi arc surface states, the reality plays a nontrivial role. In particular, the surface Fermi arcs are of richer band structure that preserves the reality. In addition, physical systems of such PT symmetry have been recently predicted in real materials [23–26] and designed in cold atom experiments [27] for the realization of PT-invariant nodal line Dirac semimetals. Starting with such Dirac semimetals, we provide a general recipe for model construction of real Dirac semimetals, which may pave the way for their future experimental realization. Before moving to the detailed study, we note that throughout this work only the combined PT symmetry is required, while individual $P$ and $T$ symmetries may be violated separately.

As real monopoles in momentum space.—We first recall some basics of a Weyl point described by

$$\text{some equation}$$

$$\text{some equation}$$

$$\text{some equation}$$

$$\text{some equation}$$

$$\text{some equation}$$

$$\text{some equation}$$
The gapless point at the origin is associated with a topologically nontrivial $U(1)$ Berry bundle on the sphere $S^2$ enclosing the gapless point, and therefore can be regarded as a monopole of $U(1)$ group. To identify the topological nature of Eq. (1), we analyze the Berry bundle by choosing stereographic coordinates for the north and south hemisphere, respectively, where without loss of generality we choose the unit sphere $|n|=1$ [See the Supplemental Material (SM) for details [28]]. The Hamiltonian (1) has eigenstates $|+,z\rangle^N = (1, z)\hat{T}$ ($(|+,z\rangle^N = (1, 1)$ and $|−,z\rangle^N = (1, −1)$) on the north (south) hemisphere with positive and negative unit energy, respectively. The “$T$” denotes vector or matrix transposition. Negative states $|−,z\rangle^N$ on the two patches of coordinates are sections of the $U(1)$ principle fiber bundle given by the projector $P = [1 - \text{sgn}(H_W(|S^1|))] / 2$, which determines the transition function on the intersection $S^1$ of the two patches.

$$|−,\beta\rangle^S = g_{SN}^C(\phi)|−,\beta\rangle^N,$$  

which makes a Berry bundle a real vector bundle associated to a $O(N)$ principle bundle. Note that there is no Kramers degeneracy for $(\hat{\mathcal{P}} \hat{T})^2 = 1$ in contrast to the case of $(\hat{\mathcal{P}} \hat{T})^2 = −1$, which guarantees that the reality condition (3) holds for each band.

For convenience and without loss of generality, we choose $\hat{\mathcal{P}} \hat{T} = \hat{\mathcal{K}}$, which simply means that the Hamiltonian is real, and a real Dirac point may be represented as

$$\mathcal{H}_{RD}(k) = (\hat{\tau}_1 j_\phi + \hat{\tau}_2 \hat{\tau}_3 \hat{\sigma}_3) \hat{\tau}_0, \quad (4)$$

as suggested by the Clifford algebra theory. It is noted that the real Hamiltonian density of Eq. (4) is actually that of Majorana fermion, but the spinor $\psi_D$ here is a Dirac spinor, rather than a Majorana one, namely, they do not agree with each other at second quantization. The topologically nontrivial real vector bundle given by the projector $P = [1 - \text{sgn}(\mathcal{H}_{RD}(|S^2|))] / 2$ is actually the generator of the reduced orthogonal $K$ group, $\hat{\mathcal{K}} O(S^2) \cong \mathbb{Z}_2$, and thus, we may call the gapless point of Eq. (4) the unit real monopole in momentum space. To identify the $O(N)$ monopole charge of Eq. (4), we still adopt the stereographic coordinates for the unit sphere $S^2$ enclosing the gapless point, similar to the case of Weyl point (See the Supplemental Material for details [28]). Now it is a four-band theory, and the real eigenstates on the north hemisphere are $|+,1\rangle^N = (1, x, 0, y)\hat{T}$, $|+,2\rangle^N = (0, y, 1, x)\hat{T}$ for positive energy and $|−,1\rangle^N = (x, −1, y, 0)\hat{T}$, $|−,2\rangle^N = (−y, 0, x, −1)\hat{T}$ for negative energy, while on the south hemisphere, $|+,1\rangle^S = (x, 1, y, 0)\hat{T}$, $|+,2\rangle^S = (−y, 0, x, 1)\hat{T}$, $|−,1\rangle^S = (1, −x, 0, y)\hat{T}$, and $|−,2\rangle^S = (0, −y, 1, −x)\hat{T}$, which satisfy Eq. (3). On the intersection $S^1$ of the two hemispheres, eigenstates for negative energy are given explicitly as $|−,1\rangle^N = (\cos \phi, −1, \sin \phi, 0)\hat{T}$, $|−,2\rangle^N = (−\sin \phi, 0, \cos \phi, −1)\hat{T}$, $|−,1\rangle^S = (1, −\cos \phi, 0, \sin \phi)\hat{T}$, and $|−,2\rangle^S = (0, −\sin \phi, 1, −\cos \phi)\hat{T}$. Then the transition function $g_{SN}^C(\phi) \in O(2)$ with $\phi \in S^1$, which gives the relation

$$|−,\alpha\rangle^S = [g_{SN}^C(\phi)] \langle −,\beta \rangle^N,$$

is calculated as

$$g_{SN}^C(\phi) = \left( \begin{array}{cc} \cos \phi & −\sin \phi \\ \sin \phi & \cos \phi \end{array} \right). \quad (6)$$

It is transparent that the transition function as a map from $S^1$ to $O(2)$ has a unit winding number, which verifies that the real Berry bundle of Eq. (4) generates $\hat{\mathcal{K}} O(S^2) \cong \mathbb{Z}_2$.

Recall that the topological invariant for a Weyl point is given by the famous Chern number or TKNN invariant $\nu = (1/2\pi i) \int_{S^1} \text{tr} \mathcal{F}$, where the Berry curvature is derived from complex valence eigenvectors on the gapped
$S^2$ enclosing the Weyl point. To formulate a topological invariant for the real monopole, similarly, we may use the real Berry curvature $F_R$, which is derived from the real connection $A_R^\alpha = \langle \alpha, k | d|\beta, k \rangle$ with the real eigenstates $|\alpha, k \rangle$ satisfying Eq. (3). The corresponding topological invariant for two valence bands is given by

$$\nu_R = -\frac{1}{4\pi} \int_{S^2} \text{tr}(IF_R) \mod 2, \quad (7)$$

where $F_R$ is the curvature for the real Berry bundle and $I = -i\lambda_2$ is the generator of the $SO(2)$ group with $\lambda_2$ being the second Pauli matrix. We may call Eq. (7) the real Chern number, and the “mod 2” should be understood according to the following paragraph.

For higher-dimensional real vector bundles on $S^2$, the transition functions on the equator are maps from $S^1$ to $O(N)$, which are classified by $\pi_1(O(N)) \cong Z_2$, and nontrivial transition functions correspond to nontrivial real vector bundles. A transition function from $S^1$ to $O(N)$ can always be continuously deformed to be a map from $S^1$ to $O(2) \subset O(N)$, and the parity of the winding number gives the homotopy class of the transition function, namely, $Z_2 \cong Z/2Z$, though two-dimensional real vector bundles have a $Z_2$ classification since $\pi_1(O(2)) \cong Z$. In this sense, the Berry bundle of the Majorana Hamiltonian, Eq. (4), is a generator of the reduce orthogonal $K$ theory, $KO(S^2) \cong Z_2$.

No-go theorem.—Following almost the same topological arguments for the NN No-go theorem of Weyl semimetals [6,7], we can show that real Dirac semimetals of the $PT$ symmetry satisfy the No-go theorem that real Dirac points always exist in pairs for a lattice model. Since a Brillouin zone as a tori is a closed orientable manifold, when one chooses an oriented $S^2$ enclosing any real monopole, the $S^2$ also encloses the rest of real monopoles with the opposite orientation, which implies the topological charge of the real monopole is equal to the inverse of the sum of topological charges for the rest, namely,

$$\sum_a \nu_R^a = 0, \quad (8)$$

where $a$ labels real monopoles in a real Dirac semimetal [28]. On the other hand, in contrast to Weyl points, real Dirac points of positive or negative unit charge are topologically indistinguishable as real monopoles due to the $Z_2$ nature. This implies there is no canonical dipole pairing of real Dirac points in the whole Brillouin zone, which is different from the situation of Weyl semimetals.

Considering copious topological features of Weyl semimetals originated from the monopole charges, particularly, the surface Fermi arc, it is also intriguing to investigate the real counterpart of them for $PT$ invariant semimetals.

Minimal semimetal models as real dipoles.—Now we shall construct a minimal model of real Dirac semimetals, which contains a single dipole of real Dirac points in the Brillouin zone. To motivate our model, we construct it from $PT$ symmetric Dirac nodal line semimetals, which have recently been predicted in several quantum materials and cold atom experimental proposals [26,27]. Let us start with the Bloch Hamiltonian in a proposed cold atom experiment for $PT$ symmetric Dirac nodal line semimetals, $H(k) = 2t\sin k_x \sigma_2 + |m - f(k)|\sigma_3$ with $f(k) = \alpha_+ \cos k_x + \alpha_\perp \cos k_z$, where the Pauli matrices $\sigma_i$ operate in every two-level lattice site, and the $PT$ operator is represented as $\hat{P} \hat{T} = \sigma_3 \hat{K}$ with $(\hat{P} \hat{T})^2 = 1$. Note that only the algebraic properties of the $PT$ operator are needed for the $PT$ symmetric topological classification, while resultant topological features are independent of concrete representations [33]. Since a real Dirac point has a four dimensional internal space, we first double the Dirac nodal line model, $H \otimes \tau_0$, and then introduce appropriate $PT$ invariant couplings between them which collapses a pair of overlapping nodal lines into two real Dirac points. For such couplings, the only $PT$ invariant gamma matrix is $\sigma_1 \otimes \tau_2$, and a possible term is therefore $\sin k_x \sigma_1 \otimes \tau_2$, which leads to our real Dirac semimetal model,

$$H_{RSM}(k) = 2t_x \sin k_x \sigma_1 \otimes \tau_2 + 2t_y \sin k_x \sigma_2 \otimes \tau_0 + |m - f(k)|\sigma_3 \otimes \tau_0. \quad (9)$$

See the Supplemental Material [28] for the corresponding tight-binding model in real space. Since the real Dirac semimetal model is constructed from a cold atom system [27], in principle, it is experimentally realizable and all parameters in Eq. (9) are independently tunable. For concreteness, we would like to work in a parameter region where only a single pair of Dirac points separating in the $k_y$ axis with linear dispersion relations for their coarse-gained effective theories. The parameter region may be identified with the region, $t_x \neq 0$, and $0 < (m - \alpha_+)/2\alpha_+ < 1$ and $(m + \alpha_+)/2\alpha_+ > 1$, where for the Dirac nodal line model only a single Dirac nodal line lies in the $k_x - k_z$ sub-Brillouin zone with $k_y = 0$ centered at the origin.

To prepare the discussions about the topology of the real Dirac semimetal, it is helpful to have a look at the topological features of the Dirac nodal line model in this parameter region. Since the nodal circle has nontrivial $Z_2$ topological charge [28], topological invariant for one-dimensional systems along $k_y$ is nontrivial either inside the circle or outside, but not both. To identify the nontrivial region, it is sufficient to check the one at the origin of the $k_x - k_z$ plane, which is given by $h(k_y) = 2t \sin k_x \sigma_2 + |m - 2\alpha_+ - \alpha_\perp \cos k_z|\sigma_3$. Assuming that $2t = \alpha_+$, we find the condition for nontrivial topological invariant is given by $|m - 2\alpha_+| < \alpha_\perp$, which is always satisfied in the aforementioned parameter region. Accordingly, when boundaries are open properly perpendicular to $y$ direction, there are drum-head states filling all the inside of the nodal loop.
Nodal loops overlaps degenerately in the $k_z$ direction, for which we may associate the same topological invariant related to the PT symmetry as for the real monopole, namely, Eq. (7) with $S^2$ being replaced by the BZ. Since the two real Dirac points have unit topological charges, a two-dimensional subsystem has its topological invariant jumped by one when passing through a Dirac point, which means the topological invariant of subsystems with $k_z \in (k_z^-, k_z^+)$ is differentiated by one from that of subsystems with $k_z \notin [k_z^-, k_z^+]$ ($k_z^\pm$ are $k_z$ coordinates of two real Dirac points). For a specific $k_y$, the subsystem of our model (9) has the Hamiltonian, $\mathcal{H}^{2D}_k(k_x, k_y) = \mathcal{H}_{RSM}(k_x, k_y, k_z)$.

Instead of computing its the topological invariant directly, we may infer it from the construction of our model. Considering the boundaries perpendicular to $\hat{y}$ direction, we start with vanishing $\sigma_1 \otimes \tau_2$ term, namely, $t_2 = 0$ in Eq. (9), where for the bulk spectrum two nodal loops overlaps degenerately in the $k_y-k_z$ plane with $k_z = 0$, so do two copies of drum-head states (see the previous paragraph). Then turning on the term $2t_e \sin k_y \sigma_1 \otimes \tau_2$ with an infinitesimal $t_e$, degeneracies in the spectrum should be sensitive to the perturbation in contrast to the gapped regions that are insensitive. Consequently, on the surface spectrum, the region outside the drum head may be still gapped, while only a degenerate line segment survives connecting two remaining real Dirac points in the bulk under the perturbation, since it is known from above discussions that two-dimensional systems parametrized by $k_z$ are topologically trivial or nontrivial separated by two real Dirac points. Thus, we infer that the Hamiltonian $\mathcal{H}^{2D}_k(k_x, k_y)$ is topologically nontrivial when $k_z \in (k_z^-, k_z^+)$.

FIG. 2. The windings of Wilson loops around $k_y$ for 2D subsystems with given $k_z$. Each Wilson loop is computed along a large circle parametrized by $k_y$ for fixed $k_x$ and $k_z$. Panel (a) for the 2D subsystem with $k_z = 0$ has unit winding number, but (b) for $k_z = \pi$ has zero winding number.

Alternately, one may work out the transition function of the real bundles for 2D subsystems through the Wilson loops

$$W(k_y) = P \exp \left( \int d k_y A^\theta(k_x, k_y) \right) \in O(N)$$

(with $P$ indicating the path order) along large circles coordinated by $k_y$, and then check the winding number around the large circle parametrized by $k_x$ [28]. The numeric results are illustrated in Fig. 2.

The spectrum of Eq. (9) with $\hat{y}$ direction being confined as a slab under natural boundary conditions respecting the PT symmetry has been shown in Figs. 1(b), 1(c), and 1(d). It is observed that for a surface the Fermi arc spectrum consists of two inclined planes with different angles connecting conduction and valence bands, where their crossing line links two real Dirac points in the bulk, as

FIG. 3. Spectra with PT being violated. (a) and (b) are gapped after breaking the PT symmetry by removing ($\uparrow A$) and ($\downarrow B$) on the top layer and ($\downarrow A$) and ($\uparrow B$) on the bottom layer, compared with (c) and (d) in Fig. 1 in the main text, respectively.
illustrated by Figs. 1(b) and 1(d), in contrast to the Fermi arc of a Weyl semimetal, which contains only one inclined plane [12]. As low-energy degrees of freedom, Fermi arcs on both surfaces have to preserve the reality of PT as a whole, while either cannot do it alone, since P maps one surface to the other and T is an internal transformation. However, it is noted that the topological protection of the surface Fermi arcs requires that the boundary conditions have to respect the PT symmetry, since the reality of the whole system has to be preserved. For instance, the gapless boundary Fermi arcs are eliminated as shown in Fig. 3, after removing degrees of (↑A) and (↓B) on the top layer and (↓A) and (↑B) on the bottom layer of the slab [σµ and τµ act on (↑, ↓) and (AB), respectively], for which the PT symmetry is broken, since \( \hat{P} = \sigma_3 \Gamma \) maps (↑A) [(↓B)] on the bottom (top) layer to (↑A) [(↓B)] on the top (bottom) layer.

**Summary.**—The real Dirac point has been identified as the real monopole of Z2 type in the real Berry bundle with the reality being endowed by the PT symmetry. The real Dirac semimetal is studied with the essential role of reality being elucidated for global topology in the whole BZ and the nonchiral surface Fermi arcs.

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[29] The inverse of the generator can be obtained by a reflection, so that the direct sum of the original bundle on the surrounding \( S^2 \) and the one from reflection is a trivial bundle that serves as the zero element.
[30] For \( O(N) \) \( [U(N)] \) Berry bundles, the clutching functions of the corresponding principle bundles are valued in the orthogonal (unitary) group \( O(N) \) \( [U(N)] \), or equivalently, the Berry connections are valued in the Lie algebra \( o(N) \) \( [u(N)] \) [31].
[33] Explicitly, the PT operator \( \hat{P} \hat{T} = \sigma_3 \hat{K} \) is related to \( \hat{K} \) by the unitary transformation given by \( e^{i \alpha \pi / 4} \).