Protection against a spin gap in two-dimensional insulating antiferromagnets with a Chern-Simons term

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We propose a mechanism for the protection against spin gapped states in doped antiferromagnets. It requires the presence of a Chern-Simons term that can be generated by a coupling between spin and an insulator. We first demonstrate that in the presence of this term the vortex loop excitations of the spin sector behave as anyons with fractional statistics. To generate such a term, the fermions should have a massive Dirac spectrum coupled to the emergent spin field of the spin sector. The Dirac spectrum can be realized by a planar spin configuration arising as the lowest-energy configuration of a square lattice antiferromagnet Hamiltonian involving a Dzyaloshinskii-Moriya interaction. The mass is provided by a combination of dimerization and staggered chemical potential. We finally show that for realistic parameters, anyonic vortex loop condensation will likely never occur and thus the spin gapped state is prevented. We also propose real magnetic materials for an experimental verification of our theory.

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I. INTRODUCTION

The discovery of high- T_c superconductivity in cuprates led to a flurry of new ideas such as the idea of spin liquid state as one of the possible explanations for the emergence of superconductivity from Mott insulating parent compounds [1] and the importance of the physics of doped antiferromagnets [2]. An exotic spin liquid state from Heisenberg types of spin models involving breaking of discrete symmetries such as parity or time-reversal symmetry, the so-called chiral spin liquid, has been proposed [3]. The effective low-energy theory of such chiral spin state normally involves a topological term called Chern-Simons term [4]. It was originally studied in particle physics [5] and mimics the fractional quantum Hall effect (FQHE) [6] where such term appears as the low-energy effective theory in the bulk [7].

The Chern-Simons term can be generated by a fermion-gauge field coupling when the fermion is integrated out. One obtains a fermion determinant which gives the Chern-Simons term as the action for gauge field A [8],

$$S_{\rm CS}[A] = i \frac{N_f}{2} \frac{e^2}{4\pi} \frac{m_{\psi}}{|m_{\psi}|} \int d^3x \epsilon^{\mu\nu\lambda} A_{\mu} \partial_{\nu} A_{\lambda}, \qquad (1)$$

in Euclidean space-time upon perturbative expansion [9], where N_f is the number of fermion flavors, e is the gauge charge, m_{ψ} is the Dirac fermion mass, and $\epsilon^{\mu\nu\lambda}$ is a totally antisymmetric tensor. The Chern-Simons action depends only on the overall sign of the mass rather than its magnitude even though the mass must be nonzero for the expansion to make sense.

In quantum magnetism, the magnetization curve may show the presence of plateaus. It is well understood that the plateau state corresponds to a state with gapped magnetic excitations while outside the plateau, the spin sector is gapless [10]. In doped antiferromagnets [11], this leads to distinct natures of interaction between the fermions [12]. A state with gapped magnetic excitations is normally associated with preserved continuous symmetry with no long-range magnetic order in the direction transverse to the applied magnetic field [10]. In this work, we show that in the presence of a Chern-Simons term, one can have either one of two scenarios: (i) a spin

gapped state occurs as an analog of the fractional quantum Hall effect (FQHE) with its chiral edge states and (ii) the system is protected against such a spin gapped state. In this paper, we consider 2D antiferromagnets on square lattice, as we did in our previous work [12].

In this paper, we also propose a way to realize a chiral theory as an effective low-energy theory of spin systems explicitly rather than spontaneously by considering doped antiferromagnets on the square lattice. The fermions hop on the lattice on top of a pre-existing magnetic background. Because of a strong Hund coupling, the spin of the electron must be parallel to that of the local spin within the adiabatic approximation. The resulting dynamics is well-described by an effective tight-binding Hamiltonian [13]:

$$H = -t \sum_{\mathbf{r}, \mathbf{r}'} \langle \Omega_{\mathbf{r}} | \Omega_{\mathbf{r}'} \rangle c_{\mathbf{r}'}^{\dagger} c_{\mathbf{r}} + \text{H.c.}, \qquad (2)$$

where the spin sector modifies the hopping integral via the overlap of spin coherent states $|\Omega_{\bf r}\rangle$ between nearest-neighbor sites, a mechanism well established from the studies of anomalous Hall effect [14] and doped antiferromagnets [15]. This coherent state overlap has two effects: providing the background flux for the fermions and the effective spin field that couples to the fermions. The spin sector should take an appropriate classical lowest-energy configuration which produces a staggered $\pi\text{-flux}$ state [16] with dimerization and staggered chemical potential needed to get a massive Dirac fermion spectrum.

II. FIELD THEORY

We use the semiclassical spin path integral approach [10] and start with the Euclidean effective action of a 2D doped antiferromagnet,

$$S_{\phi} = \int d^2x \int d\tau \frac{K_{\tau}}{2} (\partial_{\tau}\phi)^2 + \frac{K_r}{2} (\nabla\phi)^2 + i \left(\frac{S}{a^2}\right) \partial_{\tau}\phi, \tag{3}$$

$$S_{\overline{\psi},\psi} = \int d^2x \int d\tau \overline{\psi} [\gamma^{\mu}(-i\partial_{\mu} - eA_{\mu}) + m_{\psi}]\psi, \quad (4)$$

taking the form of an XY model describing quantum fluctuations around a classical lowest-energy configuration specified by $\mathbf{S} = S(\sin \theta_0 \cos \phi_0, \sin \theta_0 \sin \phi_0, \cos \theta_0)$ plus a Berry phase action whose form follows from the spin configuration we will describe. The ϕ is the quantum fluctuating phase angle field around ϕ_0 [10] and is conjugate to the momentum operator Π , $[\phi_{\mathbf{r}}, \Pi_{\mathbf{r}'}] = i\delta_{\mathbf{r}\mathbf{r}'}$, where Π can be defined in terms of θ [11]. The K_{τ}, K_{r} are stiffness coefficients of the spin sector. The $\overline{\psi}$ and ψ are Dirac spinors representing electrons that couple to the spin sector's ϕ field via the pseudogauge field given as $A_{\mu} = \partial_{\mu} \phi, \mu = \tau, x, y$. Here, two comments are in order. First, the definition of the vector field A_{μ} looks as if it were a pure gauge. However, one has to keep in mind that the field ϕ may have vorticity, which turns out to be at the origin of a nonzero flux for A_{μ} . Second, the action is not gauge invariant and as such the theory is not a genuine gauge theory. We use nevertheless the terminology of pseudogauge field for A_{μ} , which we also refer to as spin field, because of the way it couples to the charge degrees of freedom. When the fermions are integrated out, we get the Chern-Simons term due to the fermion-spin field coupling given in Eq. (1). A similar final action but without a Berry phase has been studied in a different context [17].

Integrating out the massive Dirac fermions in Eq. (4), we obtain a Chern-Simons term in terms of the phase field ϕ :

$$\mathcal{L}_{\text{CS}} = i \frac{\kappa}{2\pi} \epsilon^{\mu\nu\lambda} \partial_{\mu} \phi \partial_{\nu} \partial_{\lambda} \phi \tag{5}$$

with $\kappa = e^2/4$ ($N_f = 1$). We first investigate the effect of this Chern-Simons term added to the action (3) by applying a duality transformation which introduces a Hubbard-Stratonovich auxiliary vector field J_{μ} [10] and re-expresses the full action as

$$\mathcal{L} = \frac{J^{\mu}J_{\mu}}{2K_{\mu}} + i\left(J^{\mu} + \frac{S}{a^{2}}\delta^{\mu}_{\tau}\right)\partial_{\mu}\phi + i\kappa\partial_{\mu}\phi J^{\mu}_{\mathcal{V}}.$$
 (6)

In this case, we have decomposed the phase field into vortexful and regular parts $\phi = \phi_{\mathcal{V}} + \phi_R$ and defined the vortex loop current $J_{\mathcal{V}}^{\lambda} = (1/(2\pi))\epsilon^{\lambda\mu\nu}\partial_{\mu}\partial_{\nu}\phi_{\mathcal{V}}$. Following the boson-vortex duality transformation [10] and working in Euclidean space-time, we obtain the following effective Lagrangian for $J_{\mathcal{V}}^{\mu}$:

$$L[J_{\mathcal{V}}] = \int_{k} J_{\mathcal{V}}^{\mu}(k) \frac{1}{k^{2}} \left[\left(\delta_{\mu\nu} - \frac{k_{\mu}k_{\nu}}{k^{2}} \right) - \kappa \pi \epsilon_{\mu\nu\alpha} k^{\alpha} \right] J_{\mathcal{V}}^{\nu}(-k)$$
$$+ i2\pi \int_{x} \overline{b}_{\mu}(x) J_{\mathcal{V}}^{\mu}(x), \tag{7}$$

where $\overline{b}_{\mu} = \frac{1}{2} (\frac{s}{a^2}) \epsilon_{\tau \mu \nu} x_{\nu}$. As in Ref. [10], the above result was obtained under the assumption that only configurations with closed vortex loops do contribute to the low-energy physics. The effective magnetic field through a vortex loop is given by

$$\epsilon^{\mu\nu\lambda}\partial_{\nu}\hat{b}_{\lambda} = \frac{S}{a^2}\delta^{\mu}_{\tau} + \kappa J^{\mu}_{\mathcal{V}}. \tag{8}$$

We note that we get an extra term $\kappa J_{\mathcal{V}}^{\mu}$ to the effective magnetic field coming from the Chern-Simons term. The important consequence of this is that the resulting effective Berry phase for a vortex loop now has a total contribution coming from an ordinary Berry phase term and a Chern-Simons term and is

given by

$$\Phi_{\text{vortex loop}} = 2\pi S \frac{A}{a^2} q + 2\pi q \kappa \Phi_{\text{other vortex loop}}, \qquad (9)$$

where A is the area of a vortex loop is an integer multiple of a^2 , q is the vorticity of the vortex loop, and $\Phi_{\text{other vortex loop}}$ is the flux of other vortex loops, which is nonzero only if that other vortex loop is knotted across the first vortex loop. Comparing Eqs. (8) and (9) suggests that the Chern-Simons contribution is nonzero only if the two vortex loops are linked.

We verify the above proposition as follows. For two nonintersecting curves γ_1 and γ_2 , which act as mapping from manifold \mathbb{S}_1 (circle) to 3D Cartesian space \mathbb{R}^3 (the 3D Euclidean space in our problem), the linking number is given by

$$\mathcal{N}_{\text{linking}} = \frac{1}{4\pi} \oint_{\gamma_1} \oint_{\gamma_2} \frac{(\mathbf{r}_1 - \mathbf{r}_2) \cdot (d\mathbf{r}_1 \times d\mathbf{r}_2)}{|\mathbf{r}_1 - \mathbf{r}_2|^3}.$$
 (10)

Taking the inverse Fourier transform, with the vortex current given by $\mathbf{J}_{\mathcal{V}} = q d\mathbf{r}$ and endowing the above expression with an overall coefficient κ , the above expression corresponds to vortex current-vortex current interaction,

$$L_{\text{linking}}[J_{\mathcal{V}}] = \int_{L} J_{\mathcal{V}}^{\mu}(k) \frac{1}{k^{2}} (-\kappa \pi \epsilon_{\mu\nu\alpha} k^{\alpha}) J_{\mathcal{V}}^{\nu}(-k), \qquad (11)$$

in Fourier space. This is precisely equal to the topological part of Eq. (7). This implies that the Chern-Simons term leads to the linking between vortex loops, resulting in a nonzero linking number, precisely as implied by Eqs. (8) and (9).

III. ANYON PHYSICS

We now show how the field theory and the results in the previous section lead to the anyonic fractional statistics of the vortex-loop excitations of the 2D XY model (3). Without the Chern-Simons term, the XY model (3) has topological excitations in the form of vortex loops which, in the absence of destructive interference due to the Berry phase, can proliferate. The vortex loops can be considered as world lines of Bose particles and the proliferation of vortex loops corresponds to a condensation of the bosons. This vortex proliferation implies disordering of the dual (ϕ) field, which gives rise to gapped magnetic excitations. In the presence of the Chern-Simons term, vortex loops become anyons and a number of vortex loops must collect together in order to form a boson, and then be able to condense. One then can have either one of the two scenarios: (i) the anyonic vortex loops are able to form bosons and condense, giving rise to a state with gapped magnetic excitations as the analog of FQHE state and (ii) the anyons fail to condense, for which a state with gapped magnetic excitations will not occur.

To show this technically in a simple picture, consider a system of two vortex loops with vorticities q_1 and q_2 , respectively, knotted to each other once ($\mathcal{N}_{\text{linking}} = 1$). The partition function of this system has a contribution of the form

$$e^{i4\pi\kappa q_1q_2 + q_1q_2E_{\text{Coulomb}}}. (12)$$

Apart from the Coulomb interaction, the contribution to the partition function given in Eq. (12) resembles that of linked word lines of an anyon system with its fractional statistics [18]. The first term in the exponent $4\pi \kappa q_1 q_2$ equals twice the statistical angle Θ under the exchange of two anyons realized by vortex loops with vorticities q_1 and q_2 , giving $\Theta = 2\pi \kappa q_1 q_2$. In 3D Euclidean space-time, there exists nontrivial braiding statistics between loops [19]. Our system thus manifests anyonic loop statistics in 3D. With $q_1 = q_2 = 1$ for an elementary vortex loop, we obtain $\Theta = 2\pi\kappa$ and therefore need $\mathcal{N}_{\text{anyon}} = 2\pi/\Theta = 1/\kappa$ anyons to get a boson before the anyons can condense. The occurrence of anyon condensation (and that of spin gapped state) thus depends crucially on the magnitude of the Chern-Simons coupling κ . The Coulomb interaction then determines the critical value of the parameter that characterizes any possible condensation transition for this interacting system. We will show a scenario from a microscopic model where κ takes small enough values, leading to large \mathcal{N}_{anyon} and as a consequence prevents anyon condensation and the occurrence of a state with gapped magnetic excitations.

IV. MICROSCOPIC REALIZATION

In this work, we propose a way to generate a Chern-Simons term explicitly in a spin system from the fermion-spin field coupling. This topological term is obtained upon integrating out the fermions, as summarized in Eq. (1). The key ingredient is to have massive Dirac fermions coupled to the spin field from the spin sector for which we propose to consider the staggered π flux state [16] and we will now describe the scheme to obtain it.

According to Eq. (2), the bare hopping integral is modified by the overlap of spin coherent states $t \to t \langle \Omega_{\mathbf{r}} | \Omega_{\mathbf{r}'} \rangle$. This spin coherent state overlap generally takes complex values, which gives an exponential phase factor where the argument of the exponent, to be referred to as link spin field, contains a static background flux plus the fluctuating part, representing a dynamical spin field. In order to get a Dirac spectrum, the static background flux has to take an appropriate configuration, which can be chosen to be a staggered π flux configuration [16]. In order to attain such a staggered π flux configuration, in turn, the spin sector should take a particular spin configuration accordingly. In Appendix A, we first derive the expression for the link spin field from evaluating the spin coherent state overlap explicitly. Then, from that, we determine the corresponding spin configuration to obtain the staggered π flux state. We then propose a microscopic spin Hamiltonian that stabilizes the required spin configuration. We summarize the results as follows.

Representing the spin as a classical vector $\mathbf{S} = S(\sin\theta\cos\phi, \sin\theta\sin\phi, \cos\theta)$ and evaluating the spin coherent state overlap in Eq. (2) gives the following result for the link spin field [22]:

$$a_{\mathbf{r}\mathbf{r}'} = -\tan^{-1} \left[\frac{\sin(\phi_{\mathbf{r}'} - \phi_{\mathbf{r}})\sin\frac{\theta_{\mathbf{r}}}{2}\sin\frac{\theta_{\mathbf{r}'}}{2}}{\cos\frac{\theta_{\mathbf{r}}}{2}\cos\frac{\theta_{\mathbf{r}'}}{2} + \cos(\phi_{\mathbf{r}'} - \phi_{\mathbf{r}})\sin\frac{\theta_{\mathbf{r}}}{2}\sin\frac{\theta_{\mathbf{r}'}}{2}} \right].$$
(13)

The flux Φ_{\square} through a square plaquette is then given by $\Phi_{\square} = \sum_{\mathbf{rr'} \in \square} a_{\mathbf{rr'}}$ where the link $\mathbf{rr'}$ is taken to be such that the four

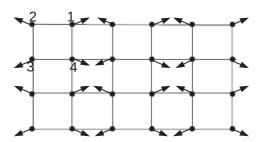


FIG. 1. The planar spin configuration on a square lattice with the labeling of the four sublattices.

links traverse the square in counter-clockwise direction. The "mean-field part" of the link spin field needed to give π -flux state is obtained by assigning $\theta_{\bf r},\phi_{\bf r}$ their classical lowest-energy configuration values. It can be checked from Eq. (13), that a $\pm\pi$ flux through square plaquette can be realized by planar spin configuration where $\theta=\pi/2$ at all sites, while the azimuthal angle takes values such that the four spins on a square rotate by a total of $\pm 2\pi$ as illustrated in Fig. 1. The need for such planar spin configuration fully agrees with the intuition that the total solid angle Φ swept by the four spins around the square should equal 2π for them to produce net flux $\Phi_S = S\Phi$ equals π with S = 1/2.

Similar planar spin configurations can be realized using $J_1 - J_2$ model with impurity [20]. Here, we propose an alternative method. We find that such kind of spin configurations can be stabilized as a classical lowest-energy configuration of the following Hamiltonian, involving Dzyaloshinskii-Moriya (DM) interactions [21]:

$$H = J \sum_{ij} \mathbf{S}_i \cdot \mathbf{S}_j + D \sum_i (S_i^z)^2 + \sum_{ij} \mathbf{D}_{DM}^{ij} \cdot (\mathbf{S}_i \times \mathbf{S}_j).$$
(14)

The easy-plane anisotropy (D>0) favors an in-plane Néel state. To retain global U(1) symmetry, only the component of $\mathbf{D}_{\mathrm{DM}}^{ij}$ along z direction can be allowed to be nonzero. DM interaction usually occurs in lattice systems with low degree of symmetries. Here, we consider a staggered DM interaction where $D_{\mathrm{DM}i,i\pm\hat{x}}^z = -D_{\mathrm{DM}i,i\pm\hat{y}}^z$, while $D_{\mathrm{DM}ij}^x = D_{\mathrm{DM}ij}^y = 0$ [22]. According to Moriya's symmetry analysis [21], this DM pattern is, in principle, allowed as long as the system has no bond-centered inversion symmetry but has mirror planes perpendicular to the bonds and passing through the bond centers.

To find the lowest-energy spin configuration, we rewrite the Hamiltonian (14) in terms of angular variables θ, ϕ as before and minimize H with respect to these variables [22]. The lowest energy configuration of H in Eq. (14) is found to be a planar spin configuration with azimuthal angles $\Delta \phi_{ij} = \phi_j - \phi_i$ between nearest-neighbor spins given by

$$\Delta \phi_{ij} = \tan^{-1} \left(\frac{D_{\rm DM}^{ij}}{J} \right) \tag{15}$$

as shown in Fig. 1. Our calculation confirms that the planar spin configuration is indeed the lowest-energy configuration of *H* at least within a finite regime of the phase diagram defined in

parameter space $(J-D-D_{\rm DM}\ {\rm space})$ and therefore requires no fine tuning.

We then derive from Eq. (2) the low-energy effective theory of fermions around the Dirac points coupled to the spin quantum fluctuations around the lowest energy spin configuration. This is done by expanding the spin vector around its lowest energy orientation and expanding the fermion field operator using gradient expansion in real space and expansion around a Dirac point in momentum space. The details are given in Appendix B. The resulting scalar theory takes the form given in Eq. (3) for the spin sector with $\phi \equiv \phi_G = (\sum_{i=1}^{N_{\text{unitcell}}} \phi_i)/N_{\text{unitcell}}$ (where ϕ_i is the ϕ of the *i*th sublattice) representing the massless Goldstone mode. Here, as the other fields are gapped, gapped magnetic excitations imply a spin gap. The Berry phase term gives a contribution $\exp(i2\pi S)$ to the partition function in imaginary time. For integer S, this equals unity corresponding to constructive interference and one can thus expect the presence of a spin gapped state. The fermion-spin field theory is given by Eq. (4) with $\psi(\mathbf{r}) = (c_D^1(\mathbf{r}), c_D^2(\mathbf{r}), c_D^3(\mathbf{r}), c_D^4(\mathbf{r}))^T$. Here, $c_D^i, i = 1, \dots, 4$ represents the Dirac fermion operator for the i^{th} sublattice, while $\gamma_{\mu}, \mu = 0,1,2$ are 4 × 4 matrices satisfying the Clifford algebra $\{\gamma_{\mu}, \gamma_{\nu}\} = 2g_{\mu\nu}$ and $\overline{\psi} = \psi^{\dagger}\gamma_{0}$ [22]. We take the unit hopping integral t = 1 and unit lattice spacing a = 1 as units of energy and length, respectively. One important result is that the gauge charge is found to be

$$e = -\frac{1}{2\sqrt{2}} \left(1 - \frac{J}{\sqrt{J^2 + D_{DM}^2}} \right)^{\frac{1}{2}},\tag{16}$$

where $\alpha = e^2/(4\pi)$ gives the dimensionless "fine structure constant" of the U(1) gauge theory. This gauge charge is obtained directly as the coupling constant of the derived fermion-spin field theory. It is to be noted that in realistic situations, $|D_{\rm DM}| \ll J$ and as a result, α becomes a small parameter, as is normally the case in quantum field theory. It will be shown in the next section that this fact will play a key role in the mechanism of protection against spin gap that we propose in this work. Strictly speaking, the derivation of our fermion-spin field theory is justified only in the large spin S limit. While we have performed explicit calculation to obtain the quantitative result for the gauge charge e as above using S = 1/2 [22], as an example, our main result regarding a mechanism for protection against spin gap that follows soon will still hold qualitatively and apply to the general spin S.

Now we show how to get the Dirac mass for the massive Dirac fermion spectrum. We consider a perturbation on the ideal system in terms of a dimerization of the strength of the hopping integral and a staggerization of the sublattice chemical potential. Since we aim for a Chern-Simons effective theory, which intrinsically breaks discrete symmetries, the full system consisting of the spin and fermion sectors including the perturbation should break the symmetry under parity and time reversal times any lattice translation. Equation (1) also suggests that Dirac fermion masses from different fermion flavors must necessarily have the same sign in order to ensure nonzero net Chern-Simons term, as masses of opposite signs

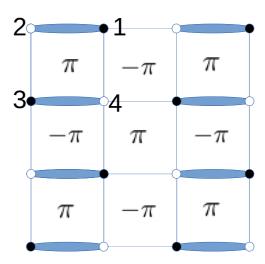


FIG. 2. (Color online) The columnar dimerized plus staggered chemical potential added to the staggered π -flux state.

will give no net parity-breaking effect [23]. We find that this can be achieved by considering a combination of a columnar dimerization and a staggered sublattice chemical potential with profiles shown in Fig. 2. It gives a Dirac mass Hamiltonian

$$H_{m} = \int d^{2}x (\overline{\psi}'_{1} m_{\psi} \psi'_{1} + \overline{\psi}'_{2} m_{\psi} \psi'_{2}), \tag{17}$$

where $\overline{\psi}'_{1(2)}, \psi'_{1(2)}$ represent the first (second) pair of sublattice 1-sublattice-2 spinor $\psi'_{1(2)} = (c'_{1(3)}, c'_{2(4)})^T$ in diagonalized basis, with $\gamma^0 = \tau^z$, the Pauli matrix in this sublattice space. The resulting mass is found to be $m_{\psi} = \sqrt{m_{\text{dimer}}^2 + m_s^2}$ where $m_{\text{dimer}} = \eta - 1$ with $\eta > 1$ is the strength of a dimerized bond with respect to the normal bond and m_s is the staggered sublattice chemical potential. The above perturbation is an example but in principle, any perturbations breaking the same discrete symmetries are expected to give rise to the same phenomena.

V. DISCUSSION

A Chern-Simons term is known to give rise to a mass to the gauge field in the Maxwell-Chern-Simons theory. What we have here is an XY-Chern-Simons theory plus the ever important Berry phase term. Plateaus in XY model without a Chern-Simons term occur when the spin sector is gapped and vortex-loops proliferate corresponding to condensation of bosons. In the presence of the Chern-Simons term, vortex loops are anyon worldlines as we noted from Eq. (12), with a statistical phase $\Theta = 2\kappa\pi = \pi e^2/2$. If the vortex loops could ever condense, we would realize the analog of a fractional quantum Hall state (FQHE) with its chiral edge states. For our microscopic model with gauge charge e as given in Eq. (16), we obtain an upper bound $\Theta = \pi/16$ from the limit $|D_{\rm DM}|/J \to \infty$, which means that in order to obtain condensation of anyons, we need at minimum $\mathcal{N}_{anyon} = 32$ vortex loops to form a boson first before they can ever condense and give rise to a spin gapped state, taking the fact that for bosons, the statistical angle is $\Theta=2\pi$. Realistic situations where $|D_{\rm DM}|/J\ll 1$ require much larger number of anyons for them to condense.

Based on the analogy with the FQHE, condensation of anyons and thus a Chern-Simons-induced spin-gapped state in doped antiferromagnets can occur only when the microscopic parameters give rise to a Chern-Simons term with coupling κ such that $v \equiv 2\kappa = \Theta/\pi = e^2/2 = P/Q$ satisfies a continued fraction expansion condition, where ν is the filling factor of typical values $1/\nu \le 9$ for Laughlin's states [24]. With a minimum of 32 vortex loops needed to form a boson (equivalent to $1/\nu \gg 16$ for the realistic case $|D_{\rm DM}|/J \ll 1$) and the strict condition on the Chern-Simons coupling, it is thus in general very unlikely to form such a vortex loops condensate and the associated spin gapped state. We can therefore conclude that the spin system (14) is protected from being in a spin gapped state, due to the Chern-Simons term induced by the fermion-spin field coupling in doped antiferromagnets with a massive Dirac fermion spectrum. In contrast, the conventional XY model with the Berry phase term would mandate the occurrence of a spin gapped state. On the other hand, if the net Chern-Simons term vanishes, then the protection effect is inactive and one can get back a conventional spin gapped state that occurs for integer spin S. It is to be noted that this result on the novel mechanism for protection against spin gap is valid for general spin S because the applicability of the mechanism is determined more by the value of coupling constants of the spin model (J and D_{DM}) rather than the spin S itself.

It would be interesting to find other microscopic spin models that can generate a Chern-Simons term via similar fermion-spin field coupling as we proposed here and yet are able to induce anyonic vortex loop condensation, giving rise to a spin gapped state, and the analog of the FQHE with chiral edge states in spin systems. Our choice for a square lattice is because a discrete Chern-Simons gauge theory with precisely the same physics as the continuum one can be constructed consistently on this lattice [25] as well as on the kagome lattice [26], due to the presence of one-to-one face-vertex correspondence on these lattices [27]. We would like to propose a study on compounds La₂CuO₄ [28] and LaMnO₃ [29], which are effectively 2D square lattice antiferromagnets as candidate materials to test our theory. Beyond these particular materials, we claim that this novel scenario of spin gap protection is general enough to be expected anytime a Chern-like charge insulator is at play as the high-energy sector. The Kondo lattice model in the triangular lattice may be another possible laboratory [30].

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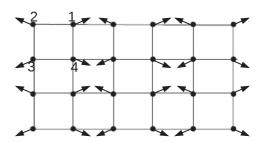


FIG. 3. The (commensurate) planar spin configuration on a square plaquette of the square lattice giving rise to a staggered π -flux state of Affleck-Marston ansatz.

APPENDIX A: THE CLASSICAL LOWEST-ENERGY SPIN CONFIGURATION

A key requirement in our proposal to generate a Chern-Simons term in a spin system is to have a Dirac spectrum for the fermions that are to be doped into the system. In order to get a Dirac spectrum on the square lattice, we need a π -flux state of the type as first proposed by Affleck-Marston and G. Kotliar [16]. The flux per square plaquette is a staggered π and $-\pi$ configuration, where the flux alternates from π to $-\pi$ between adjacent squares. Here, we will show that such a staggered π flux state can be realized by a planar spin configuration illustrated in Fig. 3. The derivation to arrive at such a spin configuration is as follows.

The spin sector couples to the fermion by providing the spin field in the form of Goldstone modes in addition to providing the background staggered π -flux state. This is well described by the following tight-binding Hamiltonian:

$$H = -t \sum_{\mathbf{r}, \mathbf{r}'} \langle \Omega_{\mathbf{r}} | \Omega_{\mathbf{r}'} \rangle c_{\mathbf{r}'}^{\dagger} c_{\mathbf{r}} + \text{H.c.}$$
 (A1)

In spherical coordinates, the classical spin vector is written as $\mathbf{S} = S(\sin\theta\cos\phi, \sin\theta\sin\phi, \cos\theta)$. The spin coherent state for general spin S is given by

$$\begin{split} |\Omega_{\mathbf{r}}(\theta(\mathbf{r}),\phi(\mathbf{r}))\rangle &= e^{iSb_{\mathbf{r}}} \sqrt{(2S)!} \sum_{m_z=-S}^{S} \\ &\times \frac{\left(\cos\frac{\theta_{\mathbf{r}}}{2}e^{i\frac{\phi_{\mathbf{r}}}{2}}\right)^{S+m_z} \left(\sin\frac{\theta_{\mathbf{r}}}{2}e^{-i\frac{\phi_{\mathbf{r}}}{2}}\right)^{S-m_z}}{\sqrt{(S+m_z)!(S-m_z)!}} |S,m_z\rangle, \end{split} \tag{A2}$$

which for S = 1/2 allows us to write

$$|\Omega_{\mathbf{r}}\rangle = e^{i\frac{1}{2}b_{\mathbf{r}}} \left(e^{i\frac{\phi_{\mathbf{r}}}{2}} \cos\frac{\theta_{\mathbf{r}}}{2}, e^{-i\frac{\phi_{\mathbf{r}}}{2}} \sin\frac{\theta_{\mathbf{r}}}{2} \right)^{T},$$
 (A3)

where $b_{\mathbf{r}}$ is a pure gauge function which we can set to a constant function $b_{\mathbf{r}} = b_0$ in the simplest case. The coherent state overlaps for electrons where S = 1/2 is given by

$$\begin{split} \langle \Omega_{\mathbf{r}} | \Omega_{\mathbf{r}'} \rangle &= e^{\frac{1}{2}i(b_{\mathbf{r}'} + \phi_{\mathbf{r}'} - b_{\mathbf{r}} - \phi_{\mathbf{r}})} \\ &\times \left(\cos \frac{\theta_{\mathbf{r}}}{2} \cos \frac{\theta_{\mathbf{r}'}}{2} + e^{i(\phi_{\mathbf{r}} - \phi_{\mathbf{r}'})} \sin \frac{\theta_{\mathbf{r}}}{2} \sin \frac{\theta_{\mathbf{r}'}}{2} \right) . \end{split} \tag{A4}$$

This spin coherent state overlap gives rise to a spin field as follows. Taking the gauge function to be $b_{\mathbf{r}} = -\phi_{\mathbf{r}}$,

$$a_{\mathbf{r}\mathbf{r}'} = -\tan^{-1} \left[\frac{\sin(\phi_{\mathbf{r}'} - \phi_{\mathbf{r}})\sin\frac{\theta_{\mathbf{r}}}{2}\sin\frac{\theta_{\mathbf{r}'}}{2}}{\cos\frac{\theta_{\mathbf{r}}}{2}\cos\frac{\theta_{\mathbf{r}'}}{2} + \cos(\phi_{\mathbf{r}'} - \phi_{\mathbf{r}})\sin\frac{\theta_{\mathbf{r}}}{2}\sin\frac{\theta_{\mathbf{r}'}}{2}} \right]. \tag{A5}$$

We will consider the Dirac spectrum obtained from the staggered π flux state configuration of Affleck-Marston. It can be checked from Eq. (A5) that for $\theta_{\bf r}=\pi/2$ at all sites ${\bf r},{\bf r}'$, we will always get $\sum_{{\bf r}{\bf r}'\in\square}a_{{\bf r}{\bf r}'}=\pm\pi$, as the link spin field is then given by $a_{{\bf r}{\bf r}'}=-\Delta\phi_{{\bf r}{\bf r}'}/2$ where $\Delta\phi_{{\bf r}{\bf r}'}=\phi_{{\bf r}'}-\phi_{{\bf r}}$. This gives rise to the needed π flux per square plaquette. This remarkable result will give rise to the planar spin configuration shown in Fig. 3 for an appropriate choice of Hamiltonian.

We will prove here that this planar spin configuration can be obtained as the lowest-energy configuration of the following Hamiltonian:

$$H = J \sum_{ij} \mathbf{S}_i \cdot \mathbf{S}_j + D \sum_i (S_i^z)^2 + \sum_{ij} \mathbf{D}_{DM}^{ij} \cdot (\mathbf{S}_i \times \mathbf{S}_j),$$
(A6)

consisting of a Heisenberg antiferromagnetic coupling (J > 0), easy-plane anisotropy $(D \ge 0)$, and a Dzyaloshinskii-Moriya terms. To retain global U(1) symmetry, only one (x, y, or z) component of D_{DM}^{ij} normal to the easy plane can be allowed to be nonzero; for concreteness we choose the z component. It will be shown here that in order to get the planar spin configuration, Fig. 3, we must take D_{DM}^{ij} z > 0 on x links and D_{DM}^{ij} z < 0 on y links (of equal magnitude but opposite in sign) or vice versa.

In spherical coordinates, the full Hamiltonian is given by

$$H = JS^{2} \sum_{ij} [\sin \theta_{i} \sin \theta_{j} \cos(\phi_{j} - \phi_{i}) + \cos \theta_{i} \cos \theta_{j}]$$

$$+ DS^{2} \sum_{i} \cos^{2} \theta_{i} + \sum_{ij} D_{DMz}^{ij} S^{2} \sin \theta_{i} \sin \theta_{j}$$

$$\times \sin(\phi_{j} - \phi_{i}). \tag{A7}$$

To find the lowest-energy configuration, we take $\partial H/\partial \phi_i = 0$, $\partial H/\partial \theta_j = 0$ and verify whether $\partial^2 H/\partial \phi_i \partial \phi_j > 0$, $\partial^2 H/\partial \theta_i \partial \theta_j > 0$, $\partial^2 H/\partial \phi_i \partial \theta_j = 0$. We find the lowest-energy configuration solution where the nearest-neighbor spins have polar angles as $\theta_i = \pi/2$ at all sites i and azimuthal angle $\Delta \phi_{ij} = \phi_j - \phi_i$ between nearest-neighbor spins given by

$$\Delta \phi_{ij} = \tan^{-1} \left(\frac{D_{\rm DM}^{ij}}{J} \right), \tag{A8}$$

whereas next-nearest-neighbor spins across the diagonal of square have $\Delta\phi_{ij}=\phi_j-\phi_i=\pm\pi$ as shown in Fig. 3. We verify that the conditions on the second derivatives are indeed satisfied. The above calculation guarantees that such planar spin state indeed exists within a finite regime of the phase diagram defined in the parameter space $(J-D-D_{\rm DM})$ space) and therefore requires no fine tuning.

Now we can describe the precise resulting lowestenergy spin configuration for a particular pattern of the Dzyaloshinskii-Moriya interaction coupling $\mathbf{D}_{\mathrm{DM}}^{ij}$ of relevance to our purpose. For case 1, we take $\mathbf{D}_{\mathrm{DM}}^{ij} = \mathbf{D}_{\mathrm{DM}}$ to be uniform equal to a constant vector at all links, and without loss of generality we take it to point to the z direction for simplicity. It can be checked from Eq. (A8) that unless $|\mathbf{D}_{\rm DM}|/J=1$ giving $\Delta \phi_{ij} = \pi/4.5\pi/4$, one cannot get π flux on a square plaquette for realistic case where $|\mathbf{D}_{\rm DM}|/J \ll 1$. The four spins just cannot be arranged to rotate in total by 2π and at the same satisfy Eq. (A8) from link to link on a square. What this means is that we have to consider a plaquatte involving further distanced spins to get π flux. Thus, in general, we will get an incommensurate state for case 1. For case 2, we consider $\mathbf{D}_{\mathrm{DM}i,i\pm\hat{x}}=(0,\pm d_1,0)$ on each "column" and alternating $\mathbf{D}_{\mathrm{DM}i,i\pm\hat{\mathbf{y}}}=(0,\pm d_2,0)$ on each "row" of square lattice, where $|d_1| = |d_2|$ and so $d_1 = \pm d_2$. It can be easily checked using Eq. (A8) that one can arrange the four spins to rotate by 2π on a square but when we tile up the pattern on a lattice, it will take a relatively large cluster of squares with their spins to form a unit cell from which the full lattice spin configuration can be formed. We, however, get a commensurate staggered π -flux state. This DM pattern occurs in real materials, e.g., La₂CuO₄ [28] but such large unit cell complicates the derivation of a Dirac fermion field theory as the spinor and matrix sizes grow with unit cell size. We then consider the model case 3 where we have $D_{\mathrm{DM}i,i\pm\hat{x}}^z = -D_{\mathrm{DM}i,i\pm\hat{y}}^z$ while $D_{\mathrm{DM}ij}^x = D_{\mathrm{DM}ij}^y = 0$. It can be checked using Eq. (A8) that the four spins on a square will take a configuration as shown in Fig. 3, and this can be nicely repeated over the whole square lattice and gives what we call (commensurate) a planar spin configuration with the resulting commensurate staggered π -flux state with just four sublattices (four sites per unit cell). The needed DM pattern is, in principle, allowed by symmetry according to Moriya's original consideration. We take this latter case as an example but our final results apply to the real materials mentioned above.

We label the four sublattices as shown in Fig. 4. It is to be noted that in such a planar spin configuration, each sublattice has its own θ and ϕ . We therefore have four sets of $(\phi^{(i)}, \theta^{(i)})$ fields: $(\phi^{(1)}, \theta^{(1)}), (\phi^{(2)}, \theta^{(2)}), (\phi^{(3)}, \theta^{(3)}), (\phi^{(4)}, \theta^{(4)})$. From these, we can define four orthogonal fields:

$$\phi_G = \frac{1}{4}(\phi_1 + \phi_2 + \phi_3 + \phi_4), \quad \phi_{t1} = \frac{1}{4}(\phi_1 + \phi_2 - \phi_3 - \phi_4),$$

$$\phi_{t2} = \frac{1}{4}(\phi_1 - \phi_2 + \phi_3 - \phi_4), \quad \phi_{t3} = \frac{1}{4}(\phi_1 - \phi_2 - \phi_3 + \phi_4),$$
(A9)

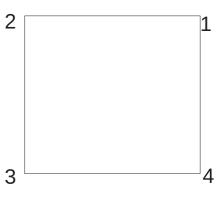


FIG. 4. Labeling of the four sublattices is as shown in the square. The labeling of the four Dirac points in the Brillouin zone also follows the same numbering.

$$\theta_{G} = \frac{1}{4}(\theta_{1} + \theta_{2} + \theta_{3} + \theta_{4}), \quad \theta_{t1} = \frac{1}{4}(\theta_{1} + \theta_{2} - \theta_{3} - \theta_{4}),$$

$$\theta_{t2} = \frac{1}{4}(\theta_{1} - \theta_{2} + \theta_{3} - \theta_{4}), \quad \theta_{t3} = \frac{1}{4}(\theta_{1} - \theta_{2} - \theta_{3} + \theta_{4})$$
(A10)

and the inverse mappings

$$\phi_1 = (\phi_G + \phi_{t1} + \phi_{t2} + \phi_{t3}), \quad \phi_2 = (\phi_G + \phi_{t1} - \phi_{t2} - \phi_{t3}),$$

$$\phi_3 = (\phi_G - \phi_{t1} + \phi_{t2} - \phi_{t3}), \quad \phi_4 = (\phi_G - \phi_{t1} - \phi_{t2} + \phi_{t3}),$$

(A11)

$$\theta_{1} = (\theta_{G} + \theta_{t1} + \theta_{t2} + \theta_{t3}), \quad \theta_{2} = (\theta_{G} + \theta_{t1} - \theta_{t2} - \theta_{t3}),$$

$$\theta_{3} = (\theta_{G} - \theta_{t1} + \theta_{t2} - \theta_{t3}), \quad \theta_{4} = (\theta_{G} - \theta_{t1} - \theta_{t2} + \theta_{t3}).$$
(A12)

The effective low-energy theory of the spin sector from the Hamiltonian (A6) will give rise to a gapless Lagrangian for ϕ_G and gapped ones for ϕ_{ti} , θ_{ti} , i=1,2,3, so that these latter fields can all be integrated out in the low-energy physics [11]. We therefore have only the ϕ_G remaining in the low-energy physics. This ϕ_G is nothing but the Goldstone modes of the rotational symmetry-breaking classical lowest-energy configuration that we have obtained in the beginning: the planar spin configuration. The effective low-energy description of the spin sector obtained upon integrating out all massive fields take the form

$$\mathcal{L} = \frac{K_{\mu}}{2} \partial^{\mu} \phi_{G} \partial_{\mu} \phi_{G} + i \frac{S}{a^{2}} \partial_{\tau} \phi_{G}, \tag{A13}$$

and it is to be noted that the Berry phase term is obtained from summing over the Berry phases of the four sublattices, which finally involves only the diagonal field ϕ_G (Goldstone mode) with overall prefactor S. Since upon integration over τ Berry phase gives the phase factor $\exp(i2\pi S)$, it equals unity for an integer spin S and thus has no effect in such a case. To simplify the notation, in the rest of this calculation, ϕ,θ represent ϕ_G,θ_G .

APPENDIX B: DERIVATION OF THE EFFECTIVE FERMION-SPIN FIELD ACTION OF AN INSULATING ANTIFERROMAGNET WITH A MASSIVE DIRAC SPECTRUM

We have shown in the previous section the microscopic spin model to obtain (gapless) the Dirac spectrum for the fermions doped into the system. We have to verify that the resulting low-energy effective theory is indeed that of Dirac fermions. Furthermore, in order to generate the Chern-Simons term properly, the Dirac fermions must be massive rather than massless. We will therefore also demonstrate in this section the way to generate a massive Dirac fermion spectrum.

To derive the low-energy effective theory of fermions coupled to the spin field, we start from Eq. (A1). A fermion in the staggered π -flux state of Affleck-Marston has the Dirac spectrum $E_k = \pm 2\chi_1 \sqrt{\cos^2 k_x + \cos^2 k_y}$ with four Dirac points at $\mathbf{k} = (\pm \pi/2, \pm \pi/2)$ as shown in Fig. 5 with their numbering. In the four-sublattice description of Affleck-Marston π -flux state, the new first Brillouin zone is defined by the square with $k_x, k_y = \pm \pi/2$.

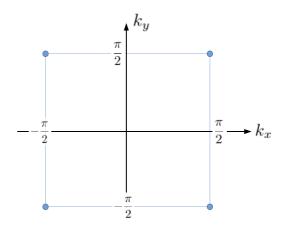


FIG. 5. (Color online) The first Brillouin zone of the staggered π -flux state in the four-sublattice description and the Dirac points (solid circles).

We first expand the fermion field operator around these four Dirac points:

$$c(\mathbf{r}) = \int \frac{d^2k}{(2\pi)^2} c_{\mathbf{k}} e^{i\mathbf{k}\cdot\mathbf{r}} \approx \sum_{i=1}^4 \int_{\mathbf{k}\approx\mathbf{k}_{Di}} \frac{d^2k}{(2\pi)^2} c_{\mathbf{k}} e^{i\mathbf{k}\cdot\mathbf{r}}$$
$$= \sum_{a=1}^4 e^{i\mathbf{k}_{Da}\cdot\mathbf{r}} c_{Da}(\mathbf{r}), \tag{B1}$$

where the slow-fermion field operator $c_{Da}(\mathbf{r})$ is given by

$$c_{Da}(\mathbf{r}) = \int_{\mathbf{k} \approx \mathbf{k}_{Da}} \frac{d^2k}{(2\pi)^2} c_{\mathbf{k}} e^{i(\mathbf{k} - \mathbf{k}_{Da}) \cdot \mathbf{r}}$$

$$\approx \int_{|\mathbf{q}_a| \leqslant q_c} \frac{d^2q}{(2\pi)^2} c_{\mathbf{k}_{Da} + \mathbf{q}_a} e^{i\mathbf{q}_a \cdot \mathbf{r}}$$
(B2)

with $\mathbf{q}_a = \mathbf{k} - \mathbf{k}_{Da}$ and \mathbf{k}_{Da} are the momentum of the a^{th} Dirac points $\mathbf{k}_{Da} = (\pm \pi/2, \pm \pi/2)$ and q_c is some appropriate UV cutoff.

We then perform a gradient expansion and in doing so, we expand the spin coherent state parameterized in terms of these two angles around the respective classical lowest-energy configuration value θ_0 and ϕ_0 at each site (sublattice). Eventually, only the Goldstone mode ϕ_G survives and in the following, the sublattice index is omitted and in the final results ϕ, θ represent ϕ_G, θ_G :

$$\begin{split} |\Omega_{\mathbf{r}}(\theta_{\mathbf{r}},\phi_{\mathbf{r}})\rangle &= \left|\Omega_{\mathbf{r}}(\theta_{\mathbf{r}}^{0},\phi_{\mathbf{r}}^{0})\right\rangle + \delta\theta_{\mathbf{r}}\partial_{\theta}\left|\Omega_{\mathbf{r}}(\theta_{\mathbf{r}}^{0},\phi_{\mathbf{r}}^{0})\right\rangle \\ &+ \delta\phi_{\mathbf{r}}\partial_{\phi}\left|\Omega_{\mathbf{r}}(\theta_{\mathbf{r}}^{0},\phi_{\mathbf{r}}^{0})\right\rangle + \mathcal{O}(\delta\theta^{2},\delta\phi^{2},\delta\theta\delta\phi), \end{split} \tag{B3}$$

$$\begin{split} |\Omega_{\mathbf{r}'}(\theta_{\mathbf{r}'},\phi_{\mathbf{r}'})\rangle &= \left|\Omega_{\mathbf{r}'}\left(\theta_{\mathbf{r}'}^{0},\phi_{\mathbf{r}'}^{0}\right)\right\rangle + \delta\theta_{\mathbf{r}'}\partial_{\theta}\left|\Omega_{\mathbf{r}'}\left(\theta_{\mathbf{r}'}^{0},\phi_{\mathbf{r}'}^{0}\right)\right\rangle \\ &+ \delta\phi_{\mathbf{r}'}\partial_{\phi}\left|\Omega_{\mathbf{r}'}\left(\theta_{\mathbf{r}'}^{0},\phi_{\mathbf{r}'}^{0}\right)\right\rangle + \mathcal{O}(\delta\theta^{2},\delta\phi^{2},\delta\theta\delta\phi), \end{split} \tag{B4}$$

$$c_{Da}^{\dagger}(\mathbf{r}') = c_{Da}^{\dagger}(\mathbf{r}) + \Delta \mathbf{r} \cdot \nabla c_{Da}^{\dagger} + \mathcal{O}(\Delta \mathbf{r}^2),$$
 (B5)

where $\Delta \mathbf{r} = \mathbf{r}' - \mathbf{r}$, we obtain from Eq. (A1),

$$H = -t \sum_{\mathbf{r}, \mathbf{r}'} \sum_{a,b=1}^{4} e^{i(\mathbf{k}_{Db} \cdot \mathbf{r} - \mathbf{k}_{Da} \cdot \mathbf{r}')} \left[c_{Da}^{\dagger}(\mathbf{r}) \langle \Omega_{\mathbf{r}}^{0} | \Omega_{\mathbf{r}'}^{0} \rangle c_{Db}(\mathbf{r}) - c_{Da}^{\dagger} i \Delta \mathbf{r} \cdot \left(-i \langle \Omega_{\mathbf{r}}^{0} | \Omega_{\mathbf{r}'}^{0} \rangle \nabla - e \mathbf{A} \right) c_{Db}(\mathbf{r}) \right] + \text{H.c.}$$

$$= -t \sum_{\mathbf{r}} \sum_{a=1}^{4} e^{-i\mathbf{k}_{Da} \cdot \Delta \mathbf{r}} \left[c_{Da}^{\dagger}(\mathbf{r}) \langle \Omega_{\mathbf{r}}^{0} | \Omega_{\mathbf{r} + \Delta \mathbf{r}}^{0} \rangle c_{Da}(\mathbf{r}) - c_{Da}^{\dagger} i \Delta \mathbf{r} \cdot \left(-i \langle \Omega_{\mathbf{r}}^{0} | \Omega_{\mathbf{r} + \Delta \mathbf{r}}^{0} \rangle \nabla - e \mathbf{A} \right) c_{Da}(\mathbf{r}) \right] + \text{H.c.}$$
(B6)

$$= -t \sum_{\mathbf{r}} \sum_{a=1}^{4} \sum_{m,n=1}^{4} e^{-i\mathbf{k}_{Da}\cdot\Delta\mathbf{r}} \left[c^{m\dagger}_{Da}(\mathbf{r}) \langle \Omega_{\mathbf{r}}^{0} | \Omega_{\mathbf{r}+\Delta\mathbf{r}}^{0} \rangle c_{Da}^{n}(\mathbf{r}) - c^{m\dagger}_{Da} i \Delta \mathbf{r} \cdot \left(-i \langle \Omega_{\mathbf{r}}^{0} | \Omega_{\mathbf{r}+\Delta\mathbf{r}}^{0} \rangle \nabla - e\mathbf{A} \right) c_{Da}^{n}(\mathbf{r}) \right] + \text{H.c.}, \quad (B7)$$

where the superscript m, n = 1, ..., 4 represents the sublattice index of the spin sector and Da = 1, ..., 4 the Dirac point index. In what remains, we take the unit hopping integral t = 1.0 and the unit lattice spacing a = 1.0 as the units of energy and length, respectively. The spin field is given by

$$e\mathbf{A} = -i\left(\delta\theta_{\mathbf{r}'}\langle\Omega_{\mathbf{r}}^{0}|\left(\partial_{\theta}|\Omega_{\mathbf{r}'}^{0}\right)\right) + \delta\theta_{\mathbf{r}}\left(\partial_{\theta}\langle\Omega_{\mathbf{r}}^{0}|\right)|\Omega_{\mathbf{r}'}^{0}\rangle + \delta\phi_{\mathbf{r}'}\langle\Omega_{\mathbf{r}}^{0}|\left(\partial_{\phi}|\Omega_{\mathbf{r}'}^{0}\right)\right) + \delta\phi_{\mathbf{r}}\left(\partial_{\phi}\langle\Omega_{\mathbf{r}}^{0}|\right)|\Omega_{\mathbf{r}'}^{0}\rangle,\tag{B8}$$

where in the continuum limit, we do a gradient expansion by defining $\nabla \theta = \delta \theta_{\mathbf{r'}} - \delta \theta_{\mathbf{r}}$, $\nabla \phi = \delta \phi_{\mathbf{r'}} - \delta \phi_{\mathbf{r}}$. We have included the gauge charge e coupling the spin field to the fermions. The expression (B8) suggests that this spin field is a Berry connection type of pseudogauge field. The actual contribution to the spin field \mathbf{A} in Eq. (B8) only comes from the ϕ field part, corresponding to a Goldstone mode representing a symmetric combination of the sublattice phase fields $\phi_G = (\phi_1 + \phi_2 + \phi_3 + \phi_4)/4$, whereas the terms containing $\delta \theta_{\mathbf{r}}$, $\delta \theta_{\mathbf{r'}}$ give rise to a massive momentum operator and correspond to the massive θ field, which can therefore be integrated out in the low-energy physics.

For S = 1/2, we get the result for the spin field:

$$e\mathbf{A}(\mathbf{r}) = -\frac{1}{2}\sin\theta_0\sin\frac{\Delta\phi_{\mathbf{r},\mathbf{r}+\Delta\mathbf{r}}}{2}\nabla\phi(\mathbf{r}),\tag{B9}$$

where $\theta_0 = \pi/2$. The above result can be derived in a more direct and straightforward way by evaluating the coherent state overlap and matching the complex phase factor exponent with the spin field

$$\langle \Omega_{\mathbf{r}} | \Omega_{\mathbf{r}'} \rangle \sim e^{ia_{\mathbf{r}\mathbf{r}'}}$$
 (B10)

and as one takes the continuum limit, one has $a_{rr'} \to A(r)$. We obtain

$$i\left(\frac{\phi_{\mathbf{r}'}+b_{\mathbf{r}'}-\phi_{\mathbf{r}}-b_{\mathbf{r}}}{2}\right)-i\tan^{-1}\left[\frac{\sin\Delta\phi_{\mathbf{r}\mathbf{r}'}^{0}\sin\frac{\theta_{\mathbf{r}}^{0}}{2}\sin\frac{\theta_{\mathbf{r}'}^{0}}{2}}{\frac{\theta_{\mathbf{r}}^{0}}{2}\cos\frac{\theta_{\mathbf{r}'}^{0}}{2}}\right]-i\frac{1}{2}\sin\frac{\Delta\phi_{\mathbf{r}\mathbf{r}'}^{0}}{2}\sin\theta_{0}\nabla\phi+i\frac{a}{\sin^{2}\theta_{0}}\tan\frac{\Delta\phi_{\mathbf{r}\mathbf{r}'}^{0}}{2}(\Pi_{\mathbf{r}}+\Pi_{\mathbf{r}'})$$

$$\langle\Omega_{\mathbf{r}}|\Omega_{\mathbf{r}'}\rangle=\left|\left\langle\Omega_{\mathbf{r}}^{0}|\Omega_{\mathbf{r}'}^{0}\right\rangle\right|e\qquad ,\qquad (B11)$$

where $\Delta\phi_{\mathbf{r}\mathbf{r}'}^0 = \phi_{\mathbf{r}'} - \phi_{\mathbf{r}}$ and

$$\langle \Omega_{\mathbf{r}}^{0} | \Omega_{\mathbf{r}'}^{0} \rangle = \sin \theta_{0} \cos \frac{\Delta \phi_{\mathbf{r}\mathbf{r}'}}{2},$$
 (B12)

where $\theta_0 = \pi/2$ for our planar spin configuration. We can immediately identify

$$e\mathbf{A}(\mathbf{r}) = -\frac{1}{2}\sin\theta_0\sin\frac{\Delta\phi_{\mathbf{r},\mathbf{r}+\Delta\mathbf{r}}^0}{2}\nabla\phi(\mathbf{r})$$
 (B13)

in complete agreement with Eq. (B9), while the last term in Eq. (B11) is dependent on the momentum operator:

$$\Pi_{\mathbf{r}} = -\frac{1}{2} \left(\delta \theta_{\mathbf{r}} \sin \theta_{\mathbf{r}} + \frac{1}{2} \delta \theta_{\mathbf{r}}^2 \cos \theta_{\mathbf{r}} \right), \tag{B14}$$

which is massive and therefore in the low-energy physics can be integrated out to give simply a renormalization correction to the coefficients (couplings) in the effective fermion-spin field action without changing the physics. It is to be noted that the expression for spin field (B13) is proportional to the gradient of the phase field, as we expected. Following our definition $A_{\mu} = \partial_{\mu} \phi$ and using Eqs. (A8) and (B13), the above result gives as the effective gauge coupling:

$$e = -\frac{1}{2\sqrt{2}} \left[\left(1 - \frac{J}{\sqrt{J^2 + D_{\rm DM}^2}} \right) \right]^{\frac{1}{2}}.$$
 (B15)

We note that if we set the Dzyaloshinskii-Moriya interaction to zero $D_{\rm DM} \to 0$, the whole picture breaks down and the gauge charge equals zero. So, DM interaction is crucially needed here.

As can be seen from Fig. 5, the Dirac points $\mathbf{k}_{Da} = (\pm \pi/2, \pm \pi/2)$ are located precisely at the corners of the first Brillouin zone. The implication of this is that the four Dirac points are now equivalent to each other and it suffices to consider only one Dirac point, e.g., the $\mathbf{k}_{D1} = (\pi/2, \pi/2)$ Dirac point. In the low-energy limit, we define a Dirac 4-spinor associated with this representative Dirac point $\psi(\mathbf{r}) = (c_D^1(\mathbf{r}), c_D^2(\mathbf{r}), c_D^3(\mathbf{r}), c_D^4(\mathbf{r}))^T$. Here, $c_D^i(\mathbf{r}), i = 1, \dots, 4$ represents the Dirac fermion annihilation operator for the i^{th} sublattice of the spin sector in the planar configuration Fig. 3.

The low-energy Hamiltonian is

$$H = \int d^2x \overline{\psi} \gamma^{\mu} (-i \partial_{\mu} - e A_{\mu}) \psi, \qquad (B16)$$

where $\mu = t, x, y \equiv 0, 1, 2$, and $\overline{\psi} = \psi^{\dagger} \gamma^{0}$.

The next task is to find the γ matrices. We obtain the following result:

$$\gamma^{x} = i\gamma^{t} \begin{pmatrix} 0 & \gamma_{12}^{x} & 0 & 0\\ \gamma_{21}^{x} & 0 & 0 & 0\\ 0 & 0 & 0 & \gamma_{34}^{x}\\ 0 & 0 & \gamma_{43}^{x} & 0 \end{pmatrix},$$
(B17)

$$\gamma^{y} = i\gamma^{t} \begin{pmatrix} 0 & 0 & 0 & \gamma_{14}^{y} \\ 0 & 0 & \gamma_{23}^{y} & 0 \\ 0 & \gamma_{32}^{y} & 0 & 0 \\ \gamma_{41}^{y} & 0 & 0 & 0 \end{pmatrix},$$
(B18)

where the matrix elements are given as follows:

$$\gamma_{12}^{x} = ite^{-ik_{D1x}} \langle \Omega_{\mathbf{r}2}^{0} | \Omega_{\mathbf{r}'1}^{0} \rangle = (\gamma_{21}^{x})^{\dagger},
\gamma_{14}^{y} = ite^{-ik_{D1y}} \langle \Omega_{\mathbf{r}4}^{0} | \Omega_{\mathbf{r}'1}^{0} \rangle = (\gamma_{41}^{y})^{\dagger},
\gamma_{23}^{y} = ite^{-ik_{D1y}} \langle \Omega_{\mathbf{r}3}^{0} | \Omega_{\mathbf{r}'2}^{0} \rangle = (\gamma_{32}^{y})^{\dagger},
\gamma_{34}^{x} = ite^{ik_{D1x}} \langle \Omega_{\mathbf{r}4}^{0} | \Omega_{\mathbf{r}'3}^{0} \rangle = (\gamma_{43}^{x})^{\dagger}.$$
(B19)

These matrix elements need to be "normalized" in such a way that each of them is of unit magnitude. We have to find the appropriate $\gamma^0 \equiv \gamma^t$. It needs to be checked that the γ matrices satisfy Dirac algebra $\{\gamma^\mu, \gamma^\nu\} = 2\delta^{\mu\nu}$.

Taking t = 1, we obtain the following gamma matrices valid for a spin system in a planar configuration on the square lattice:

$$\gamma^{x} = i\gamma^{t} \begin{pmatrix} 0 & 1 & 0 & 0 \\ 1 & 0 & 0 & 0 \\ 0 & 0 & 0 & -1 \\ 0 & 0 & -1 & 0 \end{pmatrix} = \begin{pmatrix} 0 & i & 0 & 0 \\ -i & 0 & 0 & 0 \\ 0 & 0 & 0 & -i \\ 0 & 0 & i & 0 \end{pmatrix},$$
(B20)

$$\gamma^{y} = i\gamma^{t} \begin{pmatrix} 0 & 0 & 0 & 1\\ 0 & 0 & 1 & 0\\ 0 & 1 & 0 & 0\\ 1 & 0 & 0 & 0 \end{pmatrix} = \begin{pmatrix} 0 & 0 & 0 & i\\ 0 & 0 & -i & 0\\ 0 & i & 0 & 0\\ -i & 0 & 0 & 0 \end{pmatrix}$$
(B21)

with

$$\gamma^{t} = \begin{pmatrix} 1 & 0 & 0 & 0 \\ 0 & -1 & 0 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & -1 \end{pmatrix} = (\gamma^{t})^{-1}.$$
 (B22)

We aim this Dirac fermion system to produce a Chern-Simons term. For that, we need the spectrum to be massive. We need to make sure that the Dirac masses from the first two sublattices and the second two sublattices have the same sign, otherwise the net Chern-Simons term vanishes. Such mass term can be obtained in several different ways. First, using a staggered sublattice chemical potential, which represents

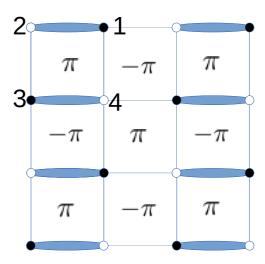


FIG. 6. (Color online) The columnar dimerized π -flux state of Affleck-Marston ansatz.

some internal sublattice or microscopic degree of freedom, which gives rise to an effective chemical potential alternating in sign between the four sublattices of the square lattice. In this case, the mass term from a staggered chemical potential takes the form $H_M = \int d^2x \psi^{\dagger} [\mathrm{diag}(m_s, -m_s, m_s, -m_s)] \psi$, where $[\mathrm{diag}(\cdots)]$ represents a diagonal matrix with the diagonal elements as given.

The second method is to consider a dimerization of the strength of the hopping integral t in the tight-binding Hamiltonian (A1). Dimerization has been known to produce a mass for the Dirac fermions. We consider one of the simplest dimerization patterns as shown in Fig. 6. The strength of the dimerized bonds is larger than the undimerized bonds. The link spin field and the flux per plaquatte remain the same as those of the original Affleck-Marston π -flux state. This dimerization does open up a gap on the Dirac points in the 4×4 form of the Dirac theory, which for the dimerization pattern shown in Fig. 6 takes the form

$$\mathcal{H}_{M}^{\text{dimer}} = \gamma^{t} \begin{pmatrix} 0 & m_{0} & 0 & 0\\ m_{0}^{*} & 0 & 0 & 0\\ 0 & 0 & 0 & m_{0}^{s}\\ 0 & 0 & m_{0}^{s*} & 0 \end{pmatrix}, \quad (B23)$$

where $m_0 = -(\eta - 1) \exp(-i\pi/4 - ik_{Dx}), m_0^s = -(\eta - 1) \exp(-i\pi/4 + ik_{Dx})$ with $\eta > 1$ is the relative strength of a dimerized bond with respect to an undimerized bond of unit strength. Using $k_{Dx} = \pi/2$ from the location of the Dirac point, we obtain $m_0^s = -m_0 = (\eta - 1) \exp(-3i\pi/4)$. So, the Dirac mass Hamiltonian induced by dimerization, Fig. 6, can be written as

$$H_{M} = \int d^{2}x \overline{\psi} \mathcal{H}_{M}^{\text{dimer}} \psi = \int d^{2}x \overline{\psi} \gamma^{t} \frac{(\eta - 1)}{\sqrt{2}} (\tau^{x} - \tau^{y}) \sigma^{z} \psi,$$
(B24)

where τ is the Pauli matrix defined within each of the 1-2 and 3-4 sublattice pairs while σ is the Pauli matrix connecting the two pairs. In the rest of this derivation, we always

imply

$$H_M = \int d^2x \overline{\psi} \mathcal{H}_M \psi = \int d^2x \psi^{\dagger} \mathbb{H}_M \psi, \tag{B25}$$

where $\mathbb{H}_M = (\gamma^t)^{-1} \mathcal{H}_M$ is the Hermitian mass matrix derived directly from the tight-binding Hamiltonian.

Evaluating the mass Hamiltonian matrix (B23) at the representative Dirac point, we obtain the following result for mass matrix:

$$\mathcal{H}_{M}^{\text{dimer}} = \gamma^{t} \begin{pmatrix} 0 & \frac{(\eta - 1)}{\sqrt{2}}(1 + i) & 0 & 0\\ \frac{(\eta - 1)}{\sqrt{2}}(1 - i) & 0 & 0 & 0\\ 0 & 0 & 0 & -\frac{(\eta - 1)}{\sqrt{2}}(1 + i)\\ 0 & 0 & -\frac{(\eta - 1)}{\sqrt{2}}(1 - i) & 0 \end{pmatrix}.$$
(B26)

The above result, however, suggests that the masses are opposite in sign between the two (two-sublattice) pairs of the four-sublattice theory (1-2 sublattice pair versus 3-4 sublattice pair). The net Chern-Simons term will be zero in this case. To remedy this, we combine columnar or staggered dimerization with a staggered sublattice potential where the staggerization takes a two-sublattice structure; the chemical potentials of sublattices 1 and 3 are equal but opposite to the chemical potentials of sublattices 2 and 4. The resulting mass matrix is

$$\mathcal{H}_{M}^{\text{dimer+stagg.chem.pot}} = \gamma^{t} \begin{pmatrix} m_{s} & -\frac{(\eta-1)}{\sqrt{2}}(1+i) & 0 & 0\\ -\frac{(\eta-1)}{\sqrt{2}}(1-i) & -m_{s} & 0 & 0\\ 0 & 0 & m_{s} & \frac{(\eta-1)}{\sqrt{2}}(1+i)\\ 0 & 0 & \frac{(\eta-1)}{\sqrt{2}}(1-i) & -m_{s} \end{pmatrix}, \tag{B27}$$

where m_s is the mass from the staggered sublattice chemical potential. It can readily be seen that the above mass matrix will give rise to a nonzero net Chern-Simons term since the mass block matrix of the 1-2 sublattice pair is not opposite of that of the 3-4 sublattice pair as the staggered sublattice potential-induced mass term matrix has precisely the same form and sign in both sectors. If upon diagonalization the masses are of the same sign in both sectors, which is indeed the case at least for certain range of parameters m_s , the net Chern-Simons term is nonzero. Performing this 4×4 matrix diagonalization of the corresponding \mathbb{H}_M , we find that the Dirac mass Hamiltonian can be written as

$$H_m = \int d^2x (\overline{\psi}'_1 m_{\psi} \psi'_1 + \overline{\psi}'_2 m_{\psi} \psi'_2),$$
 (B28)

where $\overline{\psi}'_{1(2)}, \psi'_{1(2)}$ represent the first (second) pair of sublattice 1-sublattice-2 spinor $\psi'_1 = (c'_1, c'_2)^T$ in a diagonalized basis, with $\gamma^0 = \tau^z$ and mass $m_\psi = \sqrt{m_{\rm dimer}^2 + m_s^2}$, where $m_{\rm dimer} = \eta - 1$ for both pairs, indicating that the two pairs have identical mass and thus a nonzero net Chern-Simons term eventually. This result is nicely consistent with symmetry considerations because the combination of columnar or staggered dimerization plus staggered chemical potential completely breaks parity. It can also be checked that this dimerization plus staggered chemical potential breaks the invariance under time reversal times any lattice translation and this guarantees a nonzero net Chern-Simons term.

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