

Ferromagnetism and Spin-Valley liquid states in Moiré Correlated Insulators

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Motivated by the recent observation of evidence of ferromagnetic and correlated insulating states in systems of Moiré superlattices, we study a two-orbital quantum antiferromagnetic model on the triangular lattice, where the two orbitals physically correspond to the two valleys of the original graphene sheet. For simplicity, the model is a $Sp(2)^s \otimes Sp(2)^v$ symmetry, where the two $Sp(2)$ symmetries correspond to the rotational invariance in the spin and valley space respectively. Through analytical argument, Schrieffer boson analysis and also Monte Carlo simulations, we find that even though all the couplings in the Hamiltonian are antiferromagnetic, there is still a region in the phase diagram with fully polarized ferromagnetic order. We argue that a Zeeman field can drive metal-insulator transition in our picture, as observed experimentally. We also construct spin liquids and topological ordered phases at various limits of the model. After doping the model with extra charge carriers, the system most likely becomes spin-triplet/valley-singlet $d+id$ topological superconductor as predicted previously.

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I. INTRODUCTION

Recently, a series of surprising correlated physics such as superconductivity and insulators at commensurate fractional fillings have been discovered in the systems of Moiré superlattices¹⁻⁹. These discoveries have motivated experimental and theoretical studies¹⁰⁻³⁹. For different reasons, these systems can have an exotic electronic band structure with charge neutrality, i.e. certain interaction effects are significantly enhanced. A consensus of the evidence for the observed insulator and superconductivity has not yet been reached. In the two-orbital extended Hubbard model on the triangular lattice, as proposed in Ref. 10 (physically, the orbital space corresponds to the space of two Dirac valleys in the original Billoin zone of graphene), which at least describes the bilayer graphene and the general bilayer structure (MBL)^{1,5}, as well as the twisted double bilayer graphene (DBL)⁷⁻⁹, the certain twisted angle and out-of-plane electric field (displacement field), since in these cases there is no symmetry protected band touching below the Fermi level, and the isolated bands behave as trivial quantum valley topological states^{11,37,38,44-46}. This two-orbital model could naturally predict either a spin-triplet¹⁰ or spin-singlet¹² $d+id$ topological superconductor, depending on the sign of the on-site Hund's coupling.

Evidence of spin-triplet pairing is predicted previously¹⁰ and recently found in DBL^{7,9}. Evidence of ferromagnetic correlated insulator at a filling away from charge neutrality was discovered in the same systems⁷⁻⁹. In DBL, besides clear ferromagnetic signature observed at the $1/2$ filling insulator⁷⁻⁹, it was also observed in correlated insulators at $1/4$ and $3/4$ fillings to generate a displacement field⁹, whose an effect is likely just a spin polarizing Zeeman effect. This observation relies on that in DBL, at $1/4$ and $3/4$ fillings are close to a ferromagnetic correlated insulator, and a Zeeman

field could drive a metal-insulator transition. Evidence of ferromagnetism at the insulating phase in $3/4$ filling away from the charge neutrality was also found in a one-dimensional system⁶.

Motivated by these experiments, in this work we investigate quantum spin-valley model on the triangular lattice in the one-orbital site, which corresponds to either $1/4$ filling or $3/4$ filling in the Moiré superlattice. The Hamiltonian of this model reads

$$H = \sum_{\langle i,j \rangle} \sum_{a,b=1}^3 J T_i^{ab} T_j^{ab} + J^s \sigma_i^a \sigma_j^a + J^v \tau_i^b \tau_j^b, \quad (1)$$

where σ^a and τ^b are Pauli operators in the spin and valley spaces, and $T^{ab} = \sigma^a \otimes \tau^b$. Here $J^s = J^v = J$, this model becomes the $Sp(4)$ quantum antiferromagnetic model in the dual representation on each site. The $Sp(4)$ symmetry is broken by the Hund's coupling¹⁰, which in general cases $J^v > J > J^s$, if we choose the standard sign of the Hund's coupling which favors large spin on each site. But we assume that the $Sp(4)$ breaking effect is not strong enough to change the sign of J^s , J^v and J , i.e. we keep all three coupling constants positive, i.e. antiferromagnetic. Indeed, since the Hund's coupling originates from the exchange coupling which involves overlap between wave functions at the two valleys, the Hund's coupling should be a relatively weak effect since the inter-valley wave function overlap is expected to be small because large momentum transfer between the two valleys is suppressed by the long wavelength modulation of the Moiré superlattice. For simplicity, we ignore other effects such as breaking the $Sp(4)$ symmetry, such as valley-dependent pairing¹¹, in the spin-valley model Eq. 1 in the valley space as its own $Sp(2)^v$ symmetry.

II. DERIVATION OF THE SPIN-VALLEY MODEL

The model Eq. 1 can be derived in the standard perturbation theory starting from a Hubbard model plus a σ -site σ -d's coupling on the triangular lattice. As we introduced in the introduction, this model at least applies to Py/Bi ^{1,5} and Dy ⁷⁻⁹ in certain twisted angle and displacement field, since in these cases there is a symmetry protected band touching below the Fermi level, and there are band with correlated physics as trivial unitary valley topological behavior (although the exact valley topological behavior is debated in the models used in the literature)^{11,37,38,44-46}.

$$H = H_t + H_J + H_K + \dots \quad (2)$$

$$H_t = -t \sum_{\langle i,j \rangle} \sum_{\alpha=1}^4 \left(c_{i,\alpha}^\dagger c_{j,\alpha} + \text{c.c.} \right), \quad (3)$$

$$H_J = U \sum_j (n_j - 1)^2, \quad (4)$$

$$H_K = -V \sum_j \left(\vec{\sigma}_j \right)^2 + V \sum_j \left(\vec{\tau}_j \right)^2, \quad (5)$$

where $c_{j,\alpha}^\dagger, c_{j,\alpha}$ are electron creation/annihilation operators in the j th site and α th spin/orbital degree of freedom. $n_j = \sum_{\alpha} c_{j,\alpha}^\dagger c_{j,\alpha}$ is the total particle number per site, and $\frac{1}{2} \hat{\sigma}_j^a = \frac{1}{2} c_{j,\alpha}^\dagger \sigma^a c_{j,\alpha}$ is the total σ -site spin operator. The H_J term is the σ -site on-site interaction, and the associated natural sign of the σ -d's coupling is $V > 0$, as a result of charge interaction. We treat the spin term H_t perturbatively. This amounts to integrating out the charge degree of freedom to obtain an effective spin-valley model in the correlated insulator phase.

We follow the standard approach of degenerate perturbation theory. At the unperturbed level, the ground state of $H_J + H_K$ has precisely one electron per site, and the projection of operator to the ground state manifold reads

$$\mathcal{P} = \prod_j (-1)^{\frac{1}{6} n_j (n_j - 2) (n_j - 3) (n_j - 4)}. \quad (6)$$

Considering a pair of nearest neighbor sites on the Py/Bi sublattice, the ground state manifold can be further divided into four sectors with correspond to singlet/triplet and valley-singlet/triplet states. We can write

$$\mathcal{P} = \mathcal{P}_{ss} + \mathcal{P}_{st} + \mathcal{P}_{ts} + \mathcal{P}_{tt}, \quad (7)$$

where (for example) $\mathcal{P}_{st}$ means the projection to singlet/valley-triplet states. The effective Hamiltonian can be calculated as

$$H_{\text{eff}} = \mathcal{P} H_t \frac{1}{E_0 - H_J - H_K} H_t \mathcal{P}, \quad (8)$$

where E_0 is the ground state energy for the two-site problem. A detailed analysis of the intermediate states considering the actual hopping process can be found in Ref. 10. In addition, $\mathcal{P}_{st}$ and $\mathcal{P}_{ts}$ contribute to the effective Hamiltonian which takes a diagonal form in this basis

$$H_{\text{eff}} = -\frac{2t^2}{U+4V} \mathcal{P}_{st} - \frac{2t^2}{U-4V} \mathcal{P}_{ts}. \quad (9)$$

Regarding the σ -d's of the Py/Bi generators on the nearest neighbor sites, the effective Hamiltonian is equivalent to the spin-valley model Eq. 1 with the coupling constants given by

$$\begin{aligned} J^s &= J - \frac{t^2}{U} \left(\frac{2V}{U} + O\left(\frac{V}{U}\right)^2 \right), \\ J^v &= J + \frac{t^2}{U} \left(\frac{2V}{U} + O\left(\frac{V}{U}\right)^2 \right), \\ J &= \frac{t^2}{4U} \left(1 + O\left(\frac{V}{U}\right)^2 \right). \end{aligned} \quad (10)$$

There is a Z_2 symmetry regarding the sign of the σ -d's coupling. The coupling constants transform as $J^s/J \leftrightarrow J^v/J$ when $V \leftrightarrow -V$, as is actually expected from the form of the σ -d's coupling Eq. 5.

III. THE Py/Bi $\otimes 120^\circ$ STATE

At least in certain limit, i.e. $J^v \gg J \gg J^s > 0$, it is fairly easy to see that the σ -d's agentic could emerge in model Eq. 1 with all the σ -d's agentic coupling constants. First of all, the following state will always be an eigen state of the σ -d's Hamiltonian:

$$|\text{FM}\rangle = \left(\prod_i |\sigma_i^z = +1\rangle \right) \otimes |\text{FM}\rangle \text{ of } \vec{\tau}_{\{i\}}. \quad (11)$$

This state is a direct product of two parts: the first part is a fully polarized σ -d's agentic state of the σ -d's $\vec{\sigma}_i$ space; the second part is the ground state of the nearest-neighbor σ -d's agentic unitary $\vec{\tau}_{\{i\}}$ on the triangular lattice in the $\vec{\tau}_i$ space. Although we cannot write down the explicit form of the exact microscopic wavefunction $|\text{FM}\rangle$ of $\vec{\tau}_{\{i\}}$, we do know that this state is as a 120° antiferromagnetic order with reduced symmetry due to the unitary fluctuations and geometric frustration. This state Eq. 11 is always the eigen state of Eq. 1 because a fully polarized σ -d's agentic state is the eigen state of operator $\vec{\sigma}_i \cdot \vec{\sigma}_j$ on $i, j < i, j >$. When in the limit of $J^v \gg J \gg J^s > 0$, this eigen state $|\text{FM}\rangle$ is also the ground state, because in turn, each σ -d's $\vec{\sigma}_i$ will see a background “effective” σ -d's agentic coupling

$$J_{\text{eff}} = J^s + J(\vec{\tau}_i \cdot \vec{\tau}_j). \quad (12)$$

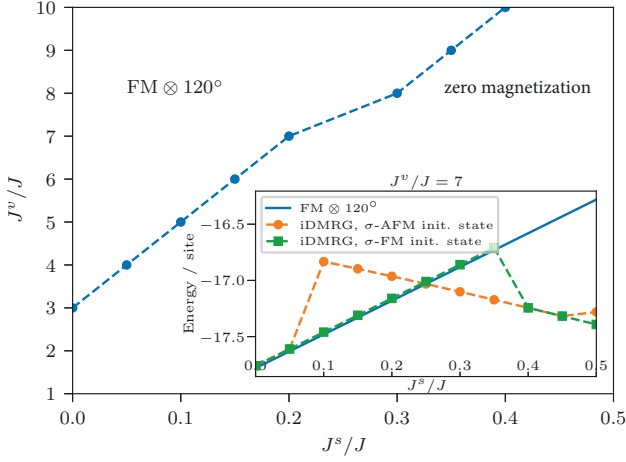


Fig. 1: Phase boundary of the $\mathbb{A}_1$ 120 state obtained using iDMRG on finite cylinders with circumference $L_y = 6$. The inset shows the energy per site obtained for $J^v/J = 7$ as a function of J^s/J . Green squares (orange circles) indicate the energies obtained using iDMRG if the spins are initialized in a $\mathbb{A}_1$ ($\mathbb{A}_2$) product state. The solid blue line indicates the energy expected for the $\mathbb{A}_1$ 120 state.

Because $\phi_i \phi_j < 0$ for the 120 state of ϕ_i , for large enough J the spins will see an effective ferromagnetic coupling, even though in the original model Eq. 1 all the couplings are antiferromagnetic.

For a fixed large J^v , while increasing J^s , eventually $J^s \phi_i \phi_j$ will not be strong enough to overcome the antiferromagnetic coupling J_s , hence we expect to see a transition from the $\mathbb{A}_1 \gg 120$ state to another state without ferromagnetic order. Numerically⁴⁷, $\phi_i \phi_j$ is found to be ~ 0.73 for the triangular lattice, using an antiferromagnet. If we evaluate the energy of FM in Eq. 11, while increasing J_s/J , this state is no longer the ground state for $J_s/J > 0.73$. Hence intuitive arguments suggest an upper bound for the transition point of J^s/J .

DMRG simulation of the spin-valley model

Our previous numerical evidence for the fact that the $\mathbb{A}_1 \gg 120$ state is the ground state in the J^v/J vs J^s/J plot and obtain the phase boundary of this state. To this end we use the density matrix renormalization group (DMRG) method^{48,49}. We note that in finite systems, bond indices can introduce strong oscillations in the expectation value of $\phi_i \phi_j$ on nearest neighbor bonds (which are expected to be uniform in the valley degree of freedom in the 120 state), and thus a careful effective coupling seen by the spins. To avoid subleading effects we use infinite DMRG⁵⁰. We observe uniform bond expectation values in the valley degree of freedom in the 120 state (indeed enough cylinders are to be considered). We relax our analysis on cylinders of circumference $L_y = 6$, for which we obtain an upper bound expectation value $\phi_i \phi_j \approx 0.7$, consistent with Ref. 47, in all simulations below the error bar.

For our numerical simulations we use the Kohn-Sham⁵¹. We assume a 3-site unit cell along the chain to allow for the formation of a 120 state in the valley and/or spin degrees of freedom. The valley degree of freedom is initialized in the total $\phi^z = 0$ sector and ϕ^x and ϕ^y are both conserved. The spin degree of freedom is initialized either in the fully polarized state, or a classical antiferromagnetic state with total $S^z = 0$. The maximal bond dimension in our simulations is $M = 1000$.

We find that at large J^v and for all J^s , the system indeed converges to a fully polarized spin $\mathbb{A}_1$ and a 120-valley ordered state, independent of the initial conditions. At larger J^s we observe a state without any net magnetization. At this stage we cannot conclude as to the nature of the effective region with zero magnetization (for larger J^s and $J^s < J$, for the total $S^z = 0$ sector and going to bond dimensions of up to 4,000 we have not identified a clear order for the spin degree of freedom), but in the next section we will explore some possibilities for testing liquid states and topological orders for this region of the phase diagram. The ground state energy obtained using iDMRG for a fixed $J^v/J = 7$ as a function of J^s/J , for the two initial states, is shown in the inset of Fig. 1. The solid blue line in the inset indicates the energy expected for the $\mathbb{A}_1 \gg 120$ state, calculated using a uniform bond expectation value $\phi_i \phi_j \approx 0.7$ that we obtain for the 120 state in finite cylinders of circumference $L_y = 6$ as mentioned above. We estimate the position of the phase boundary for each J^v/J to be at the J^s/J for which the lowest energy obtained using iDMRG drops below the one expected for the $\mathbb{A}_1 \gg 120$ state. The resulting phase boundary as a function of J^v/J and J^s/J is shown in the main plot of Fig. 1. Our results for the ground state energy and the magnetization across the phase boundary both suggest that the transition between the ferromagnetic order and the paramagnet is a first order level-crossing.

Schwinger boson analysis

We can also construct the $\mathbb{A}_1 \gg 120$ state using the Schwinger boson formalism. We first derive a four-component Schwinger boson b_j to represent the four states of the total representation under both the spin and valley $SU(2)$ symmetries, and also a four-component representation of the enlarged $SU(4)$ symmetries. The Schwinger boson Hilbert space is subject to a local constraint

$$\sum_{i=1}^4 b_j^\dagger b_j = \tau \quad (13)$$

Physically $\tau = 1$, but in the Schwinger boson mean field calculation τ is often treated as a tunable parameter. The Hamiltonian Eq. 1 can be reorganized to the following form:

$$H = J_{ts} \sum_{ij} b_{ij}^\dagger b_{ij} + J_{ss} \sum_{ij} b_{ij}^\dagger b_{ij} + J_{st} \sum_{ij} b_{ij}^\dagger b_{ij} + J_{tt} \sum_{ij} b_{ij}^\dagger b_{ij} \quad (14)$$

The operators $\vec{\Delta}_{ij}^{ts}$ and $\vec{\Delta}_{ij}^{st}$ are the spin-triplet/quartet-singlet, and spin-singlet/quartet-triplet raising operators between Su₂ gauge bosons b_α at site i, j :

$$(\vec{\Delta}_{ij}^{ts}, \vec{\Delta}_{ij}^{st}) = b_i^\dagger (\sigma^{32}, \sigma^{02}, \sigma^{12}, \sigma^{23}, \sigma^{20}, \sigma^{21}) b_j, \quad (15)$$

where $\sigma^{ab} = \sigma^a \otimes \tau^b$, $\sigma^0 = \tau^0 = 1_{2 \times 2}$. The Δ^{ss} and Δ^{tt} are the singlet/singlet, and triplet/triplet raising operators, for example $\Delta_{ij}^{ss} = b_i^\dagger \sigma^{22} b_j$.
In Eq. A,

$$J_{ts} = -\frac{1}{4}(3J_v + 3J - J_s), \quad J_{st} = -\frac{1}{4}(3J + 3J_s - J_v),$$

$$J_{ss} = \frac{3}{4}(3J - J_s - J_v), \quad J_{tt} = \frac{1}{4}(J + J_s + J_v). \quad (16)$$

In the most natural parameter regime $J_v > J > J_s > 0$, J_{ts} is always negative, and it corresponds to the staggered enfield in real, while one of the other parameters are guaranteed to be negative (for example J_{tt} is always positive). Thus for the real case of enfield analysis, we will just refer to the first two of Eq. A, and ignore all the rest three terms of Eq. A. The enfield calculation reads

$$H_{\text{MF}} = \sum_{ij} J_{ts} (\vec{\phi} \cdot \vec{\Delta}_{ij}^{ts} + H.c.) - J_{ts} |\vec{\phi}|^2$$

$$+ \sum_j \mu \left(\sum_{\alpha=1}^4 b_{j,\alpha}^\dagger b_{j,\alpha} - \kappa \right), \quad (17)$$

where $\vec{\phi}$ is a complex vector under $\text{SU}(2)^s$. We choose a minimal basis of $\vec{\phi}$ on the cubic lattice, and all the basis i, j are included in the sum with the constraint $j = i + \hat{e}$ with $\hat{e} = (1, 0), (-1/2, \pm\sqrt{3}/2)$, i.e. the enfield basis explicitly preserves the translation invariance by $2\pi/3$ every of the triangular lattice, while all the crystal symmetries are preserved as projected symmetry group ($\text{P}32_1$). μ is an on-site variational parameter of the enfield calculation which guarantees that the filling of Su₂ gauge bosons is fixed at κ on every site.

The $\text{SU}(2) \otimes 120^\circ$ ordered state corresponds to the enfield basis with $\vec{\phi} = \vec{\phi}_1 + i\vec{\phi}_2$, and the real vectors $\vec{\phi}_1$ and $\vec{\phi}_2$ orthogonal with each other. For example, where $\vec{\phi} = \phi(1, i, 0)$, only the spin-up ($\sigma^3 = +1$) Su₂ gauge bosons participate in this enfield analysis. The corresponding ordered parameter corresponds to the following gauge-invariant quantity:

$$\vec{M} \sim i\vec{\phi} \times \vec{\phi}^* \sim \vec{\phi}_1 \times \vec{\phi}_2. \quad (18)$$

In only spin-up Su₂ gauge bosons, the enfield calculation reduces precisely to the $\text{SU}(2)$ spin-1/2 Ising model on the triangular lattice⁵² with Ising coupling $J_{ij} = -4J_{ts}$ (the $\text{SU}(N)$ Ising model defined in Ref. 52 is as follows: $H = \sum_{i,j} -\frac{1}{2} J_{ij} \Delta_{ij}^\dagger \Delta_{ij}$ where Δ_{ij} is the $\text{SU}(N)$ singlet raising between Su₂ gauge bosons at sites i, j):

$$H_{\text{MF}} = \sum_{ij} J_{ts} (2\phi b_{i,\uparrow}^\dagger \tau^2 b_{j,\uparrow} + H.c.) - 2J_{ts} \phi^2$$

$$+ \sum_j \mu \left(\sum_{\alpha=1}^2 b_{j,\uparrow,\alpha}^\dagger b_{j,\uparrow,\alpha} - \kappa \right). \quad (19)$$

Because the spin-down Su₂ gauge bosons do not contribute to the enfield decomposition with $\vec{\phi} \sim (1, i, 0)$, we replace the constraint in Eq. 13 by $\sum_{\alpha=1}^2 b_{j,\uparrow,\alpha}^\dagger b_{j,\uparrow,\alpha} = \kappa$ in Eq. 19. Physically, the enfield in Eq. 19 corresponds to the “zero-field state” in Ref. 53, which has lower enfield energy than other enfield basis⁵³ for this nearest-neighbor model, and it appears in the a of the Su₂ gauge boson band structure locate at the corner of the Brillouin zone $\vec{Q} = (\pm\pi/3, 0)$. The enfield solution gives $\mu > 0$, which is consistent with the fact that the set $\sum_{\alpha=1}^2 b_{j,\uparrow,\alpha}^\dagger b_{j,\uparrow,\alpha} = 0$. And at the enfield level, the filling of the Su₂ gauge bosons κ is greater than 0.3⁵², b_α condenses, which leads to a fully polarized state in the spin space, and also 120° state in the valley space.

If the enfield value of $\vec{\phi}$ is real (or equivalently $\vec{\phi}_1$ is parallel to $\vec{\phi}_2$, for example, $\vec{\phi} \sim \phi(0, 0, 1)$), both spin-up and spin-down Su₂ gauge bosons participate in the enfield analysis, and the enfield analysis is technically equivalent to the calculation in Ref. 52 for the $\text{SU}(2) \sim \text{SO}(5)$ antiferromagnetic and the triangular lattice also with Ising coupling $J_{ij} = -4J_{ts}$, because $(\vec{\Delta}_{ij}^{ts}, \vec{\Delta}_{ij}^{st})$ together form a $\text{SO}(6)$ vector, and condense in each corner of the vector hexagon the $\text{SO}(6)$ down to $\text{SO}(5) \sim \text{SU}(2)$. Each corner of $\vec{\Delta}_{ts}$ can be regarded as the $\text{SU}(2)$ singlet introduced in Ref. 52:

$$H_{\text{MF}} = \sum_{ij} J_{ts} (\phi b_i^\dagger \sigma^{12} b_j + H.c.) - J_{ts} \phi^2$$

$$+ \sum_j \mu \left(\sum_{\alpha=1}^4 b_{j,\alpha}^\dagger b_{j,\alpha} - \kappa \right), \quad (20)$$

Quoting the results in Ref. 52, the $\text{SU}(2) \otimes 120^\circ$ state with the previous enfield basis with $\phi_1 \perp \phi_2$ as a lower enfield ground state energy density, which is consistent with our numerical observation and also our numerical simulation.

—Zeeman field driven Metal-Insulator transition

Since the insulator has a fully polarized ferromagnetic order, its energy can be tuned by an external Zeeman field. In the real magnetic field, the observed effect is the Zeeman coupling can drive first order metal-insulator transition (a level-crossing) between the polarized metal and the fully polarized ferromagnetic insulator, as was observed experimentally at the 1 Å and 3 Å filling of LiFePO_4 .^{7,9}

There is another possible evidence of metal-insulator transition driven by a Zeeman field. At the metallic side at the transition, the system is well described by a $t-J$ model with a similar J, J^s, J^v terms as Eq. 1. The Zeeman field tends to polarize the spin, which effectively increases the antiferromagnetic coupling in the valley space

$J_{eff}^v = J^v + J(\vec{\sigma}_i \cdot \vec{\sigma}_j)$. At a certain temperature, the magnitude of the 120° order in the alloy space is tunable and determined by an external Zeeman field. If the insulating behavior of the system is a consequence of the frustration in the alloy order with folds the Brillouin zone and actually gaps out the Fermi surface, as the increasing magnitude of the 120° order in the alloy space can gap out larger portion of the Fermi surface, decrease the charge carrier density, and hence eventually drive a continuous metal-insulator transition.

IV. LIQUIDS AND TOPOLOGICAL PHASES

Since $J^v \sim J^s \sim J$, it could be rather difficult for the system to form a semiclassical order due to “double frustration”: the J^s and J^v terms of Eq. 1 are both already frustrated due to the geometry of the triangular lattice, while the J term further frustrates/disorders the simultaneous 120° semiclassical order of $\vec{\sigma}_i$ and $\vec{\tau}_i$. Since there is an obvious Lieb-Schultz-Mattis theorem which forbids a completely trivial disordered phase, we expect this “double frustration” effect to lead to either a completely disordered spin-valley liquid state, or a partially ordered state with certain topological order. In this section we explore several possible spin-valley liquids or topological orders in the region $J^v \sim J^s \sim J$.

—Spin nematic Z_2 topological phase

Let us get back to the emergent Hamiltonian Eq. 17. As we discussed before, if the emergent field value of $\vec{\phi}$ is real (or equivalently if $\vec{\phi}_1$ is parallel to $\vec{\phi}_2$, for example, $\vec{\phi} \sim (0, 0, 1)$), both spin-up and spin-down Stoner gap bosons participate in the emergent analysis, and the emergent analysis is technically equivalent to the calculations in Ref. 52 for the $SO(4)$ antiferromagnet on the triangular lattice. And with large spin symmetry, the quantum fluctuations poses little difficulty for b_α to condense. If b_α is not condensed, the emergent order parameter $\vec{\phi}$ already breaks the $SO(2)^s$, and also breaks the $U(1)$ gauge symmetry down to Z_2 gauge degree of freedom.

The nature of the state with condensed $\vec{\phi}$ but uncondensed b_α depends on the nature of $\vec{\phi}$ under the exchange. The transformation of b_α under the exchange can be inferred by the fact that $\vec{\sigma} \rightarrow -\vec{\sigma}$, $(\tau^1, \tau^2, \tau^3) \rightarrow (\tau^1, \tau^2, -\tau^3)$:

$$\mathcal{T}: b_j \rightarrow i\sigma^{21} b_j, \quad \vec{\Delta}_{ij}^{ts} \rightarrow \vec{\Delta}_{ij}^{ts}, \quad (21)$$

as long as $\vec{\phi}$ is a real vector (or $\vec{\phi}_1$ parallel with $\vec{\phi}_2$), the exchange is preserved, and this state is a spin nematic Z_2 topological order. By contrast, if $\vec{\phi}_1 \perp \vec{\phi}_2$ the exchange is broken.

— $Z_2 \times Z_2$ spin-valley liquid

All the states can be constructed by introducing two flavors of Stoner gap bosons $b_{j,\alpha}^s$ and $b_{j,\alpha}^v$ for the spin and valley space on each site respectively, which are subject

to the constraint

$$\sum_{\alpha=1,2} b_{j,\alpha}^{s,\dagger} b_{j,\alpha}^s = b_{j,\alpha}^{v,\dagger} b_{j,\alpha}^v = 1. \quad (22)$$

These two constraints introduces two $U(1)$ gauge symmetries. It is fairly straightforward to construct the $SO(6) \times 120^\circ$ state using this type of Stoner gap bosons: b_α^s condenses at zero temperature, while simultaneously b_α^v condenses at the corner of the Brillouin zone.

In fact, due to the “double frustration” effect, both the spin and valley space can form a Z_2 topological order (overall speaking the system is in a $Z_2 \times Z_2$ spin-valley liquid state), whose elementary excitations are the deconfined representations of $SO(2)^s$ and $SO(2)^v$ respectively, as long as either $b_{j,\alpha}^s$ or $b_{j,\alpha}^v$ is introduced in Eq. 22 condenses in the emergent field parameters break both $U(1)$ gauge symmetries down to Z_2 .

Starting from the $Z_2 \times Z_2$ spin-valley liquid state, one can also construct a spin-valley liquid with only one Z_2 topological order. This can be for all obtained by forming bound state of the “visons” (the magnetic excitations) of both Z_2 topological orders, and condense the bound state. This condensate will condense b_α^s and b_α^v separately, but the bound state is still deconfined, and becomes the elementary of the emergent Z_2 topological order. This minimal Z_2 topological order preserves all the symmetries of the system, and it can also be constructed using the emergent field for instance as Eq. 4, as long as one condenses the spin-singlet/valley-singlet Dirac spin gap Δ_{ij}^{ss} in Eq. 4.

— $U(1) \times U(1)$ Dirac spin-valley liquid

All the exotic spin-valley liquid states can be constructed by introducing fermionic slave particles $f_{j,\alpha}^s$ and $f_{j,\alpha}^v$ which are subject to the constraints

$$\sum_{\alpha=1,2} f_{j,\alpha}^{s,\dagger} f_{j,\alpha}^s = f_{j,\alpha}^{v,\dagger} f_{j,\alpha}^v = 1. \quad (23)$$

In Ref. 4, a Dirac spin liquid with $U(1)$ gauge field and $N_f = 4$ flavors of Dirac fermions was constructed for spin-1/2 systems on the triangular lattice. And this Dirac spin liquid is the relevant state of both the 120° ordered state and the valence bond solid state^{54–56}, and it could be a deconfined quantum critical point between these two different ordered states⁵⁷.

In our case, both spin and valley space can form the Dirac liquid phases mentioned above, due to the double frustration effect. We use total there are eight flavors of Dirac fermions and two $U(1)$ gauge fields.

—The $SU(4)$ point

At the relevant $J^v = J^s = J$, this model as a $SO(6) \sim SU(4)$ symmetry. Although semiclassical approach suitable as the exact model were studied before for $SO(N)$ antiferromagnet with other representations⁵⁸, the fundamental representations on each site, this model as an obvious semiclassical limit to start with, and it is expected to be an exotic spin liquid or topological order. At this point, it is most convenient to define a four correlated Stoner gap boson $b_{j,\alpha}$ on each site, which

can simultaneously represent under both the spin and valley $SO(2)$ symmetry, and there is a constraint $\sum_{\alpha=1}^4 b_{j,\alpha}^\dagger b_{j,\alpha} = 1$.

If it is a $SO(2)$ spin system, we can require that the $SO(4)$ rotation cannot be a fully symmetric Z_2 spin liquid phase if particles are b_α slave particles. The reason is that at all the local excitations can be written as $b_{j,\alpha}^\dagger b_{j,\beta}$ with different $\alpha, \beta = 1 \cdots 4$, hence all the local excitations are invariant under the Z_4 center of the $SO(4)$ group. If a Z_2 topological order, two of the particles would be assigned to a local excitations, while two b_α slave particles cannot fuse to a representation that is invariant under the Z_4 center. This argument also requires that a Z_2 topological order phase particle is a $SO(6)$ vector is allowed.

On the other hand, using the slave particle b_α we can construct a Z_2 topological order with certain symmetric $SO(4)$ symmetry breaking. If the $SO(4)$ rotation, the model Eq. 1 can be written as

$$H = \sum_{ij} J \left(-\frac{5}{4} (\vec{\Delta}_{ij}^\dagger) \cdot (\vec{\Delta}_{ij}) + \cdots \right), \quad (2)$$

where $\vec{\Delta}_{ij}$ is a six-component vector relating between b_α . We can introduce a six-component vector $SO(6)$ vector mean field operators $\vec{\phi}$:

$$H_{MF} = \sum_{ij} J \left(-\frac{5}{4} \vec{\phi} \cdot \vec{\Delta}_{ij} + H.c. \right) + \frac{5}{4} J |\vec{\phi}|^2. \quad (25)$$

The complex vector $\vec{\phi} = \vec{\phi}_1 + i\vec{\phi}_2$, where its real and imaginary parts $\vec{\phi}_1$ and $\vec{\phi}_2$ can be either parallel or orthogonal to each other. If the $SO(6)$ gauge boson does not condense, both mean field theories could lead to a Z_2 topological order or top of the symmetric $SO(4)$ symmetry breaking.

Our DMRG simulation actually suggest that the $SO(4)$ rotation of the spin-valley model is a spin-valley liquid state with a Fermi surface of fermionic slave particles, which will be presented in detail in another work.

V. CONCLUSION

In this work we demonstrated both analytically and numerically that at a quantum spin-valley model, the all-fermion agnetic interaction can have a fully polarized ferromagnetic order in its phase diagram. We propose possible evidence for an in-plane Zeeman field to drive an insulator transition, as was observed experimentally at the 14 and 34 filling of KBG . We also discussed various possible critical spin-valley liquid state and topological order of this model.

We would like to acknowledge several previous theoretical works that studied the ferromagnetism in AB bilayer systems using different approaches and different models^{29,35,36,39}. For example, in Ref. 35, a spin-valley model with ferromagnetic couplings on a hexagonal lattice was derived for the twisted bilayer graphene system. While our work (with this manuscript) is a different bilayer system, i.e. the twisted double bilayer graphene) demonstrated that ferromagnetic order can emerge in the spin-valley model on a triangular lattice with fully antiferromagnetic interaction.

Finally, our work also suggests, under doing, again the system is likely described by the $t - J$ model with the similar J, J^s, J^v terms as Eq. 1. But the analysis in Ref. 10 still applies: the spin-triplet/valley-singlet pairing channel between electrons could become the strongest pairing channel. Due to a strong on-site Hubbard interaction, the system could still prefer to become a $d + id$ topological superconductor in the spin-triplet pairing.

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