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Superfluid Weight Bounds from Symmetry and Quantum Geometry in Flat Bands

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Flat-band superconductivity has theoretically demonstrated the importance of band topology to correlated phases. In two dimensions, the superfluid weight, which determines the critical temperature through the Berezinskii-Kosterlitz-Thouless criteria, is bounded by the Fubini-Study metric at zero temperature. We show this bound is nonzero within flat bands whose Wannier centers are obstructed from the atoms — even when they have identically zero Berry curvature. Next, we derive general lower bounds for the superfluid weight in terms of momentum space irreps in all 2D space groups, extending the reach of topological quantum chemistry to superconducting states. We find that the bounds can be naturally expressed using the formalism of real space invariants (RSIs) that highlight the separation between electronic and atomic degrees of freedom. Finally, using exact Monte Carlo simulations on a model with perfectly flat bands and strictly local obstructed Wannier functions, we find that an attractive Hubbard interaction results in superconductivity as predicted by the RSI bound beyond mean-field. Hence, obstructed bands are distinguished from trivial bands in the presence of interactions by the nonzero lower bound imposed on their superfluid weight.

Introduction. In a topological insulator, the ground state Wannier functions face an obstruction to exponential localization^{1–5}. This real-space picture connects bulk topological invariants computed from the bands in momentum space to the local chemistry of electronic states. Topological states can be either stable or fragile, and are classified by their symmetry properties in momentum space^{6–9}. Fragile states can be trivialized by mixing with non-topological bands^{10–12}, while stable states cannot. Moreover, stable topological phases are distinguished by their gapless edge states^{13,14}, while fragile phases have anomalous boundary signatures exposed by twisted boundary conditions^{15,16}, magnetic flux¹⁷, or defects¹⁸. Although our understanding of non-interacting topological bands is nearly exhaustive,^{19–22}, this is not the case for interactions within topological bands.

Superconductivity in topological bands^{23,24} is of interest since its discovery within the fragile flat bands^{25–34} of twisted bilayer graphene^{35–41}. Discovered in Ref. 42, superconducting order in flat bands, specifically the superfluid weight, originates from quantum geometry characterized by the Fubini-Study metric and has gathered much excitement^{43–49}. The quantum metric, though distinct from the band topology, is bounded by the Chern or (Euler) winding numbers^{23,50,51}. These early results suggest that topological quantum chemistry could provide general lower bounds, making contact with materials databases^{52–54}. Our work affirms this suggestion, yielding nonzero bounds in phases without winding numbers. Our bounds are given by another quantized number, the real space invariant (RSI)¹⁵, which uses symmetries to characterize topological, obstructed Wannier centers (OWCs), and trivial bands.

From a materials perspective, although topological bands are abundant within real crystals, a significant por-

tion are topologically trivial at the Fermi level. Topologically trivial bands with space group symmetries have a finer classification which divides them into trivial atomic bands, where electrons are exponentially localized at the atomic sites, and bands with OWCs which, while exponentially localized, are necessarily centered off the atoms^{55–57}. Remarkably, we show that, like topological bands, OWCs have a nonzero, lower-bounded Fubini-Study metric even without Berry curvature^{58,59}. When flat OWC bands are partially filled under attractive interactions, they possess a superconducting instability.

The zero-temperature superfluid weight $[D_s]_{ij}$ of an isolated flat band within BCS theory⁵⁹ is⁶⁰

$$[D_s]_{ij} = 2|\Delta|\sqrt{\nu(1-\nu)} \int \frac{d^2k}{(2\pi)^2} g_{ij}(\mathbf{k}) \quad (1)$$

where Δ is the superconducting gap, ν is the filling fraction of the flat bands, \mathbf{k} is a momentum in the Brillouin zone (BZ) with area $(2\pi)^2/\Omega_c$ (Ω_c is the unit cell area), and g_{ij} is the Fubini-Study quantum metric. A nonzero superfluid weight implies a finite critical superconducting temperature^{61,62} and a supercurrent $\mathbf{J} = -4D_s\mathbf{A}$, where \mathbf{A} is the vector potential in the London gauge. In a Hamiltonian with N_{orb} orbitals and N_{occ} occupied bands, we define the (abelian) quantum geometric tensor⁶³

$$\text{Tr } \mathcal{G}_{ij} = \text{Tr } P \partial_i P \partial_j P = g_{ij} + \frac{i}{2} f_{ij}, \quad (2)$$

where $P(\mathbf{k})$ is the $N_{\text{orb}} \times N_{\text{orb}}$ gauge-invariant projection matrix onto the occupied bands, ∂_i is a momentum-space derivative, and the trace is over the matrix indices. The abelian Berry curvature $f_{ij} = -f_{ji}$ is well studied while the positive semi-definite quantum metric $g_{ij} = g_{ji}$ is an object of more recent interest^{63–83}. With spatial rotation

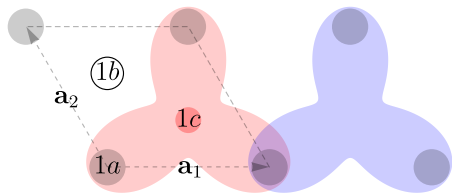


FIG. 1. Wannier basis at $1c = \frac{2}{3}\mathbf{a}_1 + \frac{1}{3}\mathbf{a}_2$. The Wannier state $|A_{1c}\rangle$ (red) centered at $1c$ is supported only on the neighboring atomic sites (grey) with A , 1E , and 2E orbitals. Only one site overlaps with neighboring Wannier states (blue).

symmetry, $[D_s]_{ij}$ is determined by the trace of g_{ij} ⁸⁴

$$G = \frac{1}{2} \int \frac{d^2k}{(2\pi)^2} \text{Tr} \nabla P \cdot \nabla P \geq 0, \quad (3)$$

which is coordinate invariant, dimensionless, and importantly is quadratic in $P(\mathbf{k})$ ⁶⁰. We give an efficient numerical discretization formula in Ref. 60.

Flat Band Model. We begin by constructing an OWC model in the space group $p3$ generated by spinless C_3 symmetry and translations along the lattice vectors $\mathbf{a}_1 = (1, 0)$, $\mathbf{a}_2 = C_3\mathbf{a}_1$. At the origin (the $1a$ position), we place electrons in the A , 1E , 2E irreps. These orbitals induce band representations^{85,86} with irreps defined by

$$\begin{aligned} A_{1a} \uparrow p3 &= \Gamma_1 + K_1 + K'_1, & C_3 &= +1 \\ {}^1E_{1a} \uparrow p3 &= \Gamma_2 + K_2 + K'_3, & C_3 &= e^{-\frac{2\pi i}{3}} \\ {}^2E_{1a} \uparrow p3 &= \Gamma_3 + K_3 + K'_2, & C_3 &= e^{\frac{2\pi i}{3}} \end{aligned} \quad (4)$$

where $\Gamma = (0, 0)^T$, $K = \frac{2\pi}{3}(\mathbf{b}_1 + \mathbf{b}_2)$, $K' = -\frac{2\pi}{3}(\mathbf{b}_1 + \mathbf{b}_2)$ are the high symmetry points, and $\mathbf{a}_i \cdot \mathbf{b}_j = \delta_{ij}$. The full group theory data can be found on the Bilbao Crystallographic server⁸⁷⁻⁸⁹. To construct a flat band OWC from these orbitals, we will use a Wannier basis centered at the $1c$ position *off* the atomic sites at $1a$ as in Fig. 1. We form the Wannier states

$$|\mathbf{R}, A_{1c}\rangle = \frac{1}{3} T_{\mathbf{R}} \sum_{j=0}^2 \tilde{C}_3^j (|0, A\rangle + |0, {}^1E\rangle + |0, {}^2E\rangle), \quad (5)$$

where $T_{\mathbf{R}}$ is the translation operator by \mathbf{R} , \tilde{C}_3 is the rotation operator about the $1c$ position, and $|0, \rho\rangle$ are the ρ orbitals in unit cell 0. Taking $\tilde{C}_3 \rightarrow e^{\mp \frac{2\pi i}{3}} \tilde{C}_3$ in Eq. (5) yields 1E and 2E states at $1c$. It is easy to check that the states $|\mathbf{R}, A_{1c}\rangle$ are orthonormal: $\langle \mathbf{R}, A_{1c} | \mathbf{R}', A_{1c} \rangle$ is nonzero only if \mathbf{R} and \mathbf{R}' are nearest neighbors, and in this case the only overlap is on a single site which vanishes due to \tilde{C}_3 eigenvalues of the orbitals. Fourier transforming Eq. (5) to obtain the eigenstate $|\mathbf{k}, A_{1c}\rangle$ yields the eigenvector $U_{\alpha}(\mathbf{k}) = \langle \mathbf{k}, \alpha | \mathbf{k}, A_{1c} \rangle$, $\alpha = A, {}^1E, {}^2E$:

$$U(\mathbf{k}) = \frac{1}{3} \begin{pmatrix} 1 \\ 1 \\ 1 \end{pmatrix} + \frac{1}{3} \begin{pmatrix} 1 \\ e^{\frac{4\pi i}{3}} \\ e^{\frac{2\pi i}{3}} \end{pmatrix} e^{i\mathbf{k} \cdot (\mathbf{a}_1 + \mathbf{a}_2)} + \frac{1}{3} \begin{pmatrix} 1 \\ e^{\frac{2\pi i}{3}} \\ e^{\frac{4\pi i}{3}} \end{pmatrix} e^{i\mathbf{k} \cdot \mathbf{a}_2}, \quad (6)$$

and the local momentum-space Hamiltonian

$$h(\mathbf{k}) = -|t|U(\mathbf{k})U^\dagger(\mathbf{k}) \equiv -|t|P(\mathbf{k}) \quad (7)$$

which has three exactly flat bands: the A_{1c} band at energy $-|t|$ and the degenerate ${}^1E_{1c}$ and ${}^2E_{1c}$ bands at zero energy. At filling $1/3$, $h(\mathbf{k})$ has the band representation

$$\Gamma_1 + K_3 + K'_3 = A_{1c} \uparrow p3, \quad (8)$$

confirming our construction in real space. We also calculate the Berry connection in crystalline coordinates,

$$A_i(\mathbf{k}) = U^\dagger(\mathbf{k}) i \partial_i U(\mathbf{k}) = (-1/3, -2/3)_i, \quad (9)$$

which is the expectation value of the lattice position operator in the occupied bands. Noting that the lattice position operator is only defined mod 1, Eq. (9) confirms that the states are located at the $1c$ position. Because $A_i(\mathbf{k})$ is independent of \mathbf{k} (up to a gauge choice), the Wilson loop bands are perfectly flat⁹⁰ and the Berry curvature is identically zero. Topologically, the model is therefore trivial. However, we calculate the quantum metric in cartesian coordinates (a is the lattice constant):

$$g_{ij}(\mathbf{k}) = \frac{1}{2} \text{Tr} \partial_i P(\mathbf{k}) \partial_j P(\mathbf{k}) = a^2 \delta_{ij} / 6, \quad (10)$$

so the mean-field superfluid weight in Eq. (1) is nonzero despite the model being topologically trivial and having compact Wannier functions (zero correlation length)⁵⁷.

It is natural to ask what indices describe these compact OWC phases. By definition, they are induced by off-site atomic orbitals, so topological quantum chemistry can identify them because their symmetry data does not match any of the orbitals present in the lattice. This is different than the stable and fragile indices which are independent of the basis orbitals.

Wilson loops can also identify OWCs. A useful reference is the SSH chain⁹¹ where an eigenvalue of π of the Wilson loop operator identifies the off-site states^{92,93}. Lastly, OWCs can be most naturally defined using the RSI formalism developed in Ref. 15. RSIs are local quantum numbers which are well-defined in fragile and OWC phases where they supply lower bounds on the number of states at the high symmetry Wyckoff positions. By definition, RSIs are invariant under symmetry-preserving adiabatic deformations. Since the Wannier states in OWCs cannot be moved to atomic sites without closing a gap, they are characterized by a nonzero RSI off an atomic site. In space group $p3$, the RSIs at a C_3 -symmetric Wyckoff positions are

$$\delta_1 = m({}^1E) - m(A), \quad \delta_2 = m({}^2E) - m(A) \quad (11)$$

and $m(\rho)$ is the number of ρ irreps. The RSIs can be conveniently calculated from the momentum space symmetry data¹⁵. In our model, the only nonzero RSIs are off the atomic sites at the $1c$ position, $(\delta_{1c,1}, \delta_{1c,2}) = (-1, -1)$. We will now show that, in generality, these off-site RSIs are responsible for a bounded superfluid weight.

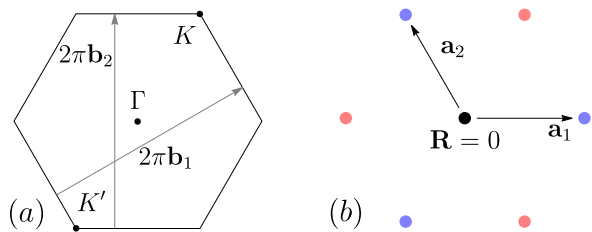


FIG. 2. C_3 Bounds. (a) We label the C_3 -invariant points Γ , K , and K' in the BZ. (b) By taking linear combinations of $P(\mathbf{k})$, we find lower bounds for the harmonics at $|\mathbf{R}| = 1$, shown in red and blue. Higher harmonics are not shown.

Lower Bounds in Real Space. We derive general lower bounds for the superfluid weight in terms of RSIs and orbital positions to show that nonzero superfluid weight is a generic feature of partially filled OWC bands, as has been shown for Chern insulators and Euler insulators^{23,50}. Our bounds also apply to all 2D topological bands⁶⁰ including fragile bands. Here, for simplicity, we prove a lower bound for our model with C_3 and all orbitals at the 1a position. Our starting point is a real space expression for $P(\mathbf{k})$, the projector onto the occupied A_{1c} band. If all orbitals are at the 1a position, then $h(\mathbf{k})$ and $P(\mathbf{k})$ are periodic under $\mathbf{k} \rightarrow \mathbf{k} + 2\pi\mathbf{b}_i$. Thus there is a Fourier representation

$$P(\mathbf{k}) = \sum_{\mathbf{R}} e^{-i\mathbf{R}\cdot\mathbf{k}} p(\mathbf{R}), \quad p(\mathbf{R}) = \int \frac{dk^1 dk^2}{(2\pi)^2} e^{i\mathbf{R}\cdot\mathbf{k}} P(\mathbf{k}), \quad (12)$$

defined in terms of the harmonics $p(\mathbf{R})$, which are $N_{\text{orb}} \times N_{\text{orb}}$ matrices, the lattice vectors \mathbf{R} , and the dimensionless crystal momenta k^i . Note that $P(\mathbf{k})$ is the momentum space Green's function⁹⁴, so $p(\mathbf{R})$ is the real space correlation function. The harmonics obey a normalization condition

$$\sum_{\mathbf{R}} \|p(\mathbf{R})\|^2 = \int \frac{dk^1 dk^2}{(2\pi)^2} \text{Tr} P(\mathbf{k}) = N_{\text{occ}} = 1, \quad (13)$$

where $\|A\|^2 = \text{Tr} A^\dagger A$ is the squared Frobenius norm⁶⁰. Rewriting Eq. (3) in real space, we find

$$G = \frac{1}{2} \int \frac{d^2k}{(2\pi)^2} \text{Tr} \nabla P \cdot \nabla P = \frac{1}{2\Omega_c} \sum_{\mathbf{R}} |\mathbf{R}|^2 \|p(\mathbf{R})\|^2. \quad (14)$$

We will use symmetry eigenvalues to show that $\|p(\mathbf{a}_1)\|$ and symmetry-related terms are bounded below. This immediately gives a bound for G because $G \geq \frac{1}{2\Omega_c} |\mathbf{a}_1|^2 \|p(\mathbf{a}_1)\|^2$ since all terms in Eq. (14) are positive semi-definite. Indeed, all are *positive* definite except for the zero mode $p(0)$, the constant mode of $P(\mathbf{k})$.

The momentum space irreps consist of the C_3 eigenvalues in the occupied bands at Γ, K, K' (see Fig. 2a). The irrep multiplicities $m(\rho)$ obey⁶⁰

$$m(\Gamma_1) + e^{\frac{2\pi i}{3}} m(\Gamma_2) + e^{-\frac{2\pi i}{3}} m(\Gamma_3) = \text{Tr} D[C_3] P(\Gamma), \quad (15)$$

and similarly for K and K' . Here $D[C_3] = \text{diag}(1, e^{-\frac{2\pi i}{3}}, e^{\frac{2\pi i}{3}})$ is the representation matrix of C_3 on the orbitals. Thus the irrep multiplicities give information about $P(\mathbf{k})$. Writing out and summing Eq. (12) at each high-symmetry momentum gives

$$P(\Gamma) + e^{\frac{2\pi i}{3}} P(K) + e^{\frac{4\pi i}{3}} P(K') = 3 \sum_{n=1}^3 p(-C_3^n \mathbf{a}_1) + \dots, \quad (16)$$

where the dots represent higher harmonics $p(\mathbf{R})$ for $|\mathbf{R}| > |\mathbf{a}_i|$ (see Fig. 2b). Crucially, the roots of unity cancel $p(0)$, so only harmonics at $\mathbf{R} \neq 0$ appear in Eq. (16). We now bound Eq. (16) on both sides. To manipulate the momentum space side, we use an elementary inequality⁹⁵

$$\|A\|^2 \geq |\text{Tr} SA|^2 / \text{Rk}(A) \quad \forall S \text{ unitary}, \quad (17)$$

proven in Ref. 60. Choosing $A = P(\Gamma) + e^{\frac{2\pi i}{3}} P(K) + e^{-\frac{2\pi i}{3}} P(K')$, we see $1/\text{Rk}(A) \geq 1/3$ because A is a 3×3 matrix, and with $S = D[C_3]$, we find with Eq. (15):

$$\begin{aligned} |\text{Tr} SA|^2 &= \frac{9}{4} (m(K_3) + m(K'_3) - m(\Gamma_2) - m(\Gamma_3))^2 + \\ &\quad \frac{3}{4} (m(\Gamma_2) - m(\Gamma_3) + m(K_1) - m(K_2) - m(K'_1) + m(K'_2))^2 \\ &= 9(\delta_{1c,1}^2 - \delta_{1c,1}\delta_{1c,2} + \delta_{1c,2}^2), \end{aligned} \quad (18)$$

where we first used Eq. (15) to write the trace in terms of momentum space irreps, and then used the tables in Ref. 15 to rewrite them in terms of the RSIs in Eq. (11).

Taking the Frobenius norm of Eq. (16) and applying the triangle inequality to the real space side gives

$$\|A\| \leq 3(\|p(\mathbf{a}_1)\| + \|p(C_3\mathbf{a}_1)\| + \|p(C_3^2\mathbf{a}_1)\| + \dots), \quad (19)$$

where the dots are higher harmonics and we used $\|p(\mathbf{R})\| = \|p(-\mathbf{R})\|$ which follows from Eq. (12). We check explicitly that Eq. (19) is not a tight inequality (by a factor of 3) and prevents our bound from being saturated by this model. All other inequalities are tight.

We next use C_3 symmetry which ensures $\|p(\mathbf{R})\| = \|p(C_3\mathbf{R})\|$ ⁶⁰. Because we have a lower bound for $\|A\|^2$, Eq. (19) proves that $\|p(\mathbf{R})\| \neq 0$ for some $\mathbf{R} \neq 0$. We now employ an optimization argument:

$$\frac{1}{2\Omega_c} \sum_{\mathbf{R}} |\mathbf{R}|^2 \|p(\mathbf{R})\|^2 \geq \min_{\psi_{\mathbf{R}}} \frac{1}{2\Omega_c} \sum_{\mathbf{R}} |\mathbf{R}|^2 |\psi_{\mathbf{R}}|^2, \quad (20)$$

where the minimization is taken over all $\psi_{\mathbf{R}} \in \mathbb{R}$ obeying

$$\sum_{\mathbf{R}} |\psi_{\mathbf{R}}|^2 = 1, \quad |\psi_{\mathbf{R}}| = |\psi_{-\mathbf{R}}|, \quad \|A\| = 9|\psi_{\mathbf{a}_1}| + \dots, \quad (21)$$

and the dots denote terms depending on $\psi_{|\mathbf{R}| > |\mathbf{a}_1|}$. As such, the space of admissible $|\psi_{\mathbf{R}}|$ described by Eq. (21) includes the choice where $|\psi_{\mathbf{R}}| = \|p(\mathbf{R})\|$. By keeping only the constraints in Eq. (21), and not the restriction that $\psi_{\mathbf{R}}$ be the Fourier transform of a projection matrix, we can perform the minimization in Eq. (20) directly.

A lemma we prove in Ref. 60 shows that the minimum occurs when $\|A\| = 9|\psi_{\mathbf{a}_1}|$, i.e., when Eq. (19) is saturated with the lowest harmonics possible. This is expected because higher harmonics have larger $|\mathbf{R}|^2$ weights. Adding up the contributions from the inner six harmonics $\|p(\mathbf{R})\|^2$ in Fig. 2, we find⁶⁰

$$\frac{1}{2\Omega_c} \sum_{\mathbf{R}} |\mathbf{R}|^2 \|p(\mathbf{R})\|^2 \geq \frac{a^2}{9\Omega_c} (\delta_{1c,1}^2 - \delta_{1c,1}\delta_{1c,2} + \delta_{1c,2}^2). \quad (22)$$

Plugging in $\delta_{1c,1} = \delta_{1c,2} = -1$ from Eq. (11), we obtain $G \geq a^2/9\Omega_c = 2/9\sqrt{3}$, a factor of 3 below the exact calculation in Eq. (10). The RSIs in Eq. (22) show that states off the atomic positions (1a in this case), which define OWCs, enforce a nonzero superfluid weight. We obtain bounds for all 2D space groups in Ref. 60.

Hubbard Model. We have shown that single-particle OWCs and fragile states have a nonzero superfluid weight at $T = 0$. However, $[D_s]_{ij}$ in Eq. (1) is obtained from mean-field BCS theory, which may seem unsuitable to treat flat band systems lacking a well-defined Fermi surface. We resort to exact numerical simulations to check its validity at finite temperature.

Using the Hamiltonian $h(\mathbf{k})$ defined in Eq. 7, we form a spinful Hamiltonian with $h_{\uparrow}(\mathbf{k}) = h(\mathbf{k})$ and $h_{\downarrow}(\mathbf{k}) = \mathcal{T}h_{\uparrow}(\mathbf{k})\mathcal{T}^{-1} = h^*(-\mathbf{k})$ which preserves time-reversal \mathcal{T} . Here \uparrow, \downarrow label the spins. Including an attractive Hubbard term with strength $|U|$, the full Hamiltonian is

$$H = -|t| \sum_{\mathbf{R}, \sigma} w_{\mathbf{R}\sigma}^{\dagger} w_{\mathbf{R}\sigma} - |U| \sum_{\mathbf{R}\alpha} c_{\mathbf{R}\alpha\uparrow}^{\dagger} c_{\mathbf{R}\alpha\downarrow}^{\dagger} c_{\mathbf{R}\alpha\downarrow} c_{\mathbf{R}\alpha\uparrow}, \quad (23)$$

where $w_{\mathbf{R}\uparrow}^{\dagger}$ creates the Wannier state in Eq. (5), $w_{\mathbf{R}\downarrow}^{\dagger} = \mathcal{T}w_{\mathbf{R}\uparrow}^{\dagger}\mathcal{T}^{-1}$, $c_{\mathbf{R}\alpha\sigma}^{\dagger}$ is the creation operator in unit cell \mathbf{R} , orbital α , and spin $\sigma = \{\uparrow, \downarrow\}$. The attractive Hubbard model does not suffer from the fermionic sign problem, and lends itself to auxiliary-field quantum Monte Carlo methods^{96,97}. We perform finite-temperature simulations in the grand canonical ensemble and tune the chemical potential $\mu(T)$ to half fill the A_{1c} band. We consider a range of Hubbard interactions $|U|$ smaller than the single-particle gap $|t|$ above the A_{1c} band: $|U| = 3, 4, 5$, with $|t| = 6$. These parameters set us away from the isolated flat band regime $|U| \ll |t|$. We focus on a system with 6×6 unit cells and periodic boundary conditions.

We can directly extract the finite-temperature superfluid weight $D_s(T)$ from the Monte Carlo results⁶⁰. The transition temperature T_c is determined by the Nelson-Kosterlitz criterion⁶²: $T_c = \pi D_s^-/2$, where D_s^- is the superfluid weight at the critical temperature approached from below. In Fig. 3, we plot $D_s(T)$ for different $|U|$ as a function of $T/|U|$, finding the curves collapse on top of each other. This confirms $T_c \propto |U|^{100}$. Our results prove that a coherent superconductor emerges upon inclusion of an attractive Hubbard interaction in the OWC flat bands, as in topological bands^{26,100,101}. Ref. 26 discusses the contrasting case of trivial atomic bands.

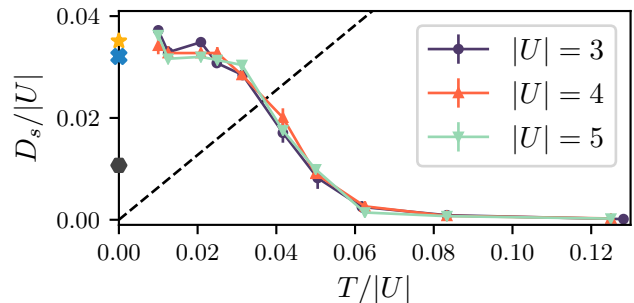


FIG. 3. Monte Carlo. The superfluid weight D_s as a function of temperature T is computed from Monte Carlo simulations on H in Eq. (23)^{98,99}. We consider $|U| = 3, 4, 5$ with $|t| = 6$ in a system with 6×6 unit cells. The crossing of D_s with the dashed line $2T/\pi$ indicates the Berezinskii-Kosterlitz-Thouless superconducting transition. The yellow star/ blue cross indicate the mean-field $D_s(T = 0)$ obtained from a multi-band/ isolated flat band calculation (Eq. (10)). The gray hexagon shows the RSI bound on $D_s(T = 0)$.

We can compare the results of our Monte Carlo simulations to the zero temperature predictions of BCS theory. In particular, we recall in Ref. 60 that the BCS wavefunction is an exact zero-temperature ground state of the attractive Hubbard model projected into the flat bands⁵¹, as follows from the equal weight of the flat band's Wannier function over all orbitals in the unit cell⁴². The blue cross in Fig. 3 shows the result of the analytical mean-field calculation after projection into the flat band. Alternatively, we solve the multi-band mean-field theory numerically in Ref. 60. The result is shown with the yellow star in Fig. 3. The agreement between our finite-temperature Monte Carlo simulations and the zero temperature mean-field calculations justify the use of the BCS result in Eq. (1), showing that our lower bounds successfully describe the many-body physics.

Discussion. We have shown that the RSIs characterizing the quantum geometry have a profound influence on the interacting groundstate when the flat bands are partially filled. Our lower bound for the superfluid weight is nontrivial in OWCs where the Wannier charge centers are obstructed from the atoms. Our bounds are not saturated by the Hamiltonian in Eq. (7), but we hope that future work can improve these bounds to be tight. Our RSI bounds also apply to OWCs with corner states, as well as stable and fragile topological phases^{102,103}. Conceptually, the gauge-invariant expression Eq. (14) in terms of the correlation function shows that long-ranged Wannier functions are *not* essential to the lower bound. Any Wannier function which is supported over multiple unit cells^{44,57}, as can be enforced by symmetry in a OWC state, produces a quantized RSI lower bound. Our derivation is general for arbitrary bands and arbitrary symmetries. Although we studied the problem in 2D, our method is generalizable to 3D where flat band OWCs have been exhaustively identified⁵⁶.

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