Detecting edge degeneracy in interacting topological insulators through entanglement entropy

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(Received 4 June 2014; revised manuscript received 24 February 2015; published 10 March 2015)

The existence of degenerate or gapless edge states is a characteristic feature of topological insulators, but is difficult to detect in the presence of interactions. We propose a new method to obtain the degeneracy of the edge states from the perspective of entanglement entropy, which is very useful to identify interacting topological states. Employing the determinant quantum Monte Carlo technique, we investigate the interaction effect on two representative models of fermionic topological insulators in one and two dimensions, respectively. In the two topologically nontrivial phases, the edge degeneracies are reduced by interactions but remain to be nontrivial.

DOI: 10.1103/PhysRevB.91.115118 PACS number(s): 71.10.Fd, 02.70.Ss, 03.65.Ud, 03.65.Vf

I. INTRODUCTION

Topologically nontrivial states of matter are a central topic of condensed-matter physics, which are classified to two categories according to their ground-state entanglement properties. The long-range entangled topological state, often named topologically ordered state, is characterized by the ground-state degeneracy on a closed manifold [1]; the short-range entangled topological insulator can be characterized by its edge degeneracy on an open boundary. (Below we use the term topological insulator to represent the general short-range entangled topological states protected by certain symmetries [2], not just for the time-reversal invariant Z2 topological insulators [3].) For a noninteracting topological insulator, edge degeneracy comes directly from the zero-energy edge mode, which is protected by its bulk topological property through the bulk-edge correspondence [4–6].

However, the single-particle picture of the edge zero modes does not apply in interacting systems. The usual bulk-edge correspondence should be understood as the relation between bulk topological property and the many-body, ground-state degeneracy on the edge [9]. The concept of edge degeneracy originates from the study of critical quantum systems [11]. Recently, it was also generalized to topologically ordered systems [12] and has been widely used for the classification of interacting topological insulators [13,14]. In this article, we apply this concept to the systems of interacting topological insulators.

Edge properties are important for the study of interaction effects in topological insulators. The first problem studied is the edge stability in the time-reversal invariant topological insulators in the presence of strong interactions [15]. Because the edge is gapless, interaction effects on the edge are more prominent than those in the bulk, which can lead to edge instabilities while maintaining the time-reversal invariance in the bulk. The above picture has also been confirmed in quantum Monte Carlo (QMC) simulations [16,17]. In recent years, interacting topological insulators have been intensively studied [15–19], which have also been classified according to different symmetries [13,14,20]. For the time-reversal invariant topological insulators, the Z2 index can be formulated in terms of the single-particle Green’s functions [21], which have been calculated based on both analytic and numeric methods [22–24].

On the other hand, quantum entanglement provides a particular perspective to investigate quantum many-body physics [25]. Entanglement entropy measures nonlocal correlations between part A and the rest of the system, denoted as part B. Entanglement entropy can be defined as the von Neumann entropy,

\[ S_e = -\text{Tr}[\rho_A \ln \rho_A], \]  

based on the reduced density matrix \( \rho_A = \text{Tr}_B(\rho_{A:B}) \). For systems characterized by short-range entanglement, entanglement entropy obeys an area law: It is proportional to the area/length of the boundary, i.e., the entanglement cut. However, in the quantum critical region, entanglement entropy shows a logarithmic dependence on the subsystem size due to the divergence of coherence length [26]. In topologically ordered systems with long-range entanglements, a negative subleading term appears, termed as topological entanglement entropy [27], which depends on the degeneracy of ground states. Topological entanglement entropy has been used to identify different topological orders in quantum spin-liquid systems [28,29]. For short-range entangled topological insulators, the single-particle entanglement spectrum [30] is found to exhibit a “zero mode”-like behavior [31,32] in the noninteracting case. However, interaction invalidates such a single-particle picture, and thus a more delicate method is required to describe interacting topological insulators.

In this article, we propose a method to determine the edge degeneracy using entanglement entropy. A quantity termed edge entanglement entropy \( S_{n,\text{edge}} \) [defined in Eq. (5)] is employed to measure edge degeneracy for both noninteracting and interacting topological insulators. This work is motivated by a recently developed algorithm using the fermionic determinant QMC to calculate the Renyi entanglement entropy [33],

\[ S_n = -\frac{1}{n-1} \ln \text{Tr}[\rho_A^n]. \]  

We employ this algorithm to study edge degeneracy of fermionic interacting topological insulators by measuring \( S_{2,\text{edge}} \) in both one- and two-dimensional systems. Our methodology is first explained by using the Su-Schrieffer-Heeger-Hubbard (SSHH) model [34]. In the topologically nontrivial phase, the Hubbard \( U \) reduces \( S_{2,\text{edge}} \) from 2 ln 2...
to \(\ln 2\), corresponding to reducing edge degeneracy from 4 to 2 in the thermodynamic limit. In two dimensions (2D), \(S_{2,\text{edge}}\) also contributes a subleading term to the entanglement entropy area law, as we observed in the Kane-Mele-Hubbard (KMH) model \cite{35} for a cylindrical geometry. Moreover, \(S_{2,\text{edge}}\) shows even-odd dependence on the system size along the entanglement cut, in agreement with the helical liquid behavior on the edge.

II. THE SSHH MODEL

The 1D SSHH model is defined as

\[
H_{\text{SSHH}} = -\sum_{i=1}^{2L} [t + \delta t (-1)^i] c_{i\sigma}^\dagger c_{i+1,\sigma} + \text{H.c.} \\
+ \sum_{i=1}^{2L} \frac{U}{2} (n_i - 1)^2,
\]

where \(\sigma = \uparrow, \downarrow\), \(n_i = \sum_\sigma c_{i\sigma}^\dagger c_{i\sigma}\), \(\delta t\) controls the hopping dimerization strength, \(t\) is set 1 below, and \(U\) is the Hubbard interaction.

At \(U = 0\), this model is well known, exhibiting two topologically distinct ground states, characterized by Berry phase 0 (\(\delta t > 0\)) and \(\pi\) (\(\delta t < 0\)), respectively. We use the convention that two sites \(2i - 1\) (odd) and \(2i\) (even) are combined into one unit cell. The \(\pi\)-valued Berry phase guarantees the existence of one zero-energy mode for each spin on each end \cite{31} (inset of Fig. 1(a)) at \(\delta t < 0\). In order to study the entanglement entropy, the chain is divided into two parts, A and B: Using the truncated correlation matrix \(\tilde{C}_{ij,\sigma} = (c_{i\sigma}^\dagger c_{j\sigma})\) with \(i, j \in A\), the Renyi entanglement entropy can be obtained \cite{36} as

\[
S_n = -\frac{1}{n-1} \sum \ln f_{i\sigma}^n,
\]

where \(f_{i\sigma}\) are eigenvalues of \(C_{ij,\sigma}\). \(S_2\) is calculated under both the periodic boundary condition (PBC) and the open boundary condition (OBC), as plotted in Fig. 1(a). In practice, for \(\delta t > 0\), we choose \([1; L/2]\) as subsystem A and \([L/2 + 1; L]\) as sublattice B for both OBC and PBC. In this case, all the cuts are at weak bonds. On the other hand, for \(\delta t < 0\), we use a different partition method such that the cuts are still at weak bonds. For the case of OBC, we choose \([1; L/2 + 1]\) and \([L/2 + 2; L]\) as subsystems A and B, respectively, while for the case of PBC, we choose \([2; L/2 + 1]\) and \([L/2 + 2; 1]\) as subsystems A and B, respectively, such that again all the cuts are at weak bonds. Figure 1(a) shows that neither PBC nor OBC gives quantized entanglement entropy because of the short-range entanglement near the cut to divide the system \cite{31}. Of course, we can also choose the cuts on strong bonds and define the edge entanglement.

To extract the entanglement between two edges, we define the edge entanglement entropy as

\[
S_{n,\text{edge}} = S_{n,\text{obc}} - \frac{S_{n,\text{pbc}}}{2},
\]

where half of \(S_{n,\text{pbc}}\) is subtracted because there are two cuts for defining entanglement in the case of PBC but only one in the case of OBC. This definition also applies for the interacting case. Equation (5) measures the nonlocal entanglement between the edges. Although \(n\) can be any integer number, we only consider the case of \(n = 2\) below because of the numerical convenience by QMC. Certainly, we can also choose cuts on strong bonds and define \(S_{n,\text{edge}}\) by subtraction accordingly; the results of the quantization of \(S_{n,\text{edge}}\) remain robust.

The edge entanglement entropy exhibits a quantized behavior in two gapped phases. At \(U = 0\), Fig. 1(a) shows \(S_{2,\text{edge}} = 2 \ln 2 = \ln 4\) at \(\delta t < 0\), while \(S_{2,\text{edge}} = 0\) at \(\delta t > 0\). This result...
can be understood as follows. For each spin component $\sigma$, two zero modes $\gamma^i_{L,R\sigma}$ at two ends are coupled through an effective hopping $t_{\text{eff}} \sim \exp(-\delta t L)$, and then the bonding state, $\frac{1}{\sqrt{2}}(\gamma^i_{L\sigma} + \gamma^i_{R\sigma})|0\rangle$, contributes a $\ln 2$ to $S_{n,\text{edge}}$ in each spin component. More explicitly, these bonding edge states are occupied by both spin components; i.e., the two-particle edge states can be written as

$$\frac{1}{\sqrt{2}}(\gamma^i_{L\uparrow} \gamma^i_{R\downarrow} + \gamma^i_{L\downarrow} \gamma^i_{R\uparrow} + \gamma^{i*}_{L\uparrow} \gamma^{i*}_{R\downarrow} + \gamma^{i*}_{L\downarrow} \gamma^{i*}_{R\uparrow})|0\rangle,$$

which clearly exhibit the $\ln 4$ contribution to the edge entanglement entropy.

Now let us consider taking the limit of $L \to +\infty$ in which $t_{\text{eff}}$ approaches to zero and the edge modes become exactly zero modes. Then each term in Eq. (6) corresponds to a zero mode for either edge. Say, after tracing out the degree of freedom on the right edge, we arrive at the zero modes at the left edge as $\gamma^{i}_{L\downarrow} \gamma^{i}_{R\downarrow}|0\rangle, |0\rangle, \gamma^{i}_{L\uparrow} \gamma^{i}_{R\uparrow}|0\rangle, |0\rangle$. Thus, the above-defined $S_{2,\text{edge}}$ can be used as a topological index, which corresponds to the thermodynamic entropy at zero temperature of one edge. This explains the relation between entanglement entropy and the ground-state degeneracy $D$ on one open end as

$$\ln D = \lim_{L \to \infty} S_{n,\text{edge}}(L),$$

which converges to the same value independent of $n$. It is sufficient to calculate $S_2$ to determine edge degeneracy. A similar quantity to Eq. (5) was used as a topological invariant to study a 1D noninteracting $\pi$-wave superconductor [37]. The physical meaning of this topological invariant becomes clear in our approach: It represents entanglement between two edges for a finite lattice size and converges to the thermodynamic limit very quickly. Note that at the critical point $\delta t = 0$, $S_{2,\text{edge}}$ is negative and unquantized, in agreement with the “noninteger” edge degeneracy in critical quantum systems [11].

Now let us turn on the Hubbard interaction $U$. We combine the zero temperature projector QMC [38] with the new developed algorithm to calculate $S_2$ [33] under PBC and OBC, respectively. Due to the particle-hole symmetry, the half-filled SSHH model is free of the sign problem, and thus the QMC simulation can be performed in a controllable way. The results of $S_2$ vs $\delta t$ at $L = 40$ and $U = 1$ are calculated and plotted in Fig. 1(b). The behavior of $S_{2,\text{edge}}$ is similar to the case of $U = 0$ in Fig. 1(a), except that its quantized value becomes $\ln 2$ when $\delta t < 0$. At large values of $L, U \gg t_{\text{eff}}$, and thus the singlet ground state changes to

$$\frac{1}{\sqrt{2}}(\gamma^i_{L\downarrow} \gamma^i_{R\downarrow} - \gamma^i_{L\uparrow} \gamma^i_{R\uparrow})|0\rangle,$$

leads to $S_{2,\text{edge}} = \ln 2$. Again in the limit of $L \to +\infty$, the edge modes become exactly zero modes. If we trace out the right edge, the zero modes left at the left edge are $\gamma^i_{L\sigma}|0\rangle$ and $\gamma^{i*}_{L\sigma}|0\rangle$, which means that edge degeneracy $D$ is reduced from 4 to 2 by the Hubbard $U$; i.e., the double and empty occupations of the edge states are projected out. Due to the exponential decay of $t_{\text{eff}}$, the finite-size effect of $S_{2,\text{edge}}$ is weak. It converges to $\ln D$ quickly even before $L$ goes large.

We have also calculated $S_{2,\text{pbc}}, S_{2,\text{OBC}}$, and $S_{2,\text{edge}}$ at different values of $U$ ranging from 0.5 to 2 as shown in Fig. 1(c). $S_{2,\text{pbc}}$ and $S_{2,\text{OBC}}$ are nonquantized, which decreases as increasing $U$ due to the suppression of charge fluctuations across the cuts. Nevertheless, $S_{2,\text{edge}}$ is pinned at $\ln 2$ regardless of different values of $U$ due to the exponential suppression of $t_{\text{eff}}$. In Fig. 2, we set up the phase diagram of the SSHH model using $S_{n,\text{edge}}$ and edge degeneracy. A similar phase diagram has been obtained by calculating the bulk topological number $Z = \langle 0 | U \sigma_\uparrow \sigma_\downarrow | 0 \rangle$ for $\delta t > 0 (< 0)$ using Green’s functions extracted from the density matrix renormalization group [22]. Our study here further indicates the edge behavior: In the topologically nontrivial region, edge degeneracy is reduced from 4 to 2 by the Hubbard interaction [14] at half filling.

When $U$ is large, the low-energy physics of the SSHH model is described by the spin-$\frac{1}{2}$ Heisenberg-Peierls model $H = \sum_{i,j} J_i \delta_i \delta_j \delta_{i+1}$, where $J_1 = J$ or $J'$ for the odd or even bond, respectively [39]. Our study shows that the cases of $J' < J$ and $J' > J$ belong to topologically distinct phases. At $J' > J$, there is a free local moment at one end, resulting in a double-edge degeneracy. The transition occurs at $J = J'$ consistent with the critical behavior of the spin-$\frac{1}{2}$ Heisenberg model [40,41].

For the above results, the particle-hole symmetry gives rise to zero-energy edges in noninteracting cases. The finite size effect couples the two edge states together and contributes to $S_{n,\text{edge}}$. Even in the interacting case, our numeric simulations show that $S_{n,\text{edge}}$ remains robust. When the particle-hole symmetry is gone, for the 1D case, both edge states are not at zero energy. They are either both occupied or empty, and thus $S_{n,\text{edge}}$ will be reduced to zero. Nevertheless, our method still applies to the 2D KMH model because the chemical potential crosses the 1D band of edge states. The single-particle states right at the chemical potential play the role of zero-energy states.

### III. THE KMH MODEL

Next we move to 2D and investigate the KMH model on a honeycomb lattice defined as

$$H_{\text{KMH}} = -\sum_{\langle i,j \rangle,\sigma} t_{\langle i,j \rangle,\sigma} c^\dagger_{i,\sigma} c_{j,\sigma} + \sum_{\langle i,j \rangle,\sigma} i\lambda c^\dagger_{i,\sigma} \sigma c_{j,\sigma}$$

$$+ \sum_i \frac{U}{2} (n_i - 1)^2,$$

where $c^\dagger_{i,\sigma}$ and $c_{i,\sigma}$ is the creation and annihilation operator at site $i$ and spin $\sigma$, respectively. $\lambda$ is the hopping parameter in the Kondo model, and $U$ is the Hubbard interaction.
where $\lambda$ is the next-nearest-neighbor, spin-orbit coupling, $\sigma = \uparrow, \downarrow$, and again $\tau$ is set 1. This model is free of the sign problem and has been investigated by the determinant QMC [16,17]. Along the y direction (zigzag), the PBC is applied, and along the x direction (armchair), both of the PBC and the OBC are applied. The PBC and OBC correspond to the toric and cylindrical geometries, respectively. The lattice is divided into the subsystem $A$ with $1 \leq x \leq L/2$ and the environment $B$ with $L/2 + 1 \leq x \leq L$ for the study of entanglement entropy.

The QMC results of $S_{2,\text{pbc}}$ and $S_{2,\text{obc}}$ for the KMH model are shown in Fig. 3. $S_{2,\text{pbc}}$ exhibits a standard area law, i.e., $S_{2,\text{pbc}} \propto L_y$, while $S_{2,\text{obc}}$ shows an even-odd oscillating behavior. Then $S_{2,\text{edge}} = \ln 2$ and 0 for even and odd values of $L_y$, respectively, as shown in the inset. On the other hand, $S_{2,\text{edge}}$ can also be obtained by extrapolating $S_{2,\text{obc}}(L_y)$ for even values of $L_y$, in which $S_{2,\text{edge}}$ appears as the subleading term of the area law as

$$S_{2,\text{obc}}(L_y) \approx \alpha L_y + S_{2,\text{edge}}. \quad (10)$$

Such a subleading term is an analogy to the topological entanglement entropy in the long-range entangled topological orders [27–29]. We propose to use $S_{2,\text{edge}}$ to characterize the short-range entangled topological insulators in 2D. Both topological entanglement entropy and $S_{2,\text{edge}}$ are related to the ground-state degeneracy, but account for bulk and edge states, respectively.

Next we explain the origin of the nonzero $S_{2,\text{edge}}$ by analyzing the edge degeneracy. At $U = 0$, such a behavior is a direct consequence of the zero-energy edge states, which has also been found in the Kitaev model [42] and noninteracting triplet topological superconductors [43]. In Fig. 4(a), the energy spectrum with the open edges is plotted as a function of $k_y$, which is conserved due to the PBC along the y direction. The edge zero mode is located at $k_y = \pi$, which is accessible for even values of $L_y$, but not for odd $L_y$; thus, the many-body ground-state degeneracy varies between 4 and 1.

At $U > 0$, the above single-particle picture does not hold any more. Interaction effects have to be fully taken into account to investigate the many-body edge degeneracy. We use an effective edge helical liquid defined in momentum space [15],

$$H_{hl} = \sum_{k,\sigma} \sigma v_F (k - \pi) c_{k,\sigma}^\dagger c_{k,\sigma} + \sum_{k'kq} U c_{k+q,\uparrow}^\dagger c_{k,\downarrow} c_{k'-q,\downarrow}^\dagger c_{k',\uparrow}. \quad (11)$$

where $c_{k,\sigma}$ are creation operators of noninteracting edge states. The $y$-direction momenta $k_y = \frac{2\pi}{L_y} n (n = 0,1, \ldots, L_y - 1)$ are chosen only for edge states and satisfy $|v_F (k - \pi)| \lesssim \Lambda$, where $\Lambda$ is the energy cutoff. Since edge modes with different values of $k$ have different localization lengths, rigorously speaking, the interaction matrix elements for the edge modes are $k$ dependent even for the case of the Hubbard model. Nevertheless, for simplicity, we neglect this dependence. In real calculations, we choose $\Lambda = \pi v_F / 2 = 1$ without loss of generality. The number of momentum points for edge states within the cutoff is denoted as $N_k$.

Exact diagonalization method is employed to numerically solve the many-body edge energy levels at $L_y \leq 15$ ($N_k \leq 8$), and the energy gaps are plotted in Fig. 4(b). For even $L_y$'s, $\Delta = 0$ corresponds to a double degeneracy in agreement with the QMC result $S_{2,\text{edge}} = \ln 2$. For odd $L_y$'s, the ground state behavior is ...
has no degeneracy; thus, $S_{2,\text{edge}} = 0$. Nevertheless, the gap decreases to zero as increasing $L_y$, as shown in the inset. This gapless behavior in the thermodynamic limit has been obtained from the bosonization analysis [15], which shows that forward scattering does not open a gap in a helical liquid in the weak interacting regime.

IV. SUMMARY

We propose a quantized quantity of the edge entanglement entropy $S_{\text{edge}}$ to determine edge degeneracy in topological insulators in the presence of interactions. Using the fermionic QMC algorithm, $S_{n,\text{edge}}$ is calculated for both the interacting 1D SSHH model and 2D KMH model. In topologically nontrivial phases of these models, the Hubbard $U$ suppresses the quantized values of $S_{2,\text{edge}}$ from 2 in the noninteracting cases to its half value ln 2. In 2D, such a nonzero $S_{\text{edge}}$ also contributes a subleading term in the entanglement entropy area law for a cylindrical geometry.

Before closing this paper, some remarks are in order. (I) Our QMC calculations are only performed at small and medium $U$. When $U$ goes large, the QMC numeric error of entanglement entropy increases significantly [33]. Significant numeric efforts are needed to obtain reliable entanglement entropy. Very recently, a new QMC algorithm using the replica technique was proposed for fermionic systems [44], which is more stable in the large $U$ regime and can be helpful to study the Mott transition regime in the future. (II) If the third-nearest-neighbor hopping is added to the KMH model, two Dirac nodes are produced at $k_y = 0$ and $\pi$, respectively, on an edge [24]. In this case, any small $U$ will gap out the edge states due to the Umklapp scattering [15]. Therefore, $S_{2,\text{edge}} = 0$ is expected to be consistent with the physical implication of the $Z_2$ topological insulator. (III) We have seen that the above edge state entanglement is built up through the effective coupling $\sim \exp(-L_y/\xi)$ between two edges, in which $L_y$ is the width of the system and $\xi$ is the typical localization length of the edge modes. On the other hand, for the 2D case, the length $L_y$ along the edge direction also gives another energy scale for the low-energy edge excitations, which is $\sim 1/L_y$. When $\exp(-L_y/\xi) \ll 1/L_y$, the edge modes not right located at the Fermi energy, we can neglect their contributions to the total entanglement entropy (EE) and only need to consider the zero-energy mode right at the Fermi energy. In this regime, our method applies.

ACKNOWLEDGMENTS

D.W. thanks Zhusen Huang and Xiao Chen for helpful discussions. This work is supported by NSF Grant No. DMR-1410375 and AFOSR Grant No. FA9550-14-1-0168. Y.W. and C.W. acknowledge the financial support from the National Natural Science Foundation of China under Grant No. 11328403 and the Fundamental Research Funds for the Central Universities. C.W. also acknowledges the support from the Presidents Research Catalyst Awards of University of California. Part of the computational resources required for this work were accessed via the GlideinWMS [45] on the Open Science Grid [46].

[6] In some special cases, there is no zero mode on an open boundary even in a topologically nontrivial state, e.g., Refs. [7, 8]. Then the bulk-edge correspondence should be generalized by introducing twisted boundary condition [5].
[9] There is a subtlety when we say the degeneracy of a “gapless” system. We emphasize that it depends on the sequence of two limits: zero temperature and infinite lattice size [10]. In this article, we take zero temperature limit first.


