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Measurement of the Entanglement Spectrum of a Symmetry-Protected Topological State Using the IBM Quantum Computer

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Entanglement properties are routinely used to characterize phases of quantum matter in theoretical computations. For example, the spectrum of the reduced density matrix, or so-called “entanglement spectrum”, has become a widely used diagnostic for universal topological properties of quantum phases. However, while being convenient to calculate theoretically, it is notoriously hard to measure in experiments. Here, we use the IBM quantum computer to make the first ever measurement of the entanglement spectrum of a symmetry-protected topological state. We are able to distinguish its entanglement spectrum from those we measure for trivial and long-range ordered states.

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Introduction.—The patterns of entanglement between local degrees of freedom (d.o.f.) are a fingerprint of many-body quantum phases [1–3]. Various measures for quantum entanglement capture different universal aspects of quantum states and allow for their classification. Of these measures, the entanglement spectrum, obtained from the spectrum of the reduced density matrix of a bipartitioned quantum system, reveals the most detailed information. It was first used in this way by Li and Haldane [4], who characterized the topological order of a fractional quantum Hall system by matching its low-lying entanglement level counting to the universal level structures of conformal field theories.

Entanglement spectroscopy further became an important tool in identifying symmetry-protected topological states of matter (SPTs) [1,2,5–8]. The defining property of a SPT is that it cannot be connected to a trivial product state via a finite-depth quantum circuit of local symmetry preserving unitary operations. SPTs are not topologically ordered in the sense of the fractional quantum Hall effect, but still feature a topological bulk-boundary correspondence: (in one dimension) they always support gapless boundary excitations in an open geometry, as long as the boundary does not break the protecting symmetry. This bulk-boundary correspondence is also manifested in the entanglement spectrum: the bipartitioning used to define the reduced density matrix introduces a “virtual” boundary between parts A and B of the system, and in a SPT topological excitations are associated with this virtual boundary leading to protected degeneracies in the entanglement spectrum [9].

While entanglement spectra conveniently characterize many-body quantum states in numerical simulations, they are intrinsically hard to measure in condensed matter experiments. It is only recently with the improvements in trapped ions and superconducting quantum simulators that the entanglement entropy [10] or density matrix [11] of many-spin systems have been accessed [12,13]. The recent release of the IBM cloud quantum computing service allows one to test various conceptual ideas on an actually existing quantum simulator [14]. Several recent works characterize the entanglement properties of the IBM devices [15] and use them to test various quantum algorithms [16–22] and to solve physical demonstration problems [23,24]. Here, we show that the high level of control that can be reached on the IBM digital quantum computer allows one to both prepare a SPT state and to measure its entanglement spectrum.

Topology in quantum paramagnets.—SPT phases are gapped short-ranged entangled states with a protecting symmetry. In this Letter, we are interested in SPT phases of one-dimensional paramagnets arising in a chain of spin-1/2 d.o.f. with local quantum states $|\uparrow\rangle$ and $|\downarrow\rangle$ at each site (corresponding to a chain of qubits on the IBM quantum computer). According to the classification of SPT phases [1], such one-dimensional bosonic systems support a topologically nontrivial phase protected by time-reversal symmetry $T$ with $T^2 = +1$. We choose the representation $T = K \prod_i \sigma_{\mu,i}$, where $K$ is complex conjugation and $\sigma_{\mu,i}$ is the Pauli operator with $\mu = 0, x, y, z$ acting on site $i$ of the chain ($\sigma_0$ being the identity operator).

A trivial paramagnetic state invariant under $T$ is given by the ground state of

$$H_{\text{triv}} = -\sum_i \sigma_{x,i}, \quad (1)$$

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which, for the example of a chain of length $N$, is a trivial product state,
\[ |PM\rangle = \frac{1}{\sqrt{2^N}} \sum_r |r\rangle = |+\rangle^N, \]  
(2)

where $r$ represents all possible binary strings of length $N$ and $|+\rangle$ is the eigenstate of $\sigma_x$ with eigenvalue $+1$.

On the other hand, a topologically nontrivial SPT state protected by $T$ is given by the ground state of the stabilizer Hamiltonian [25],
\[ H_{\text{topo}} = -\sum_i \sigma_{z,i-1} \sigma_{x,i} \sigma_{z,i+1}, \]  
(3)

with periodic boundary conditions imposed. Since all terms in $H_{\text{topo}}$ commute, its unique ground state $|\text{SPT}\rangle$ is defined by
\[ \sigma_{z,i-1} \sigma_{x,i} \sigma_{z,i+1} |\text{SPT}\rangle = |\text{SPT}\rangle, \quad \forall \ i. \]  
(4)

This state is the so-called graph state, which can be used as a key resource to achieve one-way quantum computation [26,27]. It is also topologically equivalent to the well-known Affleck-Kennedy-Lieb-Tasaki (AKLT) state [28,29].

We now discuss how to distinguish the trivial Eq. (2) and nontrivial SPT ground states Eq. (4). The classification of one-dimensional SPT phases in Ref. [1] is based on the fact that in one dimension any paramagnetic or short-range correlated state can be transformed into a trivial product state via a local unitary (LU) transformation. However, if the paramagnetic state is a nontrivial SPT, some of the local operations comprising such a LU transformation necessarily break the protecting symmetry. LU transformations are well approximated by finite-depth quantum circuits [30]. To define a quantum circuit, we first introduce the piecewise local unitary operators,
\[ \hat{U}_P = \prod_{i \in \mathcal{P}} U_i, \]  
(5)

where $U_i$ is a set of unitary operations acting on non-overlapping local regions listed in the set $\mathcal{P}$. For the IBM quantum computer the only available multiqubit gate is a CNOT gate; that is, $U_i$ acts on sites $i$ and $i+1$. The full LU transformation is then given by a finite product of $M$ piecewise unitaries of the form of Eq. (5):
\[ U_{\text{LU}}^M = \hat{U}_{\mathcal{P}_1} \cdots \hat{U}_{\mathcal{P}_M}. \]  
(6)

A LU transformation that transforms the SPT state (SPT) [Eq. (4)] into the trivial product state $|PM\rangle$ [Eq. (2)] is obtained by choosing
\[ U_i = \frac{1}{2} \sigma_{0,i}(\sigma_{0,i+1} + \sigma_{z,i+1}) + \frac{1}{2} \sigma_{z,i}(\sigma_{0,i+1} - \sigma_{z,i+1}), \]  
(7)

which represents a controlled-Z gate between spin $i$ and $i+1$. Further using $M = 2$ with $\mathcal{P}_1$ and $\mathcal{P}_2$ containing all even and all odd sites, respectively, one obtains
\[ |PM\rangle = \prod_i U_i |\text{SPT}\rangle \]  
(8)
as the LU transformation. Since $U_i^2 = 1$, the same LU transformation can be used to construct the state (SPT) from (PM). The quantum circuit corresponding to $\prod_i U_i$ is shown in Fig. 1(d). It is straightforward to see that none of the $U_i$ commute with $T$, as expected for a nontrivial SPT state (SPT) and a trivial product state (PM).

Another characterization of the SPT state is that with open boundary conditions, Hamiltonian Eq. (3) supports gapless excitations localized at the chain end. To see this, notice that the operators $\sigma_{x,1} \sigma_{z,2}$, $\sigma_{y,1} \sigma_{z,2}$, and $\sigma_{z,1}$ all commute with the Hamiltonian and form a Pauli algebra. They then enforce a twofold degeneracy of the ground state associated with a localized excitation at the boundary. Since

\[ \begin{array}{cccc}
(a) & (b) & (c) & (d) \\
Q0 & Q1 & Q2 & Q7 \\
Q15 & Q14 & Q13 & Q12 \\
Q11 & Q10 & Q9 & Q8 \\
Q1 & Q0 & Q1 & Q0 \\
\end{array} \quad \begin{array}{cccc}
|q1\rangle = |0\rangle & - & |q1\rangle = |0\rangle & - \\
|q2\rangle = |0\rangle & - & |q2\rangle = |0\rangle & - \\
|q3\rangle = |0\rangle & - & |q3\rangle = |0\rangle & - \\
|q4\rangle = |0\rangle & - & |q4\rangle = |0\rangle & - \\
\end{array} \quad \begin{array}{cccc}
(d) & (e) \\
\text{CNOT} & \text{Hadamard} \\
\end{array} \]

FIG. 1. (a) Layout of the qubits and two-qubit gates (arrows) in the IBM quantum computer ibmqx5. Blue and red qubits were used to measure entanglement spectra and formed subsystem $A$ and $B$ of the simulated quantum spin chain (with periodic boundary conditions), respectively. (b)–(d) Circuits used for constructing the quantum states, where $H$ stands for a Hadamard and + for a CNOT gate, respectively. (b) Constructs the ground state of the trivial paramagnet. (c) Constructs the eight qubit cat state (also “GHZ state”). (d) Constructs the ground state of a topological paramagnet (also “graph state”). Hadamard gates that appear next to each other are automatically removed. (e) Symbols for the various logic gates.
all these operators are odd under $\mathcal{T}$, they cannot be added as perturbations to the Hamiltonian and we conclude that the gapless topological end excitation is protected by $\mathcal{T}$.

Entanglement spectrum.—For any quantum state $|\psi\rangle$, the reduced density matrix of subsystem $A$ (some subset of the lattice) is given by the partial trace of the full density matrix $\rho = |\psi\rangle\langle\psi|$ over the complementary subsystem $B = A^c$, i.e.,

$$\rho_A = \text{Tr}_B \rho.$$

(9)

The entanglement spectrum is then the set of numbers $\lambda_i$, $i = 1, \ldots, N_A$, where $e^{-\lambda_i}$ are the eigenvalues of the operator $\rho_A$ and $N_A$ is the dimension of the Hilbert space describing all d.o.f. in subsystem $A$. The entanglement spectrum can reveal information about universal, in particular, topological, properties of the state $|\psi\rangle$. For example, if the bipartitioning is realized by a cut separating a linear chain into a left and a right half, it has been argued [6] that nontrivial SPT phases have entanglement spectra whose low energy spectra are in one-to-one correspondence with the spectrum of the system in the presence of a boundary. This statement can be proved straightforwardly for noninteracting free-fermion systems [31]. For the case at hand, we study a chain with periodic boundary conditions, so that a an entanglement cut introduces two virtual boundaries between parts $A$ and $B$. If each cut comes with a twofold degeneracy, we expect a $2 \times 2$ or fourfold degenerate entanglement spectrum of the SPT. In contrast, no degeneracies are expected in the low-lying entanglement spectrum of the trivial PM phase, as it has no protected boundary modes. The degeneracy is thus a useful diagnostic for identifying topological phases.

We obtain the entanglement spectrum by measuring the reduced density matrix of the subsystem. Given any reduced density matrix $\rho_A$ we can decompose it into a sum of Pauli matrices. For a $n$-spin subsystem at sites $1, \ldots, n$, we have

$$\rho_A = \sum_{a_1, \ldots, a_n} \frac{1}{2^n} \sigma_{a_1,1} \sigma_{a_2,2} \cdots \sigma_{a_n,n},$$

(10)

where $1/2^n$ is a normalization factor and the indices $a_i$, $i = 1, \ldots, n$, run over the set $\{0, x, y, z\}$. The aim then is to obtain the coefficients $c_{a_1, \ldots, a_n}$ given that we have access to the full density matrix. Since the Pauli matrices are orthogonal under the trace norm, the coefficients are given by

$$c_{a_1, \ldots, a_n} = \text{Tr}_A[\sigma_{a_1,1}\sigma_{a_2,2}\cdots\sigma_{a_n,n}\rho_A] = \text{Tr}_A[\sigma_{a_1,1}\sigma_{a_2,2}\cdots\sigma_{a_n,n}\text{Tr}_B(\rho)] = \text{Tr}_A\text{Tr}_B[\sigma_{a_1,1}\sigma_{a_2,2}\cdots\sigma_{a_n,n}\rho] = \text{Tr}[\sigma_{a_1,1}\sigma_{a_2,2}\cdots\sigma_{a_n,n}\rho].$$

(11)

This means that in order to obtain the coefficients $c_{a_1, \ldots, a_n}$, we need only measure all the Pauli operators corresponding to subsystem $A$. For 4 spins, and ignoring the identity operator, this gives $4^4 - 1 = 255$ measurements.

After obtaining the coefficients, we reconstruct the reduced density matrix according to Eq. (10) and we numerically diagonalize $\rho_A$ to obtain the entanglement spectrum.

Results.—We consider the different quantum states on a system of 8 spins (qubits): the trivial paramagnet (PM) from Eq. (2), the cat state,

$$|\text{cat}\rangle = \frac{1}{\sqrt{2}} (|00000000\rangle + |11111111\rangle),$$

(12)

and the SPT state (graph state) as defined in Eq. (4). We construct these states on the IBM 16 qubit quantum computer ibmqx5 using the quantum circuits shown in Fig. 1.

To obtain the reduced density matrix, we have to obtain the expectation values of the Pauli operators according to Eq. (10). Since each measurement gives only a 0 or 1, the experiment consisting of state preparation and measurement is repeated multiple times in order to estimate the expectation value. The results we show are taken with 1024 repetitions all executed within about 15 min on the quantum computer. See the Supplemental Material for performance characteristics of the quantum computer (as well as further information about the measurements) [32].

There are several sources of error which prevent us from exactly reproducing the density matrix. These errors could lead to unphysical density matrices with negative eigenvalues. The standard way to avoid this is to use a maximum likelihood estimator [33] to obtain the closest physical density matrix based on the estimated one. All entanglement spectra obtained is shown in Fig. 2 (in the Supplemental Material, we provide the calibration data of the ibmqx5 for the day where the measurement of Fig. 2(c) was taken. We also show in Fig. 1 of the Supplemental Material [32] the entanglement spectra obtained another day and thus with another calibration set. While there is some variation between measurements on different days, the qualitative picture remains unchanged). We now list the various sources of error.

The repeated measurement gives rise to a statistical noise. For example, if the observable $\sigma_z$ in the state $|+\rangle = (1/\sqrt{2})(|0\rangle + |1\rangle)$ is measured 100 times, within 1 standard deviation, one may obtain 55 times “0” and 45 times “1”, and would conclude that $\langle \sigma_z \rangle = 0.1$ instead of the exact 0. To estimate the scale of the latter error, we computed the probabilistic outcome of 1024 $\times$ 256 measurements on the exact states for each of the three cases [PM], $|\text{cat}\rangle$, [SPT]. The entanglement spectra computed from the resulting estimate of the density matrix are plotted in Fig. 2 in the column “Ideal”.

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Additionally, errors in the real quantum computer arise from a finite coherence time of the qubits, as well as errors involved in implementing the gates and in reading out the qubits. They account for the difference between the columns “Ideal” and “IBM” (the latter obtained from the actual measurement) in Fig. 2. To quantify the statistical noise in the latter, we have generated a distribution of eigenvalues of reduced density matrices drawn from their distribution (see Supplemental Material for details [32]). The distributions are shown in the respective right-hand panels in Fig. 2. They show a robust entanglement gap in each of the cases [PM], [cat], [SPT], with the expected one, two, and four states above the gap. The fourfold degeneracy of the entanglement spectrum of [SPT] is only approximate in the measurement due to a combination of statistical noise and more importantly also the errors present in the quantum computer.

Using the measured reduced density matrices, we can further compute various entanglement measures. In Sec. IV of the Supplemental Material [32], we give the von Neumann entropy $S_{1,A} = -\text{Tr}_A[\rho_A \log \rho_A]$ and the second Rényi entropy $S_{2,A} = -\log \text{Tr}_A[\rho_A^2]$. The entanglement structure of gapped ground states of local Hamiltonians in one-dimensional systems is short-ranged. This property can be studied via the mutual information

$$I(A, B) = S_{1,A} + S_{1,B} - S_{1,AB},$$

where $A$ and $B$ are two subsystems. For the 8-site chain that we study, $A, B$ consist of two adjacent sites each. $A, B$ can then be a distance $d = 0, 1, 2$ sites apart, owing to the periodic boundary conditions. We measured the density matrix on the four sites comprising subsystems $A, B$ for $d = 0, 1, 2$ in each of the three prepared states [PM], [cat], and [SPT]. The results, summarized in Fig. 3, are qualitatively consistent with the following theoretical facts. (i) The mutual information of [cat] cannot be a gapped ground state of a local Hamiltonian, as its mutual information does not decay with distance, and (ii) [SPT] has only local entanglement that decays with distance. Furthermore, (iii) [PM] is a local product state with vanishing correlations, as even for $d = 0$ the mutual information vanishes; in our data, this mutual information does not vanish, but it is certainly smaller compared to the two other states. The

![FIG. 3. (a) Measured mutual information between two regions of the 8-site spin chain. The results are obtained from the measured reduced density matrices, measured in the same way as for Fig. 2. The theoretically exact values are $I(A, B) = 0$ and $I(A, B) = \log 2$ for all $d$ in the case of [PM] and [cat], respectively, and $I(A, B) = 2 \log 2, \log 2, 0$ for [SPT] at distances $d = 0, 1, 2$, respectively. (b) Schematic depiction of subsystem $A$ (blue) and subsystem $B$ (red) for the various distances $d = 0, 1, 2$.](image-url)
systematic discrepancy between the exact values and the measured values (irrespective of the states) is purely due to decoherence from qubit and gate errors.

Summary.—We measured the entanglement spectrum, a universal fingerprint of topological phases of matter, using the IBM quantum computer. Although the method as presented in this Letter is not suitable for obtaining the entanglement spectra of large subsystems due to the exponential scaling, the spectrum of a small subregion is already a useful probe when the region size exceeds the characteristic correlation length set by the gap. This is exactly the case for the states studied in this Letter and is more generally true for the ground states of gapped Hamiltonians. Nevertheless, if we wish to obtain the entanglement spectra of extensive subsystems, there are more sophisticated schemes available to reduce the exponential complexity of the problem. For example, the one presented in Ref. [13] outlines a quantum algorithm to obtain the $p$ largest eigenvalues ($\lambda_1 > \ldots > \lambda_p$) of the reduced density matrix, requiring only a parallel circuit of depth $O(p(\lambda_1/\lambda_p)^p)$ and $O(pL)$ qubits for a system of size $L$. With a small quantum computer of a few hundred qubits, this scheme would already allow one to obtain at least part of the entanglement spectrum of a system which cannot be computed classically [34].

However, with the improving gate fidelities and increasing number of qubits [35,36], it seems only a matter of time before quantum simulators or quantum computers will be able to implement more efficient algorithms and make more precise measurements of the entanglement spectrum of larger systems, thus providing a useful tool in understanding topological phases of matter. Further, since “topological software,” for instance the surface (toric) code [37], is envisioned as one venue to turn an analogue quantum computer into a digital (error-corrected) quantum computer, the characterization of quantum states in terms of the entanglement spectrum will be indispensable.

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