Quantum Entanglement as a Diagnostic of Phase Transitions in Disordered Fractional Quantum Hall Liquids

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We investigate the disorder-driven phase transition from a fractional quantum Hall state to an Anderson insulator using quantum entanglement methods. We find that the transition is signaled by a sharp increase in the sensitivity of a suitably averaged entanglement entropy with respect to disorder—the magnitude of its disorder derivative appears to diverge in the thermodynamic limit. We also study the level statistics of the entanglement spectrum as a function of disorder. However, unlike the dramatic phase-transition signal in the entanglement entropy derivative, we find a gradual reduction of level repulsion only deep in the Anderson insulating phase.

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Introduction.—Following the advances in understanding the fascinatingly complex phase diagram of two-dimensional electron systems in a strong perpendicular magnetic field in the fractional quantum Hall (FQH) regime [1–12], there has been intense interest in phase transitions in topological systems [13–17]. Disorder is a ubiquitous ingredient that may affect, even drive such phase transitions. In fact, it has been understood since the original explanation of the integer QH effect [18] that disorder plays an essential role in providing a finite domain for various QH states. Its presence allows a detailed study of QH plateau transitions in experiments [19–22], and in numerical studies of noninteracting models [23–27].

By contrast, there have been relatively few numerical studies of transitions from FQH states driven by disorder. The breaking of spatial symmetries by random disorder makes numerics even more challenging for this many-body problem with an exponentially large Hilbert space. One study [28,29] over a decade ago examined the disorder-driven transition from the \( f = 1/3 \) filling FQH state to the insulator using the ground-state Chern number as a diagnostic. It was shown that with increased disorder, the gap characterizing the FQH state collapsed, leading to an insulating phase at large disorder. Further, the calculated disorder dependence of the gap agreed with experiments.

In recent years, topological phases have been characterized by the underlying patterns of quantum entanglement [30–32]. Besides the vast theoretical literature, entanglement has been the subject of recent experiments [33] and related proposals [34]. Though many studies have been done for clean FQH states using the concepts of entanglement entropy and the entanglement spectrum (see Refs. [35–51] and references therein), there is relatively little corresponding effort [52,53] in the presence of disorder. In this work, we fill this void by studying the ground-state entanglement properties in a disordered \( f = 1/3 \) FQH system, as the disorder strength \( W \) is increased. We find that the magnitude of the derivative of a suitably defined entanglement entropy \( S \) with respect to \( W \) exhibits a sharp peak at a characteristic value of \( W = W_c \). The peak increases with system size, and is consistent with a divergence in the thermodynamic limit. We identify this behavior as signaling the FQH-insulator phase transition. Besides being a completely different diagnostic of the phase transition, our method (which only uses periodic boundary conditions) is significantly faster than calculating Chern numbers [28,29], which requires integrating over boundary conditions. We also analyze the entanglement spectrum (ES), and find that its level statistics undergoes an evolution which differs from that of highly excited states studied very recently in the context of many-body localization [54–56]. In our case, the low-energy part of the ES obeys Gaussian Unitary Ensemble (GUE) statistics at low \( W \), and changes to Poisson only at very large \( W \).

When the entanglement entropy changes its scaling behavior at a phase transition, e.g., from a volume law to an area law, one can define an “order parameter” by dividing by the volume. One thus obtains a zero order parameter on one side of the transition, and finite on the other, just like symmetry-breaking transitions. In our case, however, the entanglement entropy follows an area law on both sides [35,36,57], so this method does not work. Instead, we analyze our data using finite-size scaling (FSS) ideas (see, e.g., Ref. [58]), and do obtain a good FSS collapse of the data [59]. However, the exponent \( \nu \) that we obtain for the diverging length scale violates the expected inequality \( \nu \geq 2/d \) for \( d \)-dimensional disordered systems [60–62], a violation also seen in recent studies of many-body localization [63,64]. We provide possible explanations of this result in our concluding remarks.
We consider \( N \) interacting electrons in a two-dimensional random potential on a square with periodic boundary conditions (torus geometry). Landau levels are formed in the presence of a perpendicular external magnetic field. We assume that the energy scales of both interaction and disorder are small compared with the Landau level spacing, so we can focus on the lowest Landau level (LLL). The Hamiltonian of the system is

\[
H = \mathcal{P}_{\text{LLL}} \left( \sum_{i<j} V(\mathbf{r}_i - \mathbf{r}_j) + \sum_{i} U(\mathbf{r}_i) \right),
\]

where \( \mathcal{P}_{\text{LLL}} \) is the projector to the LLL, and \( V(\mathbf{r}) \) and \( U(\mathbf{r}) \) are the interaction and the random potentials, respectively. We further suppose that electrons interact via Haldaneopotential \cite{65} \( V = \text{Landau level spacing} \), so we can focus on the lowest interaction and disorder are small compared with the magnetic field. We assume that the energy scales of both boundary conditions (torus geometry). Landau levels are

\[
\text{exact}
\]

the ground states are exact\cite{24}. Since all single-particle states exist at the center of the LLL band\cite{24}. Since all single-particle states below the Fermi level at \( f = 1/3 \) are localized, the ground state is an Anderson insulator. In the following, we monitor the ground-state entanglement properties as a function of \( W \)\cite{68}.

Ground-state entanglement entropy.---The entanglement in a bipartite system can be measured by the von Neumann entropy of one subsystem. We make two cuts at orbital 0 and \([N_\phi/2] - 1\), respectively, where \([x]\) is the integer part of \( x \). This procedure divides \( N_\phi \) LLL orbitals into sub-systems A and B with boundary length \( L = 2 \times \sqrt{2\pi N_\phi} \) (in units of the magnetic length \( \ell_B \)), consisting of orbitals \( 0, \ldots, \left[ N_\phi/2 \right] - 1 \) and \([N_\phi/2], \ldots, N_\phi - 1\), respectively. In clean samples, there are three exactly degenerate Laughlin states. Disorder splits them. However, in the topological phase, the degeneracy is recovered in the thermodynamic limit\cite{69}. Such a topological degeneracy motivates us to consider a ground-state manifold containing the lowest three eigenstates \( |\Psi_{i=1,2,3}\rangle \) of the Hamiltonian at any \( W \) for consistency, rather than a single eigenstate. Similar to the situation of a single ground state, we define the ground-state entanglement entropy in our case as

\[
\rho = -\text{Tr}\rho_A \ln \rho_A, \quad \rho = \text{Tr}_{\mathcal{B}} \rho \text{ is the reduced density matrix of part A, and } \rho \text{ is the density matrix describing the ground-state manifold. There are several options to choose } \rho, \text{ for example, by using either } S(\bar{\rho}) \text{ with } \bar{\rho} = \frac{1}{3} \sum_{i=1}^{3} |\Psi_i\rangle \langle \Psi_i| \text{ or } \bar{\rho} = \frac{3}{4} \sum_{i=1}^{3} S(|\Psi_i\rangle). \text{ Both of them include the contributions of all } |\Psi_i\rangle's, \text{ thus minimizing the finite-size effect. Moreover, they give similar results} \cite{70}, \text{ so we adopt the first one in what follows.}
\]

At each \( W \), we implement different disorder configurations and compute the sample-averaged entanglement entropy \( \langle S(\bar{\rho}) \rangle \). The evolution of \( \langle S(\bar{\rho}) \rangle \) as a function of \( W \) is shown in Fig. 1(a) for various system sizes. For a fixed system size, \( \langle S(\bar{\rho}) \rangle \) decreases with \( W \). However, it increases with the system size at a fixed \( W \). \( \langle S(\bar{\rho}) \rangle \), although likely suffering from finite-size effects for our small sizes, agrees with an area law variation \( \langle S(\bar{\rho}) \rangle = aL - \beta [30-32] \) at all \( W \)'s[Fig. 1(b)]. The extracted entanglement density \( a \) starts to drop at \( W \approx 0.4 \), providing a rough estimate for the collapse of the Laughlin phase.

A more precise location of the phase transition can be obtained from the derivative of the entanglement entropy with respect to the disorder strength, i.e., \( dS(\bar{\rho})/dW \). To compute the sample-averaged derivative \( dS(\bar{\rho})/dW \), we average the derivative for each disorder configuration, approximated by

\[
dS(\bar{\rho})/dW = [S(\bar{\rho})]_{W+\Delta W} - [S(\bar{\rho})]_{W}/\Delta W \text{[71]. Figure 2(a) shows } dS(\bar{\rho})/dW \text{ as a function of } W \text{ for various system sizes. All curves exhibit a pronounced minimum at } W_c \approx 0.6, \text{ which gets deeper and sharper with increasing system size. A double logarithmic plot of its magnitude } h = \text{[min}(dS(\bar{\rho})/dW)] \text{ versus } N \text{[Fig. 2(b)] shows that } h \propto N^{1.35} \text{ consistent with a divergence in the thermodynamic limit. We consider this divergence to be a convincing signature of a phase transition between two phases with different entanglement properties.}
\]

Since the entanglement entropy obeys an area law in both phases, for a continuous phase transition, we may expect a scaling behavior for large size \( N \) of the form

\[
|\Psi_{i=1,2,3}\rangle \text{ is the density matrix describing the ground-state manifold. There are several options to choose } \rho, \text{ for example, by using either } S(\bar{\rho}) \text{ with } \bar{\rho} = \frac{1}{3} \sum_{i=1}^{3} |\Psi_i\rangle \langle \Psi_i| \text{ or } \bar{\rho} = \frac{3}{4} \sum_{i=1}^{3} S(|\Psi_i\rangle). \text{ Both of them include the contributions of all } |\Psi_i\rangle's, \text{ thus minimizing the finite-size effect. Moreover, they give similar results} \cite{70}, \text{ so we adopt the first one in what follows.}
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\]
the translational symmetry; therefore, we expect that versus their momenta, match the combination of two edge
in the ground-state manifold. For all superpositions obtained from the minimally entangled states is the same
in the parameter space spanned by ($\theta_i, \theta_2, \phi_1, \phi_2$) with $\theta_1, \theta_2 \in [0, \pi/2]$ and $\phi_1, \phi_2 \in [0, 2\pi]$, the local minima of $S([\Psi])$ in the parameter space spanned by ($\theta_1, \theta_2, \phi_1, \phi_2$) correspond to the MES. In the presence of topological degeneracy between $|\Psi_i\rangle$’s, MES are essential for the extraction of the topological entropy [72,73] and modular matrices [73–75]. We numerically search all minimally entangled states $|\Psi_{\text{min}}^i\rangle, i = 1, \ldots, M$ in each sample [76], and compute their average entanglement entropy $S_{\text{min}} = (1/M) \sum_{i=1}^{M} S(|\Psi_{\text{min}}^i\rangle)$). In Fig. 3, we show the sample-averaged entropy ($S_{\text{min}}$) as a function of $W$. At small disorder, ($S_{\text{min}}$) is almost a constant, indicating that the ground-state topological properties are the same as those of Laughlin states in clean samples. ($S_{\text{min}}$) starts to drop at $W \approx 0.6$, signifying a phase transition. Strikingly, the critical $W$ obtained from the minimally entangled states is the same as that suggested by the singular behavior of $dS(\rho)/dW$.

Ground-state entanglement spectrum (ES).—The spectrum of the reduced density matrix $\rho_A$ usually contains more information than the entanglement entropy, which is a single number obtained from the whole spectrum. Therefore, we now consider the ground-state ES $\{\xi_n\}$—the spectrum of $-\ln \rho_A$. Since the number of electrons $N_A$ in part $A$ is still conserved in the presence of disorder, we focus on the ES with $N_A = \lfloor N/2 \rfloor$, which has the largest rank.

In clean torus samples, the ES levels, when plotted versus their momenta, match the combination of two edge modes of pertinent FQH states [39,77]. Disorder breaks the translational symmetry; therefore, we expect that ground-state properties will be revealed by the statistics of the ES [78]. We compute the distribution $P(s)$ of the normalized level spacing $s_n/s_n$, where $s_n = \xi_n - \xi_{n-1}$ with $\xi_n$ sorted in ascending order, as a diagnostic of ES level statistics. To allow for variation between different parts of the ES, we examine level statistics in different windows of $\xi$ in the density of states (DOS) $D(\xi)$. Results from three individual ground states are quite similar [70], so we consider $D(\xi)$ and $P(s)$ averaged over three ground states below. This gives us more statistics in the evaluation. With increasing $W$, ES levels with large $\xi$ suffer from machine precision issues [70]. Consequently, we limit our discussion to those with $\xi \leq 40$, which are reliable.

With increasing $W$, we observe the development of localization in the ES. Localization first appears among low ES levels, then expands to higher $-\xi$ levels at larger $W$. In Fig. 4, we show the ES level statistics in three windows [79] for $N = 9$ electrons at various disorder strengths. At $W = 0.4$, we find strong level repulsion governed by GUE in all three windows; thus, the level spacing of the entire ES follows GUE [Fig. 4(a)]. The same ES level statistics is found at $W = 0.6$ [Fig. 4(b)] and $W = 1$ [Fig. 4(c)]. Only at significantly higher $W = 10$ is there a significant change in the level statistics—while $P(s)$ in the two higher $-\xi$ windows still obeys GUE, the level repulsion in $\xi \in (0, 10)$ is weaker, where GUE is replaced by a semi-Poisson distribution [Fig. 4(d)]. At $W = 100$, the spacing of the low ES levels becomes almost Poissonian without level repulsion, as for a localized system. Simultaneously, the level statistics in the middle window has evolved to semi-Poisson, while GUE is still preserved in the third window [Fig. 4(e)]. Finally, in the noninteracting limit with $W = \infty$, we observe the Poisson distribution for all ES levels below $\xi = 40$ [Fig. 4(f)]. It should be emphasized that this evolution of the ES takes place at significantly higher values of $W$ than the ground-state phase transition at $W \approx 0.6$ indicated by the entanglement entropy derivative.

We also track the number of reliable ES levels below $\xi = 40$ to get more insights about the ES evolution. We find that more and more ES levels move to the $\xi > 40$ region with increasing disorder [80]. At $W = \infty$, only a small
fraction of ES levels (18% for \( N = 9 \) and 8% for \( N = 10 \) electrons) forms one peak below \( \xi = 40 \) [Fig. 4(f)], where the Poisson distribution dominates the level statistics. The majority of ES levels have moved to the \( \xi > 40 \) region, which we cannot track due to the machine precision problem. Since ES levels in the rightmost part of the DOS always display GUE [Figs. 4(a)–4(e)], we might expect that the level statistics in the \( \xi > 40 \) region at \( W = \infty \) also follows GUE. Since \( \langle \mathcal{D}(\xi) \rangle \) is almost zero at \( \xi = 40 \), this would imply that there is an entanglement gap separating the Poisson part and GUE part for \( W = \infty \).

Summarizing the analysis above, we can obtain the following picture of the ES evolution. The entire ES follows GUE at small disorder [Figs. 4(a)–4(c)]. With the increase of \( W \), localization is first activated among low ES levels, then propagates towards the higher \( -\xi \) region, characterized by the evolution of level statistics from GUE to semi-Poisson, then to Poisson distribution [Figs. 4(d)–4(f)]. However, ES levels with GUE still exist, which are located at higher and higher \( \xi \) with the expansion of the DOS [Figs. 4(d), 4(e)]. Finally, in the noninteracting limit, the Poisson part forms a wide and small peak in the DOS [Fig. 4(f)], where only a small fraction of ES levels are located. The majority of ES levels belong to the GUE part with quite high \( \xi \) and is separated from the Poisson part by an entanglement gap. We notice that another two-component structure was also observed in the ES of highly excited energy eigenstates, but with different level statistics [54,55], reflecting the different evolution of the ground state and highly excited states with disorder.

Discussion.—In summary, through extensive exact-diagonalization studies of quantum entanglement properties of the \( f = 1/3 \) FQH system with \( N = 4–9 \) electrons on a torus geometry as a function of disorder, we find that the system undergoes a transition from the topological Laughlin state to a (topologically trivial) Anderson insulating state. The phase transition is signaled by a sharp peak (at a characteristic disorder \( W_c \approx 0.6 \)) in the magnitude of the disorder derivative of the entanglement entropy, \( dS/dW \), which appears to diverge in the thermodynamic limit (much as thermal transitions are often characterized by a singularity of the specific heat, which is proportional to the temperature derivative of the thermal entropy). From the scaling of the magnitude of the peak with size, as well as the entire \( dS/dW \) curve as a function of disorder in the vicinity of the transition, we obtain \( \nu \approx 0.6 \), which is very different from the exponent \( \nu \approx 1.5–2.6 \) characterizing plateau transitions [19–27]. Further, it violates the conventionally accepted bound [60] for nontopological transitions of disordered systems in two dimensions, which the result for integer quantum Hall transitions of noninteracting electrons does obey. A possible explanation for this anomalous result could be corrections to finite-size scaling, due to subleading terms in the entanglement entropy. However, in that case, the good finite-size scaling of the data will have to be accepted as being fortuitous. Violation of the bound was also found in recent numerical studies of many-body localization in one dimension [63,64].

Whatever the resolution of this scaling, the location of the transition appears to be quite robust. We find similar results for transitions of Coulomb interacting electrons from other FQH states to the insulator, to be reported in a subsequent publication [81]. Our results provide motivation for experimental studies of the length exponent characterizing the transition to the insulating state from various FQH states. A study of the universality of the diagonal conductivity at such a transition [82] was motivated by theoretical [83] and numerical [84] studies. In contrast to the entanglement entropy, the evolution of the level spacing of the entanglement spectrum is found to occur at much larger disorder, over an order of magnitude larger than the critical disorder, showing evidence of level repulsion in the entanglement spectrum until deep in the Anderson insulating phase.

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We normalize the pseudopotential such that the energy scale of the two-particle problem is unity.

In this work, the largest system size that we can reach by exact diagonalization is $N = 10$ electrons, whose Hilbert space dimension is $30045015$. This system size is smaller than the limit in clean systems due to the breaking of translation invariance by disorder. Considering that we often need at least hundreds of samples to do the ensemble average, we focus on $N \leq 9$ electrons (with Hilbert space dimension up to $4686825$) in most of our calculations to finish the computation in a reasonable time.

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See the Supplemental Material at http://link.aps.org/supplemental/10.1103/PhysRevLett.117.206801 for more numerical data of the ground-state entanglement entropy and entanglement spectrum.

The disorder configuration is fixed when we compute $S(\langle|W+h\rangle|W\rangle)$ and $S(\langle|W\rangle|W\rangle)$. Only the magnitude of $W$ is changed by a small percentage. We use $\Delta W = 0.001W$, though $\Delta W = 0.01W$ gives almost the same results.

In clean samples ($W = 0$) with translation invariance, there are three minimally entangled states with different momenta, corresponding to the three types of quasiparticles of the Laughlin state.

We compare the number of reliable ES levels below $\xi = 40$, i.e., $\lambda_{\xi<40} = \sum_{\xi<40} \langle\mathcal{D}(\xi)\rangle d\xi$, with the total number of ES levels $n_{\text{tot}} = \sum_{\xi} \langle\mathcal{D}(\xi)\rangle d\xi$. For $N = 9$ electrons, we have $n_{\text{tot}} = 715$ and $n_{\xi<40} = 715, 715, 715, 683, 468, 468, 126$ at $W = 0.4, 0.6, 1, 10, 100, \infty$, respectively.

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