INTRODUCTION

A main focus of modern practical quantum information research is on the generation of large-scale quantum entanglement involving many particles with the goal of achieving real applications of quantum technologies. \cite{1,2} Recent quantum experiments dealing with a large number of particles, such as optical lattice simulations involving $10^3$–$10^4$ atoms, \cite{3–4} experiments with hundreds of trapped ions \cite{5} or thousands of qubits in D-Wave systems (See: https://www.dwavesys.com/), show the real potential for applications of quantum technologies in the near future. An important instance of this challenge is the verification problem, i.e., how to reliably detect the presence of quantum resources, in particular quantum entanglement. The plausibility of standard verification schemes is questionable, since they require repeated measurement on large ensemble of identically prepared copies, which is highly demanding to achieve in practice.

One way of detecting quantum entanglement is to perform full quantum state tomography \cite{6–7} from which one can extract full information about the quantum state preparation (one can recover the entire density matrix). However, the full tomography becomes an unrealistic task already for a moderate size of quantum systems as the number of required measurement settings grows exponentially fast with the size of system. Luckily, in many cases, the knowledge of the entire quantum state is not needed, i.e., one can witness the presence of entanglement by measuring the mean values or the higher moments of a moderate number of identically prepared copies such that the sample averages are close to the true mean values. However, as any practical situation deals with a finite amount of data generated by a quantum measurement, the presence of entanglement can be verified only with a certain level of confidence, thus an adequate statistical analysis is necessary. \cite{8–10} Moreover, it is very hard to fulfill these requirements when dealing with large-scale entangled systems as only a limited (rather low) number of instances of a given quantum resource is available, due to various technical challenges, such as the lack of a good control and manipulation. As an example, one can take a recent experiment done with single photons \cite{11} where a 10-photon coincidence was registered every 5 min in average. With the same technology, that is by using the parametric down-conversion and postselection techniques, every additional photon pair (e.g., 12, 14, … photons) would reduce the count rate by at least one order of magnitude, consequently making the duration of the experiment (to verify entanglement) months or even years longer. In such situations, where only a limited number of resources is available, the main question arises whether it is still possible to reliably and efficiently certify the presence of quantum entanglement? This question is not only interesting from the theoretical point of view, it is also of great importance for practical quantum information.

In recent years we have seen several works that go beyond i.i.d. scenario, in the context of quantum state tomography \cite{12–14} and reliable entanglement verification. \cite{15–16} Although the techniques
and methods developed there are quiet generic, they still require a large sample sizes to verify entanglement with high confidence. On the other hand, in situations where only a low number of instances of a given quantum resource are available it appears natural to employ random sampling techniques for reliable detection. The advantage of such methods comes from simplicity of the data analysis, since minimal prior knowledge of the global population is needed. In the quantum scenario, random sampling has been proven very useful for quantum communication complexity, tomography via compressed sensing, fidelity estimation, self-testing methods, quantum state certification, quantum secret sharing, and verification of quantum computing. Some of these methods can be used to verify entanglement probabilistically as demonstrated in. In the same spirit, we shall incorporate random sampling methods together with techniques of quantum communication complexity to propose entanglement verification scheme in the form of a quantum information task. Unlike focusing on a large ensemble of i.i.d. copies, our main target here is a single experimental run, i.e., the central quantity for entanglement detection is the probability of success to perform certain binary task, given that the state was entangled/separable. Therefore, our scheme is designed to detect entanglement probabilistically. This framework has two main advantages as compared to conventional detection schemes:

a. it promises a dramatic reduction of the resources needed for reliable verification in large quantum systems, and
b. it provides a simple tool for reliable statistical analysis of errors and confidence intervals.

Most importantly, we show that in many situations the probability of success (to accomplish certain binary task) decreases exponentially fast with the size of system for all separable input states, whereas it approaches certainty if a particular entangled state (the target state) was prepared. Thus, even a single experimental run (single copy) can reveal the presence of entanglement with high accuracy. To our best knowledge, this is the first demonstration of entanglement detection in a single-copy regime (apart from the well-studied example of an i.i.d. state ρS24, see example of k-producible state below). We explicitly construct the detection procedure for k-producible states and cluster states. We show that, for example, one copy of a 16-qubit k-producible state or a single-copy of 24-qubit linear cluster state suffices to certify entanglement with more than 95% confidence. Thus, our method is applicable for quantum experiments involving tens of entangled qubits with the promise of a dramatic reduction of the resources needed for reliable entanglement detection (as compared to standard methods). Furthermore, the method developed for k-producible states can be used to naturally embed the standard techniques based on entanglement witnesses into our framework, meaning the statistical analysis of confidence intervals and errors simple and straightforward. Finally, we develop a general method for entanglement detection in ground states of local Hamiltonians (that exhibit the so-called entanglement gap). These include many important classes of quantum states, such as the matrix product states and projected-entangled pair states as they can be seen as unique ground states of the so-called parent Hamiltonians. At the end, we analyze the noise effect and we show that our probabilistic detection is very robust against the noise modeled by an arbitrary separable state.

RESULTS

Detection framework

Our method relies on a decision procedure where entanglement is seen as the ability of quantum systems to answer certain “yes-no questions”. The main figure of merit is the probability of success that a certain binary cost function F evaluates to 1, i.e., P(F = 1). The main goal is to provide examples of quantum states where P decreases exponentially fast to zero as the size of system grows for all separable states, whereas it approaches certainty (P = 1) if a particular entangled state (the target state) was prepared. Therefore, one can verify the presence of quantum entanglement with high probability even by measuring a single copy of a large quantum system.

In order to explain how our scheme works we consider a quantum system composed of N subsystems each residing in a finite-dimensional Hilbert space of dimension d. We usually assume N is large, although all the formulas derived hold for general N. To each subsystem we associate a certain set of possible measurements that can be performed locally. For example, in the case of qubits we may chose to measure each of them in complementary bases (such as X and Z measurement). In the general case, we shall include the most general quantum measurements (POVMs). Thus to each subsystem we associate a set of M different measurement settings defined by the set of positive semi-definite operators E(k), where \[ \sum_{k} E(k) = 1 \] and m = 1...M. Here k labels the subsystem, m the measurement setting and l labels the measurement outcome.

Given a single-copy of an N-partite quantum system, the detection procedure consists of the following four steps (see Fig. 1):

1. A sequence of measurement settings \( \{m_1, m_2, ..., m_N\} \) is randomly generated from the probability distribution of settings \( \Pi(m_1, ..., m_N) \).
2. The measurements are locally executed on each subsystem and the set of outcomes \( \{l_1, ..., l_N\} \) is obtained.
3. A certain binary (0/1) cost function of settings and outcomes \( F_{[N]} = F^{\Pi(m_1, ..., m_N)} \) is computed,
4. If \( F_{[N]} = 0/1 \) we associate “success/failure” to the experimental run.

This is the way to establish the probabilistic framework for entanglement detection through the probability of success \( P(F_{[N]} = 1) \). Our goal here is to choose the cost functions such that the probability of success vanishes exponentially fast in N for all

![Fig. 1 Probabilistic entanglement detection. A single-copy of N-partite quantum state is prepared. The sequence of measurement settings \( \{m_1, ..., m_N\} \) is randomly drawn from distribution \( \Pi(m_1, ..., m_N) \). Each \( m_k \) is locally executed on kth subsystem and the set of outcomes \( \{l_1, ..., l_N\} \) is obtained. The value of binary cost function \( F_{[N]} = F^{\Pi(m_1, ..., m_N)} \) prescribes either “success” \( (F_{[N]} = 1) \) or “failure” \( (F_{[N]} = 0) \) to the experimental run.](image)
separable states $\rho_{\text{sep}}$

$$P_{\text{sep}}[F_{\text{sep}} = 1] \leq \exp[-Mc],$$

(1)

where $c > 0$ is some constant. On the other hand, the cost function $F_{\text{sep}}$ is chosen such that there is an entangled state for which $P_{\text{sep}}[F_{\text{sep}} = 1] \approx 1$, meaning that whenever the state $\rho_{\text{ent}}$ (target state) has been prepared, the detection scheme works even in a single-copy scenario. An explicit bound on the probability of success for concrete examples will be derived.

Example of k-productive state

A good example to start with is that of the k-productive entangled state, i.e. $|\psi_1\rangle|\psi_2\rangle \cdots |\psi_k\rangle$, where the products $|\psi_i\rangle$ involve at most k parties. For simplicity, we chose the target state to be the product of quantum singlets $|\psi_0\rangle = (|0\rangle - |1\rangle)\otimes N$, where $|\psi_\delta\rangle = \frac{1}{\sqrt{3}}(|01\rangle - |10\rangle)$. The example and task presented here is quiet similar to “non-local” quantum games (see refs. 36,37). Nevertheless, we believe it is a good starting point for introducing a more delicate examples that follow in subsequent sections. We consider the set of $\{X, Y, Z\}$ measurement settings for each qubit, meaning that the measurement is performed in the eigenbasis of Pauli operators with the set of binary outcomes $i = 0, 1$. The quantum singlet is the unique state for which $X \otimes X = Y \otimes Y = Z \otimes Z = -1$, meaning that the measurement of $X \otimes X, Y \otimes Y,$ and $Z \otimes Z$ reveals perfect anticorrelations. Let us introduce the projectors on the outcome $-1$ for the correlation measurement

$$Q = \frac{1 - X \otimes X}{2},$$

(2)

$$W = \frac{1 - Y \otimes Y}{2},$$

(3)

$$R = \frac{1 - Z \otimes Z}{2}.$$  

(4)

To these projectors we associate three measurement settings $S = \{XX, YY, ZZ\}$. Although, the projectors are commutative, there is no separable state for which the measurement reveals $Q = W = R = 1$ simultaneously (this is a property of the singlet state only). Therefore, if we pick one of the settings from $S$ randomly (with probability $1/3$), there is a chance of at most $2/3$ to get the outcome $1$ for all separable inputs. More precisely, in such a case, the probability of success

$$P_{\text{sep}} = \frac{1}{3}(Q + W + R) \leq \frac{2}{3},$$

(5)

for all separable two-qubit states $\rho_{\text{sep}}$. Here $(\cdot) = \text{Tr}(\cdot)\rho$ denotes the mean value. This observation clearly suggests a detection scheme. We shall divide the set of $2N$ qubits into consecutive pairs and for every individual pair we pick one of the settings from $S$ randomly (with probability $1/3$) and perform the corresponding correlation measurement $(2), (3)$ and $(4)$. For separable inputs, the bound $(5)$ suggests that the relative frequency of the outcome 1 cannot significantly exceed the value of $2/3$ (provided that $N$ is large). Formally, we define the frequency $R_{\text{sep}} = \sum_{i=1}^{N} F_k$, where $F_k$ is the outcome of the correlation measurement on individual pairs

$$F_k = \frac{1}{2} \left( 1 - (-1)^{i+j} \right).$$

(6)

Here $i, j = 0, 1$ label the single-qubit measurement outcomes for the $k$th pair. The cost function is defined as

$$F_{\text{sep}} = \begin{cases} 1, & R_{\text{sep}} \geq \frac{2}{3} \delta N; \\ 0, & R_{\text{sep}} < \frac{2}{3} \delta N, \end{cases}$$

(7)

where $\delta > 0$ is some constant we keep at the moment as a free parameter. In other words, we associate "success" to the run if the number of local successes $F_k$ exceeds certain threshold of $(\frac{2}{3} + \delta)N$. The overall probability of success reads

$$P_{\text{sep}}[F_{\text{sep}} = 1] = P_F + \cdots + P_\delta \geq \frac{2}{3} \delta N,$$

(8)

and we recognize in the last equation the probability that the sum of random variables $F_1 + \cdots + F_\delta$ exceeds the value of $(\frac{2}{3} + \delta)N$. If the input state is a product state $\rho_{\text{prod}} = \rho_1 \otimes \cdots \otimes \rho_{2N}$, the random variables $F_k$ are independent with $F_k \leq \frac{\delta}{2}$. For such a case the bound on $(8)$ is well studied in classical probability theory and the results are known as the Chernoff bounds.\(^\text{37}\) We show in the Appendix that

$$P_{\text{prod}}[F_{\text{sep}} = 1] \leq \exp[-D(X\|\mu)],$$

(9)

where $D(X\|\mu) = x \log x + (1 - x) \log (1 - x) \geq 0$ is the Kullback-Leibler divergence. Furthermore, if the bound holds for all product states, it also holds for their mixtures, i.e., it holds for all separable states. We see that the probability of success vanishes exponentially fast in $N$, for all $\delta > 0$. This is quite convenient, as we do not have to fix $\delta$ in advance. Once the experiment has been performed, we can calculate directly from the experimental data $F_1, F_2, \ldots, F_N$ how much the frequency deviates from $2N/3$, i.e., we set $\delta = (F_1 + \cdots + F_N)/N - 2/3$, and consequently calculate the bound on probability of success for separable states by using $(9)$. For the case of $|\psi_\delta\rangle = \frac{1}{\sqrt{3}}(|01\rangle - |10\rangle)$ input state, each local cost function $F_k = 1$ deterministically, thus we get $\delta = 1/3$. The bound $(9)$ reduces to

$$P_{\text{sep}}[F_{\text{sep}} = 1] \leq \left( \frac{2}{3} \right)^N.$$

(10)

If $N$ is sufficiently large, a single-copy of $|\psi_\delta\rangle$ suffices to certify presence of entanglement with high probability. For example, if we want to have a detection probability of at least 95% (i.e., we want to be sure that no separable state has a probability of success more than 5%) in a single-shot experiment, we get the minimal number $N_{\text{min}} = 8$, which is a remarkably low number.

The present example shows how the standard detection of entanglement in quantum singlet based on the witness operator $W = Q + W + R$ can be naturally embedded in our framework. Conventionally, one has to measure $Q, W$, and $R$ in three separate experiments by using an i.i.d. ensemble of qubit pairs $\rho^{(i)} \otimes \rho^{(j)}$ in order to estimate the mean values $\langle Q \rangle, \langle W \rangle$, and $\langle R \rangle$. However, the i.i.d. assumption is difficult to justify operationally, and consequently the statistical analysis involving many experiments is non-trivial. Furthermore, if the number of singlet pairs is low, it not clear how to actually pursue the detection scheme. For example, imagine a situation where only $N = 8$ pairs are available. The question is how to divide the pairs and perform the corresponding measurements. We may use the first three copies to measure $Q$, the second three to measure $W$, and the last two for the measurement of $R$. However, if the order is known and fixed in advance, than the following product state $\frac{1}{\sqrt{3}}(|x\rangle + |y\rangle + |z\rangle)$ gives exactly the same result as the i.i.d. state $|\psi_\delta\rangle$. Thus, we cannot conclude the presence of entanglement or we may even wrongly claim its presence. Certainly, a correct statement requires a proper statistical analysis. On the other hand, one of the key procedures in our method is the random sampling of measurement settings, which provides us a simple tool to analyze the errors and confidence intervals trough the probability of success. Therefore, there is a clear separation between the state $|\psi_\delta\rangle$ and the product state given above, as the latter has only the chance of $(2/3)^8 = 0.039$ to reveal the result $F_1 + \cdots + F_8 = 8$.

In general, any detection based on witness operator can be incorporated in our framework, with the goal to achieve more resource-efficient entanglement detection. For a witness $W = \sum W_i$, one has to sample the measurements of $W_i$ randomly every single experimental run. The bound similar to $(9)$ can be easily derived. Nevertheless, one may observe that in
such a case, one still requires many copies for reliable detection (i.e., $N$ copies of $k$-particle state $|\psi\rangle$ folded into a single multipartite copy $|\psi^{[N]}\rangle$). In the next examples we will unambiguously show that, indeed, one can certify entanglement by measuring only a single-copy.

Example of cluster states

Another example we present here is that of cluster states. In contrast to the previous example of $k$-producible states, cluster states contain genuine multipartite entanglement and are known to be a universal resource for measurement-based quantum computation. For the sake of simplicity, we shall explain how the single-copy detection scheme works for the case of linear cluster states (LCS). The generalization to higher dimensions is straightforward and we briefly discuss it at the end of this section. The $N$-qubit LCS is uniquely defined by the set of $2^N$ stabilizers, i.e.,

$$G_{q_1\ldots q_N}|LCS\rangle = G_{q_1}^\dagger \cdots G_{q_N}^\dagger |LCS\rangle = +1|LCS\rangle,$$

(11)

where $G_{q_1}$ are $Z$-type operators $Z_{q_1} \ldots Z_{q_{k-1}}$ and $q_k = 0, 1$. Here $\{X, Y, Z\}$ represent the set of standard Pauli matrices acting on $k$th qubit and, for simplicity, we have chosen the cluster state with periodic boundaries, i.e., $X_{N+1} = X_1$ and $X_{N+1} = X_1$. We consider the set of $\{X, Y, Z\}$ measurement settings for each qubit, meaning that the measurement is performed in the eigenbasis of Pauli operators with the set of binary outcomes $i = 0, 1$. We start with a simple analysis by considering a small subset (cluster) of four qubits, let say (1, 2, 3, 4) with the corresponding stabilizers $G_1 = Z_1 Z_2 Z_3$, $G_2 = Z_1 Z_2 Z_3$, and $G_3 = Z_1 Z_2 Z_3$. All three stabilizers are commutative, they are not locally compatible (in a sense, there is no local measurement of all three of them simultaneously), therefore there is no product (separable) state for which all three observables can take the same value, i.e., $G_1 G_2 G_3 = +1$ simultaneously. For that reason, if we randomly chose (with probability 1/3) to measure one of the stabilizers there is only a chance of 2/3 to get the result +1, for all separable inputs (similar to the previous example of singlet state). This is the key observation that enables our detection method to work. Our main idea is to show that if we pick a random partition of the set of $N$ qubits into 4-qubit clusters and subsequently on each of them randomly measure one of the corresponding stabilizers, the relative frequency of the outcome +1 cannot significantly exceed the value of 2/3. More formally, we start by introducing partitions of $N$ qubits into 4-qubit clusters $\{c_1, c_2, \ldots, c_{\ell}\}$, where $c_\ell$ is the cluster involving the sequence of four qubits $c_\ell = \{t_1, t_1 + 1, t_1 + 2, t_1 + 3\}$. Furthermore, we see from the previous analysis that the border qubits in each cluster are always measured in the $Z$-basis (when measuring the corresponding stabilizer). Thus, we shall allow for possible overlap between neighbouring clusters on border qubits. More precisely, we say that the partition is regular if the neighbouring clusters overlap on at most one (border) qubit, i.e., $t_\ell - t_{\ell-1} \geq 3$. For example, the partition $\{\ldots, 7, 8, 9, 10, \ldots\}$ is considered regular, whereas $\{\ldots, 7, 8, 9, 10, 9, 10, 11, 12\}$ is irregular, as the two clusters in partition overlap on qubits 9 and 10 (see SI for more examples). We denote the set of all numbers of qubits of size $L$ by $\mathcal{C}_L$. We shall think of $L$ as being large, e.g., on the same scale as $O(N)$ with the number of qubits and, at the same time, we shall choose $L$ such that the set $\mathcal{C}_L$ is large in size as well. The clusters in the partition serve as the building-blocks for the construction of the cost function $F_{\text{prod}}$. For every cluster $c_\ell$ in the partition there are three stabilizers associated to it: $G_{c_\ell+1} = Z_{t_1} Z_{t_1 + 1} Z_{t_1 + 2}$, $G_{c_\ell+2} = Z_{t_1 + 1} Z_{t_1 + 2} Z_{t_1 + 3}$ and $G_{c_\ell+1 t_1 + 2} = G_{c_\ell+1} G_{c_\ell+2} = Z_{t_1} Y_{t_1 + 1} Y_{t_1 + 2} Z_{t_1 + 3}$. To each of them we associate three projectors

$$Q_{c_\ell} = \frac{1 + G_{c_\ell+1}}{2},$$

(12)

$$W_{c_\ell} = \frac{1 + G_{c_\ell+2}}{2},$$

(13)

$$R_{c_\ell} = \frac{1 + G_{c_\ell+1} G_{c_\ell+2}}{2},$$

(14)

that project on the +1 outcome. We associate the following measurement settings with each projector

$\{ZXXZ, ZZZZ, ZYYZ\}$,

(15)

and we assign “success” to the cluster measurement only if the outcome +1 (for the value of measured stabilizer) occurs. Formally speaking, for every cluster we define the following local cost function

$$F_m = F_m^{b/h} = \frac{1}{2} + \frac{1}{2} \left\{ \begin{array}{ll}
-1 & m = ZXXZ; \\
-1 & m = ZZZZ; \\
-1 & m = ZYYZ;
\end{array} \right.$$

(16)

where $s = 1 \ldots L$. Finally, for a given partition $\{c_1, c_2, \ldots, c_L\}$ the overall cost function is defined in the following way

$$F_{\text{prod}}[s] = \left\{ \begin{array}{ll}
1 & F_1 + \ldots + F_L \geq \left( \frac{2}{3} + \delta \right) L; \\
0 & F_1 + \ldots + F_L \leq \left( \frac{2}{3} + \delta \right) L.
\end{array} \right.$$  

(17)

Now, we will show that the probability of success vanishes exponentially fast for all separable states as the number of qubits grows. Firstly, for a fixed partition $\{c_1, c_2, \ldots, c_L\}$ it is clear that product states fail to satisfy $F_{\text{prod}} = 1$ for all three settings $\{ZXXZ, ZZZZ, ZYYZ\}$, because $XZ, ZX, YY$ are locally incompatible on a second and third cluster qubit. Thus, if the settings are uniformly distributed (with probability 1/3), one can easily show that the probability of success for individual clusters

$$P_{\text{prod}}[s = 1] = \langle F_1 \rangle = \frac{1}{3} \left( Q_1 + W_1 + R_1 \right) \leq \frac{2}{3},$$

(18)

for all product states $\rho_{\text{prod}} = \rho_1 \otimes \ldots \otimes \rho_N$. Furthermore, if the input state is a product state, the local cost functions $F_i$ can be seen as independent binary (0/1) random variables with $\langle F_i \rangle \leq 2/3$ for all $s = 1 \ldots L$. The overall probability of success reads

$$P_{\text{prod}}[F_{\text{prod}} = 1] = P_{\text{prod}} \left[ F_1 + \ldots + F_L \geq \left( \frac{2}{3} + \delta \right) L \right].$$

(19)

and we recognize in the last equation the probability that the sum of independent random variables $F_1 + \ldots + F_L$ exceeds the value of $\left( \frac{2}{3} + \delta \right) L$. As $\langle F_i \rangle \leq 2/3$ we expect that the sum $F_1 + \ldots + F_L$ cannot exceed $2/3L$ significantly. Similar to the previous example (of singlet state), the Chernoff bound holds (see Appendix for the proof), i.e.,

$$P_{\text{prod}}[F_{\text{prod}} = 1] \leq e^{-D(\frac{2}{3} + \delta | \frac{2}{3}) L},$$

(20)

where $D(x|y)$ is the Kullback–Leibler divergence. Furthermore, if the bound holds for all product states, it also holds for their mixtures, i.e., it holds for all separable states. Thus, as long as $L$
grows with $N$ (for example we can set $L = [N/2]$, where $[\cdot]$ denotes the integer part) the probability of success vanishes exponentially fast, for all $\delta > 0$. As before, we do not have to fix $\delta$ in advance. Once the experiment has been performed, we can calculate directly from the experimental data $F_1, F_2, \ldots$ how much the sum of results deviates from $2L/3$, i.e., we set $\delta = (F_1 + \ldots + F_l)/L - 2/3$, and consequently calculate the bound on probability of success for separable states by using (20). For the case of cluster state preparation $|\text{LCS}\rangle$, each local cost function $F_l = 1$ deterministically, thus we get $\delta = 1/3$. The bound (20) reduces to

$$P_{\rho_\text{L}}[F_1 = 1] \leq \left( \frac{2}{3} \right)^L.$$  \hspace{1cm} (21)

If the number of qubits is sufficiently even, even a single-copy of LCS suffices to certify presence of entanglement with high probability. For example, if we want to have a detection probability of at least 95% (i.e., we want to be sure that no separable state has a probability of success more than 5%) in a single-shot experiment, we get the minimal number of clusters $L_{\text{min}} = 8$. The lowest number of qubits with such support is $N = 24$. Nevertheless, in such a case the set of all partitions $C_s$ reduces to three only (see discussion at the end of this section), one may want to have $|C_s|$ significantly larger. For example, already $N = 25$ has $|C_s| = 25$, for $N = 26$ we get $|C_s| = 117$ etc. The scheme can be used not only to detect entanglement, it can be also used to certify the presence of LCS. To see this, note that $P_{\rho}[F_1 = 1] = \text{Tr} P \rho$, where

$$\Pi = \frac{1}{|C_1|} \sum_{\{c_1, \ldots, c_N\} \in C_1} \prod_{k=1}^{N} \frac{1}{3} (Q_{c_k} + W_{c_k} + R_k).$$  \hspace{1cm} (22)

Clearly LCS is the eigenstate $\Pi |\text{LCS}\rangle = |\text{LCS}\rangle$ for the maximal eigenvalue. Now, the operator $\Pi$ can be expanded in terms of stabilizers $G_{c_1 \ldots c_N}$ defined by the Eq. (11). If the set $C_1$ is sufficiently large the expansion will include all $2^N$ stabilizers. Since the LCS is the only state with $G_{c_1 \ldots c_N} |\text{LCS}\rangle = |\text{LCS}\rangle$ for all stabilizers, we conclude that LCS is a unique eigenstate of $\Pi$ for eigenvalue 1. Thus, the LCS state is the only state with the maximal probability of success, i.e., $P_{\rho}[F_1 = 1] = 1$.

We shall briefly comment on the type of entanglement certified by the single-copy detection scheme. Firstly, if we are willing to detect multipartite entanglement, it is very important that the set $C_1$ is large in size. Recall, that the bound (20) holds for arbitrary partition from the set $C_1$ for all separable states. Therefore, if the partition $\{c_1, \ldots, c_N\}$ is fixed and known in advance, the bound (20) still holds. Nevertheless, for such a case, the following 4-producible state $|\psi\rangle = |\psi_1\rangle |\psi_2\rangle |\psi_3\rangle |\psi_4\rangle$, where $|\psi_i\rangle$ is the common eigenstate for all three projectors $Q_{c_i}, W_{c_i}$ and $R_i$ (for eigenvalue 1), reveals $F_1 = 1$ deterministically for every cluster. Consequently we have $P_{\rho}[F_1 = 1] = 1$ for $|\psi\rangle$ being the input state. Such a state contains localized entanglement on individual clusters (blocks of entanglement). To prevent $|\psi\rangle$ maximizing the probability of success, a random choice of partition from a large set $C_1$ is necessary. For example, already including additional partition $\{c_1, \ldots, c_{N+1}\}$ obtained by shifting one qubit to the right, prevents $|\psi\rangle$ for the reason explained below (see discussion at the end of this section), one may want to have $|C_1|$ significantly larger. For example, already $N = 25$ has $|C_1| = 25$, for $N = 26$ we get $|C_1| = 117$ etc. The scheme can be used not only to detect entanglement, it can be also used to certify the presence of LCS.

Finally, let us briefly explain the generalization to the higher dimensional case. Take an example of a 2D cluster state (known to be universal for quantum computation). Here, one can introduce partitions into $4 \times 4$ qubit clusters with the corresponding stabilizer projectors (in analogy to $Q_{c_i}, W_{c_i}$ and $R_i$ for LCS) and define the local cost functions. In complete analogy to the 1D case, the 2D detection scheme consists of drawing a random partition followed by a random measurement of local projectors on individual clusters. The separable bound similar to (20) can be derived. On the other hand, if the 2D cluster state has been prepared, the probability of success is 1.

Ground states of local Hamiltonians

One of the reasons that single-copy entanglement detection works for cluster states can be associated to the robustness of entanglement with respect to local perturbations. For example, if we measure one or even a group of localized qubits in cluster state, entanglement remains present between the rest qubits. Ground states of local Hamiltonians are believed to share this property (robustness of entanglement),\(^5\) therefore we can expect that they are also amenable to single-copy verification. Here we show that indeed this is the case.

Consider a $L$-local Hamiltonian on some graph of $N$ particles $H = \sum_{k=1}^{N} H^{(k)}$, where $H^{(k)}$ acts on at most $L$ subsystems ($L$ is fixed and independent of $N$). For simplicity reason, we assume the number of local terms $h$ of $H^{(k)}$ to match the number of particles $N$ (this is common for many physical situations, see for example\(^6\)). One may consider a more general case where the number of local terms grows as a polynomial function of $N$. Nevertheless, the detection scheme shall work the same way. Let $|\psi_0\rangle$ is the ground state $H|\psi_0\rangle = \mathcal{N}|\psi_0\rangle$, where $\mathcal{N} = \text{cst}$ is the ground-state energy. We are interested in Hamiltonians that exhibit the so-called entanglement gap $\delta = \epsilon_1 - \cdots - \epsilon_D > 0$, which is the minimal achievable energy (per particle) by a separable state. Furthermore, we assume $\mathcal{N}$ to be finite and non-zero in the thermodynamical limit, i.e., $\mathcal{N} = \text{const}$. To summarize, we are interested in Hamiltonians where the average energy $\langle H \rangle$, $\langle H \rangle = \mathcal{N}$, for LCS) and $\langle H \rangle \geq N\mathcal{N} = N\mathcal{N}$ for all separable states, whereas at least the ground state violates this bound.

We shall develop a general scheme that works for arbitrary local Hamiltonian. For that reason we introduce a set of tomographically complete measurement for each particle. For example, in the case of qubits, a natural choice are the three complementary measurements defined by $X, Y$, and $Z$ Pauli operators. Thus, the set of measurement operators $E_{(k)}^m$ forms a complete basis in the space of observables, i.e., any observable $A^{(k)}$ acting on $k$th subsystem can be decomposed as $A^{(k)} = \sum_{m=1}^{N} E_{(k)}^m$. Here $m = 1 \cdots M$ and $i = 1 \cdots D$, where $M$ is the number of settings and $D$ is the number of outcomes. Furthermore, in order to simplify the notation, we introduce a new variable $x_i = (m_i, i_j)$ which labels a pair of measurement setting and outcome, hence $E_{x_i}^m$ refers to $E_{(k)}^m$. Note that $\sum_{x_i} E_{x_i}^m = M^{(k)}$.

For a given local Hamiltonian $H = \sum_{k=1}^{N} H^{(k)}$, operator $H^{(k)}$ acts only on at most $L$ neighbouring subsystems (neighbours of $k$ including $k$ itself). It is convenient to introduce the $N \times L$ “neighbouring” matrix $N_{x_k}^{(k)}$, where $N_{x_k}^{(k_i)} \ldots N_{x_k}^{(k_L)}$ is a sequence of integers labeling all the neighbours of $k$th subsystem (including $k$th subsystem itself) on which the local operator $H^{(k)}$ acts. The “neighbouring” matrix can be seen as the list of neighbourhoods $(N(1), \ldots, N(N))$, where $N(k)$ denotes the set of all neighbours of $k$. For example, the notation $(n_{x_3}, n_{x_2}, n_{x_1}) \in \{2, 3, 4\}$ means that $H^{(3)}$ acts on subsystems 2, 3, and 4. Because the set of measurement operators is tomographically complete, each $H^{(k)}$ can be decomposed into the sum of products of local measurement operators

$$H^{(k)} = \sum_{x_1, \ldots, x_k} \prod_{i=1}^{L} E_{x_i}^{(n_{x_i}^{(k)})} \cdots E_{x_k}^{(n_{x_k}^{(k)})}.$$  \hspace{1cm} (23)

The operator $H^{(k)}$ can be completely identified with the tensor $h^{(k)} = h^{(k)}_{x_1, \ldots, x_k}$. Similarly, the full Hamiltonian $H$ reads

$$H = \sum_{x_1, \ldots, x_N} \prod_{i=1}^{L} E_{x_i}^{(n_{x_i}^{(k)})} \cdots E_{x_N}^{(n_{x_N}^{(k)})},$$  \hspace{1cm} (24)

where we set $M^{N-L} H_{x_1, \ldots, x_N} = \sum_{k=1}^{N} h^{(k)}_{x_1, \ldots, x_k}$ (the factor $M^{N-L}$
comes because of the normalization $\sum_{i} E_i = M1$. Now we can set-up the detection procedure. Firstly, we pick measurement settings for individual subsystems randomly (i.e., with probability $1/M$) and generate the sequence $(m_1, ..., m_N)$. The measurements are executed on local subsystems and the set of outcomes $(i_1, ..., i_N)$ is obtained. Equivalently, we say that the sequence of random variables $(x_1, ..., x_N)$ is generated, where $x_i = (m_i, i)$. Now, we shall define the cost function $F_{\text{HI}}$. It is convenient to define $H_{\text{HI}} = M^{1/2} (x_1, ..., x_N) = M^{1/2} \sum_{i=1}^{N} h^{(i)}$. A straightforward inspection shows $(h^{(i)}(\mathbf{H})) = \text{Tr} \, \rho \, h^{(i)}(\mathbf{H})$, thus the classical random variable $H_{\text{HI}}$ can serve to extract the mean value of Hamiltonian $(\mathbf{H})$. Since $(\mathbf{H}) \geq N \epsilon$, holds for all separable states, it is natural to choose the following cost function

$$F_{\text{HI}} = \begin{cases} 1, & H_{\text{HI}} \leq N(\epsilon_1 - \delta); \\ 0, & H_{\text{HI}} > N(\epsilon_1 - \delta), \end{cases}$$

(25)

where $0 < \delta < \epsilon_1 - \epsilon_0 = g_0$ is constant we keep at the moment as a free parameter. Since the random variable $H_{\text{HI}}$ completely captures properties of Hamiltonian, we expect $H_{\text{HI}}$ not to precede the separable bound $N \epsilon_1$ significantly in a single-shot experiment (provided that $N$ is large). Indeed, in the Appendix we show that for all separable states $\rho_{\text{sep}}$ the following bound holds

$$P_{\rho_{\text{sep}}} \{ F_{\text{HI}} = 1 \} \leq \exp \left[-N \kappa^2 \delta^2 \right],$$

(26)

where $\kappa > 0$ is constant. Thus, the probability of success vanishes exponentially fast with $N$ for all separable inputs. On the other hand, if the ground state $|\psi_0\rangle$ is prepared, we show in the Appendix that the probability of success reaches 1 in the thermodynamical limit, i.e.,

$$P_{\psi_0} \{ F_{\text{HI}} = 1 \} \geq 1 - \frac{\beta^2}{N(g_2 - \delta)^2},$$

(27)

where $\beta > 0$ is constant. In other words, if $N$ is sufficiently large, the probability of success is close to 1.

There are several points worth of mentioning here. Firstly, the previous example of cluster states can be incorporated in the present scheme, since cluster states can be seen as unique ground-states of local Hamiltonians.$^{61}$ Nevertheless, the detection scheme introduced in the previous section is more resource-efficient for cluster states for the following reasons: a) the bound (20) is more tight than (26), and b) the probability of success evaluates to 1 (for the cluster-state input), in contrast to (27) which reaches 1 asymptotically. On the other hand, the ground-state detection method has certain practical advantages. Namely, one of the crucial elements for detection is the use of tomographically complete set of measurements. In principle, they can be substituted by a single informationally complete POVM (ICPOVM).$^{62}$ More precisely, instead of a set of tomographically complete measurements, a single POVM with the measurement operators $E_i$, forming a complete basis in the space of observables, can be used. Thus, an $N$-partite Hamiltonian can be expressed as $H = \sum_{i_1, ..., i_N} h_{i_1, ..., i_N} E_{i_1(1), ..., i_N(N)}$, where $i_k = 1, ..., D$ labels the measurement outcome. The properties of Hamiltonian are fully captured by the classical random variable $H_{\text{HI}} = h_{i_1, ..., i_N}$ (function of the measurement outcomes $i_1, ..., i_N$). Now, there is no random sampling of measurement settings, there is only one measurement (ICPOVM) for each particle. The variable $H_{\text{HI}}$ is calculated from the set of measurement outcomes $(i_1, ..., i_N)$. The cost function is defined as (25), and derivation of bounds (20) and (26) is essentially the same as before. Formally, both methods are equivalent. Nevertheless, practical advantage of using ICPOVM compared to random sampling of measurement settings can be significant in certain cases, conditioned on the physical implementation of POVM. For example, if ICPOVM is implemented by using additional degrees of freedom (e.g., for the case of single-photons by combining the path and polarization degree of freedom$^{63}$), then, the same, single measurement setting is applied on every local subsystem. This is very convenient when dealing with large-scale quantum systems, for which full manipulation and addressability of individual particles is demanding to achieve.

Discussion

We introduced a probabilistic technique for resource-efficient entanglement detection in large-scale multiparticle quantum systems. We have shown that for variety of quantum states, probability to detect entanglement (as quantified by the
probability of success) approaches one exponentially fast with the size of system, implying that even a single copy suffices to verify entanglement with high probability. Our method promises a dramatic reduction of the resource needed for reliable entanglement verification, therefore it has great potential for practical applications in current and near future experiments aiming at generation and manipulation of massive entanglement.

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AUTHOR CONTRIBUTIONS
B.D. supervised the project, provided theoretical analysis and wrote the manuscript. A.D. provided theoretical analysis and wrote the manuscript.

ADDITIONAL INFORMATION
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