

Probability-density-function characterization of multipartite entanglement

P. Facchi,^{1,2} G. Florio,^{3,2} and S. Pascazio^{3,2}

¹*Dipartimento di Matematica, Università di Bari, I-70125 Bari, Italy*

²*INFN, Sezione di Bari, I-70126 Bari, Italy*

³*Dipartimento di Fisica, Università di Bari, I-70126 Bari, Italy*

(Received 31 March 2006; published 26 October 2006)

We propose a method to characterize and quantify multipartite entanglement for pure states. The method hinges upon the study of the probability density function of bipartite entanglement and is tested on an ensemble of qubits in a variety of situations. This characterization is also compared to several measures of multipartite entanglement.

DOI: [10.1103/PhysRevA.74.042331](https://doi.org/10.1103/PhysRevA.74.042331)

PACS number(s): 03.67.Mn, 03.65.Ud

I. INTRODUCTION

Entanglement is one of the most intriguing features of quantum mechanics. Although it is widely used in quantum communication and information processing and plays a key role in quantum computation, it is not fully understood. It is deeply rooted in the linearity of quantum theory and in the superposition principle and basically consists (for pure states) in the impossibility of factorizing the state of the total system in terms of states of its constituents.

The quantification of entanglement is an open and challenging problem. It is possible to give a good definition of *bipartite* entanglement in terms of the von Neumann entropy and the entanglement of formation [1]. The problem of defining *multipartite* entanglement is more difficult [2] and no unique definition exists: different measures capture in general different aspects of the problem [3]. Attempts to quantify the degree of quantum entanglement are usually formulated in terms of its behavior under local operations/actions that can be performed on different (possibly remote) parts of the total system. Some recent work has focused on clarifying the dependence of entanglement on disorder and its interplay with chaos [4,5], or its behavior across a phase transition [6,7].

The work described here is motivated by the observation that as the size of the system increases, the number of measures (i.e., real numbers) needed to quantify multipartite entanglement grows exponentially. A good definition of multipartite entanglement should therefore hinge upon some statistical information about the system. We shall look at the distribution of the purity of a subsystem over all possible bipartitions of the total system. As a characterization of multipartite entanglement, we will not take a single real *number*, but rather a whole *function*: the probability density of bipartite entanglement between two parts of the total system. The idea that complicated phenomena cannot be “summarized” in a single (or a few) number(s) stems from studies on complex systems [8] and has been considered also in the context of quantum entanglement [9]. In short, we expect that multipartite entanglement will be large when bipartite entanglement is large *and* does not depend on the bipartition, namely when its probability density is a narrow function centered at a large value. This characterization of entanglement will be tested on several classes of states and will be compared with several measures of multipartite entanglement.

II. THE SYSTEM

We shall focus on a collection of n qubits. The dimension of the Hilbert space is $N=2^n$ and the two partitions A and B are made up of n_A and n_B spins ($n_A+n_B=n$), respectively, where the total Hilbert space reads $\mathcal{H}=\mathcal{H}_A\otimes\mathcal{H}_B$ and the Hilbert spaces \mathcal{H}_A and \mathcal{H}_B have dimensions $N_A=2^{n_A}$ and $N_B=2^{n_B}$, respectively ($N_A N_B=N$). We shall consider only pure states

$$|\psi\rangle = \sum_{k=0}^{N-1} z_k |k\rangle, \quad (1)$$

where $|k\rangle=|j_A\rangle\otimes|l_B\rangle$, with a bijection between k and (j_A, l_B) , $0\leq j_A\leq N_A-1$, and $0\leq l_B\leq N_B-1$. As a measure of bipartite entanglement between A and B , we consider the participation number

$$N_{AB} = \pi_{AB}^{-1}, \quad \pi_{AB} = \text{Tr}_A \rho_A^2, \quad \rho_A = \text{Tr}_B \rho, \quad (2)$$

where $\rho=|\psi\rangle\langle\psi|$, and Tr_A (Tr_B) is the partial trace over the degrees of freedom of subsystem A (B). N_{AB} can be viewed as the relevant number of terms in the Schmidt decomposition of $|\psi\rangle$ [10]. The quantity $n_{AB}=\log_2 N_{AB}$ represents the effective number of entangled spins. Clearly, for a completely separable state, $\text{Tr}_A \rho_A^2=1$ for all possible bipartitions, yielding $N_{AB}=1$ and $n_{AB}=0$. In this sense, the participation number can distinguish between entangled and separable states. Moreover, π_{AB} is directly related to the linear entropy $S_L=1-\pi_{AB}$, which is an entanglement monotone, i.e., it is nonincreasing under local operations [11] and classical communication. In general, the quantity N_{AB} will depend on the bipartition, as generally entanglement will be distributed in a different way among all possible bipartitions. Therefore, its distribution $p(N_{AB})$ will yield information about multipartite entanglement: its mean will be a measure of the amount of entanglement in the system, while its variance will measure how well such entanglement is distributed, a smaller variance corresponding to a higher insensitivity to the particular choice of the partition.

We will show that for a large class of pure states, statistically sampled over the unit sphere, $p(N_{AB})$ is very narrow and has a very weak dependence on the bipartition: thus entanglement is uniformly distributed among all possible bipartitions. Moreover, $p(N_{AB})$ will be centered at a large

value. These are both signatures of a very high degree of *multipartite* entanglement.

By plugging Eq. (1) into Eq. (2), one gets

$$\pi_{AB} = \sum_{j,j'=0}^{N_A-1} \sum_{l,l'=0}^{N_B-1} z_j \bar{z}_{j'} \bar{z}_{l'} z_{l'} \bar{z}_{j'} \bar{z}_{j'} \bar{z}_{l'}. \quad (3)$$

We note that $\pi_{AB} = \text{Tr}_A \rho_A^2 = \text{Tr}_B \rho_B^2$ and $1/N_A \leq \text{Tr}_A \rho_A^2 \leq 1$, with the minimum (maximum) value attained for a completely mixed (pure) state ρ_A . Therefore,

$$1 \leq N_{AB} = N_{BA} \leq \min\{N_A, N_B\}. \quad (4)$$

A larger value of N_{AB} corresponds to a more entangled bipartition (A, B) , the maximum value being attainable for a *balanced* bipartition, i.e., when $n_A = \lfloor n/2 \rfloor$ (and $n_B = \lfloor (n+1)/2 \rfloor$), where $\lfloor x \rfloor$ is the integer part of the real x , i.e., the largest integer not exceeding x , and the maximum possible entanglement is $N_{AB} = N_A = 2^{n_A} = \sqrt{N}$ ($= \sqrt{N/2}$) for an even (odd) number of qubits. As anticipated, as a characterization of multipartite entanglement, we will consider the distribution of N_{AB} over all possible balanced bipartitions.

III. MEASURING MULTIPARTITE ENTANGLEMENT: SOME EXAMPLES

Let us illustrate this approach on the simplest nontrivial situation, that of three entangled qubits. If the pure state is fully factorized, say

$$|\psi\rangle = |k\rangle \quad (5)$$

for a given $0 \leq k \leq 7$, then the reduced density matrix ρ_A of every qubit is a pure state, from which

$$p(N_{AB}) = \delta_{N_{AB},1}. \quad (6)$$

there is no entanglement. On the other hand, for a maximally entangled state

$$|\psi\rangle = \frac{1}{\sqrt{2}}(|000_2\rangle + |111_2\rangle), \quad (7)$$

one gets a completely mixed state for every partition, namely $\rho_A = I_2/2$, and thus

$$p(N_{AB}) = \delta_{N_{AB},2}, \quad (8)$$

with maximum average and zero variance: there is maximum multipartite entanglement, fully distributed among the three qubits. The above probability distributions should be compared with an intermediate case like

$$|\psi\rangle = \frac{1}{\sqrt{2}}(|000_2\rangle + |110_2\rangle), \quad (9)$$

where the first couple of qubits are maximally entangled (Bell state) while the third one is completely factorized. In such a situation, one gets $\rho_1 = \rho_2 = I_2/2$ while $\rho_3 = |1\rangle\langle 1|$, from which

$$p(N_{AB}) = \delta_{N_{AB},1}/3 + 2\delta_{N_{AB},2}/3. \quad (10)$$

This simple application discloses the rationale behind the quantity $p(N_{AB})$ as a measure of multipartite entanglement.

When the system becomes larger, the natural extension is toward larger (balanced) bipartitions. We stress that, besides the comment that follows Eq. (4), the use of balanced bipartitions is simply motivated by the fact that, in the thermodynamic limit, the unbalanced ones give a small contribution, from the statistical point of view: this can be easily understood if one considers that for n large and $n_A \ll n$, the binomial coefficients

$$\binom{n}{n/2} \gg \binom{n}{n_A}, \quad (11)$$

so that our characterization of multipartite entanglement will be largely dominated by balanced bipartitions. Notice also that very unbalanced bipartitions of large systems yield negligible average entanglement [12,20]. For all these reasons, if one considers the distribution over all bipartitions, the contribution from the balanced bipartitions will dominate due to Eq. (11). By contrast, if only *unbalanced* bipartitions are considered, the results will be in general very different.

It is interesting to study the features of the characterization of entanglement proposed in Sec. II when applied to particular classes of states. For the Greenberger-Horne-Zeilinger (GHZ) states [13], we find

$$N_{AB}(\text{GHZ}) = 2 \quad (12)$$

for all possible bipartitions (both balanced and unbalanced) and for an arbitrary number of qubits. Clearly, the width of the distribution is 0, i.e., $p(N_{AB}) = \delta_{N_{AB},2}$.

For the W states [14], we obtain

$$N_{AB}(W) = \frac{n^2}{n_A^2 + n_B^2}. \quad (13)$$

This value depends only on the relative size of the two partitions, i.e., also in this case the width of the distribution of bipartite entanglement is 0. Notice that, if n is even, $N_{AB}(W) = 2$ for balanced bipartitions (and in this case a discrimination between W and GHZ states would require the analysis of unbalanced bipartitions). Moreover, in the large- n limit, $N_{AB}(W) \approx 2$ also for n odd.

These results indicate that, for n large, the amount of (multipartite) entanglement is limited both for GHZ and W states. These states essentially share the same amount of entanglement when n is large. They can be distinguished only by considering less relevant (from the statistical point of view) bipartitions. Moreover, for n large, $N_{AB}(W) \neq 1$ for balanced bipartitions. This means that also in the thermodynamic limit the W states retain some entanglement.

IV. TYPICAL STATES

Let us now study the typical form of our characterization of multipartite entanglement $p(N_{AB})$ for a very large class of pure states of the form (1), sampled according to a given statistical law. Several features of these random states are already known in the literature [5,16,17], but we shall focus on those quantities that are relevant for our purpose. We write

$$|\psi\rangle = \sum_{k=0}^{N-1} r_k e^{i\phi_k} |k\rangle, \quad (14)$$

where ϕ_k are independent random variables with expectation

$$E[e^{i\phi_k}] = 0 \quad (15)$$

and $\mathbf{r}=(r_1, \dots, r_N)$ is a random point with a given symmetric distribution $p(\mathbf{r})$ on the hypersphere $S^{N-1}=\{\mathbf{r} \in \mathbb{R}^N | r^2=1\}$. The features of these random states are readily evaluated: one first splits π_{AB} in two parts,

$$\pi_{AB} = X_{AB} + M_{AB}, \quad (16)$$

where

$$X_{AB} = \sum'_{j,j'} \sum'_{l,l'} r_{jl} r_{j'l'} r_{j'l} r_{j'l'}, \quad (17)$$

$$M_{AB} = \sum'_{j,j'} \sum'_{l,l'} r_{jl}^2 r_{j'l}^2 + \sum'_{j,l,l'} r_{jl}^2 r_{j'l}^2 + \sum'_{j,l} r_{jl}^4, \quad (18)$$

with $j, j'=0, \dots, N_A-1$, $l, l'=0, \dots, N_B-1$, and primes banning equal indices in the sums.

We note that the expectation value $E[r_{jl}^2]=O(1/N)$, thus X_{AB} and M_{AB} are sums of at most N^2 terms of order $1/N^2$. By the central limit theorem, for large N , π_{AB} tends to a Gaussian random variable with mean and variance

$$\begin{aligned} \mu_{AB} &= E[\pi_{AB}], \\ \sigma_{AB}^2 &= E[\pi_{AB}^2] - \mu_{AB}^2, \end{aligned} \quad (19)$$

respectively, namely it is distributed as

$$f(\pi_{AB}) = \frac{1}{(2\pi\sigma_{AB}^2)^{1/2}} \exp\left(-\frac{(\pi_{AB} - \mu_{AB})^2}{2\sigma_{AB}^2}\right). \quad (20)$$

From $E[X_{AB}]=0$ and the independence between phases ϕ_k and moduli r_k , we get

$$\mu_{AB} = E[M_{AB}] = N(N_A + N_B - 2)E[r_1^2 r_2^2] + NE[r_1^4] \quad (21)$$

and

$$\sigma_{AB}^2 = E[X_{AB}^2] + E[M_{AB}^2] - \mu_{AB}^2, \quad (22)$$

where

$$E[X_{AB}^2] = 2N(N_A - 1)(N_B - 1)E[r_1^2 r_2^2 r_3^2 r_4^2] \quad (23)$$

and

$$\begin{aligned} E[M_{AB}^2] &= N(N_A + N_B - 2) \\ &\times [(N_A + N_B)(N - 4) - 2(N - 5)]E[r_1^2 r_2^2 r_3^2 r_4^2] \\ &+ 2N(N_A + N_B - 2)(N + 2N_A + 2N_B - 8)E[r_1^2 r_2^2 r_3^4] \\ &+ N(N + 2N_A + 2N_B - 5)E[r_1^4 r_2^4] + 4N(N_A + N_B \\ &- 2)E[r_1^2 r_2^6] + NE[r_1^8], \end{aligned} \quad (24)$$

where we used $E[r_1^\alpha r_2^\beta r_3^\gamma r_4^\delta] = E[r_i^\alpha r_j^\beta r_l^\gamma r_k^\delta]$ with i, j, l, k all distinct. Notice that the above results do not depend on the particular distribution of ϕ_k , as far as the condition (15) is

satisfied [otherwise the analysis is still valid, but Eqs. (21)–(24) become more involved]. Our results particularize for the case of a typical pure state (1), sampled according to the unitarily invariant Haar measure, where each $z_k \in \mathbb{C}$ is chosen from an ensemble that is uniformly distributed over the projective Hilbert space $\sum_k |z_k|^2 = 1$. In such a case, in Eq. (14), $\phi_k \in [0, 2\pi]$ are independent uniformly distributed random variables and $\mathbf{r}=(r_1, \dots, r_N)$ is a random point uniformly distributed on the hypersphere S^{N-1} , with distribution function

$$p(\mathbf{r}) = \frac{2^N}{\pi^{N/2}} \Gamma\left(\frac{N}{2}\right) \delta(1 - r^2), \quad (25)$$

the prefactor being twice the inverse area of the hyperoctant $\{r_i > 0\}$, with $\Gamma(x)$ the Gamma function.

The explicit expressions of Eqs. (21)–(24) can be computed through Eq. (25), recovering the values of mean and variance obtained by different approaches [5,16,17]. However, one can easily estimate them for large N by the following reasoning. For large N , the marginal distributions of the amplitudes r_k become normal,

$$\begin{aligned} p(r_k) &= \frac{2}{\sqrt{\pi}} \frac{\Gamma(N/2)}{\Gamma((N-1)/2)} (1 - r_k^2)^{(N-3)/2} \\ &\sim 2 \sqrt{\frac{N}{2\pi}} \exp\left(-\frac{N}{2} r_k^2\right) \quad (\forall k), \end{aligned} \quad (26)$$

with variance $1/N$. One can convince oneself of the correctness of the above expression just by recalling the asymptotic behavior of the Γ function and expanding $(1 - r_k^2)^{N/2}$. Moreover, it is not difficult to show that the r_k 's become uncorrelated, hence independent. Therefore, the expectation of products factorizes and $E[r_1^{2m}] = (2m-1)!!/N^m$, yielding

$$\mu_{AB} = \frac{N_A + N_B - 1}{N}, \quad \sigma_{AB}^2 = \frac{2}{N^2}. \quad (27)$$

It is important to notice that when $N \gg 1$, we can effectively replace r_k with its mean-square root value, $r_k = 1/\sqrt{N}$, from which Eq. (27) immediately follows. In the simulation plotted in Fig. 1, we used the above substitution. The fact that for Haar distributed states the average (27) is concentrated around a large value was already recognized by other authors [5,16,17].

The quantity of interest is N_{AB} defined in Eq. (2). From Eq. (20), its probability density reads

$$p(N_{AB}) = \frac{1}{N_{AB}^2 (2\pi\sigma_{AB}^2)^{1/2}} \exp\left(-\frac{(N_{AB}^{-1} - \mu_{AB})^2}{2\sigma_{AB}^2}\right). \quad (28)$$

It is interesting to compare the features of the random states with those of other states studied in the literature. Table I displays the average value of N_{AB} (evaluated for $n=5-12$) for GHZ states [13], W states [14], the generic states (14), and one-dimensional cluster states [18] defined as

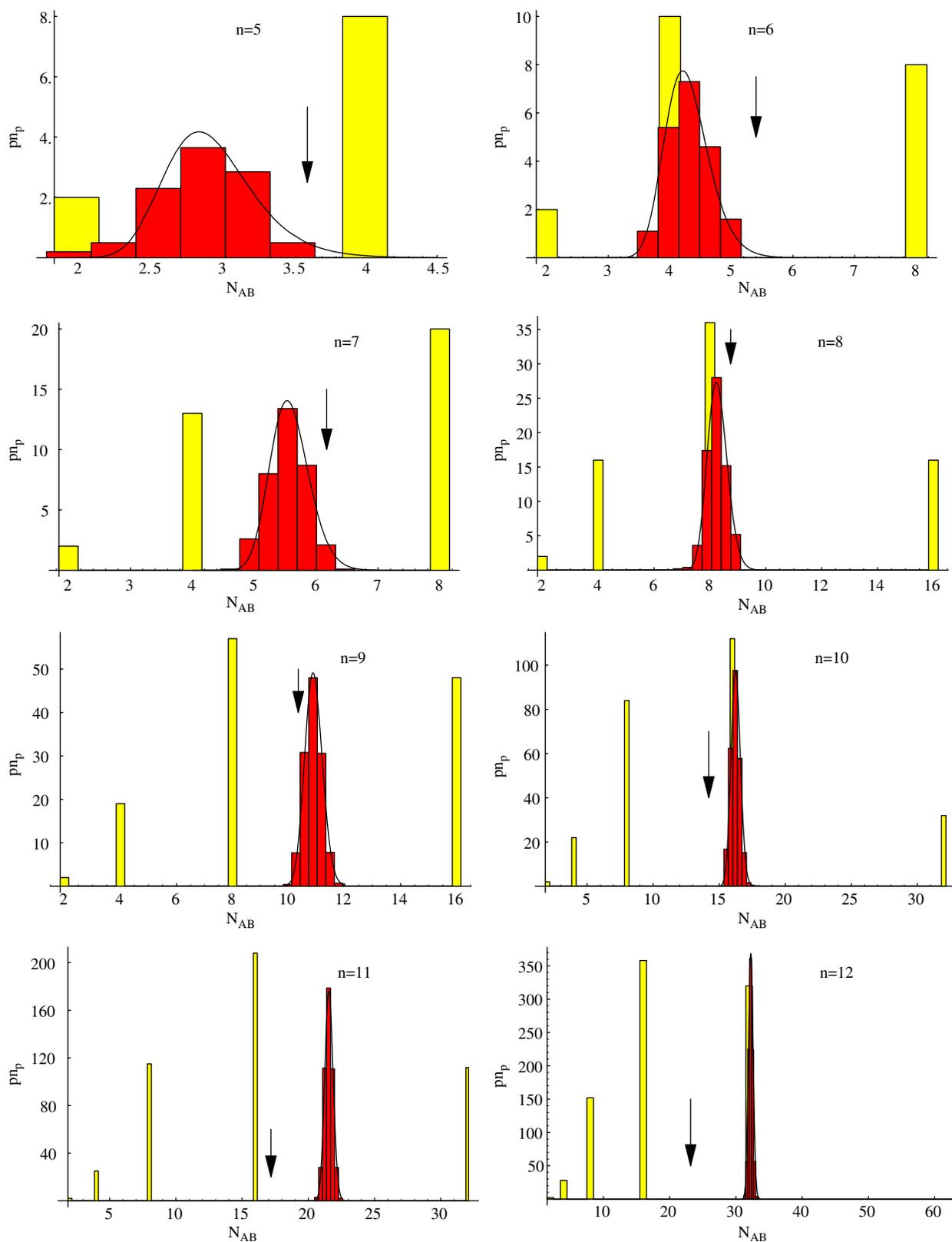


FIG. 1. (Color online) Number of balanced bipartitions vs N_{AB} ; p is the probability density, $n_p = n! / n_A! n_B!$ is the number of bipartitions. The yellow bars represent one-dimensional cluster states [see Eq. (29)], the red ones random states; the solid line is the distribution (27) and (28); the black arrows indicate the average $\langle N_{AB} \rangle_{\text{cluster}}$. For even n ($n=12$ in particular), the distribution of the random state partially hides a peak of the corresponding cluster state distribution, centered at $N_{AB} = 2^{n_A-1} = 2^{\lfloor n/2 \rfloor - 1}$.

TABLE I. Mean bipartite entanglement $E[N_{AB}]$, analytically evaluated according to Eqs. (12), (13), and (27). The values for the cluster state were computed by inserting Eq. (29) into the definitions (2) and (3).

n	GHZ	W	cluster	random
5	2	1.923	3.6	2.909
6	2	2	5.4	4.267
7	2	1.96	6.171	5.565
8	2	2	8.743	8.258
9	2	1.976	10.349	10.894
10	2	2	14.206	16.254
11	2	1.984	17.176	21.558
12	2	2	23.156	32.252

$$|\phi_n\rangle = \frac{1}{\sqrt{2^n}} \otimes_{k=1}^n (|0\rangle_k \sigma_z^{(k+1)} + |1\rangle_k), \quad (29)$$

where σ_z is the third Pauli matrix and the convention $\sigma_z^{(n+1)} = 1$ is applied. While the entanglement of the GHZ and W states is essentially independent of n [see Eqs. (12) and (13)], the situation is drastically different for cluster and random states. In both cases, the average entanglement increases with n ; for $n > 8$, the average entanglement is higher for random states. However, it is now clear that the average $E[N_{AB}]$ yields poor information on multipartite entanglement. For this reason, it is useful to analyze the distribution of bipartite entanglement over all possible balanced bipartitions. The results for the cluster and random states are shown in Fig. 1, for $n=5-12$, where the product of the probability density p times the number of bipartitions $n_p = n! / n_A! n_B!$ is plotted versus N_{AB} . Notice that the distribution function of the random state is always peaked around μ_{AB}^{-1} in Eq. (27) and becomes narrower for larger n , in agreement with σ_{AB}^2 in Eq. (27). Notice also that the cluster state can reach higher values of N_{AB} (the maximum possible value being $2^{\lfloor n/2 \rfloor}$), however the fraction of bipartitions giving this result becomes smaller for higher n . This is immediately understood if one realizes that cluster states are designed for optimized applications and therefore perform better in terms of specific bipartitions. On the other hand, according to the characterization we propose, the random states (14) are characterized by a large value of multipartite entanglement that is roughly independent on the bipartition. The probability density functions (28) are displayed in Fig. 2.

A few additional comments on random states are in order. In the thermodynamic limit,

$$\frac{\sigma_{AB}}{\mu_{AB}} = \frac{\sqrt{2}}{N_A + N_B - 1} = O(1/\sqrt{N}) \quad (30)$$

and the single real number $E[N_{AB}]$ is sufficient to characterize multipartite entanglement (modulo more accurate thermodynamic considerations).

In general, for finite systems, the mean bipartite entanglement $N_{AB} \approx \mu_{AB}^{-1}$ in Eq. (27) is maximum for $N_A = N_B = \sqrt{N}$

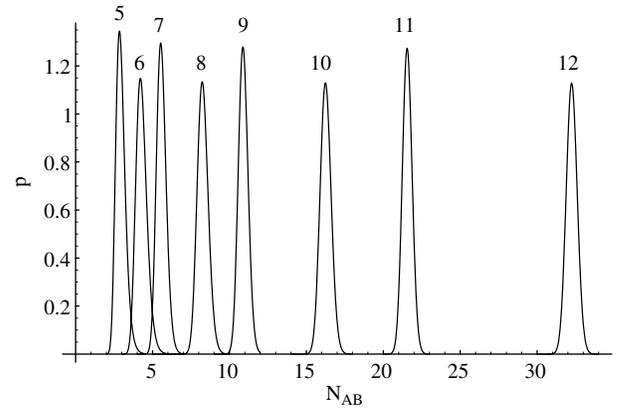


FIG. 2. Probability density functions (28) vs N_{AB} . Each curve is labeled with the corresponding value of n (number of qubits). The standard deviation of the distribution is essentially independent of n .

($N_A = N_B / 2 = \sqrt{N/2}$) for even (odd) n , namely for balanced bipartitions. Notice, however, that, as we already emphasized a number of times in this paper, although we focused on balanced bipartitions for illustrative purposes, the main results are valid when one includes also unbalanced bipartitions, as, by virtue of Eq. (11), the contribution of the balanced bipartition will be exponentially dominant.

Moreover, for large N , any (symmetric) radial distribution $p(r)$ yields the same results (27), the only relevant feature being the curvature in the projective Hilbert space, forced by the normalization $r^2 = 1$ [see, for example, Eq. (25)]. In this sense, the above analysis is of general validity, being independent of the particular choice of the ensemble.

V. COMPARISON WITH SOME MULTIPARTITE ENTANGLEMENT MEASURES

It is interesting to compare our proposed characterization of multipartite entanglement with some other entanglement measures. In general, we will find that this characterization sheds additional light on this issue and helps specify some of the global features of multipartite entanglement in a clear-cut way.

The quantity [15]

$$Q(|\psi\rangle) = 2 \left(1 - \frac{1}{n} \sum_k \text{Tr} \rho_{\{k\}}^2 \right), \quad (31)$$

where $\rho_{\{k\}}$ is the reduced density matrix of qubit k , i.e., ρ_A with $A = \{k\}$. In our language, it corresponds to the mean value of π_{AB} over maximally unbalanced bipartitions, namely

$$Q(|\psi\rangle) = 2(1 - E_{\max \text{ unbal}}[\pi_{AB}]). \quad (32)$$

For W states, this yields $Q(W) \sim 0$ for large n . This should be compared with the value $N_{AB}(W) = 2$ (exact for even n , approximate for odd n), obtained by considering balanced bipartitions of the system. As previously stressed, this means that the W states retain some entanglement even in the thermodynamic limit.

Moreover, at variance with Q , the mean value of N_{AB} can distinguish subglobal entanglement. For instance, the state $|\psi\rangle = (|0\rangle|0\rangle + |1\rangle|1\rangle) \otimes (|0\rangle|0\rangle + |1\rangle|1\rangle)/2$ cannot be distinguished from the GHZ state by using only Q . On the other hand, one gets an average $\langle N_{AB} \rangle = 3$ and a width for the distribution $\sigma = 1.55$. Another interesting point is that the distribution of N_{AB} can distinguish GHZ and cluster states (actually the average is already sufficient, as can be seen from Table I). From these results, one can argue that the probability density function of the participation number N_{AB} not only better specifies the meaning of Q but also yields additional information.

It is also interesting to recall the behavior of the pairwise entanglement (concurrence) and the tangle [3]. The former is defined (for states $\rho_{\{i,j\}}$ of two qubits i and j) as

$$C_{ij} = \max(0, \lambda_1 - \lambda_2 - \lambda_3 - \lambda_4), \quad (33)$$

where λ_k are the square roots of the eigenvalues (in decreasing order) of the matrix $\rho_{\{i,j\}} \sigma_y \otimes \sigma_y \rho_{\{i,j\}}^* \sigma_y \otimes \sigma_y$, and is therefore related to π_{AB} with $A = \{i, j\}$ (highly unbalanced bipartitions when N is large). The tangle is defined as

$$\tau_1^{(i)} = 4 \det \rho_{\{i\}} = 2(1 - \text{Tr} \rho_{\{i\}}^2), \quad (34)$$

where $\rho_{\{i\}}$ is the reduced density matrix for qubit i . Note that $\tau_1^{(i)} = 2(1 - \pi_{AB})$, with $A = \{i\}$, is nothing but the local version of Q in Eq. (31). In particular, one can consider the ratio $R^{(i)} = \tau_2^{(i)} / \tau_1^{(i)}$ [7], where $\tau_2^{(i)} = \sum_{j \neq i} C_{ij}^2$ is the sum of the squared concurrences of qubit i with qubit j . Due to the Coffman-Kundu-Wootters conjecture $\tau_1^{(i)} \geq \tau_2^{(i)}$ [3], one can take $R^{(i)}$ as a witness of multipartite entanglement: if $R^{(i)} < 1$, pairwise entanglement is less relevant than multiqubit correlations. In particular, in order to elucidate their relation with the bipartite entanglement of highly unbalanced bipartitions, it is interesting to apply these measures to typical states. We notice that, in the limit of large n , one has, on the average,

$$E[\tau_1] = Q = 1 - 1/2^{n-1} \sim 1, \\ E[\tau_2] \sim 0. \quad (35)$$

These results are interesting because they show how, in the thermodynamic limit, pairwise entanglement is negligible for typical states. At the same time, Eq. (35) does not yield much information about the very structure of multipartite entanglement: actually one can see that the same result can be ob-

tained for GHZ states (for arbitrary n). In this sense, our characterization in terms of the probability density function corroborates and better specifies the results obtained by studying the behavior of R .

VI. CONCLUSIONS

It is well known that an efficient way to generate states endowed with random features is by a chaotic dynamics [4,5], or at the onset of a quantum phase transition [6]. In particular, the random states (14) describe quite well states with support on chaotic regions of phase space, before dynamical localization has taken place. Interestingly, other ways have been recently proposed [17,19] in order to generate these states, in particular by operating on couples of qubits with random unitaries followed by CNOT gates [19]. The introduction of a probability density function as a measure of multipartite entanglement paves the way for further investigations of this intimate relation between entanglement and randomness. Work is in progress in order to clarify whether the random states can be efficiently used in quantum information processing.

In some sense, the characterization we propose quantifies the robustness of entanglement against all possible partial tracing. Clearly, it is more effective for large numbers of qubits and when relatively few moments are sufficient to specify the distribution. We stress that although we studied the distribution function of the inverse purity (linear entropy) (2), our analysis could have been performed in terms of any other measure of bipartite entanglement, such as the entropy.

Finally, we emphasize again the main motivation behind this work: as the number of subsystems increases, the number of measures (i.e., real numbers) needed to quantify multipartite entanglement grows exponentially. It is therefore not surprising if a satisfactory global characterization of entanglement requires the use of a function.

ACKNOWLEDGMENTS

This work is partly supported by the bilateral Italian-Japanese Projects II04C1AF4E on ‘‘Quantum Information, Computation and Communication’’ of the Italian Ministry of Instruction, University and Research and by the European Community through the Integrated Project EuroSQIP. G.F. acknowledges the support and kind hospitality of the Department of Physics of Waseda University, where part of this work was done.

[1] W. K. Wootters, *Quantum Inf. Comput.* **1**, 27 (2001); C. H. Bennett, D. P. DiVincenzo, J. A. Smolin, and W. K. Wootters, *Phys. Rev. A* **54**, 3824 (1996).
 [2] D. Bruss, *J. Math. Phys.* **43**, 4237 (2002).
 [3] V. Coffman, J. Kundu, and W. K. Wootters, *Phys. Rev. A* **61**, 052306 (2000); A. Wong and N. Christensen, *ibid.* **63**, 044301 (2001); D. A. Meyer and N. R. Wallach, *J. Math. Phys.* **43**, 4273 (2002).

[4] J. N. Bandyopadhyay and A. Lakshminarayan, *Phys. Rev. Lett.* **89**, 060402 (2002); S. Montangero, G. Benenti, and R. Fazio, *ibid.* **91**, 187901 (2003); S. Bettelli and D. L. Shepelyansky, *Phys. Rev. A* **67**, 054303 (2003); A. J. Scott and C. M. Caves, *J. Phys. A* **36**, 9553 (2003); L. F. Santos, G. Rigolin, and C. O. Escobar, *Phys. Rev. A* **69**, 042304 (2004); N. Lambert, C. Emary, and T. Brandes, *Phys. Rev. Lett.* **92**, 073602 (2004); C. Mejia-Monasterio, G. Benenti, G. G. Carlo, and G. Casati,

- Phys. Rev. A **71**, 062324 (2005).
- [5] A. J. Scott and C. M. Caves, J. Phys. A **36**, 9553 (2003).
- [6] A. Osterloh, L. Amico, G. Falci, and R. Fazio, Nature (London) **416**, 609 (2002); T. J. Osborne and M. A. Nielsen, Phys. Rev. A **66**, 032110 (2002); I. Bose and E. Chattopadhyay, *ibid.* **66**, 062320 (2002); G. Vidal, J. I. Latorre, E. Rico, and A. Kitaev, Phys. Rev. Lett. **90**, 227902 (2003); U. Glaser, H. Büttner, and H. Fehske, Phys. Rev. A **68**, 032318 (2003); S. J. Gu, H. Q. Lin, and Y. Q. Li, *ibid.* **68**, 042330 (2003); L. Amico, A. Osterloh, F. Plastina, R. Fazio, and G. M. Palma, *ibid.* **69**, 022304 (2004); V. E. Korepin, Phys. Rev. Lett. **92**, 096402 (2004); J. Vidal, G. Palacios, and R. Mosseri, Phys. Rev. A **69**, 022107 (2004); F. Verstraete, M. Popp, and J. I. Cirac, Phys. Rev. Lett. **92**, 027901 (2004).
- [7] T. Roscilde, P. Verrucchi, A. Fubini, S. Haas, and V. Tognetti, Phys. Rev. Lett. **93**, 167203 (2004); **94**, 147208 (2005).
- [8] G. Parisi, *Statistical Field Theory* (Addison-Wesley, New York, 1988).
- [9] V. I. Man'ko, G. Marmo, E. C. G. Sudarshan, and F. Zaccaria, J. Phys. A **35**, 7137 (2002).
- [10] R. Grobe, K. Rzażewski, and J. H. Eberly, J. Phys. B **27**, L503 (1994); J. H. Eberly, e-print quant-ph/0508019.
- [11] C. Emary, J. Phys. A **37**, 8293 (2004); A. J. Scott, Phys. Rev. A **69**, 052330 (2004).
- [12] V. M. Kendon, K. Życzkowski, and W. J. Munro, Phys. Rev. A **66**, 062310 (2002).
- [13] D. M. Greenberger, M. Horne, and A. Zeilinger, Am. J. Phys. **58**, 1131 (1990).
- [14] W. Dür, G. Vidal, and J. I. Cirac, Phys. Rev. A **62**, 062314 (2000).
- [15] K. Brennen, Sciences (N.Y.) **3**, 619 (2003).
- [16] E. Lubkin, J. Math. Phys. **19**, 1028 (1978); S. Lloyd and H. Pagels, Ann. Phys. (N.Y.) **188**, 186 (1988); K. Życzkowski, and H.-J. Sommers, J. Phys. A **34**, 7111 (2001); Y. Shimoni, D. Shapira, and O. Biham, Phys. Rev. A **69**, 062303 (2004).
- [17] J. Emerson, Y. S. Weinstein, M. Saraceno, S. Lloyd, and DG Cory, Science **302**, 2098 (2003); P. Hayden, D. W. Leung, and A. Winter, Commun. Math. Phys. **265**, 95 (2006).
- [18] H. J. Briegel and R. Raussendorf, Phys. Rev. Lett. **86**, 910 (2001).
- [19] R. Olivera, O. Dahlsten, and M. B. Plenio, e-print quant-ph/0605126.
- [20] However, particularly for small systems, but sometimes also for large systems (see later), whenever a finer resolution is needed, unbalanced bipartitions can also be considered.