Nonergodic Phases in Strongly Disordered Random Regular Graphs

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(Received 16 June 2016; published 6 October 2016)

We combine numerical diagonalization with semianalytical calculations to prove the existence of the intermediate nonergodic but delocalized phase in the Anderson model on disordered hierarchical lattices. We suggest a new generalized population dynamics that is able to detect the violation of ergodicity of the delocalized states within the Abou-Chakra, Anderson, and Thouless recursive scheme. This result is supplemented by statistics of random wave functions extracted from exact diagonalization of the Anderson model on ensemble of disordered random regular graphs (RRG) of N sites with the connectivity K = 2. By extrapolation of the results of both approaches to $N \rightarrow \infty$ we obtain the fractal dimensions $D_1(W)$ and $D_2(W)$ as well as the population dynamics exponent D(W) with the accuracy sufficient to claim that they are nontrivial in the broad interval of disorder strength $W_E < W < W_c$. The thorough analysis of the exact diagonalization results for RRG with $N > 10^5$ reveals a singularity in $D_{1,2}(W)$ dependencies which provides clear evidence for the first order transition between the two delocalized phases on RRG at $W_E \approx 10.0$. We discuss the implications of these results for quantum and classical nonintegrable and many-body systems.

DOI: 10.1103/PhysRevLett.117.156601

Introduction.-The concept of many-body localization (MBL) [1] emerged as an attempt to extend the celebrated ideas of Anderson localization (AL) [2] from one-particle eigenstates formed by a static random potential to the many-body eigenfunctions of macroscopic quantum systems. Later, the MBL in various models (XXZ spin chain subject to a random magnetic field [3,4], array of Josephson junctions [5], etc.) became a subject of intensive theoretical studies. The ideas of MBL appear naturally in discussions of applicability of the conventional Boltzmann-Gibbs statistical mechanics to isolated many-body systems. This description based on the equipartition postulate should not be valid for the localized many-body states. Moreover, in Ref. [5] it was shown that the Boltzmann-Gibbs description of the isolated Josephson arrays most likely remains invalid even in the so-called "bad metal" phase where the eigenstates are extended but not ergodic; e.g., they occupy a vanishing fraction of the Hilbert space.

Because of complexity and diversity of many-body systems it is worthwhile to exploit the MBL-AL analogy to demonstrate existence of the nonergodic extended states first in models for single-particle localization. It is known that such states do exist at the critical point of Anderson transition (AT) [6]. However, in order to be relevant for MBL they have to constitute a distinct *phase*, i.e., to exist in a finite range of parameters, such as the disorder strength. A natural candidate for a model where it can happen is the disordered Bethe lattice (BL), where the number of

resonance sites increases exponentially with distance. This increase can compensate for the exponential smallness of the transition amplitude, thus leading to an extended critical phase. There are reasons to believe [7] that the topology of Hilbert space of a generic many-body system shares (to a leading approximation) the basic properties of BL: (i) the exponential growth of the number of sites $N = K^R$ on the radius of the tree R with the branching number K and (ii) the absence of loops. The latter simplifies the problem of AL as compared to AL in finite dimensions. In the seminal paper [8] Abou-Chakra, Anderson, and Thouless developed an analytical approach to the Anderson model on an infinite BL that allowed them not only to demonstrate the existence of the AL transition but also to evaluate the critical disorder with a pretty good accuracy. More recently, some mathematically rigorous results for AL on BL were obtained [9,10].

The most interesting and the least studied aspect of AL on the BL is the statistics of extended wave functions. Recently it was suggested [11,12] that these statistics may be multifractal, i.e., extended nonergodic. A similar conclusion was reached in early analytical work [13] and a more recent numerical one [14] on "directed BL models." The contradiction with other results on the BL and sparse random matrices [15], where only ergodic states were found below AT, provoked a vigorous discussion [16–20].

Note that the mere formulation of statistics of normalized extended wave functions in a *closed* system requires

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understanding of the thermodynamic limit of a *finite-size* problem. For the BL this poses a major problem: a finite fraction of sites belong to the boundary making the results crucially dependent on the boundary conditions. A known remedy [11,12] is to consider a random regular graph (RRG) [21,22], where each of N sites is connected to a fixed number (K + 1) of other sites. Such a graph has a local tree structure similar to a BL but no boundary. In contrast to a BL, RRG has loops but the length of the smallest statistically relevant ones is macroscopically large $\sim \ln N / \ln K$.

In this Letter we reformulate the approach of Ref. [8] in a way that distinguishes extended nonergodic states from ergodic ones. A new recursive algorithm [similar to population dynamics (PD) [23]] of treatment, the Abou-Chakra-Anderson-Thouless (ACAT) equations [8], enables us to justify semianalytically the existence of the intermediate extended nonergodic phase for a BL with K = 2. This result is relevant for a broad class of systems (e.g., Refs. [13,17]) described by self-consistent ACAT equations, where the loops are absent, or rare, or irrelevant [24]. Our extensive exact diagonalization numerics on the Anderson model on RRG with N up to 128 000 brought up a strong support for such a phase too. Moreover, we discovered evidence for the first order transition between ergodic (EES) and nonergodic states (NEES) within the delocalized phase. Its position corresponds to the condition for the Lyapunov exponent $L(W, E=0) = \frac{1}{2} \ln K$ discussed in Ref. [9]. The results are summarized in Fig. 1.

The model and fractal dimensions D_q .—Below we analyze the properties of the eigenfunctions of the Anderson model described by the Hamiltonian

$$\hat{H} = \sum_{i=1}^{N} \varepsilon_i |i\rangle \langle i| + \sum_{i,j=1}^{N} t_{ij} |i\rangle \langle j|.$$
(1)

Here ε_i are random on-site energies uniformly distributed in the interval [-W/2, +W/2], the connectivity matrix t_{ij} equals to 1 for nearest neighbors, and 0 otherwise.

Let $|a\rangle$ and $\langle i|a\rangle$ be the normalized eigenstates and wave function coefficients of Hamiltonian (1) in the site representation. One can introduce the moments $I_q = \sum_i |\langle i|a\rangle|^{2q}$, which generically scale with the number of the lattice sites $N \gg 1$ as $I_q \propto N^{-\tau(q)}$. For localized states $\tau(q) = 0$, while the ergodicity implies $\tau(q) = q - 1$. Multifractal nonergodic states are characterized by the set of nontrivial fractal dimensions $0 < D_q = \tau(q)/(q-1) < 1$, e.g., $D_1 = \lim_{q \to 1} D_q$ and $D_2 = \tau(2)$. Exact diagonalization of a large RRG (see Fig. 1) suggests that the fractal dimensions experience a jump from $D_q < 1$ for $W > W_E \approx 10.0$ to $D_q = 1$ for $W < W_E$ manifesting the first order ergodic transition.

Generalized recursive algorithm for ACAT equations.— Following Ref. [8] we introduce a single-site Green function, $G_i^{(k)}(\omega) = \langle i | (\omega - \tilde{H}_k)^{-1} | i \rangle$ for a site *i* at a



FIG. 1. Fractal dimensions D_2 and D_1 for K = 2 RRG and the population dynamics exponent D as functions of disorder strength W. The W dependence of D extrapolated to $N \to \infty$ is presented by the "brush-painted" blue line where the width corresponds to the uncertainty of extrapolation. In spite of this uncertainty, D is distinctly different from 0 and 1 in a broad interval of W manifesting the nonergodic (multifractal) nature of extended wave functions. The red solid line with black data points is a "running" fractal dimension $D_2(N, W) = -d\langle \ln I_2 \rangle/d \ln N$ obtained by exact diagonalization at the maximal size N =128 000 of a disordered RRG. The fat red line is a sketch of the fractal dimension $D_2(N \to \infty, W) \equiv D_2(W)$ extrapolated to infinite N. Inset: the jump singularity in the running fractal dimensions $D_1(N = 60\,000, W)$ and $D_2(N = 128\,000, W)$ manifesting the ergodic transition at $W = W_E \approx 10.0$.

generation k of the reduced Hamiltonian H_k , obtained from \hat{H} by disconnecting generations k and k + 1. The random Green functions are characterized by distribution functions $P_k(G)$. Individual $G_i^{(k)}$ obey the ACAT recursion equation [8]

$$G_i^{(k)} = \frac{1}{\omega - \varepsilon_i - \sum_{j(i)} G_j^{(k-1)}(\omega)},$$
(2)

where j(i) are sites at the generation k - 1 connected to site *i*. These equations are ill determined: the polelike singularities in the right-hand sides have to be regularized. This is usually achieved by adding an infinitesimal imaginary part to $\omega \rightarrow \omega + i\eta$. The recursion Eq. (2) might become unstable with respect to this addition. This happens for *W* below the critical disorder of the AL transition W_c and manifests the delocalized phase. For $W > W_c$ the solution $P(G) \propto \delta(\text{Im}G)$ is stable. The two types of behavior occur generically in a broad class of Anderson models [2].

The spectrum of the Hamiltonian on a finite lattice is given by a discrete set of energies E_a , corresponding to states $|a\rangle$. Although the global density of states is a sum of delta functions $\nu(\omega) = \sum_a \delta(\omega - E_a)$ it *always* has a welldefined thermodynamic limit: one introduces an infinitesimal broadening of each delta function η , takes first the limit of the infinite number of sites $N \to \infty$, and afterwards $\eta \to 0$. As a result, $\nu(\omega)$ tends to a smooth function. In contrast, for the local density of states (LDOS) $\nu_i(\omega) = \sum_a |\langle i|a \rangle|^2 \delta(\omega - E_a)$ the result of this procedure is not always a smooth function. Indeed, in the limit $W \to \infty$ the on-site states $|i\rangle$ are exact eigenstates and $\nu_i(\omega) = \delta(\omega - \varepsilon_i)$ even for the infinite system. For finite but large W, satellite δ -like peaks appear. The total number of the peaks is infinite in the thermodynamic limit, but almost all of them have exponentially small weight. Hence, the effective number of peaks remains finite: it increases as W is decreased and becomes infinite at $W = W_c$. At this point LDOS becomes smooth provided that the limit $N \to \infty$ is taken *before* $\eta \to 0$. Note that the opposite order of limits ($\eta \to 0$ before $N \to \infty$) always leads to discrete peaks in LDOS.

At $W < W_c$ LDOS contains an extensive number M of peaks with significant weight: $M \to \infty$ as $N \to \infty$. Generally, one expects $M \propto N^D$ with some $0 < D \le 1$. For $\nu_i(\omega)$ to be smooth, the broadening η should exceed the spacing between the peaks $\delta_M \propto M^{-1} \propto N^{-D}$. Thus, the simultaneous limit $N \to \infty$, $\eta \to 0$, $N^{\gamma}\eta =$ const results in a smooth LDOS *iff* $\gamma < D$. Studying such generalized limits yields information on the scaling of the number of peaks, i.e., on the structure of the eigenfunctions. Wave functions of *ergodic* states are uniformly spread on a lattice, so that $M \propto N$, i.e., D = 1 and the LDOS is smooth for any $\gamma < 1$. We show below that in a broad interval of disorder strengths in the delocalized regime D = D(W) < 1 and equals to the critical value of γ corresponding to the transition between a smooth and a singular LDOS, $D(W) = \gamma_c(W)$.

For $W < W_c$ (delocalized regime) and an infinitesimal $\eta > 0$, ImG increases exponentially with the number of recursion steps *n* in Eq. (2) describing an *infinite* tree:

$$\mathrm{Im}G \propto \eta e^{\Lambda n}.$$
 (3)

For a *finite* RRG of size N, $n < \ln N / \ln K$ [21]. For larger n the loops terminate the exponential growth of a typical ImG limiting it by $\text{Im}G \propto \eta N^{\Lambda/\ln K}$. Thus, for $\nu_i(\omega) \sim N^{-D} \sum_a \eta / [(\omega - E_a)^2 + \eta^2] \approx \int dE_a \eta / [E_a^2 + \eta^2]$ to be smooth (and Im $G \sim 1$ independent of η) η should scale as $\eta \propto N^{-\Lambda/\ln K}$, i.e.,

$$D(W) = \Lambda(W) / \ln K.$$
(4)

Ideally, one would deal with infinitely small $\eta \rightarrow 0$ in order to determine the exponent Λ . However, the limited precision of any numerical computation makes it impossible in practice: for any realistic initial Im $G \neq 0$, the value of ImG becomes significant after few recursions. To avoid this problem we included an additional step to the recursion Eq. (2):

$$\mathrm{Im}G_i^{(k)} \to e^{-\Lambda_k}\mathrm{Im}G_i^{(k)},\tag{5}$$

so we keep the *typical* imaginary part fixed and k independent: $\exp\langle \ln \operatorname{Im} G_i^{(k)} \rangle_k = \delta$ (where $\langle \cdots \rangle_k$ denotes averaging over all sites *i* in the kth generation). As soon as the stationary distribution of *G* is reached in this recursive procedure, $\Lambda_k \to \Lambda$.

To realize this algorithm we adopted a modified *population dynamics* (PD) method [23]. In each step we used the set of N_p Green functions $G_i^{(k)}$ (population) obtained at the previous step and new on-site energies ε_i to generate N_p new Green functions $G_i^{(k+1)}$ according to Eq. (2), in which each site is connected to K randomly chosen sites of the previous population set.

In order to obtain D(W) one needs to take the limits $N_p \to \infty$, $\delta \to 0$ of $D(N_p, \delta, W)$. The convergence turns out to be slow (logarithmic), resulting in a considerable uncertainty in D(W). Luckily, $D(N_p, \delta, W)$ depends only on the combined variable $X = -1/\ln(N_p^{-1} + a\delta^b)$, with $a, b \sim 1$, rather than on $\ln N_p$ and $\ln \delta$ separately. Extrapolation of D(W, X) to X = 0 yields D(W) shown in Fig. 2. The lower inset of Fig. 2 shows the collapse of the data for several N and δ from the intervals $10^3 < N < 10^8$ and $10^{-3} < \delta < 10^{-17}$. Since $b \approx 0.5$, one needs exceptionally small δ to reach small X. This required computation with higher than usual precision.

Note that the exponent Λ is a property of an infinite BL, $N = \infty$. Therefore, Λ is free of the finite-size effects which



FIG. 2. The population dynamic exponent D(W) (blue points with gray error bars) extrapolated to $N = \infty$ and $\delta = 0$ for K = 2. The condition $D(W_c) = 0$ yields $W_c = 18.4^{+0.4}_{-0.2}$. In a broad interval of $W < W_c$ we obtained D(W) distinctly different from the ergodic limit D(W) = 1. Lower inset: The collapse of data for a fixed W and different N_p , δ to a function D(W, X) of $X = -1/\ln(N_p^{-1} + a\delta^b)$. Extrapolation to X = 0 gives the population dynamic exponent D(W). The delocalized phase corresponds to $1 \ge D(W) > 0$, whereas in the localized phase D(W) < 0. Upper inset: the finite-size critical disorder $W_c(X)$ defined as $D[W_c(X), X] = 0$ and its extrapolation to X = 0 by the power-law fit $W_c(X) = W_c - aX^{1/\nu}$ with $W_c = 18.4$, $\nu = 0.56$ (blue), and $W_c = 19.0$, $\nu = 0.7$ (red). Without extrapolation the value of W_c at maximal population size $N_p^* \sim 10^8$ is $W_c(N_p^*) \approx 17.5$.



FIG. 3. Left panel: $D_2(N, W)$ deep in the delocalized phase. The curves tend to converge to two different values of D_2 for $W = W_E + 0$ and $W_E - 0$, where $W_E \approx 10.0$. Right panel: Formation of a jump in $D_2(W)$.

dominate the moments $I_q(N)$ at $N < N_c$, where the correlation volume $N_c \sim \exp[1/\Lambda(W)]$ diverges at $W \to W_c$. The uncertainty of extrapolation of Λ to $N_p \to \infty$ and $\delta \to 0$ turns out to be small enough not to raise doubts that 0 < D < 1 at least in the interval 10 < W < 18 for K = 2. Additional support of existence of the phase with 0 < D < 1 comes from the analytical solution to Eq. (2) in the large-K limit [24]. It turns out that in this limit D(W) = 0 and D(W) = 1 correspond to the special values of the Lyapunov exponent $L = \ln K$ and $L = \frac{1}{2} \ln K$ discussed in Ref. [9].

Exact diagonalization on RRG.—While the ACAT approach is commonly believed to describe well the localized phase of RRG, its applicability in the delocalized regime requires further investigation. We performed a direct study of the Anderson model on RRG by exact diagonalization at the system sizes N up to 128 000 in the range of disorder strength 7.5 < W < 20 [25]. The main focus was on calculating the inverse participation ratio $I_2 = \sum_i |\langle i | a \rangle|^4$ and the Shannon entropy $S = -\sum_i |\langle i | a \rangle|^2 \ln(|\langle i | a \rangle|^2)$ for the eigenstates $|a\rangle$ with



FIG. 4. $D_2(N, W)$ close to the localization transition at $W = W_c$. The N dependence shows minima (red spot) for $W < W_c$ at $N_{\min} \to \infty$ as $W \to W_c$ [19]. Inset: $D_2(N_{\min}, W)$ as a function of W. Extrapolation by a second-order polynomial gives $W_c = 18.1 \pm 0.5$.

energies E_a near the band center. The expected asymptotic behavior of the typical averages at $N \to \infty$ is [12]

$$\langle \ln I_2 \rangle = -D_2 \ln N + c_2, \qquad \langle \ln S \rangle = D_1 \ln N + c_1, \quad (6)$$

where $\langle \cdots \rangle$ are the averages both over the ensemble of RRG with fixed connectivity K = 2 and over random on-site energies ε_i , $D_{1,2}$ are the multifractal dimensions, and $c_{1,2} \sim 1$. The derivatives $D_2(N, W) = -d\langle \ln I_2 \rangle/d \ln N$ and $D_1(N, W) = d\langle \ln S \rangle/d \ln N$ should saturate at D_2 and D_1 , respectively, in the limit $N \to \infty$.

We present the results for $D_2(N, W)$ deep in the delocalized phase (Fig. 3) and close to the localization transition (Fig. 4). Note two important features on these plots: (i) an abrupt change of behavior for W close to 10 and (ii) a minimum in the N dependence of $D_2(N, W)$ (recently reported in Ref. [19]) in the vicinity of AL transition: as $W \rightarrow W_c - 0$, $D_2(N_{\min}, W)$ at the minimum and $1/\ln N_{\min}$ vanish. Extrapolation of $D_2(N_{\min}, W)$ leads to $W_c = 18.1 \pm 0.5$ (see inset to Fig. 4) in agreement with PD results, Fig. 2.

A striking result of the exact diagonalization is the existence of a jump in both $D_2(N, W)$ and $D_1(N, W)$ shown in Fig. 1. A feature, which is almost invisible at small N, evolves to a more and more abrupt jump as N increases above 60 000 (see Fig. 3, right panel). Extrapolation of $D_2(N, W)$ to $N \to \infty$ for W < 10.0 gives $D_2 = D_2(N \to \infty, W) = 1$, whereas $D_2(W = 10.0) = 0.86 \pm 0.02$ [25]. We conclude that on RRG at $W = W_E \approx 10.0$ [26] there is a first-order transition from the non-ergodic delocalized phase at $W > W_E$ to the ergodic one at $W < W_E$.

Conclusion.—The existence of the nonergodic phase of the BL Anderson model together with the similarity of this model with generic many-body ones gives basis for farreaching speculations. The point is that in contrast to the conventional Anderson localization, which is the property of any wave dynamics, the MBL is a genuine quantum phenomenon. Indeed, in the classical limit, a weakly perturbed integrable system with d > 2 degrees of freedom

always demonstrates some diffusion in the phase space known as Arnold diffusion [27]. Although the celebrated Kolmogorov Arnold Moser theorem [28] guarantees the survival of the vast majority of the invariant tori, the chaotic part of the phase space is connected (unless d = 2), thus allowing the diffusion for arbitrary weak perturbation. Therefore, one should not expect MBL in the classical limit. On the other hand, the glassy states of matter without doubt exist for any \hbar including $\hbar = 0$ and are obviously not ergodic. It is safe to assume that the extended nonergodic phase of the MBL models is not qualitatively different from a classical glassy state [29]. Therefore, our arguments in favor of the existence of the delocalized nonergodic phase of the BL Anderson model and the true phase transition between the ergodic and nonergodic states can be considered as arguments in favor of glassy states being distinct states of matter and their transition to fluids being a true phase transition.

We appreciate hospitality at KITP of University of California at Santa Barbara under the NSF Grant No. NSF PHY11-25915, at KITPC in Beijing (V.E.K.) and at the Center for Theoretical Physics of Complex Systems (Daejeon, S. Korea) where the research has been carried out. E. C. thanks partial financial support by the Murcia Regional Agency of Science and Technology (Project No. 19907/GERM/15). The research of L. B. I. was partially supported by the Russian Science Foundation Grant No. 14-42-00044. We are grateful to G. Biroli, E. Bogomolny, J. T. Chalker, M. V. Feigelman, M. Foster, I. M. Khaymovich, G. Parisi, V. Ros, A. Scardicchio, M. A. Skvortsov, V. N. Smelyanskiy, K. S. Tikhonov, and S. Warzel for stimulating discussions..

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