Anderson localization on the Bethe lattice: non-ergodicity of extended states

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Statistical analysis of the eigenstates of the Anderson tight-binding model with on-site disorder on regular random graphs strongly suggests that the extended states are multi-fractal at any finite disorder. The spectrum of fractal dimensions \( f(\alpha) \) remains positive for \( \alpha \) noticeably far from 1 even when the disorder is several times weaker than the one which leads to the Anderson localization, i.e. the ergodicity can be reached only in the absence of disorder. The one-particle multifractality on the Bethe lattice signals on a possible inapplicability of the equipartition law to a generic many-body quantum system as long as it remains isolated.

\( H = - \sum_{<ij>} (c_i^+ c_j + \text{h.c.}) + \sum \varepsilon_i c_i^+ c_i, \quad (1) \)

where \( \varepsilon_i \in [-W/2, W/2] \). A normalized wave function \( \psi(i) (i = 1, ..., N) \) can be characterized by the moments \( I_q = \sum_i |\psi(i)|^{2q} \propto N^{-\tau(q)} \) \( (I_1 = 1 \text{ for the normalization}) \). Ergodicity is violated when \( \tau(q) \) differs from \( (q-1) \). If the ratio \( D_q = \tau(q)/(q-1) \) depends on \( q \), the wave function \( \psi(i) \) is called \textit{multi-fractal}. It is customary to characterize \( \psi(i) \) by the spectrum of fractal dimensions \( \text{SFD} f(\alpha) \) defined as the Legendre transform of \( \tau(q) \).

\[ f(\alpha) = \frac{1}{q-1} \frac{\partial \tau(q)}{\partial q}, \quad \tau(q) = \frac{\partial f(\alpha)}{\partial \alpha} \]

\( \text{SFD} f(\alpha) \) can be computed from the numerical diagonalization of the Hamiltonian \( \text{Eq. (1)} \) on the regular graph with finite number \( \text{of lattice sites} \) when the disorder is several times weaker than the one which leads to the Anderson localization, i.e. the critical \( W = 17.5 \) states.

\[ \psi(i) = \psi_{\text{en}}(i) \phi_{\text{osc}}(i), \quad (2) \]

Below we describe how to separate statistics of \( \psi_{\text{en}}(i) \)
and demonstrate that at all strengths $W$ of the on-site disorder the distribution function (DF) of $x = N|\psi_{en}(i)|^2$ is consistent with the multifractal ansatz [14–16], i.e. can be expressed through SFD $f(\alpha)$ as:

$$P(x) = \frac{A}{x} N^\alpha x^{-1}, \quad \alpha(x) = 1 - \ln x / \ln N,$$

where $A \sim O(N^0)$ is the normalization constant. We found that with decreasing disorder $f(\alpha)$ evolves from almost triangular shape in the insulator to a steep parabolic shape concentrated near $\alpha = 1$ (see Fig.1).

Fractal behaviour of quantum dynamics on the disordered BL have been discussed previously. Transmission from the root to the given surface point of the Cayley tree turned out to be multifractal [12]. This surface multifractality of the extended states does not necessarily mean that bulk of the BL is multifractal: it is known, e.g., that in 2D the bulk multifractality is much weaker than the surface one [12]. Our analysis of the results of exact numerical diagonalization of the Hamiltonian (1) on the RRG demonstrated that the extended wave functions are multifractal in the bulk of the BL. This statement is much stronger and more general than the analogous statement for the transmissions of an artificial model of a tree with surface sites connected to ideal leads [12].

Authors of Ref. [11] analyzed numerically the statistics of the spectra and by population dynamics the distribution of the Green functions of the model (1) and conjectured a transition between extended-ergodic and extended non-ergodic phases in addition to the Anderson transition. Contrarily, we do not see any evidence of the second transition and believe that the entire extended phase is non-ergodic.

Numerics on the BL: rectification and extrapolation.– With the exception of the deeply localized states discussed below, analytical methods to address the problem of wave function statistics on BL are yet to be developed. One can try to access these statistics numerically by diagonalization of the Anderson model Eq. (1) on a RRG. The first challenge along this route is the necessity to extract the statistics of the smooth envelope $\psi_{en}(i)$ of the wave function $\psi(i)$, Eq. (2). The short-range oscillations of $\phi_{osc}(i)$ have nothing to do with AL but dominate the numerically obtained DF of $|\psi(i)|^2$ at small $|\psi(i)|^2$. This tail of the DF thus reflects the density of the nodes of $\phi_{osc}(i)$ rather than the probability for $|\psi_{en}(i)|^2$ to be small.

Since the scales of spatial dependencies of $\psi_{en}(i)$ and $\phi_{osc}(i)$ are so different, it is natural to assume that these two functions are statistically independent and $|\phi_{osc}|^2$ is characterized by the Porter-Thomas DF of the Gaussian Orthogonal Ensemble (GOE) [17] $P_{GOE}(|\phi_{osc}|^2) = (\sqrt{2\pi} |\phi_{osc}|)^{-1} \exp (-|\phi_{osc}|^2 / 2)$. Under these assumptions $\bar{\alpha} = 1 - \ln(N|x|^2) / \ln N$, which is evaluated numerically, is a sum of two statistically independent random variables: $\bar{\alpha} = \alpha(x) + \alpha_{osc}$ where $\alpha(x)$ is given by Eq. (3) and $\alpha_{osc} = -\ln|\phi_{osc}|^2 / \ln N$. The DF of $\bar{\alpha}$ is thus a convolution of the DF $p(\alpha) = P(x(\alpha))x(\alpha)\ln N$ of $\alpha(x)$ with $P_{GOE}(\alpha_{osc}) = \ln N (2\pi N^{-\alpha_{osc}})^{-1} \exp (-N^{-\alpha_{osc}} / 2)$, i.e. $P(x)$ determined by (3) can be obtained (“rectified”) from the DF of $\bar{\alpha}$ and by the Laplace transform method (see Supplemental material for details).

Another, even bigger challenge is that Eq. (3) is expected to hold only in the limit of a sufficiently large graph, i.e. $f(\alpha)$ can be determined only from the limit of $f(\alpha,N) = f(\alpha) + \delta_N f(\alpha)$ at $N \to \infty$, where:

$$f(\alpha,N) = 2 - \alpha + \ln P(x) / \ln N, \quad x = N^{1-\alpha}.$$

The biggest number of sites accessible to us was $N = 32,000$. Increasing further $N$ does not buy much, as the computation time increases as $N^3$ while the finite-size correction $\delta_N f(\alpha)$ is small only as $1 / \ln N$. However, we found a bright side of the slowness of the convergence: in the broad interval of $\alpha$ the correction $\delta_N f(\alpha)$ turns out to be linear in $1 / \ln N$ to a surprisingly high accuracy. This allowed us to make a reliable linear extrapolation in $1 / \ln N$ well beyond the numerical data.

Fixed points in the N-dependence of $f(\alpha,N)$– We numerically diagonalized the Hamiltonian Eq. (1) for the regular graph with connectivity $K + 1 = 3$ (which does not contain boundary sites) and evaluated $f(\alpha,N)$ for $N = 32, 16, 8, 4, 2 \times 10^3$ at several disorder strengths $W$, both above the Anderson transition at $W_c = 17.5$ and below it, down to $W = 5$. Plots of $f(\alpha)$ for $W = 10$ and $W = 5$ are shown in Fig.2 and Fig.3.
An important observation is the existence of two fixed points \(\alpha_+\) and \(\alpha_-\) (shown by arrows): \(f(\alpha_+, N)\) and \(f(\alpha_-, N)\) are essentially \(N\)-independent. This rules out the possibility of \(f(\alpha)\) evolving with further increase of \(N\) into a sharp parabola at \(\alpha = 1\) (dashed line in Fig. 1). In addition to that we verified to a high degree of accuracy that \(\delta N f(\alpha)\) is linear in \(1/\ln N\), at least for \(\alpha_- < \alpha < \alpha_+\). The insert of Fig. 2 demonstrates that the values of \(f(\alpha = 1, N)\) being plotted as a function of \(1/\ln N\) for \(N = 32; 16; 8; 4; 2 \times 10^3\) (red points) form an almost ideal straight line which can be prolonged down to \(1/\ln N = 0\). This is how we obtained the extrapolated SFD \(f_{\text{ext}}(\alpha)\). It was already mentioned that the maximal value of \(f(\alpha)\) in Eq. (3) should be 1. It is not the case for \(f(\alpha, N)\) as one can see from Figs. 2, 3. After extrapolation, however, the maxima of \(f_{\text{ext}}(\alpha)\) turn out to be much closer to 1: \(f_{\text{ext}}^{\text{max}} = 0.99, 1.03, 1.01, 1.00\) for \(W = 5, 10, 17.5, 22.5\), respectively. We just conclude that the extrapolation passed an important and non-trivial test for consistency.

**Verification of the symmetry of \(f(\alpha)\)**—Another important observation providing us with the additional confidence in the validity of the extrapolation is the symmetry of SFD and DF:

\[
\begin{align*}
 f(1 + \alpha) &= f(1 - \alpha) + \alpha, \\
P(x) &= x^{-3} P(x^{-1}).
\end{align*}
\]

One can use Eq. (3) to check that Eqs. (5), (6) follow from each other. Log-normal distribution found for weakly multifractal states in 2D disordered systems \(\text{[18]}\) is one of the examples of this symmetry. A relation similar to Eq. (6) was proven for the DF of the local density of states \(\rho(\epsilon, \xi)\) in a one-dimensional chain \(\text{[19]}\) and a variety of systems (e.g. short and long disordered wires, 2D and 3D disordered systems) described by the nonlinear sigma-model \(\text{[20, 21]}\). The precise conditions of validity of Eq. (6) for the individual eigenfunctions are yet to be formulated. It does not hold for the localized eigenstates, while for weakly multifractal extended states in 2D systems it is valid \(\text{[18]}\). A vast numerical evidence of the validity of Eq. (6) for the multifractal states at the Anderson transition point in 2D and 3D systems was reported \(\text{[10]}\). In the insert of Fig. 3 we present the separate plots of \(f_{\text{ext}}(1 + \alpha)\) and \(f_{\text{ext}}(1 - \alpha) + \alpha\) for the weakest disorder we studied, \(W = 5\) (deep in the region of extended states, the fractal dimensions \(D_1\) and \(D_2\) are very close to 1). One can see that the two curves are indistinguishable in the interval \(-0.4 < \alpha < 0.4\), while \(f(1 + \alpha, N = 16 \times 10^3)\) and \(f(1 - \alpha, N = 16 \times 10^3) + \alpha\) differ noticeably.

In the localized regime \(W > W_c = 17.5\) and at the critical point \(W = W_c\) the shape of the SFD is approximately triangular (see Fig. 1):

\[
f_{\text{ext}}(\alpha) = k \alpha \theta(1 - k \alpha), \quad \theta(z) = \begin{cases} 
1 & \text{if } z > 0 \\
0 & \text{if } z < 0.
\end{cases}
\]

The slope \(k\) depends on the disorder \(k = k(W)\) with \(k(W_c)\) is very close to 1/2. Note, that the only linear \(f(\alpha)\) allowed by Eq. (5) is the one with \(k = 1/2\). Thus one concludes that the critical states with not very small amplitude of wave function (not very large \(\alpha\)) for \(W = W_c\) obey Eq. (6). At the same time, the most abundant critical states around the maximum value of \(f(\alpha)\) reached at \(\alpha = \alpha_0 \approx 2.3\) clearly violate the symmetry Eq. (6). Indeed \(f(1 - \alpha)\) is defined only for \(\alpha < 1\) and thus according to Eq. (6) \(f(1 + \alpha)\) should make no sense for \(\alpha > 1\). In the localized regime we found that \(\alpha_0\) increases with disorder, while \(k(W) \approx \alpha_0^{-1}\) decreases below 1/2. This is in a clear contradiction with Eq. (6).

**Power-law DF for strongly localized states on BL**—Our numeric-based conclusions on the regime of strong localization \(W \gg W_c\) fully agree with the analytical results which follow from the locator expansion \(\text{[6]}\) (also in Supplemental material). In the shortest path approximation \(\text{[22, 23]}\) one obtains: \(\psi^{(0)}(i) = \prod_{j \in p_{0,i}} (\epsilon_0 - \epsilon_j)^{-1}\), where \(\psi^{(0)}(i)\) is the eigenfunction of the Hamiltonian Eq. (1) on BL which at \(W = \infty\) is located on the site 0, and \(p_{0,i}\) is the shortest path connecting sites 0 and \(i\). One can show that within this approximation the DF \(P(x)\) can be represented as:

\[
P(x) = I(1,0) - N^{1-\kappa} I(1-m,\kappa).
\]

Here \(m = \ln N/\ln K\) is the BL radius, \(\kappa = \ln(W/2)/\ln K\), and

\[
I(p,q) = \frac{1}{\sqrt{N} s \pi^2} \int_{B} ds \, s^p (x N^{2q-1} s)^{1/2},
\]

The contour \(B \in (r-i\infty,r+i\infty)\) is parallel to the imaginary axis and crosses the real axis at \(s_0 < r < s_+\), where \(s_\pm\) are the larger and the smaller of the only two
tracting statistics of the smooth envelopes. However, even for $k > \kappa_c = \ln(\kappa_c \ln K) + \ln(eK)/\ln K$, which can be rewritten as

$$W \geq W_c = 2eK \ln(W_c/2) \approx 2eK \ln(eK).$$

(11)

Solution of Eq. (11) is nothing but the critical disorder of Ref. [2] (see Eq. (8)) there, see also the “upper limit critical condition” of Refs. [6, 7]). Since $x = N/|\psi(i)|^2 < N$ in the first term of R.H.S. of Eq. (8), one can deform the contour of integration in Eq. (10) to encircle the pole $s = s_\pm$. For the second term, $N^{1/2} I(1-m, \kappa)$, this can be done only provided that $x(N/K)^{2\kappa} < N$. Under this condition the two terms cancel each other and $P(x) \equiv 0$. In the opposite case $x > N(N/K)^{2\kappa}$, the integral $I(1-m, \kappa)$ in Eq. (8) is determined by the poles $s = s_-$ and $s = 0$. Within the region of validity of the shortest path approximation, $x \sim \frac{N}{W} \ln N \ll 1$ the two contributions cancel each other, i.e. $P(x) = I(1,0)$. Finally within the shortest path approximation, for $W \gg W_c$ we have:

$$P(x) = \frac{\theta(x - x_{\min})}{N^2} \left( \frac{x}{N} \right)^{(s_+ - 3)/2},$$

(12)

where

$$x_{\min} = N^{1-2\kappa} K^{2\kappa} \Rightarrow \alpha_{\max} = 2\kappa (1 - m^{-1}) \approx 2\kappa.$$

(13)

Using the definition of $\alpha(x)$ Eq. (3) one obtains from the power-law DF Eq. (12) the linear SFD $f(\alpha)$, Eq. (7) with:

$$k(W) = \frac{1}{2} (1 - s_+),$$

(14)

truncated at $\alpha > \alpha_{\max}$. Note that for $\kappa \gg 1$ (i.e. for $W \gg 2K$), Eq. (10) yields $s_+ \approx 1 - \kappa^{-1}$, so that the condition $k(W) \alpha_{\max} = 1$ encoded in Eq. (7) is satisfied at large disorder ($\kappa \gg 1$) and large system size ($m \gg 1$).

Conclusion. We developed an effective method for extracting statistics of the smooth envelopes $\psi_{\alpha}(i)$ of random eigenfunctions of the Anderson model (1) on RRG and to extrapolate these results from RRG to the BL with an infinite number of sites. Our results strongly suggests that DF of $|\psi_{\alpha}(i)|^2$ in the limit $N \rightarrow \infty$ indeed converges to the form Eq. (3) regardless to the strength of disorder. As long as the states are localized the spectrum of fractal dimensions turns out to be triangular: $f(0) = 0$ and the linear $f(\alpha)$ is well described by Eq. (7). The slope $k$ increases as the disorder $W$ decreases and reaches its maximal possible value $k_e = 1/2$ at the Anderson transition point $W = W_c = 17.5$. With further decrease of the disorder below the critical one $f(\alpha)$ gradually crosses over to the parabolic shape typical for weak multi-fractality: the two roots of $f(\alpha)$ become positive $0 < \alpha_{\min} < 1 < \alpha_{\max}$ ($\alpha_{\min} \rightarrow 0^+$ as $W \rightarrow W_c^+$). However even for $W$ several times smaller than $W_c$ both $\alpha_{\min}$ and $\alpha_{\max}$ turn out to be quite far from 1, while the ergodicity would imply that $\alpha_{\min}, \alpha_{\max} \rightarrow 1$. We conclude that the non-ergodicity and multi-fractality persist in the entire region of delocalized states $0 < W < W_c$, and the only critical point is the point of the Anderson localization transition.

It goes without saying that only RRG with not too big $N$ are accessible for the numerical analysis and one has to deal with $f(\alpha, N)$ determined by Eq. (4). However the existing data allow us to exclude the possibility that the observed non-ergodicity is a finite size effect. Our confidence is based, among other things, on the existence of two fixed points: $f(\alpha, N)$ is $N$-independent at $\alpha = \alpha_-(W) < 1$ and $\alpha = \alpha_+(W) > 1$. The extrapolation of $f(\alpha, N)$ to $N \rightarrow \infty$ in the interval $\alpha_--\leq \alpha \leq \alpha_+$ turned out to be tremendously reliable. It is thus hard to imagine how $f(\alpha, N)$ could evolve to the ergodic limit with further increase of $N$.

There is an apparent contradiction between our results and the claims [24, 25] that all of the extended eigenvectors of sparse random matrices (which manifest some similarities with the eigenstates on BL [24]) are ergodic: $I_2 = C(W)/N$ for $N \gg C(W)$ playing the role of the critical volume, $C(W) \rightarrow \infty$ when $W \rightarrow W_c$. We do not believe that this discrepancy can be attributed to a small finite size $N < C(W)$ of our ensemble of RRGs. Indeed, in Refs. [24, 25], as well as in all the critical phenomena with one single critical length, the intensive quantity like $f(\alpha, N, W)$ is either independent of $W$ ($f(\alpha, N, W) = f(\alpha, W, c)$ in the critical region $N \ll C(W)$) or it is independent of $N$ ($f(\alpha, N) = f(\alpha, W)$ for $N \gg C(W)$). The multi-fractality established in this paper is characterized by $f(\alpha, N)$ which changes rapidly with $W$ (in particular $\alpha_{\min}$ moves from 0 at $W = W_c = 17.5$ to about 0.7 at $W = 5$) and thus the system should be outside the critical region for our system sizes and $W = 10$ or $W = 5$. At the same time $f(\alpha, N)$ does not collapse to the ergodic limit $\alpha_{\min} \approx \alpha_{\max} = 1$ at any finite $W$ that we studied. Discussion of the relevance of the approach of Ref. [24, 25] to the properties of the quantum diffusion on BL goes beyond the scope of this paper.

The absence of ergodicity for the dynamics of the one-particle Anderson model on the BL, in light of the possible connection with the many-body dynamics, suggests serious implications on the statistical mechanics of isolated systems with a large number of degrees of freedom. If the same phenomenon occurs in the many-body case, the equipartition law is likely not to be valid exactly even for strongly non-integrable systems.

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