

## Scaling Theory of Few-Body Delocalization

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## Many-Body Localization: Phenomenology

Example: local charge density after a quantum quench in cold atom chain

$$
\hat{H}=-J \sum_{\langle i j\rangle \sigma} c_{i \sigma}^{\dagger} c_{j \sigma}+\sum_{i \sigma} \Delta_{i} n_{i \sigma}+U \sum_{i} n_{i \uparrow} n_{j \downarrow}
$$


initial state


Ref: Schreiber, Science 2015

## Anderson Localization

Disorder without interactions:

$$
H=-t \sum_{\substack{\langle i j\rangle \\ \text { hopping }}}\left(c_{i}^{\dagger} c_{j}+h . c .\right)+\sum_{i} \mu_{i} n_{i}
$$



No disorder: ballistic spread of wavefunction

Disorder: localization

> position

Disorder potential

In d=1 or $\mathrm{d}=2$ dimensions all wavefunctions are exponentially localized: $|\Psi(r)| \sim e^{-r / \xi}$

Occupation number of each wavefunction is a Local Integral of Motion: $H=\epsilon_{i} \tilde{n}_{i}$

## Many-Body Integrals of Motion

Add interactions to the Anderson insulator, still get LIOMs?

Yes! Dress the Anderson LIOMs:

$$
\tilde{n}_{i}=U n_{i} U^{\dagger}=n_{i}+\alpha_{i ; j k l m} c_{j}^{\dagger} c_{k}^{\dagger} c_{l} c_{m}+\ldots
$$

Serbyn et al, PRL 2013; Huse et al, PRB 2014

In d=1 with short-range interactions and strong disorder: MBL

localized

Localized $\operatorname{Tr}\left[\tilde{n}_{i} n_{j}\right] \sim e^{-\left|r_{i}-r_{j}\right| / \xi}$
Short-range interactions $V_{i j} \sim e^{-\left|r_{i}-r_{j}\right| / \xi_{V}}$

## Many-Body Delocalization?



## Few-Body States

Short-range interactions $\quad H=\sum_{i=1}^{N} \epsilon_{i} n_{i}+t \sum_{i=1}^{N-1}\left(c_{i}^{\dagger} c_{i+1}+c_{i+1}^{\dagger} c_{i}\right)+V \sum_{i=1}^{N-1} n_{i} n_{i+1}$
What are the possible states of $\mathbf{n}=\mathbf{2}$ or $\mathbf{3}$ particles?


When particles are far apart: $n$-particle state is unaffected

When particles are close: changed $n$-particle states!
Expectation: seeds of many-body delocalization

## Few-Body Greens Functions

How to quantify this? Greens functions!
One-particle: $G_{1}(x ; y ; E)=\langle 0| c_{x}(E-H)^{-1} c_{y}^{\dagger}|0\rangle$
Two-particle: $G_{2}\left(x_{1}, x_{2} ; y_{1}, y_{2} ; E\right)=\langle 0| c_{x_{2}} c_{x_{1}}(E-H)^{-1} c_{y_{1}}^{\dagger} c_{y_{2}}^{\dagger}|0\rangle$


Greens function allows for effective localization length
One-particle: $\left.\lambda_{1}^{-1}(W, L)=-\left.\frac{2}{L-1}\langle\log | G_{1}(1 ; L)\right|^{2}\right\rangle_{\text {dis }}$
Two-particle: $\left.\lambda_{2}^{-1}(W, L)=-\left.\frac{2}{L-2}\langle\log | G_{2}(1,2 ; L-1, L)\right|^{2}\right\rangle_{\text {dis }}$
Which is related to the transmission coefficient $T_{n}(W, L)=\exp \left(\frac{-2 L}{\lambda_{n}(W, L)}\right)$

## Scaling theory

Dimensionless conductance $g(L)$ is a general form of transmission coefficient

Determined by scaling $\beta(g)=d \ln g / d \ln L$


## Numerical results for scaling theory

$$
\left.\lambda^{-1}=\lim _{N \rightarrow \infty}[2(N-1)]^{-1} \ln \operatorname{Tr}|\langle 1| G(N)| N\right\rangle\left.\right|^{2} \quad \lambda(W, M) / M=f_{d}\left(\lambda_{\infty}(W) / M\right)
$$



Ref: MacKinnon PRL 1981


## Calculating few-body Greens functions

## Exact calculation of Greens function is inefficient so use a trick

Two-particle noninteracting Greens function is

$$
G_{2}^{(0)}=\sum_{m n} \frac{\phi_{x_{2} n} \phi_{x_{1} m} \phi_{y_{1} m} \phi_{y_{2} n}-\phi_{x_{2} m} \phi_{x_{1} n} \phi_{y_{1} m} \phi_{y_{2} n}}{E-\epsilon_{m}-\epsilon_{n}}
$$

Dyson's equation states $G_{2}=G_{2}^{(0)}+G_{2}^{(0)} H_{\mathrm{int}} G_{2}$
But local interactions only act on $\mathcal{O}(L)$ part of Hilbert space
Calculate the restricted Greens function $\tilde{G}_{2}=\tilde{G}_{2}^{(0)}+\tilde{G}_{2}^{(0)} H_{\mathrm{int}} \tilde{G}_{2}$
Speeds up the computation of localization length $\mathcal{O}\left(L^{6}\right) \rightarrow \mathcal{O}\left(L^{4}\right)$

## Scaling of two-body states in d=2

## 2d Bose-Hubbard

$$
\begin{aligned}
H & =t \sum_{\{i, k\}, j}|i, j\rangle\langle k, j|+t \sum_{i,\{j, l\}}|i, j\rangle\langle i, l| \\
& +\sum_{i, j}|i, j\rangle\left(\epsilon_{i}+\epsilon_{j}\right)\langle i, j|+U \equiv H_{0}+U
\end{aligned}
$$

Two-body states

$$
\left.\left.\ln \operatorname{Tr}|\widetilde{G}(l)|^{2} \equiv\left\langle\ln \sum_{i, j}\right| \widetilde{G}(1, i ; l, j)\right|^{2}\right\rangle
$$



## Scaling of few-body states in d=1

$$
\text { Scaling function } \lambda_{n}(W, L) / L=f_{n}^{ \pm}\left(\lambda_{n}^{\infty}(W) / L\right)
$$



Ref: Rademaker PRB 2021

## Beta function

The scaling function $f_{n}$ can be transformed into a beta function

$$
\begin{aligned}
& \log T_{n}(W, L)=\frac{-2}{f_{n}\left(\lambda_{n}^{\infty}(W) / L\right)} \\
& \beta(T)=\frac{d \log T_{n}}{d \log L} \\
& \quad=\log T_{n} \frac{d \log f_{n}(x)}{d \log x}
\end{aligned}
$$



## Few-body delocalization

Beta function for:



## Delocalization transition

for $n$ particles in $d$ dimensions when

$$
n+d>3
$$

## Why is this possible?

In d=2:

## ○ $\longrightarrow$ Bound state of two particles

Single particle with internal structure

Sigma models with symplectic symmetry allow for delocalization in $d=2$ (spin-orbit coupling)

Example: Spinless fermions with nearest neighbor interaction

$|R x\rangle$

$|R y\rangle$


$$
\left|\psi_{R \pm}\right\rangle=\frac{1}{2}[(|R x\rangle+|(R-x) x\rangle) \pm(|R y\rangle+|(R-y) y\rangle)]
$$

## Is it really true?

Criticism in $\boldsymbol{d}=\mathbf{2}$ :
"just finite size effects in numerics"

... but these are results at weak disorder
Ref: Stellin, Orso PRB 2020

Criticism in $\boldsymbol{d}=1$ :

Three-particle states do not have clear symplectic symmetry

... but d=1 delocalization does exist!

Ref: Evers, Mirlin RMP 2008

## On to many-body delocalization




Critical disorder for $n$-body delocalization increases with $n$ up to the critical disorder for many-body delocalization

Energy/charge transport becomes increasingly difficult when $\mathrm{W}>\mathrm{W}^{\mathrm{c}}{ }_{n}$ for large $n$

Possible mechanism for subdiffusion?
$x^{2}(t) \sim t^{2 / z}$

Ref: Luitz, Bar Lev 2017

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## Conclusion

Delocalization transition for $n$ particles in dimensions when

$$
n+d>3
$$



## Extra slides

## How to break thermalization



Calculate (Local) Integrals of Motion
Ref: Rademaker, Ortuño, PRL 2016


Scaling theory of few-body delocalization
Ref: Rademaker, PRB 2021


Dynamics of a quantum spin glass
Ref: Rademaker, Abanin, PRL 2020


The landscape of a self-generated electron glass
Ref: Mahmoudian, Rademaker, et al., PRL 2015

## Failure of perturbation theory

$$
H=\sum_{\substack{i \\ \text { Anderson LIOMs }}} \epsilon_{i} n_{i}+\frac{1}{2} \sum_{i j k l} V_{i j k l} c_{i}^{\dagger} c_{j}^{\dagger} c_{k} c_{l}
$$

Perturbative construction: dress the electrons with particle-hole excitations

$$
\begin{aligned}
c_{i} \rightarrow c_{i} & +\underbrace{\frac{V_{i j k l}}{\epsilon_{i}+\epsilon_{j}-\epsilon_{k}-\epsilon_{l}}}_{2^{\text {nd }}} c_{j}^{\dagger} c_{\text {particle-hole exturbation theory }}^{\dagger} c_{l} \\
& \text { This guy can blow up due to resonances! }
\end{aligned}
$$

## Displacement transformations

$$
H=\sum_{\substack{i \\ \text { Anderson LIOMs }}} \epsilon_{i} n_{i}+\frac{1}{2} \sum_{i j k l} V_{i j k l} c_{i}^{\dagger} c_{j}^{\dagger} c_{k} c_{l}
$$

Our solution: Consider one interaction term:

$$
\begin{aligned}
& X=c_{i}^{\dagger} c_{j}^{\dagger} c_{k} c_{l} \\
& H=\sum_{m} \epsilon_{m} n_{m}+V_{i j k l}\left(X+X^{\dagger}\right)
\end{aligned}
$$

Introduce displacement transformations

$$
\begin{aligned}
& \mathcal{D}_{\lambda}(X)=\exp \left(\lambda\left(X^{\dagger}-X\right)\right) \\
& \tan 2 \lambda=-\frac{V_{i j k l}}{\epsilon_{i}+\epsilon_{j}-\epsilon_{k}-\epsilon_{l}} \\
& \mathcal{D}_{\lambda}^{\dagger}(X) H \mathcal{D}_{\lambda}(X)=\sum_{i} \epsilon_{i} n_{i}+\sum_{i j} V_{i j} n_{i} n_{j}+\ldots
\end{aligned}
$$

The interaction term disappeared!

Compute Local Integrals of Motion
Repeated displacement transformations: $U=\mathcal{D}_{\lambda_{1}}\left(X_{1}\right) \mathcal{D}_{\lambda_{2}}\left(X_{2}\right) \cdots$
Local integrals of motion: $\tilde{n}_{i}=U n_{i} U^{\dagger}=n_{i}+\alpha_{i ; j k l m} c_{j}^{\dagger} c_{k}^{\dagger} c_{l} c_{m}+\mathbb{Q}$
Classical Hamiltonian: $H=\sum_{i} \epsilon_{i} \tilde{n}_{i}+\sum_{i j} V_{i j} \tilde{n}_{i} \tilde{n}_{j}+$. $\longleftarrow$ approximation: cut-off expansion


Study 1d Anderson insulator + NN repulsion

$$
H=\sum_{i=1}^{N} \epsilon_{i} n_{i}+t \sum_{i=1}^{N-1}\left(c_{i}^{\dagger} c_{i+1}+c_{i+1}^{\dagger} c_{i}\right)+V \sum_{i=1}^{N-1} n_{i} n_{i+1}
$$

Disorder scale: $\epsilon_{i} \in[-W / 2, W / 2]$
(4) Large system sizes ( $N$ up to 60 )

- Cannot describe ergodic transition (?!)


## Bring on the bath



How to calculate (Local) Integrals of Motion


Scaling theory of few-body delocalization


Dynamics of a quantum spin glass


The landscape of a self-generated electron glass

## Quantum spin glass

Only known one-dimensional spin-glass has long-range interactions


Kotliar et al, PRB 1983


Each spins feels an effective field from all the other spins

$$
\phi_{i} \equiv \sum_{j} \frac{J_{i j}}{|i-j|^{\alpha}} Z_{j}
$$

With transverse field, only resonant spins $\left|\phi_{i}\right|<h_{x}$ will flip

## Dynamics of quantum spin glass



At low T , the distribution of effective fields has a soft gap

Resonant spins are very scarce

Dynamics: Most spins remain frozen, only resonant spins entangle


Numerical technique: Monte Carlo to find low T state

+ Exact Diagonalization for resonant spins

Verified new phase: ergodic for resonant spins localized for other spins

## Experimental realization

Hyperfine states of Yb ions allows exactly the right Hamiltonian


But their experiment has 10 spins only
Typical distance between resonant spins is, for $h=0.05 \mathrm{~J}$,

## Infinite temperature

$$
d=60 \text { sites }
$$

## Low temperature

$d=100-500$ sites

# Forget disorder 



How to calculate (Local) Integrals of Motion


Scaling theory of few-body delocalization


## Dynamics of a quantum spin glass

$000 \because 0$.
The landscape of a self-generated electron glass $0.000 \cdot$

## Self-generated electron glass

Electrons in organic crystal $\theta$-(BEDT-TTF) 2 RbZn $(S C N)_{4}$ Ref: Kagawa, Nat Phys 2014; Sato PRB 2014
 molecules in a triangular lattice one electron per two molecules
amorphous configurations


## Monte Carlo simulations

Electrons on a triangular lattice with long-range Coulomb interactions

$$
\begin{array}{r}
H=-t \sum_{\langle i, j\rangle} c_{i}^{\dagger} c_{j}+\frac{1}{2} \sum_{i j} V_{i j}\left(n_{i}-\frac{1}{2}\right)\left(n_{j}-\frac{1}{2}\right) \\
\\
V_{i j}=\frac{V}{\left|R_{i}-R_{j}\right|}
\end{array}
$$

Ground state is stripe phase
Monte Carlo simulations slow down below at low T System doesn't reach stripes but remains amorphous


Ref: Mahmoudian, Rademaker, et al., PRL 2015; Rademaker, et al., NJP 2018

## Landscape picture

Exponentially many metastable states
Example: $24 \times 24$ lattice has $\mathbf{1 0}^{\mathbf{3 5}} \mathrm{MS}$ states


Configuration space

DOS satisfies Efros-Shklovskii bound

without marginal stability!

Self-generated glasses are different from quenched disorder glasses!

Ref: Mahmoudian, Rademaker, et al., PRL 2015; Rademaker, et al., NJP 2018

