LOCALIZATION IN INTERACTING FERMIONIC CHAINS WITH QUASI-RANDOM DISORDER

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- In 1D typically any amount of disorder produces localization, while in 3D the disorder has to be sufficiently strong and a metal to insulator transition is expected varying the strength of the random field. Still open problems in 2D and 3D.

• Many Body Localization. Evidence that localization persists in the presence of weak inter-particle interactions on the basis of perturbative analysis Basko, Aleiner, Altshuler (2006) or numerical analysis Oganesyan, Huse (2007); MBL rigorous consequence in 1d of an assumption of level attraction Imbrie (2014).

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• A proof of MBL in generality is a challenging problem (single particle description breaks down, full N-particle Schroedinger)

• Cold atoms as "quantum simulators"; Experimental evidence of MBL with quasi-random disorder and local interaction Schreiber, Hodgman, Bordia, Luschen, Fischer, Vosk, Altman, Schneider, Bloch, (2015) by monitoring the time evolution of local observables following a quench. No dependence on the sign of the interaction.

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- With no interaction very good theoretical understanding based on advanced mathematical tools; quest for understanding the role of interaction.

THE INTERACTING AUBRY-ANDRE' MODEL

 If a⁺_x, a⁻_x, x ∈ Z ≡ Λ are spinless creation or annihilation operators on the Fock space verifying {a⁺_x, a⁻_y} = δ_{x,y}, {a⁺_x, a⁺_y} = {a⁻_x, a⁻_y} = 0. The Fock space Hamiltonian is

$$H = -\varepsilon \left(\sum_{x \in \Lambda} (a_{x+1}^+ a_x + a_{x-1}^+ a_x^-) + \sum_{x \in \Lambda} u \cos(2\pi(\omega x + \theta)) a_x^+ a_x^- + U \sum_{x,y} v(x - y) a_x^+ a_x^- a_y^+ a_y^- \right)$$

with $v(x-y) = \delta_{y-x,1} + \delta_{x-y,1}$.

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• Early studies of the extended phase in Mastropietro (1999) and Giamarchi, Mohunna, Vidal (1999)

• In the non interacting case the states are obtained by the antisymmetrization (fermions) of the eigenfunctions of almost Mathieu equation

 $-\varepsilon\psi(x+1)-\varepsilon\psi(x-1)+u\cos(2\pi(\omega x+\theta)\psi(x))=E\psi(x)$

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- the spectrum is a Cantor set for all irrational ω. For almost every ω, θ the almost Mathieu operator has a)for ε/u < 1/2 exponentially decaying eigenfunctions (Anderson localization);
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b)for $\varepsilon/u > \frac{1}{2}$ purely absolutely continuous spectrum (extended quasi-Bloch waves)

• Metal insulator transition (with no interaction) seen experimentally by Roati et al (2008) • Such remarkable properties are related to a deep connection between the non interacting Aubry-Andre model and the Kolmogorov-Arnold-Moser (KAM) theorem of classical mechanics.

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- Such remarkable properties are related to a deep connection between the non interacting Aubry-Andre model and the Kolmogorov-Arnold-Moser (KAM) theorem of classical mechanics.
- KAM ensures the existence of quasi-periodic solutions of Hamiltonian systems close to integrable one, that is the stability of invariant tori. Applications to the stability of solar system.

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- KAM ensures the existence of quasi-periodic solutions of Hamiltonian systems close to integrable one, that is the stability of invariant tori. Applications to the stability of solar system.
- A crucial assumption of KAM and of the analysis of almost mathieu is that the frequency verify a number theoretical condition called Diophantine condition to deal with small divisors. It says that a number is a "good irrational" and is full mesaure.

• We impose a Diophantine condition on the frequency

$$||\omega x|| \geq C_0 |x|^{-\tau} \quad \forall x \in \mathbb{Z}/\{0\} \quad (*)$$

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 $\bullet\,$ The continued fraction representation of a number $\omega\,$

$$\omega = a_0 + rac{1}{a_1 + rac{1}{a_2 + rac{1}{a_3 + \dots}}}$$

The golden ratio $\omega = \frac{\sqrt{5}+1}{2}$ has representation 1; 1, ..1, .. and it verifies the Diophantine condition with $\tau = 1$ and $C_0 = \frac{3+\sqrt{5}}{2}$.

• The construction of all the eigenvectors of the *N*-body Schroedinger equation with almost-Mathieu potential and interaction seems at the moment out of reach, especially for infinite *N*.

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- The construction of all the eigenvectors of the *N*-body Schroedinger equation with almost-Mathieu potential and interaction seems at the moment out of reach, especially for infinite *N*.
- More modest goal. Information on the localization of the interacting ground state can be obtained by the zero temperature grand-canonical truncated correlations of local operators, whose exponential decay with the distance is a sign of localization. This allow to use exact RG methods combined with KAM (Lindstedt series).

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- For $\omega \ \theta$ verifying Diophantine conditions, for small $\frac{\varepsilon}{u}$, $\frac{U}{u}$ the fermionic zero temperature grand canonical infinite volume truncated correlations of local operators decays exponentially for large distances.

Comm Math Phys 342, 1, 217(2016); Phys. Rev. Lett. 115, 180401 (2015) , arxiv 1604.08264

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• Renormalized expansion around the anti-integrable limit

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$$H = -\varepsilon \left(\sum_{x \in \Lambda} (a_{x+1}^+ a_x + a_{x-1}^+ a_x^-) + \sum_{x \in \Lambda} u \cos(2\pi(\omega x + \theta)) a_x^+ a_x^- + U \sum_{x,y} v(x - y) a_x^+ a_x^- a_y^+ a_y^- \right)$$

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with $v(x - y) = \delta_{y-x,1} + \delta_{x-y,1}$. • If $a_x^{\pm} = e^{(H-\mu N)x_0} a_x^{\pm} e^{-(H-\mu N)x_0}$, $\mathbf{x} = (x, x_0)$, $N = \sum_x a_x^+ a_x^-$ and μ the chemical potential, the Grand-Canonical imaginary time 2-point correlation is

$$<\mathbf{T}a_{\mathbf{x}}^{-}a_{\mathbf{y}}^{+}>=\frac{Tre^{-\beta(H-\mu N)}\mathbf{T}\{a_{\mathbf{x}}^{-}a_{\mathbf{y}}^{+}\}}{Tre^{-\beta(H-\mu N)}}$$

where **T** is the time-order product and μ is the chemical potential.

 It is convenient to write the chemical potential as a function of the interaction so that the density has the same value in the free or interacting case. We introduce a counterterm ν so that the interacting chemical potential is u cos 2π(ωx̂ + θ).

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- A condition on the phase is also imposed

$$||\omega x \pm 2\theta|| \ge C_0 |x|^{-\tau} \quad \forall x \in \mathbb{Z}/\{0\} \quad (**)$$

LOCALIZED REGIME

Theorem

Under conditions (*) and (**), u = 1 $\mu = \cos 2\pi(\omega \hat{x} + \theta) + \nu$ there exists an ε_0 such that, for $|\varepsilon|, |U| \le \varepsilon_0$, it is possible to choose ν so that the limit $\beta \to \infty$

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- (**) could be replaced by a condition on the density

 Different behavior is found close to the integrable limit. Fix
 ε = 1,θ = 0, U, u small, μ = cos p_F, ||2πωn||_{2π} ≥ C|n|^{-τ}, n ≠ 0,
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- 1)If $||2p_F + 2\pi n\omega||_{2\pi} \ge C|n|^{-\tau}$ a decay of the two point function $O(|x y|^{-1-\eta})$, $\eta = aU^2 + O(U^3)$ (metallic Luttinger liquid behavior).

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- 2) If $p_F = n\omega\pi$ a faster than any power decay with rate

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with $F = O(|U| + |\lambda|)$, a_n non vanishing and $X_n = X_n(U) = 1 + bU + O(U^2)$; the decay rate is of the order of the interacting gap. Dense set of gaps.

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All gaps are renormalized via a critical exponent

• In the non interacting case there are infinitely many gaps in correspondence of quasi-momenta $2n\pi\omega \mod 2\pi$, forming a dense set, and their size is decreasing exponentially with n.

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- In the case of a Fibonacci quasi-periodic potential there is evidence that the interaction closes the smallest gaps, Giamarchi (1999), causing an insulating to metal transition.
- In the case of Aubry-Andre' potential all gaps persists instead; no quantum phase transition at small coupling.

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Close to the singularity

$$\cos(k'\pm p_F)-\mu\sim\pm\sin p_Fk'+O(k'^2)$$

linear dispersion relation.

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- The 2-point function is

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ho,k_0)\sim rac{1}{-ik_0\pm v_0(\omega x')_{\mathrm{mod}.1}}$$

• The denominator can be arbitrarily large; for $x \neq \rho \hat{x}$ by (*),(**) , $||\omega x'|| = ||\omega(x - \rho \hat{x}) + 2\delta_{\rho,-1}\theta|| \ge C|x - \rho \hat{x}|^{-\tau}$. $(\omega x')_{mod,1}$ can be very small for large x (infrared-ultraviolet mixing) The 2-point function is given by $\frac{\partial^2}{\partial \phi_{\mathbf{x}}^+ \partial \phi_{\mathbf{y}}^-} W|_0$

$$e^{W(\phi)} = \int P(d\psi) e^{-V(\psi) - \mathcal{B}(\psi,\phi)}$$

with $P(d\psi)$ a gaussian Grassmann integral with propagator $\delta_{x,y}\bar{g}(x, x_0 - y_0)$

$$\begin{split} V(\psi) &= U \int d\mathbf{x} \sum_{\alpha = \pm} \psi_{\mathbf{x}}^+ \psi_{\mathbf{x}}^- \psi_{\mathbf{x}+\alpha \mathbf{e}_1}^+ \psi_{\mathbf{x}+\alpha \mathbf{e}_1}^- \\ &+ \varepsilon \int d\mathbf{x} (\psi_{\mathbf{x}+\mathbf{e}_1}^+ \psi_{\mathbf{x}}^- + \psi_{\mathbf{x}-\mathbf{e}_1}^+ \psi_{\mathbf{x}}^-) + \nu \int d\mathbf{x} \psi_{\mathbf{x}}^+ \psi_{\mathbf{x}}^- \end{split}$$

where $\int d\mathbf{x} = \sum_{\mathbf{x} \in \Lambda} \int_{-\frac{\beta}{2}}^{\frac{\beta}{2}} dx_0$, Finally $B = \int d\mathbf{x} (\phi_{\mathbf{x}}^+ \psi_{\mathbf{x}}^- + \psi_{\mathbf{x}}^+ \phi_{\mathbf{x}}^-)$

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• In absence of many body interaction there are only chain graphs, $\alpha_i=\pm$

$$\varepsilon^{n} \sum_{x_{1}} \int dx_{0,1} \dots dx_{0,n} \bar{g}(x_{1}, x_{0} - x_{0,1}) \bar{g}(x_{1} + \sum_{i \leq n} \alpha_{i}, (x_{0,n} - y_{0}))$$
$$\prod_{i=1}^{n} \bar{g}(x_{1} + \sum_{k \leq i} \alpha_{k}, x_{0,i+1} - x_{0,i})$$

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- When $U \neq 0$ there also loops producing additional divergences, absent in classically.
- To establish localization in presence of interaction one has to prove that such small divisors are harmless. Sort of quantum KAM.
 Constructive RG approach.

• We perform an RG analysis decomposing the propagator as sum of propagators living at scale $|\phi_x - \phi_{\hat{x}}| \sim \gamma^h$, h = 0, -1, -2..., $\phi_x = \cos 2\pi(\omega x + \theta)$; this correspond to two regions, around \bar{x}_+ and \bar{x}_- .

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• This implies that the single scale propagator has the form $\sum_{\rho=\pm} g_{\rho}^{(h)}$ with $|g_{\rho}^{(h)}(\mathbf{x})| \leq \frac{C_N}{1+(\gamma^h(x_0-y_0))^N}$; the corresponding Grasmann variable is $\psi_{\mathbf{x},\rho}^{(h)}$.

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- Very similar to what is done in the integrable limit u/ε small (Aubry duality).
- We integrate the fields with decreasing scale; for instance W(0) (the partition function) can be written as

$$\int P(d\psi)e^V = \int P(d\psi^{\leq -1})\int P(d\psi)e^V = \int P(d\psi^{\leq -1})e^{V^{-1}}...$$

The effective potential V^h , sum of monomials of any order in ψ_{ρ}^{\pm} .

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- One has to distinguish among the monomials Π_i ψ_{x'_i,x_{0,i},ρ_i} in the effective potential between resonant and non resonant terms. Resonant terms; x'_i = x'. Non Resonant terms x'_i ≠ x'_j for some i,j. (In the non interacting case only two external lines are present).
- It turns out that the non resonant terms are irrelevant (even if they are relevant according to power counting).
- Roughly speaking, the idea is that if two propagators have similar (not equal) small size (*non resonant subgraphs*), then the difference of their coordinates is large and this produces a "gain" as passing from x to x + n one needs n vertices. That is if $(\omega x'_1)_{mod1} \sim (\omega x'_2)_{mod1} \sim \Lambda^{-1}$ then by the Diophantine condition

$$2\Lambda^{-1} \geq ||\omega(x_1' - x_2')|| \geq C_0 |x_1' - x_2'|^{-\tau}$$

that is $|x_1' - x_2'| \ge \bar{C} \Lambda^{ au^{-1}}$

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- Consider two vertices w_1, w_2 such that x'_{w_1} and x'_{w_2} are coordinates of the external fields, and let be c_{w_1,w_2} the path (vertices and lines) in \overline{T}_v connecting w_1 with w_2 ; we call $|c_{w_1,w_2}|$ the number of vertices in c_{w_1,w_2} . The following relation holds, if $\delta'_w = \pm 1$ it corresponds to an ε end-point and $\delta'_w = (0, \pm 1)$ is a U end-point

$$x'_{w_1} - x'_{w_2} = \bar{x}_{\rho_{w_2}} - \bar{x}_{\rho_{w_1}} + \sum_{w \in c_{w_1,w_2}} \delta^{i_w}_w$$

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• As $x_i - x_j = M \in \mathbb{Z}$ and $x'_i = x'_j$ then $(\bar{x}_{\rho_i} - \bar{x}_{\rho_j}) + M = 0$, so that $\rho_i = \rho_j$ as $\bar{x}_+ = \hat{x}$ and $\bar{x}_- = -\hat{x} - 2\theta/\omega$ and $\hat{x} \in \mathbb{Z}$.


FIG. 1: A tree \overline{T}_v with attached wiggly lines representing the external lines P_v ; the lines represent propagators with scale $\geq h_v$ connecting w_1, w_a, w_b, w_c, w_2 , representing the end-points following vin τ .

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• By the Diophantine condition a) $\rho_{w_1} = \rho_{w_2}$ the (*); b)if $\rho_{w_1} = -\rho_{w_2}$ by (**)

$$2cv_0^{-1}\gamma^{h_{\tilde{v}'}} \ge \\ ||(\omega x'_{w_1})||_1 + ||(\omega x'_{w_2})||_1 \ge ||\omega(x'_{w_1} - x'_{w_2})||_1 \ge C_0(|c_{w_2,w_1}|)^{-\tau}$$

so that $|c_{w_1,w_2}| \geq A \gamma^{\frac{-h_{\tilde{v}'}}{\tau}}$. If two external propagators are small but not exactly equal, you need a lot of hopping or interactions to produce them

• If $\bar{\varepsilon} = \max(|\varepsilon|, |U|)$ from the $\bar{\varepsilon}^n$ factor we can then extract

$$\bar{\varepsilon}^{\frac{n}{4}} \leq \prod_{v \in L} \varepsilon^{N_v 2^{h_{v'}}}$$

where N_v is the number of points in v; as $N_v \ge |c_{w_1,w_2}| \ge A\gamma^{\frac{-h_{v'}}{\tau}}$ then

$$\bar{\varepsilon}^{\frac{n}{4}} \leq \prod_{v \in L} \bar{\varepsilon}^{A\gamma^{\frac{-h_{v'}}{\tau}} 2^{h_{v'}}}$$

where L are the non resonant vertices. If $\gamma^{\frac{1}{\tau}}/2 > 1$ then $\leq C^n \prod_{v \in L} \gamma^{3h_v S_v^L}$ where S_v^L is the number of non resonant clusters in v.

• We localize the resonant terms $\mathbf{x} = x_{0,i}, x$ with all x'_i equal

$$\mathcal{L}\psi_{\mathbf{x}_{1},\rho_{1}}^{\varepsilon_{1}}...\psi_{\mathbf{x}_{n},\rho_{n}}^{\varepsilon_{n}}=\psi_{\mathbf{x}_{1},\rho_{1}}^{\varepsilon_{1}}...\psi_{\mathbf{x}_{1},\rho_{n}}^{\varepsilon_{n}}$$

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Note that one has to renormalize monomial of all orders, a potentially very dangerous situation (this is like in KAM).

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- The result can be rephrased fixing θ and changing the chemical potential.

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- Combined effect of Umklapp and the incommensurability of potential has the effect that a large momentum exchange can connect points arbitrarily close to the Fermi points.

$$\sum_{i=1}^{m} \varepsilon_{i} \rho_{i} k_{i}' = -\sum_{i=1}^{m} \varepsilon_{i} \rho_{i} p_{F} + 2n\pi\omega + 2l\pi$$

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- This is true for quasi-periodic functions with fast decaying Fourier transform; With other quasi-random noise, is believed instead that there are infinitely many rcc.

• System of fermions with quasi-random Aubry-Andre' noise and interaction.

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- Spin? Coupled chains? other eigenstates? 2 or 3 dimension?