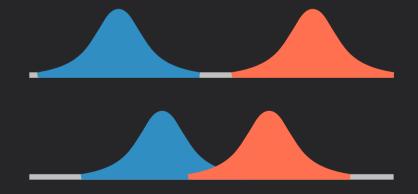
EXACTLY SOLVABLE MODEL OF CORRELATED METAL-INSULATOR TRANSITION

Insights on Non-Fermi Liquid and Mott Insulator

ABHIRUP MUKHERJEE

February 16, 2025 EPQM Seminar



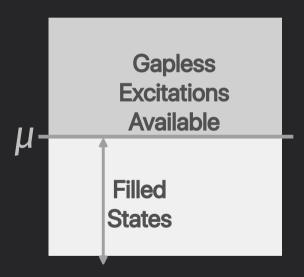
In A Nutshell

- An exactly solvable model that displays correlation-driven transition from a non-Fermi liquid to a Mott insulator.
- Analyse the non-Fermi liquid in this controlled setting to understand its features.
- Study the superconducting instability of this metal to figure out the nature of the superconductor arising from this exotic normal state.

Where Does This Fit?

Mott Insulators Are Different

Half-filled system is **metallic** in absence of interactions.

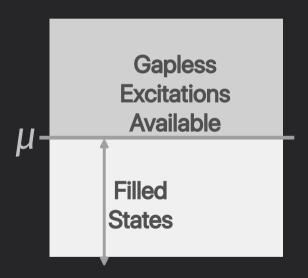


Dispersion away from band edge is **non-interacting**.

Mott (1968)

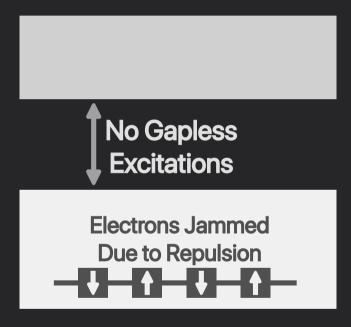
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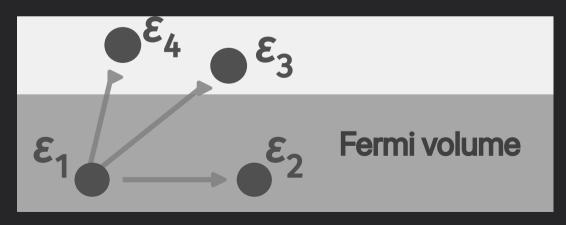
Add strong interactions - Mott Insulator!



Gap opens inside
$$\begin{pmatrix} \varepsilon & U \\ U & \varepsilon \end{pmatrix} \rightarrow \varepsilon \pm U$$
 the band:

Landau Fermi Liquid Theory (Postulates)

- Theory describing how metals arise in interacting systems
- Lack of scattering phase space at low-energies
- Fermi surface and low-lying electronic excitations survive.



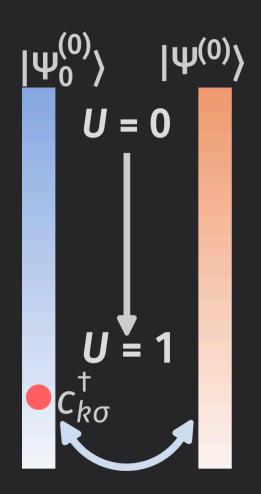
```
\Gamma
\sim \int d\varepsilon_4 d\varepsilon_3 d\varepsilon_2 \delta(\varepsilon - \varepsilon_2 - \varepsilon_3 - \varepsilon_4)
\sim \varepsilon^2
\rho \sim T^2
```

Landau Fermi Liquid Theory (Quantification)

- Self-energy $\Sigma \sim i\omega^2$. Roughly equal to scattering rate. Vanishes very fast as $\omega \rightarrow$ 0: essential for quasiparticle picture
- Quasiparticle residue: how similar are the true excitations to 1-particle excitations

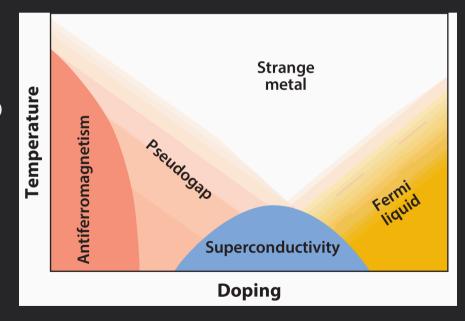
$$Z = \left\langle \Psi \middle| c_{k\sigma}^{\dagger} \middle| \Psi_{0} \right\rangle$$

■ $Z = \left(1 - \frac{\partial(\text{Re }\Sigma)}{\partial\omega}\right)^{-1}$ ~ 1. Must be non-zero for Landau Fermi liquid.



Violations Of Landau Fermi Liquid Theory

- Tomonaga-Luttinger Liquid:
 Interacting electrons in 1D
- Overscreened fixed points in Kondo models
- Strange Metal: Normal state of unconventional SCs in Cu oxides, heavy fermions, Fe pnictides, etc



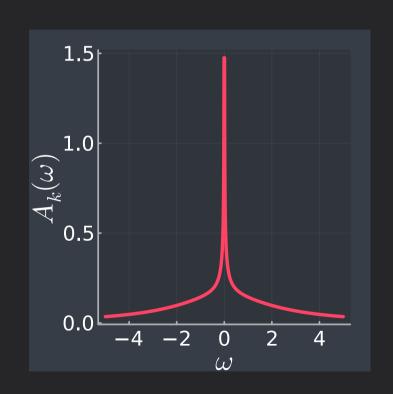
Custers et al. (2003); Doiron-Leyraud et al. (2009); Emery & Kivelson (1992); Haldane (1981); Keimer et al. (2015)

The Marginal Fermi Liquid

- Phenomenological explanation of normal state of cuprates: $Σ \sim ω log(|ω|) iπ |ω|$
- Quasiparticle residue vanishes at Fermi surface

$$Z^{-1} = 1 - \frac{\partial(\text{Re }\Sigma)}{\partial\omega} \sim -\log(\omega) \to \infty$$

 Not accessible through perturbative corrections of Landau Fermi liquid



Main Takeaways

- Landau Fermi Liquid theory requires interacting eigenstates to be adiabatically connected to non-interacting eigenstates
- Non-Fermi liquids involve vanishing quasiparticle residue, signalling that the states are in fact not adiabatically connected.
- This typically means non-perturbative approaches are required to deal with such phases.
- The qualitatively different nature of excitations means that LFL and NFL correspond to distinct fixed points in the RG sense.

An Exactly Solvable Model

The Hatsugai-Kohmoto Model

Consider long-ranged interaction in real-space.

$$H = -t \sum_{\langle i,j \rangle,\sigma} c^{\dagger}_{i,\sigma} c_{j,\sigma} + \frac{U}{L^d} \sum_{i_1,i_2,r} c^{\dagger}_{i_1+r,\uparrow} c^{\dagger}_{i_2-r,\downarrow} c_{i_2,\downarrow} c_{i_1,\uparrow}$$



Switch to momentum space, Hamiltonian becomes local!

$$c_r^{\dagger} \sim \sum_k e^{-ikr} c_k^{\dagger}; \quad H = \sum_{\vec{k}} H_{\vec{k}}; \quad H_{\vec{k}} = \varepsilon_{\vec{k}} \sum_{\sigma} n_{\vec{k},\sigma} + U n_{\vec{k}\uparrow} n_{\vec{k}\downarrow}$$

The Hatsugai-Kohmoto Model

Contrast with the Hubbard interaction.

$$H_{\text{int}} \sim \sum_{i} n_{i,\uparrow} n_{i,\downarrow} = \sum_{k_1,k_2,q} c_{k_1+q,\uparrow}^{\dagger} c_{k_2-q,\downarrow}^{\dagger} c_{k_2,\downarrow} c_{k_1,\uparrow}$$













- local in real-space, highly non-local in k -space
- HK model is q = 0, $k_1 = k_2$ (zero mode!) component of the Hubbard
- HKM is easier to solve than Hubbard (KE and PE do not commute)

$$H = \sum_{\vec{k}} H_{\vec{k}}; \quad H_{\vec{k}} = \varepsilon_{\vec{k}} \sum_{\sigma} n_{\vec{k},\sigma} + U n_{\vec{k}\uparrow} n_{\vec{k}\downarrow}$$

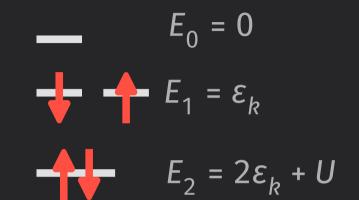
Each H_k can be diagonalised.

$$|0\rangle : E = 0, \quad |\sigma\rangle : E = \varepsilon_b, \quad |2\rangle : E = 2\varepsilon_b + U$$

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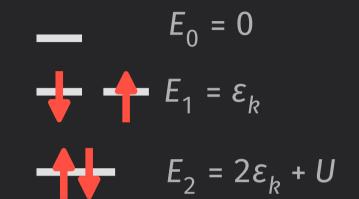
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Case Of Half-Filling

$$E(\mu) = E - \mu n_k, \quad \mu = \frac{U}{2}$$

$$E_0 = 0$$
, $E_1 = \varepsilon_k - \frac{U}{2}$, $E_2 = 2\varepsilon_k$



$$H = \sum_{\vec{k}} H_{\vec{k}}; \quad H_{\vec{k}} = \varepsilon_{\vec{k}} \sum_{\sigma} n_{\vec{k},\sigma} + U n_{\vec{k}\uparrow} n_{\vec{k}\downarrow}$$

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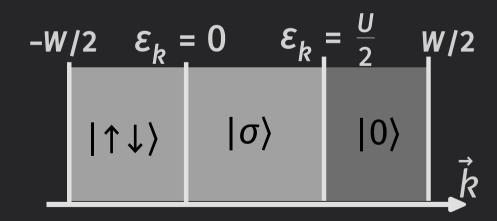
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, $E_1 = \varepsilon_k - \frac{U}{2}$, $E_2 = 2\varepsilon_k$



Introduction to Greens Functions

Nature of excitations can be studied through Greens function

$$G_{v}(t) = -i\theta(t)\langle \left\{ c_{v(t)}, c_{v}^{\dagger} \right\} \rangle$$

- Non-interacting system: $G_k(\omega + i\eta) = \frac{1}{\omega + i\eta \varepsilon_k}$
- Poles of Greens function → one-particle excitations
- **Zeroes** of Greens function \rightarrow destruction of one-particle excitations

Greens function can be calculated
$$E_0 = 0$$

as
$$G_{k\sigma}(\omega) = \frac{P_e(k\sigma)}{\omega - E_X} + \frac{P_h(k\sigma)}{\omega + E_X}$$

$$E_1 = \varepsilon_k$$

$$E_2 = 2\varepsilon_k + U$$

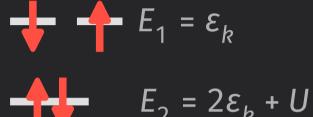
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$$E_1 = \varepsilon_k$$

$$E_2 = 2\varepsilon_k + U$$

$$E_0 = 0$$



If opposite spin is unoccupied

- Particle addition: $E_x = \varepsilon_b$
- Particle removal: $E_X = -\overline{\varepsilon_k}$

$$G \to \frac{1 - \langle n_{k\overline{o}} \rangle}{\omega - \varepsilon_b}$$

$$|0\rangle \rightarrow |\sigma\rangle$$

$$|\sigma\rangle \rightarrow |0\rangle$$

Greens function can be calculated

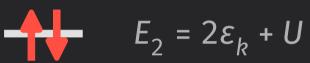
$$G_{k\sigma}(\omega) = \frac{P_e(k\sigma)}{\omega - E_X} + \frac{P_h(k\sigma)}{\omega + E_X}$$

$$E_1 = \varepsilon_k$$

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If opposite spin is occupied

- Particle addition: $E_x = \varepsilon_b + U$
- Particle removal: $E_x = -\varepsilon_k U$

$$G \to \frac{\langle n_{k\overline{o}} \rangle}{\omega - \varepsilon_b - U}$$

$$|0\rangle \rightarrow |\sigma\rangle$$

$$|\sigma\rangle \rightarrow |0\rangle$$

Greens function can be calculated
$$F_0 = 0$$

as
$$G_{k\sigma}(\omega) = \frac{P_e(k\sigma)}{\omega - E_X} + \frac{P_h(k\sigma)}{\omega + E_X}$$

$$E_0 = 0$$

$$E_1 = \varepsilon_k$$

$$E_2 = 2\varepsilon_k + U$$

Total Greens Function

$$G_{k\sigma} = \frac{1 - \left\langle n_{k\overline{\sigma}} \right\rangle}{\omega - \varepsilon_k} + \frac{\left\langle n_{k\overline{\sigma}} \right\rangle}{\omega - \varepsilon_k - U}$$
$$\left(\varepsilon_k \to \varepsilon_k - \mu\right)$$

Correlated Metal-Insulator Transition

The Case Of Half-Filling:
$$2\mu = U$$
, $\langle n_{k\sigma} \rangle = \frac{1}{2}$

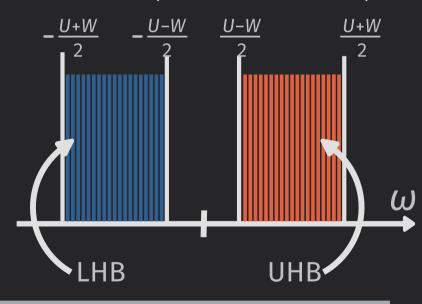
$$G_{k\sigma} = \frac{1}{2} \left[(\omega - \varepsilon_k + U/2)^{-1} + (\omega - \varepsilon_k - U/2)^{-1} \right]$$

Correlated Metal-Insulator Transition

The Case Of Half-Filling:
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, $\langle n_{k\sigma} \rangle = \frac{1}{2}$

$$G_{k\sigma} = \frac{1}{2} \left[(\omega - \varepsilon_k + U/2)^{-1} + (\omega - \varepsilon_k - U/2)^{-1} \right]$$

U > *W* (Mott Insulator)

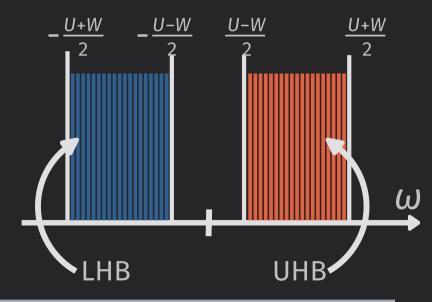


Correlated Metal-Insulator Transition

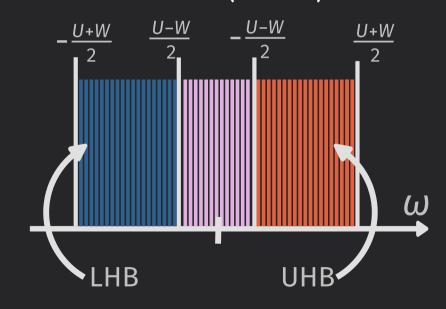
The Case Of Half-Filling:
$$2\mu = U$$
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$$G_{k\sigma} = \frac{1}{2} \left[\left(\omega - \varepsilon_k + U/2 \right)^{-1} + \left(\omega - \varepsilon_k - U/2 \right)^{-1} \right]$$

U > *W* (Mott Insulator)



U < W (Metal)

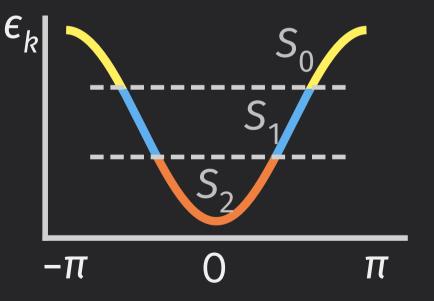


Non-Fermi Liquid Signatures In Metallic Phase

$$G_{k\sigma} = \frac{1}{2} \left[\left(\omega - \varepsilon_k + U/2 \right)^{-1} + \left(\omega - \varepsilon_k - U/2 \right)^{-1} \right]$$

Momentum states classified into three groups:

- $S_2 : \varepsilon_k < -U/2$: Both poles below $\omega = 0 : \langle n_k \rangle = 2$
- $S_1: -U/2 < \varepsilon_k < U/2$: One pole below $\omega = 0: \langle n_k \rangle = 1$
- $S_0: \varepsilon_k > U/2$: No pole below $\omega = 0: \langle n_k \rangle = 0$

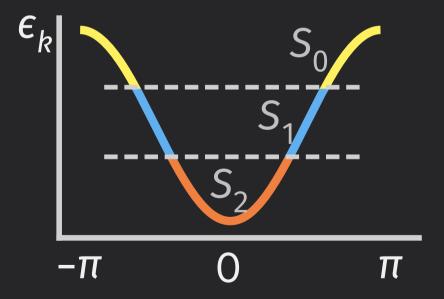


$$G_{k\sigma} = \frac{1}{2} \left[\left(\omega - \varepsilon_k + U/2 \right)^{-1} + \left(\omega - \varepsilon_k - U/2 \right)^{-1} \right]$$

Momentum states classified into three groups:

Ground state is a mixed state.

- k -states in S₂ are doublyoccupied.
- k -states in S_1 are half-filled.
- 2^{N₁}-fold degenerate.

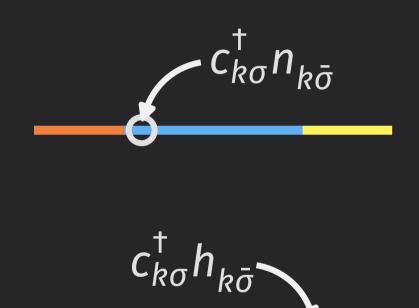


Near
$$S_2 - S_1$$
 boundary $(\varepsilon_k = -U/2)$

- Excitation operator: $c_{k\sigma}^{\dagger} n_{k\overline{\sigma}}$
- Excitation energy is $\varepsilon_k + U \mu \rightarrow 0^+$

Near
$$S_1 - S_0$$
 boundary $(\varepsilon_k = U/2)$

- Excitation operator: $c_{k\sigma}^{\dagger}(1 n_{k\overline{\sigma}})$
- Excitation energy is $\varepsilon_k \mu \rightarrow 0^+$



Projectors are needed because the other excitations are gapped.

Near
$$S_2$$
 – S_1 boundary $c_{k\sigma}^{\dagger} n_{k\overline{\sigma}}$

Near
$$S_1 - S_0$$
 boundary $c_{k\sigma}^{\dagger} (1 - n_{k\overline{\sigma}})$

Excitations are Non-Fermi Liquid In Nature!

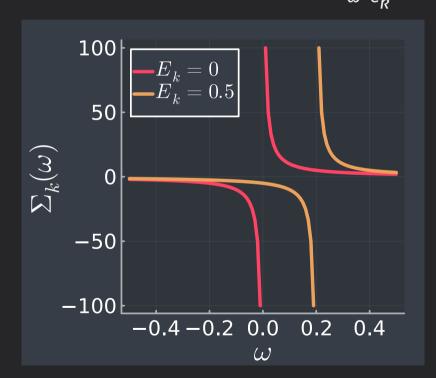
- Strong correlations lead to composite (hole/double) excitations
- Excitations are **not electronic** (do not satisfy {·} relations)
- Breakdown of quasiparticle picture, and hence of Fermi liquid theory

Signature II: Divergence of Self-Energy

Greens function at 1/2-filling can be rewritten as $G_{k\sigma} = \frac{1}{\omega - \varepsilon_k + \frac{U^2/4}{\omega - \varepsilon_k}}$.

Self-energy is
$$\Sigma(\omega) = \frac{U^2/4}{\omega - \varepsilon_k}$$

- Diverges along $\varepsilon_k = 0$ as $\omega \to 0$
- Violates Fermi Liquid Theory
- Leads to zeros of Greens function
- Death of Landau quasiparticles



Signature II: Divergence of Self-Energy

How Does A Diverging Self-Energy Leave The System Metallic?

Greens functions for composite excitations do not have self-energy!

$$d_{k\sigma}^{\dagger} = c_{k\sigma}^{\dagger} n_{k\sigma}, \quad G_{d} = \frac{1}{\omega - \varepsilon_{k} - \frac{U}{2}}$$

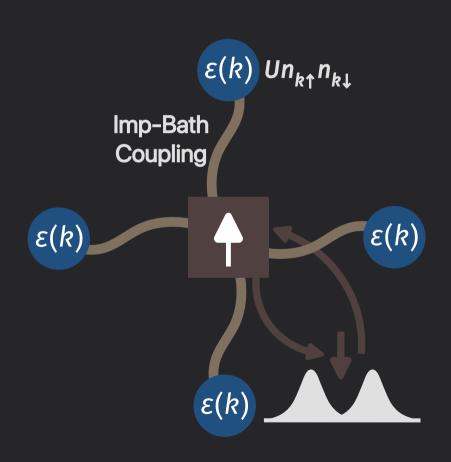
$$h_{k\sigma}^{\dagger} = c_{k\sigma}^{\dagger} (1 - n_{k\sigma}), \quad G_{h} = \frac{1}{\omega - \varepsilon_{k} + \frac{U}{2}}$$

These can therefore propagate with long lifetimes.

Summary of Main Ideas

Avenues for Futher Investigation

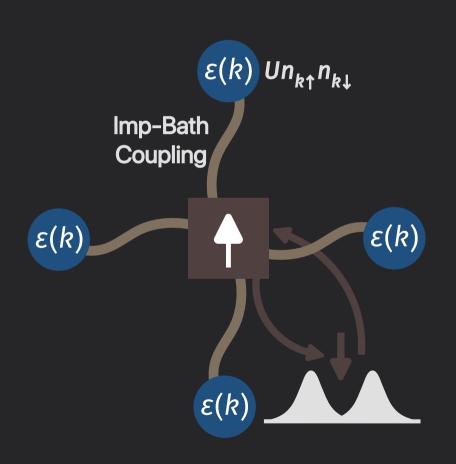
Kondo Screening in Hatsugai-Kohmoto Model



Consider local moment hybridising with HK Model

$$H = H_{Kondo} + H_{HKM}$$

Kondo Screening in Hatsugai-Kohmoto Model

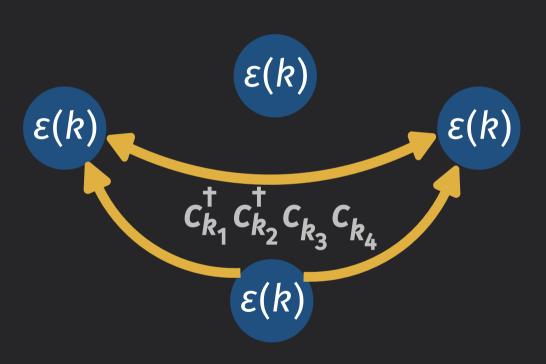


Consider local moment hybridising with HK Model

$$H = H_{Kondo} + H_{HKM}$$

How does **absence** of quasiparticles affect Kondo screening?

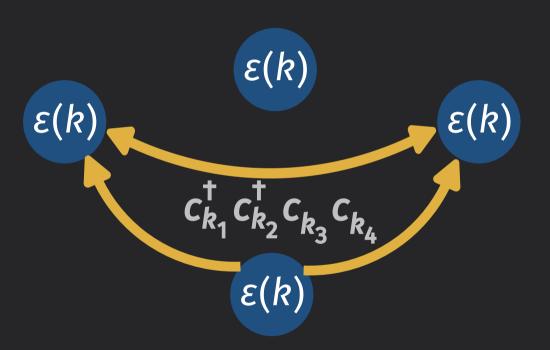
Toy Model for Thermalisation and Many-Body Scars



Consider HK Model perturbed by Hubbard interaction

$$H = H_{HKM} + P_{\nu}H_{Hub}P_{\nu}$$

Toy Model for Thermalisation and Many-Body Scars

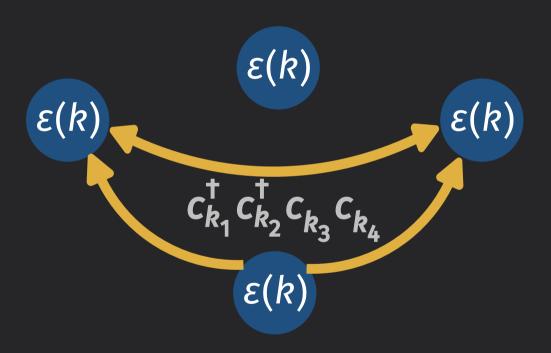


Consider HK Model perturbed by Hubbard interaction

$$H = H_{HKM} + P_v H_{Hub} P_v$$

H_{Hub} allows thermalisation of k -states

Toy Model for Thermalisation and Many-Body Scars

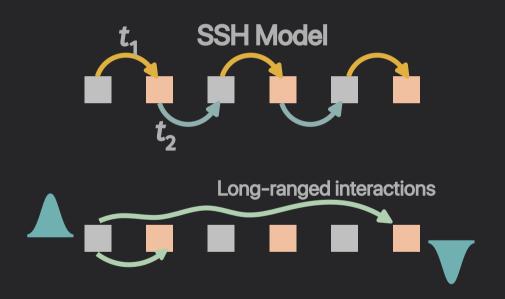


Consider HK Model perturbed by Hubbard interaction

$$H = H_{HKM} + P_v H_{Hub} P_v$$

- H_{Hub} allows thermalisation of k -states
- P_v will preserve certain sectors.
 Scars?

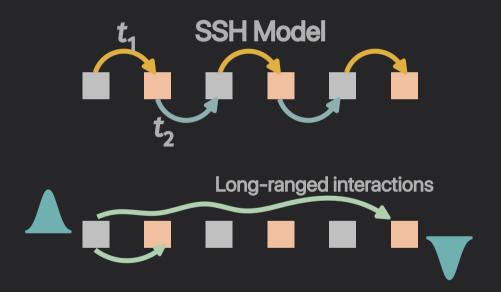
Effect of Hatsugai-Kohmoto Interaction on Edge Modes of SSH Chain



Add long-ranged HKM interactions to SSH chain.

$$H_{\text{SSH}} = -t_1 c_1^{\dagger} c_2^{} - t_2 c_2^{\dagger} c_3^{} + ...$$
 $H = H_{\text{HKM}} + H_{\text{SSH}}$

Effect of Hatsugai-Kohmoto Interaction on Edge Modes of SSH Chain



Add long-ranged HKM interactions to SSH chain.

$$H_{\text{SSH}} = -t_1 c_1^{\dagger} c_2^{} - t_2 c_2^{\dagger} c_3^{} + ...$$
 $H = H_{\text{HKM}} + H_{\text{SSH}}$

Fate of topological edge modes?

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