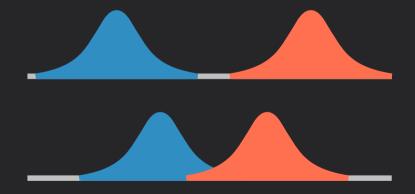
EXACTLY SOLVABLE MODEL OF CORRELATED METAL-INSULATOR TRANSITION

Insights on Non-Fermi Liquid and Mott Insulator

ABHIRUP MUKHERJEE

April 14, 2025 EPQM Seminar



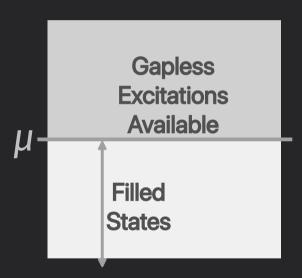
In A Nutshell

- 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 a generalisation of this model to obtain Fermi arcs

Where Do Mott Insulators and Non-Fermi Liquids Fit in the "Standard Model"?

Mott Insulators Are Different

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

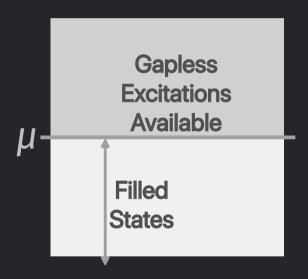


Dispersion away from band edge is **non-interacting**.

Mott (1968)

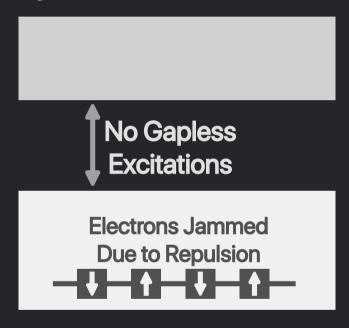
Mott Insulators Are Different

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



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

Add strong interactions - Mott Insulator!

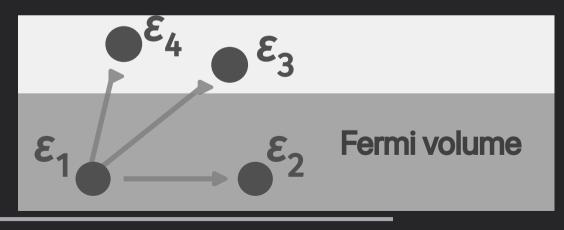


Gap opens inside
$$\begin{pmatrix} \varepsilon & U \\ U & \varepsilon \end{pmatrix} \rightarrow \varepsilon \pm U$$
 the band:

Mott (1968)

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 (quasiparticles).



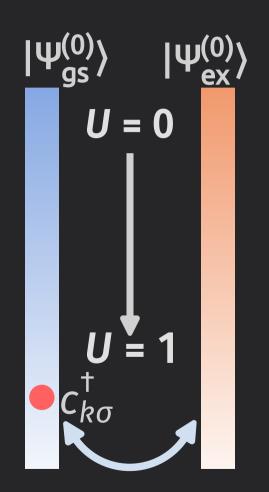
```
\Gamma
\sim \int d\varepsilon_4 d\varepsilon_3 d\varepsilon_2 \delta(\varepsilon - \varepsilon_2 - \varepsilon_3 - \varepsilon_4)
\sim \varepsilon^2
\rho \sim T^2
```

Landau Fermi Liquid Theory (Quantification)

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

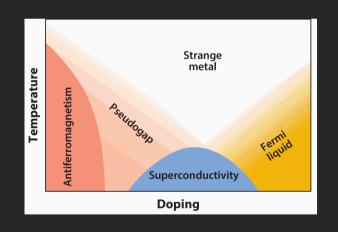
$$Z = \left\langle \Psi_{\rm ex} \middle| c_{k\sigma}^{\dagger} \middle| \Psi_{\rm gs} \right\rangle$$

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



Violations Of Landau Fermi Liquid Theory

- Tomonaga-Luttinger Liquid: Interacting electrons in 1D → spin-charge separation!
- Overscreened fixed points in Kondo models \rightarrow fractional entropy, diverging χ , C_v
- Strange Metal: Normal state of unconventional SCs in Cu oxides, heavy fermions, pnictides

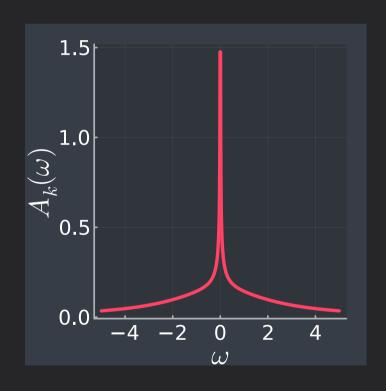


The Marginal Fermi Liquid

- Phenomenological explanation of normal state of cuprates: $\Sigma \sim \omega \log(|\omega|) i\pi |\omega|$
- 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.

$$H = -t \sum_{\langle i,j \rangle,\sigma} C_{i,\sigma}^{\dagger} C_{j,\sigma} + \frac{U}{L^{d}} \sum_{i_{1},i_{2},r} C_{i_{1}+r,\uparrow}^{\dagger} C_{i_{2}-r,\downarrow}^{\dagger} C_{i_{2},\downarrow} C_{i_{1},\uparrow}$$



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.

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













- 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)

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

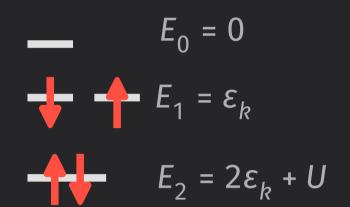
Each H_k can be diagonalised.

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

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

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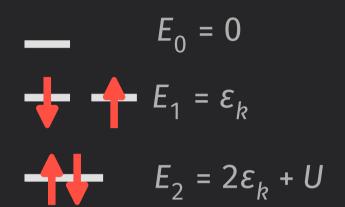
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Case Of Half-Filling

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

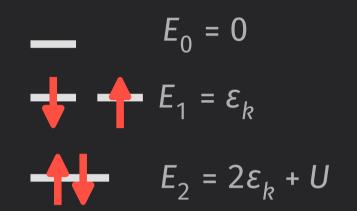
$$E_0 = 0$$
, $E_1 = \varepsilon_k - \frac{U}{2}$, $E_2 = 2\varepsilon_k$



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

Each H_k can be diagonalised.

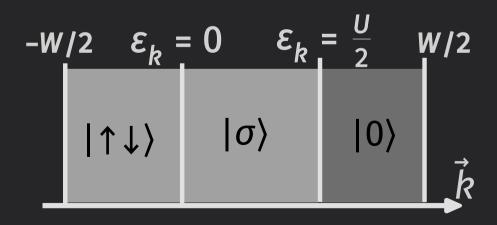
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Case Of Half-Filling

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

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



Introduction to Greens Functions

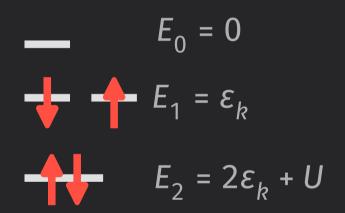
Nature of excitations can be studied through Greens function

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

- 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 as

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



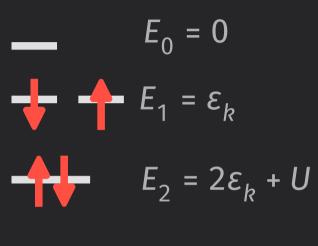
Greens function can be calculated as

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

If opposite spin is unoccupied

- Particle addition: $E_X = \varepsilon_k$
- Particle removal: $E_X = -\varepsilon_k$

$$G \to \frac{1 - \langle n_{k\overline{\sigma}} \rangle}{\omega - \varepsilon_k}$$



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

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

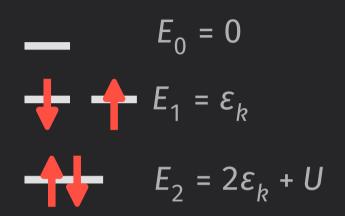
Greens function can be calculated as

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

If opposite spin is occupied

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

$$G \to \frac{\langle n_{k\overline{\sigma}} \rangle}{\omega - \varepsilon_k - U}$$

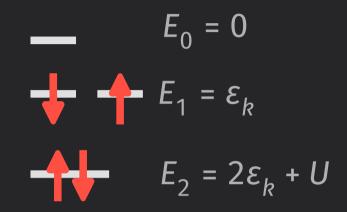


$$|-\sigma\rangle \rightarrow |\sigma, -\sigma\rangle$$

$$|\sigma, -\sigma\rangle \rightarrow |-\sigma\rangle$$

Greens function can be calculated as

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



Total Greens Function

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

Correlated Metal-Insulator Transition

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

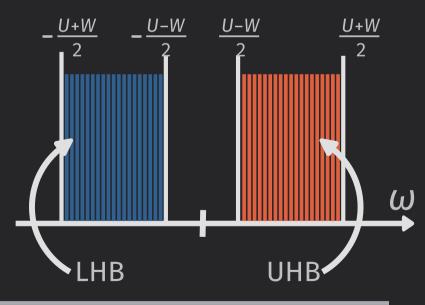
$$G_{k\sigma} = \frac{1}{2} \left[(\omega - \varepsilon_k + U/2)^{-1} + (\omega - \varepsilon_k - U/2)^{-1} \right]$$

Correlated Metal-Insulator Transition

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

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

U > W (Mott Insulator)

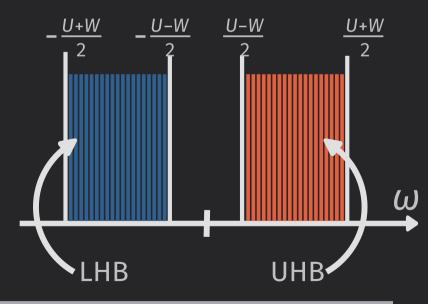


Correlated Metal-Insulator Transition

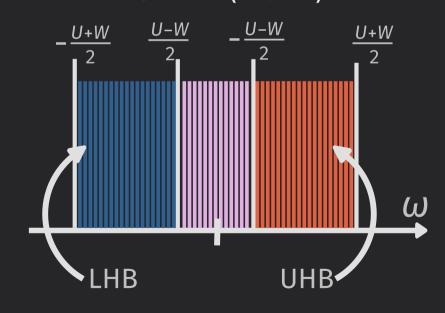
The Case Of **Half-Filling**:
$$2\mu = U$$
, $\langle n_{k\sigma} \rangle = \frac{1}{2}$

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

U > W (Mott Insulator)



U < W (Metal)

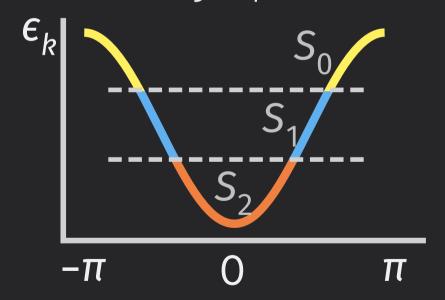


Non-Fermi Liquid Signatures In Metallic Phase

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

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 $\omega = 0: \langle n_k \rangle = 0$

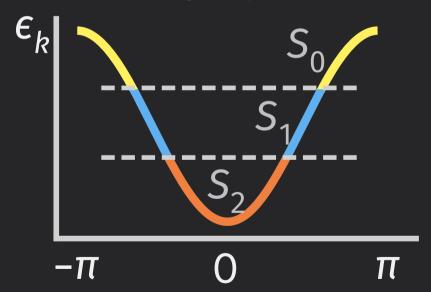


$$G_{k\sigma} = \frac{1}{2} \left[\left(\omega - \varepsilon_k + U/2 \right)^{-1} + \left(\omega - \varepsilon_k - U/2 \right)^{-1} \right]$$

Momentum states classified into three groups:

Ground state is a **mixed** state.

- k -states in S_2 are doubly-occupied.
- k –states in S_1 are half-filled.
- 2^{N_1} -fold degenerate.

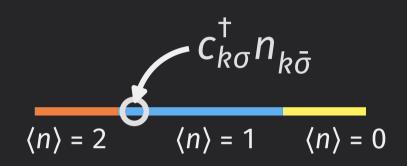


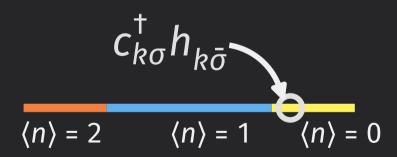
Near $S_2 - S_1$ boundary $(\varepsilon_k = -U/2)$

- Excitation operator: $c_{k\sigma}^{\dagger}n_{k\overline{\sigma}}$
- Excitation energy is $\varepsilon_k + \frac{u}{2} \to 0^+$

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

- Excitation operator: $c_{k\sigma}^{\dagger}(1-n_{k\overline{\sigma}})$
- Excitation energy is $\varepsilon_k \frac{u}{2} \to 0^+$





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

We have at 1/2-filling:

$$G_{k\sigma} = \frac{1}{\omega - \varepsilon_k + \frac{U^2/4}{\omega - \varepsilon_k}}.$$
Self-energy is $\Sigma(\omega) = \frac{U^2/4}{\omega - \varepsilon_k}$

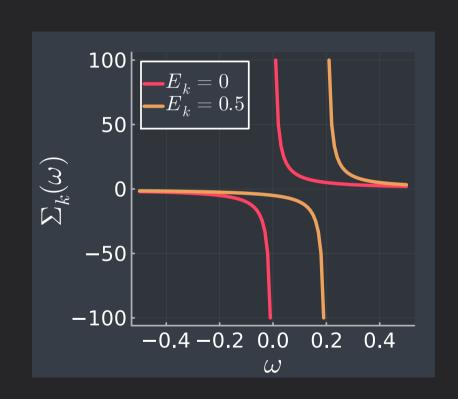
Self-energy is
$$\Sigma(\omega) = \frac{U^2/2}{\omega - \varepsilon_0}$$

We have at 1/2-filling:

$$G_{k\sigma} = \frac{1}{\omega - \varepsilon_k + \frac{U^2/4}{\omega - \varepsilon_k}}.$$

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

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



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

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.

Is This A Realistic Model of Interacting Electrons?

Stability Under Small Perturbations

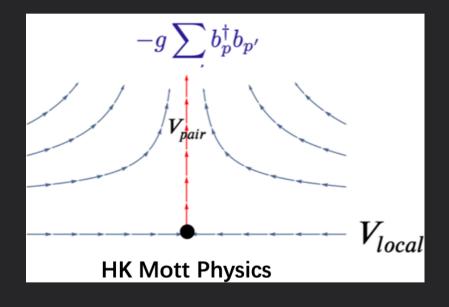
$$S = S_{FL} + U \int c_{k\uparrow}^{\dagger} c_{k\downarrow}^{\dagger} c_{k\downarrow} c_{k\uparrow} + U \int c_{k\uparrow}^{\dagger} c_{k\downarrow}^{\dagger} c_{k\downarrow} c_{k\uparrow} + U \int c_{k\downarrow}^{\dagger} c_{k\downarrow}^{\dagger} c_{k\downarrow}^{\dagger} c_{k\downarrow} c_{k\uparrow}^{\dagger} + U \int c_{k\downarrow}^{\dagger} c_{k\downarrow$$

Stability Under Small Perturbations

$$S = S_{FL} + U \int c_{k\uparrow}^{\dagger} c_{k\downarrow}^{\dagger} c_{k\downarrow} c_{k\uparrow} +$$

$$U \int c_{k_1+q,\uparrow}^{\dagger} c_{k_2-q,\downarrow}^{\dagger} c_{k_2,\downarrow} c_{k_1,\uparrow}$$

- HK interaction strongly relevant under RG flow
- Stable under local interaction terms
- Displays superconducting instability

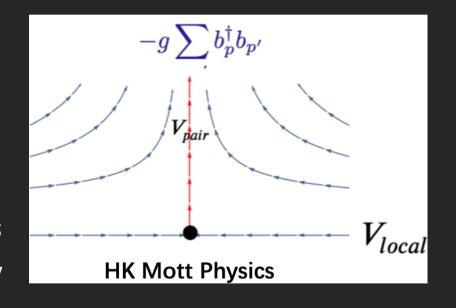


Stability Under Small Perturbations

$$S = S_{FL} + U \int c_{k\uparrow}^{\dagger} c_{k\downarrow}^{\dagger} c_{k\downarrow} c_{k\uparrow} +$$

$$U \int c_{k_1+q,\uparrow}^{\dagger} c_{k_2-q,\downarrow}^{\dagger} c_{k_2,\downarrow} c_{k_1,\uparrow}$$

- HK interaction strongly relevant under RG flow
- Stable under local interaction terms
- Displays superconducting instability



Key Point: HK interaction is the **source of Mottness** in Hubbard model.

Link With The 2D Hubbard Model

Alternative form of HK interaction:

$$U \sum_{k} n_{k\uparrow} n_{k\downarrow} \sim -J \sum_{k} \mathbf{S}_{k} \cdot \mathbf{S}_{k}$$

Baskaran Model

$$\sum_{k} \varepsilon_{k} n_{k\sigma} + \left(\sum_{k} \mathbf{S}_{k}\right)^{2} - J \sum_{k} \mathbf{S}_{k} \cdot \mathbf{S}_{k}$$

- Spin-charge separation of NFL
- Gapped spinon excitations

Link With The 2D Hubbard Model

Alternative form of HK interaction: $U \sum_{k} n_{k\uparrow} n_{k\downarrow} \sim -J \sum_{k} \mathbf{S}_{k} \cdot \mathbf{S}_{k}$

Baskaran Model

$$\sum_{k} \varepsilon_{k} n_{k\sigma} + \left(\sum_{k} S_{k}\right)^{2} - J \sum_{k} S_{k} \cdot S_{k}$$

- Spin-charge separation of NFL
- Gapped spinon excitations

Similar model emerges from 2D Hubbard Model, Mott liquid

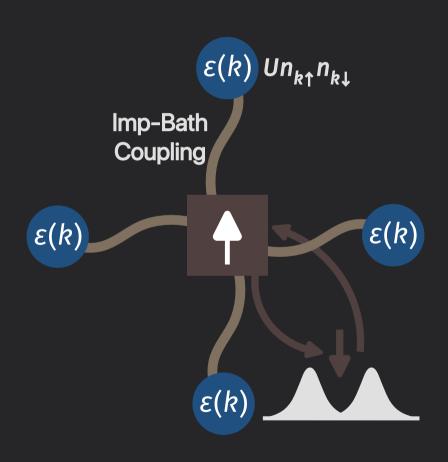
$$-\sum_{k} \mathbf{S}_{k} \cdot \mathbf{S}_{-k} + \sum_{k} \mathbf{C}_{k} \cdot \mathbf{C}_{-k}$$

$$S^z = n_{k\uparrow} - n_{k\uparrow\uparrow}, C^z = n_{k\uparrow} + n_{k\uparrow\uparrow} - 1$$



Avenues for Futher Investigation

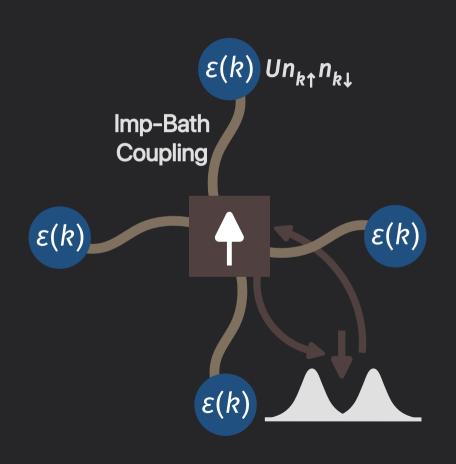
Kondo Screening in Hatsugai-Kohmoto Model



Consider local moment hybridising with HK Model

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

Kondo Screening in Hatsugai-Kohmoto Model

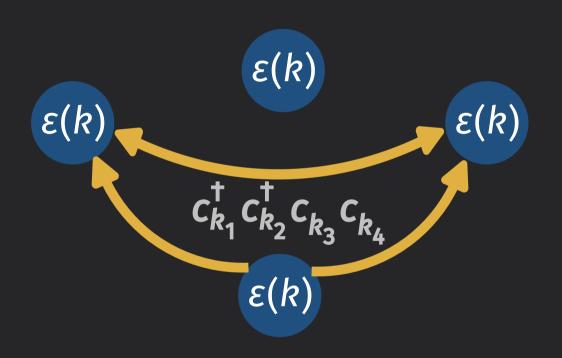


Consider local moment hybridising with HK Model

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

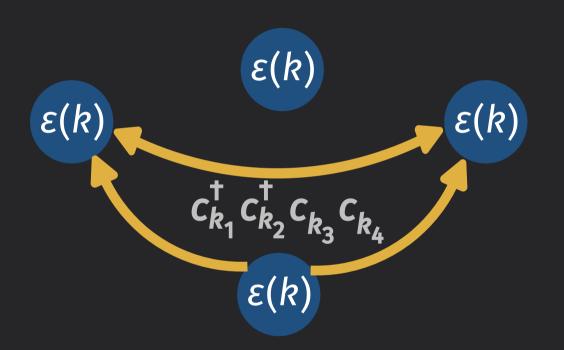
How does **absence** of quasiparticles affect Kondo screening?

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

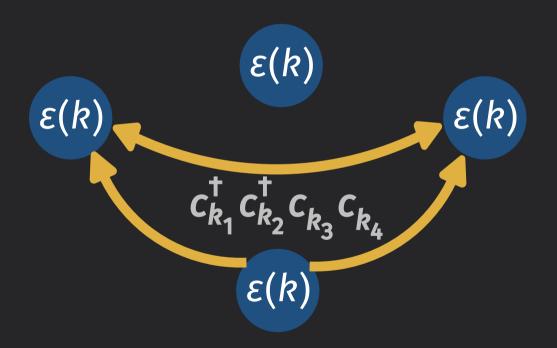


Consider HK Model perturbed by Hubbard interaction

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

• H_{Hub} allows **thermalisation** of k –states

Toy Model for Thermalisation and Many-Body Scars

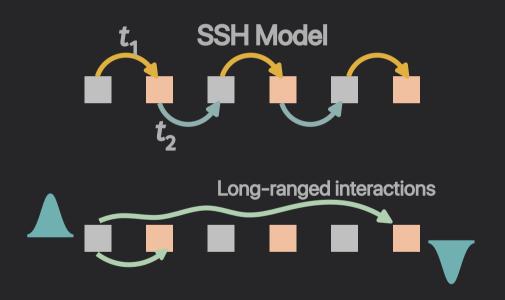


Consider HK Model perturbed by Hubbard interaction

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

- \blacksquare H_{Hub} allows **thermalisation** of k –states
- P_{ν} will preserve certain sectors. Scars?

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

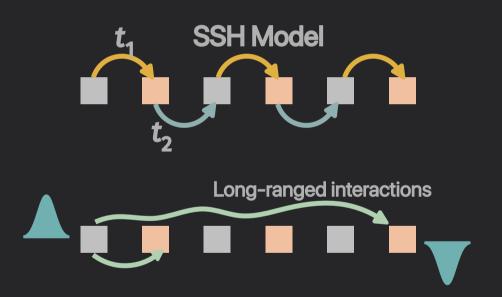


Add long-ranged HKM interactions to SSH chain.

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

 $H = H_{\text{HKM}} + H_{\text{SSH}}$

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



Add long-ranged HKM interactions to SSH chain.

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

Fate of topological edge modes?

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