Phase transitions in Quantum many-body System

Zhituo Wang

Harbin Institute of Technology

2022.07.28, The First Harbin-Moscow Conference on Analysis

 Example of phase transitions (PT) and statistical physics models.

 Example of phase transitions (PT) and statistical physics models.

► The 2-D Honeycomb-Hubbard model and Main results.

- Example of phase transitions (PT) and statistical physics models.
- ► The 2-D Honeycomb-Hubbard model and Main results.
- Sketch of the proof: fermionic cluster expansions and constructive renormalization theory.

- Example of phase transitions (PT) and statistical physics models.
- ► The 2-D Honeycomb-Hubbard model and Main results.
- Sketch of the proof: fermionic cluster expansions and constructive renormalization theory.

< □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > <

Hubbard model on the square lattice.

- Example of phase transitions (PT) and statistical physics models.
- ► The 2-D Honeycomb-Hubbard model and Main results.
- Sketch of the proof: fermionic cluster expansions and constructive renormalization theory.

- Hubbard model on the square lattice.
- Conclusions and perspectives.

- Example of phase transitions (PT) and statistical physics models.
- ► The 2-D Honeycomb-Hubbard model and Main results.
- Sketch of the proof: fermionic cluster expansions and constructive renormalization theory.

- Hubbard model on the square lattice.
- Conclusions and perspectives.
- References.

 Matters exhibit different macroscopic phases at different physical conditions.

 Matters exhibit different macroscopic phases at different physical conditions.

◆□▶ ◆□▶ ◆臣▶ ◆臣▶ 臣 の�?

• Consider the mercury (Hg) at 1 atm:

 Matters exhibit different macroscopic phases at different physical conditions.

◆□▶ ◆□▶ ◆三▶ ◆三▶ 三三 のへぐ

- Consider the mercury (Hg) at 1 atm:
 - $T \ge 356.7$ degree Celsius: Gas

 Matters exhibit different macroscopic phases at different physical conditions.

- Consider the mercury (Hg) at 1 atm:
 - $T \ge 356.7$ degree Celsius: Gas
 - ► -38.8 ≤ T ≤ 356.7; Liquid;

 Matters exhibit different macroscopic phases at different physical conditions.

- Consider the mercury (Hg) at 1 atm:
 - $T \ge 356.7$ degree Celsius: Gas
 - ► -38.8 ≤ T ≤ 356.7; Liquid;
 - ► T ≤ -38.8; Solid;

- Matters exhibit different macroscopic phases at different physical conditions.
- Consider the mercury (Hg) at 1 atm:
 - $T \ge 356.7$ degree Celsius: Gas
 - ► -38.8 ≤ T ≤ 356.7; Liquid;
 - ► T ≤ -38.8; Solid;
 - These are called the *first order* phase transitions.

- Matters exhibit different macroscopic phases at different physical conditions.
- Consider the mercury (Hg) at 1 atm:
 - $T \ge 356.7$ degree Celsius: Gas
 - ► -38.8 ≤ T ≤ 356.7; Liquid;
 - ► T ≤ -38.8; Solid;
 - These are called the *first order* phase transitions.

< □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > <

• $-268.8 \le T \le -38.8$; Normal conductors.

- Matters exhibit different macroscopic phases at different physical conditions.
- Consider the mercury (Hg) at 1 atm:
 - $T \ge 356.7$ degree Celsius: Gas
 - ► -38.8 ≤ T ≤ 356.7; Liquid;
 - ► T ≤ -38.8; Solid;
 - These are called the *first order* phase transitions.

- $-268.8 \le T \le -38.8$; Normal conductors.
- $T \leq -268.8$; Superconductor;

- Matters exhibit different macroscopic phases at different physical conditions.
- Consider the mercury (Hg) at 1 atm:
 - $T \ge 356.7$ degree Celsius: Gas
 - ► -38.8 ≤ T ≤ 356.7; Liquid;
 - ► T ≤ -38.8; Solid;
 - These are called the *first order* phase transitions.
 - $-268.8 \le T \le -38.8$; Normal conductors.
 - $T \leq -268.8$; Superconductor;
 - These are called the *second order* phase transition.

Microscopic particles are governed by the law of quantum mechanics. Study such system by solving Schrödinger equations?

- Microscopic particles are governed by the law of quantum mechanics. Study such system by solving Schrödinger equations?
 - ► Not possible. There are 10²³ such particles. Huge number of coupled Schrödinger equations.

- Microscopic particles are governed by the law of quantum mechanics. Study such system by solving Schrödinger equations?
 - Not possible. There are 10²³ such particles. Huge number of coupled Schrödinger equations.

< □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > <

• $N = 10^{23} = \infty$, and using probability theory:

- Microscopic particles are governed by the law of quantum mechanics. Study such system by solving Schrödinger equations?
 - Not possible. There are 10²³ such particles. Huge number of coupled Schrödinger equations.
- $N = 10^{23} = \infty$, and using probability theory:
 - Construct a many-body Hamiltonian on a suitable lattice (a model).

- Microscopic particles are governed by the law of quantum mechanics. Study such system by solving Schrödinger equations?
 - Not possible. There are 10²³ such particles. Huge number of coupled Schrödinger equations.
- $N = 10^{23} = \infty$, and using probability theory:
 - Construct a many-body Hamiltonian on a suitable lattice (a model).
 - ► (Formally) Define a finite measure on some suitable sample space, whose total mass is called the partition function Z_T.

- Microscopic particles are governed by the law of quantum mechanics. Study such system by solving Schrödinger equations?
 - Not possible. There are 10²³ such particles. Huge number of coupled Schrödinger equations.
- $N = 10^{23} = \infty$, and using probability theory:
 - Construct a many-body Hamiltonian on a suitable lattice (a model).
 - ► (Formally) Define a finite measure on some suitable sample space, whose total mass is called the partition function Z_T.
 - ▶ Phase transition happens when $T \leq T_c$, at which the partition function blows up.

- Microscopic particles are governed by the law of quantum mechanics. Study such system by solving Schrödinger equations?
 - Not possible. There are 10²³ such particles. Huge number of coupled Schrödinger equations.
- $N = 10^{23} = \infty$, and using probability theory:
 - Construct a many-body Hamiltonian on a suitable lattice (a model).
 - ► (Formally) Define a finite measure on some suitable sample space, whose total mass is called the partition function Z_T.
 - ▶ Phase transition happens when $T \leq T_c$, at which the partition function blows up.

Universal properties of different models.

Mathematical models in statistical physics

 The best known model is the Ising model (Ising 1924)," the drosophila (fruit flies) of statistical physics". Smirnov (2010), Duminil-Copin (2022).

Mathematical models in statistical physics

- The best known model is the Ising model (Ising 1924)," the drosophila (fruit flies) of statistical physics". Smirnov (2010), Duminil-Copin (2022).
- ► The Hubbard model (Hubbard, 1963) for correlated fermions, "the *tsetse flies* of Quantum many-body theory".

Mathematical models in statistical physics

- The best known model is the Ising model (Ising 1924)," the drosophila (fruit flies) of statistical physics". Smirnov (2010), Duminil-Copin (2022).
- ► The Hubbard model (Hubbard, 1963) for correlated fermions, "the *tsetse flies* of Quantum many-body theory".
- ► We will introduce some mathematically rigorous results in the Hubbard model in the 2-D Honeycomb lattice.

The Hubbard model on the 2-d Honeycomb lattice.



The honeycomb lattice Λ = Λ^A ∪ Λ^B is the superposition of the triangular lattice Λ^A (White dots) with Λ^B = Λ^A + δ_i (Black dots): δ₁ = (1,0), δ₂ = ½(−1, √3), δ₃ = ½(−1, −√3).

イロト 不得 トイヨト イヨト

The Hubbard model on the 2-d Honeycomb lattice.



The honeycomb lattice Λ = Λ^A ∪ Λ^B is the superposition of the triangular lattice Λ^A (White dots) with Λ^B = Λ^A + δ_i (Black dots): δ₁ = (1,0), δ₂ = ½(-1, √3), δ₃ = ½(-1, −√3).

► Let $L \in \mathbb{N}$, define the finite honeycomb lattice of side L: $\Lambda_L = \Lambda/L\Lambda$, $\lim_{L\to\infty} \Lambda_L = \Lambda$.

- The one-particle Hilbert space
 - $$\begin{split} \mathcal{H}_L &= \{ \Psi_{\mathbf{x},\alpha,\tau} : \Lambda_L \times \{A,B\} \times \{\uparrow,\downarrow\} \to \mathbb{C} \} \text{ such that } \\ \|\Psi\|_2^2 &= \sum_{\mathbf{x},\tau,\alpha} |\Psi_{\mathbf{x},\alpha,\tau}|^2 = 1, \ \Lambda_L = \Lambda/L\Lambda. \end{split}$$

The one-particle Hilbert space

$$\begin{array}{l} \mathcal{H}_{L} = \{\Psi_{\mathbf{x},\alpha,\tau} : \Lambda_{L} \times \{A,B\} \times \{\uparrow,\downarrow\} \rightarrow \mathbb{C} \} \text{ such that} \\ \|\Psi\|_{2}^{2} = \sum_{\mathbf{x},\tau,\alpha} |\Psi_{\mathbf{x},\alpha,\tau}|^{2} = 1, \ \Lambda_{L} = \Lambda/L\Lambda. \end{array}$$

• The Fermionic Fock space \mathcal{F}_L is:

$$\mathcal{F}_L = \mathbb{C} \oplus \bigoplus_{N=1}^{4L^2} \mathcal{F}_{\Lambda}^{(N)}, \quad \mathcal{F}_L^{(N)} = \bigwedge^N \mathcal{H}_L.$$

- The one-particle Hilbert space
 - $\begin{array}{l} \mathcal{H}_L = \{ \Psi_{\mathbf{x},\alpha,\tau} : \Lambda_L \times \{A,B\} \times \{\uparrow,\downarrow\} \rightarrow \mathbb{C} \ \} \text{ such that} \\ \|\Psi\|_2^2 = \sum_{\mathbf{x},\tau,\alpha} |\Psi_{\mathbf{x},\alpha,\tau}|^2 = 1, \ \Lambda_L = \Lambda/L\Lambda. \end{array}$

• The Fermionic Fock space \mathcal{F}_L is:

$$\mathcal{F}_L = \mathbb{C} \oplus \bigoplus_{N=1}^{4L^2} \mathcal{F}_{\Lambda}^{(N)}, \quad \mathcal{F}_L^{(N)} = \bigwedge^N \mathcal{H}_L.$$

• The Fermionic operators a^{\pm} , b^{\pm} on \mathcal{F}_{L} , $(\xi = (\mathbf{x}, \tau))$:

$$(a_{\mathbf{z},\tau}^{+}\Psi)^{(N)}(\xi_{1},\cdots,\xi_{N}) = \frac{1}{\sqrt{N}} \sum_{j=1}^{N} (-1)^{j} \delta_{\mathbf{z},\mathbf{x}_{j}} \delta_{\tau,\tau_{j}} \Psi^{(N-1)}(\xi_{1},\cdots,\xi_{j-1},\xi_{j+1},\cdots,\xi_{N}), (a_{\mathbf{z},\tau}^{-}\Psi)^{(N)}(\xi_{1},\cdots,\xi_{N}) = \sqrt{N+1} \Psi^{(N+1)}(\mathbf{z},\tau;\ \xi_{1},\cdots,\xi_{n}),$$

The one-particle Hilbert space

 $\begin{array}{l} \mathcal{H}_L = \{ \Psi_{\mathbf{x},\alpha,\tau} : \Lambda_L \times \{A,B\} \times \{\uparrow,\downarrow\} \rightarrow \mathbb{C} \ \} \text{ such that} \\ \|\Psi\|_2^2 = \sum_{\mathbf{x},\tau,\alpha} |\Psi_{\mathbf{x},\alpha,\tau}|^2 = 1, \ \Lambda_L = \Lambda/L\Lambda. \end{array}$

• The Fermionic Fock space \mathcal{F}_L is:

$$\mathcal{F}_L = \mathbb{C} \oplus \bigoplus_{N=1}^{4L^2} \mathcal{F}_{\Lambda}^{(N)}, \quad \mathcal{F}_L^{(N)} = \bigwedge^N \mathcal{H}_L.$$

• The Fermionic operators a^{\pm} , b^{\pm} on \mathcal{F}_L , $(\xi = (\mathbf{x}, \tau))$:

$$(a_{z,\tau}^{+}\Psi)^{(N)}(\xi_{1},\cdots,\xi_{N}) = \frac{1}{\sqrt{N}} \sum_{j=1}^{N} (-1)^{j} \delta_{z,x_{j}} \delta_{\tau,\tau_{j}} \Psi^{(N-1)}(\xi_{1},\cdots,\xi_{j-1},\xi_{j+1},\cdots,\xi_{N}),$$

$$(a_{z,\tau}^{-}\Psi)^{(N)}(\xi_{1},\cdots,\xi_{N}) = \sqrt{N+1} \Psi^{(N+1)}(z,\tau;\ \xi_{1},\cdots,\xi_{n}),$$

$$\blacktriangleright \text{ The CAR: } \{a_{x,\tau}^{+},a_{x',\tau'}^{-}\} = \delta_{x,x'} \delta_{\tau,\tau'},\ \{a_{x,\tau}^{+},a_{x',\tau'}^{+}\} = 0,$$

$$\{a_{x,\tau}^{-},a_{x',\tau'}^{-}\} = 0. \text{ The same for } b_{z,\tau}^{\pm}.$$

The Hubbard model on the honeycomb lattice

The Hubbard model Hamiltonian is:

$$\begin{aligned} H_{\Lambda_{L}} &= -t \sum_{\substack{\mathbf{x} \in \Lambda_{A} \\ i=1,2,3}} \sum_{\tau=\uparrow\downarrow} \left(a_{\mathbf{x},\tau}^{+} b_{\mathbf{x}+\vec{\delta}_{i},\tau}^{-} + b_{\mathbf{x}+\vec{\delta}_{i},\tau}^{+} a_{\mathbf{x},\tau}^{-} \right) \\ &- \mu \sum_{\mathbf{x} \in \Lambda_{A}} \sum_{\tau=\uparrow\downarrow} \left(a_{\mathbf{x},\tau}^{+} a_{\mathbf{x},\tau}^{-} + b_{\mathbf{x}+\vec{\delta}_{i},\tau}^{+} b_{\mathbf{x}+\vec{\delta}_{i},\tau}^{-} \right) \\ &+ \lambda \sum_{\mathbf{x} \in \Lambda_{A}} \left(a_{\mathbf{x},\uparrow}^{+} a_{\mathbf{x},\uparrow}^{-} a_{\mathbf{x},\downarrow}^{+} a_{\mathbf{x},\downarrow}^{-} + b_{\mathbf{x},\uparrow}^{+} b_{\mathbf{x},\uparrow}^{-} b_{\mathbf{x},\downarrow}^{+} b_{\mathbf{x},\downarrow}^{-} \right) \end{aligned}$$

• $t \in \mathbb{R}^+$ is called the hopping parameter, $\lambda \in \mathbb{R}$ is called the coupling constant, $\mu \in \mathbb{R}$ is called the chemical potential.

The Hubbard model on the honeycomb lattice

The Hubbard model Hamiltonian is:

$$\begin{aligned} H_{\Lambda_{L}} &= -t \sum_{\substack{\mathbf{x} \in \Lambda_{A} \\ i=1,2,3}} \sum_{\tau=\uparrow\downarrow} \left(a_{\mathbf{x},\tau}^{+} b_{\mathbf{x}+\vec{\delta}_{i},\tau}^{-} + b_{\mathbf{x}+\vec{\delta}_{i},\tau}^{+} a_{\mathbf{x},\tau}^{-} \right) \\ &- \mu \sum_{\mathbf{x} \in \Lambda_{A}} \sum_{\tau=\uparrow\downarrow} \left(a_{\mathbf{x},\tau}^{+} a_{\mathbf{x},\tau}^{-} + b_{\mathbf{x}+\vec{\delta}_{i},\tau}^{+} b_{\mathbf{x}+\vec{\delta}_{i},\tau}^{-} \right) \\ &+ \lambda \sum_{\mathbf{x} \in \Lambda_{A}} \left(a_{\mathbf{x},\uparrow}^{+} a_{\mathbf{x},\uparrow}^{-} a_{\mathbf{x},\downarrow}^{+} a_{\mathbf{x},\downarrow}^{-} + b_{\mathbf{x},\uparrow}^{+} b_{\mathbf{x},\uparrow}^{-} b_{\mathbf{x},\downarrow}^{+} b_{\mathbf{x},\downarrow}^{-} \right) \end{aligned}$$

t ∈ ℝ⁺ is called the hopping parameter, λ ∈ ℝ is called the coupling constant, μ ∈ ℝ is called the chemical potential.

• \vec{x} are the coordinates on the lattice, $\tau = \uparrow \downarrow$ are the spins.

The Hubbard model on the honeycomb lattice

The Hubbard model Hamiltonian is:

$$\begin{aligned} \mathcal{H}_{\Lambda_{L}} &= -t \sum_{\substack{\mathbf{x} \in \Lambda_{A} \\ i=1,2,3}} \sum_{\tau=\uparrow\downarrow} \left(a_{\mathbf{x},\tau}^{+} b_{\mathbf{x}+\vec{\delta}_{i},\tau}^{-} + b_{\mathbf{x}+\vec{\delta}_{i},\tau}^{+} a_{\mathbf{x},\tau}^{-} \right) \\ &- \mu \sum_{\mathbf{x} \in \Lambda_{A}} \sum_{\tau=\uparrow\downarrow} \left(a_{\mathbf{x},\tau}^{+} a_{\mathbf{x},\tau}^{-} + b_{\mathbf{x}+\vec{\delta}_{i},\tau}^{+} b_{\mathbf{x}+\vec{\delta}_{i},\tau}^{-} \right) \\ &+ \lambda \sum_{\mathbf{x} \in \Lambda_{A}} \left(a_{\mathbf{x},\uparrow}^{+} a_{\mathbf{x},\uparrow}^{-} a_{\mathbf{x},\downarrow}^{+} a_{\mathbf{x},\downarrow}^{-} + b_{\mathbf{x},\uparrow}^{+} b_{\mathbf{x},\uparrow}^{-} b_{\mathbf{x},\downarrow}^{+} b_{\mathbf{x},\downarrow}^{-} \right) \end{aligned}$$

- t ∈ ℝ⁺ is called the hopping parameter, λ ∈ ℝ is called the coupling constant, μ ∈ ℝ is called the chemical potential.
- \vec{x} are the coordinates on the lattice, $\tau = \uparrow \downarrow$ are the spins.
- ► When λ = 0, any fermion is only hopping to its nearest neighbor. When λ > 0, all fermions are correlated through the interaction term.

Why the honeycomb-Hubbard model?

It is easy to define yet highly nontrivial; It is mathematically challenging.

◆□▶ ◆□▶ ◆臣▶ ◆臣▶ 臣 の�?
Why the honeycomb-Hubbard model?

- It is easy to define yet highly nontrivial; It is mathematically challenging.
- ► This model captures the essence of many phenomena exhibited in the *Graphene* (Geim, Novoselov, 2004, Nobel Prize in Physics 2010), a mono-layer graphite, such as Dirac fermion, topological insulator, semi-metal, high-*T_c* superconductivity...

Time evolution and the correlation functions

► Define $\mathbf{a}_{x,\alpha}^{\pm}$, $\alpha = 1, 2$, s.t. $\mathbf{a}_{x,1}^{\pm} = \mathbf{a}_{x}^{\pm}$, $\mathbf{a}_{x,2}^{\pm} = \mathbf{b}_{x}^{\pm}$, the imaginary-time evolution of $\mathbf{a}_{x,\alpha}^{\pm}$ is: $\mathbf{a}_{x,\alpha}^{\pm} = e^{H_{\Lambda_{L}}x^{0}}\mathbf{a}_{x,\alpha}^{\pm}e^{-H_{\Lambda_{L}}x^{0}}$, $x = (x^{0}, \mathbf{x}), x_{0} \in [-\beta, \beta), \beta = 1/T$.

< □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > <

Time evolution and the correlation functions

► Define $\mathbf{a}_{x,\alpha}^{\pm}$, $\alpha = 1, 2$, s.t. $\mathbf{a}_{x,1}^{\pm} = \mathbf{a}_{x}^{\pm}$, $\mathbf{a}_{x,2}^{\pm} = \mathbf{b}_{x}^{\pm}$, the imaginary-time evolution of $\mathbf{a}_{x,\alpha}^{\pm}$ is: $\mathbf{a}_{x,\alpha}^{\pm} = e^{H_{\Lambda_{L}}x^{0}}\mathbf{a}_{x,\alpha}^{\pm}e^{-H_{\Lambda_{L}}x^{0}}$, $x = (x^{0}, \mathbf{x}), x_{0} \in [-\beta, \beta), \beta = 1/T$.

• The Gibbs state associated with the Hamiltonian H_{Λ_L} is: $\langle \cdot \rangle = \text{Tr}_{\mathcal{F}_L} \left[\cdot e^{-\beta H_{\Lambda_L}} \right] / Z_{\beta,\Lambda_L}, Z_{\beta,\Lambda_L} = \text{Tr}_{\mathcal{F}_L} e^{-\beta H_{\Lambda_L}}.$

Time evolution and the correlation functions

- ► Define $\mathbf{a}_{x,\alpha}^{\pm}$, $\alpha = 1, 2$, s.t. $\mathbf{a}_{x,1}^{\pm} = \mathbf{a}_{x}^{\pm}$, $\mathbf{a}_{x,2}^{\pm} = \mathbf{b}_{x}^{\pm}$, the imaginary-time evolution of $\mathbf{a}_{x,\alpha}^{\pm}$ is: $\mathbf{a}_{x,\alpha}^{\pm} = e^{H_{\Lambda_{L}}x^{0}}\mathbf{a}_{x,\alpha}^{\pm}e^{-H_{\Lambda_{L}}x^{0}}$, $x = (x^{0}, \mathbf{x}), x_{0} \in [-\beta, \beta), \beta = 1/T$.
- ► The Gibbs state associated with the Hamiltonian H_{Λ_L} is: $\langle \cdot \rangle = \text{Tr}_{\mathcal{F}_L} \left[\cdot e^{-\beta H_{\Lambda_L}} \right] / Z_{\beta,\Lambda_L}, Z_{\beta,\Lambda_L} = \text{Tr}_{\mathcal{F}_L} e^{-\beta H_{\Lambda_L}}.$
- the n-point Schwinger function is defined as:

$$S_{n,\beta}(\mathbf{x}_{1}, \alpha_{1}\cdots \mathbf{x}_{n}, \alpha_{n}, \lambda, \mu) = \lim_{L \to \infty} \langle \mathbf{T} \mathbf{a}_{\mathbf{x}_{1}, \alpha_{1}}^{\varepsilon_{1}} \cdots \mathbf{a}_{\mathbf{x}_{n}, \alpha_{n}}^{\varepsilon_{n}} \rangle_{\beta, L}$$
$$\mathbf{T} \mathbf{a}_{(\mathbf{x}_{1}, \mathbf{x}_{1}^{0}), \alpha_{1}}^{\varepsilon_{1}} \cdots \mathbf{a}_{(\mathbf{x}_{n}, \mathbf{x}_{n}^{0}), \alpha_{n}, \tau_{n}}^{\varepsilon_{n}}$$
$$= \operatorname{sgn}(\pi) \ \mathbf{a}_{(\mathbf{x}_{\pi(1)}, \mathbf{x}_{\pi(1)}^{0}), \alpha_{\pi(1)}}^{\varepsilon_{\pi(1)}} \cdots \mathbf{a}_{(\mathbf{x}_{\pi(n)}, \mathbf{x}_{\pi(n)}^{0}), \alpha_{\pi(n)}}^{\varepsilon_{\pi(n)}},$$

is the time-ordering operator, $sgn(\pi)$ is the sign of the permutation π , in which $x_{\pi(1)}^0 \ge x_{\pi(2)}^0 \ge \cdots \ge x_{\pi(n)}^0$.

• The most interesting quantities are:

- The most interesting quantities are:
 - The partition function: $Z_{\beta,\Lambda}(\lambda,\mu) = \lim_{L\to\infty} Z_{\beta,\Lambda_L}(\lambda,\mu)$,

(ロ)、(型)、(E)、(E)、 E) の(の)

- The most interesting quantities are:
 - The partition function: $Z_{\beta,\Lambda}(\lambda,\mu) = \lim_{L\to\infty} Z_{\beta,\Lambda_L}(\lambda,\mu)$,
 - The two-point Schwinger's function:

$$[S_{2,\beta}(x_1, x_2, \lambda, \mu)]_{\alpha_1, \alpha_2} = \lim_{L \to \infty} \langle \mathsf{Ta}_{x_1, \alpha_1, \tau_1}^{\varepsilon_1} \mathsf{a}_{x_2, \alpha_2, \tau_2}^{\varepsilon_2} \rangle_{\beta, L}$$

◆□▶ ◆□▶ ◆三▶ ◆三▶ 三三 のへぐ

- The most interesting quantities are:
 - The partition function: $Z_{\beta,\Lambda}(\lambda,\mu) = \lim_{L\to\infty} Z_{\beta,\Lambda_L}(\lambda,\mu)$,
 - The two-point Schwinger's function:

$$[S_{2,\beta}(x_1, x_2, \lambda, \mu)]_{\alpha_1, \alpha_2} = \lim_{L \to \infty} \langle \mathsf{Ta}_{x_1, \alpha_1, \tau_1}^{\varepsilon_1} \mathsf{a}_{x_2, \alpha_2, \tau_2}^{\varepsilon_2} \rangle_{\beta, L}$$

◆□▶ ◆□▶ ◆三▶ ◆三▶ 三三 のへぐ

► The connected Schwinger's function [S^c_{2,β}(λ, μ)]_{α1,α2} "cummulants of the Gibbs state"

- The most interesting quantities are:
 - The partition function: Z_{β,Λ}(λ, μ) = lim_{L→∞} Z_{β,Λ_L}(λ, μ),
 - The two-point Schwinger's function:

$$[S_{2,\beta}(x_1, x_2, \lambda, \mu)]_{\alpha_1, \alpha_2} = \lim_{L \to \infty} \langle \mathsf{Ta}_{x_1, \alpha_1, \tau_1}^{\varepsilon_1} \mathsf{a}_{x_2, \alpha_2, \tau_2}^{\varepsilon_2} \rangle_{\beta, L}$$

- ► The connected Schwinger's function [S^c_{2,β}(λ, μ)]_{α1,α2} "cummulants of the Gibbs state"
- The self-energy $[\Sigma_{2,\beta}(x_1, x_2, \lambda, \mu)]_{\alpha_1, \alpha_2}$

- The most interesting quantities are:
 - The partition function: Z_{β,Λ}(λ, μ) = lim_{L→∞} Z_{β,Λ_L}(λ, μ),
 - The two-point Schwinger's function:

$$[S_{2,\beta}(x_1, x_2, \lambda, \mu)]_{\alpha_1, \alpha_2} = \lim_{L \to \infty} \langle \mathsf{Ta}_{x_1, \alpha_1, \tau_1}^{\varepsilon_1} \mathsf{a}_{x_2, \alpha_2, \tau_2}^{\varepsilon_2} \rangle_{\beta, L}$$

- ► The connected Schwinger's function [S^c_{2,β}(λ, μ)]_{α1,α2} "cummulants of the Gibbs state"
- The self-energy $[\Sigma_{2,\beta}(x_1, x_2, \lambda, \mu)]_{\alpha_1, \alpha_2}$
- The fundamental questions are:

- The most interesting quantities are:
 - The partition function: Z_{β,Λ}(λ, μ) = lim_{L→∞} Z_{β,Λ_L}(λ, μ),
 - The two-point Schwinger's function:

$$[S_{2,\beta}(x_1, x_2, \lambda, \mu)]_{\alpha_1, \alpha_2} = \lim_{L \to \infty} \langle \mathsf{Ta}_{x_1, \alpha_1, \tau_1}^{\varepsilon_1} \mathsf{a}_{x_2, \alpha_2, \tau_2}^{\varepsilon_2} \rangle_{\beta, L}$$

- ► The connected Schwinger's function [S^c_{2,β}(λ, μ)]_{α1,α2} "cummulants of the Gibbs state"
- The self-energy $[\Sigma_{2,\beta}(x_1, x_2, \lambda, \mu)]_{\alpha_1, \alpha_2}$
- The fundamental questions are:
 - ► Is $\lim_{L\to\infty} \frac{Z_{\beta,\Lambda_L}(\lambda)}{Z_{\beta,\Lambda_L}(0)}$ a well-defined quantity, or can we rigorously define this model?

- The most interesting quantities are:
 - The partition function: Z_{β,Λ}(λ, μ) = lim_{L→∞} Z_{β,Λ_L}(λ, μ),
 - The two-point Schwinger's function:

$$[S_{2,\beta}(x_1, x_2, \lambda, \mu)]_{\alpha_1, \alpha_2} = \lim_{L \to \infty} \langle \mathsf{Ta}_{x_1, \alpha_1, \tau_1}^{\varepsilon_1} \mathsf{a}_{x_2, \alpha_2, \tau_2}^{\varepsilon_2} \rangle_{\beta, L}$$

- ► The connected Schwinger's function [S^c_{2,β}(λ, μ)]_{α1,α2} "cummulants of the Gibbs state"
- The self-energy $[\Sigma_{2,\beta}(x_1, x_2, \lambda, \mu)]_{\alpha_1, \alpha_2}$
- The fundamental questions are:
 - ► Is $\lim_{L\to\infty} \frac{Z_{\beta,\Lambda_L}(\lambda)}{Z_{\beta,\Lambda_L}(0)}$ a well-defined quantity, or can we rigorously define this model?

► The regularity of S_{2,β}(x₁, x₂, λ, μ) as a function of λ, β and the coordinates x₁, x₂.

The Hubbard model at $\mu = 0$.

Theorem (Giuliani, Mastropietro, 2010)

There exists a positive constant U such that the "pressure function" $\log \frac{Z_{\beta,\Lambda}(\lambda)}{Z_{\beta,\Lambda}(0)}$ and the connected Schwinger function $S_{2,\beta}^{c}(x_{1}, x_{2}, \lambda)$ are both analytic functions of λ when $\beta \to \infty$, for $|\lambda| \leq U$. The ground state is a fermi liquid up to T = 0.

Definition (Fermi liquid, Salmhofer, 1998)

Let $\hat{S}_{2,\beta}^{c}(k,\lambda)$ be the Fourier transform of $S_{2,\beta}^{c}(x_1, x_2, \lambda)$. The ground state of an interacting many-fermion system is said to be a Fermi liquid if

*Ŝ*_{2,β}(k, λ) is an analytic function of the coupling constant λ
 for β < ∞.
 </p>

くしゃ (雪) (雪) (雪) (雪) (雪) (

The Hubbard model at $\mu = 0$.

Theorem (Giuliani, Mastropietro, 2010)

There exists a positive constant U such that the "pressure function" $\log \frac{Z_{\beta,\Lambda}(\lambda)}{Z_{\beta,\Lambda}(0)}$ and the connected Schwinger function $S_{2,\beta}^{c}(x_{1}, x_{2}, \lambda)$ are both analytic functions of λ when $\beta \to \infty$, for $|\lambda| \leq U$. The ground state is a fermi liquid up to T = 0.

Definition (Fermi liquid, Salmhofer, 1998)

Let $\hat{S}_{2,\beta}^{c}(k,\lambda)$ be the Fourier transform of $S_{2,\beta}^{c}(x_1, x_2, \lambda)$. The ground state of an interacting many-fermion system is said to be a Fermi liquid if

- *Ŝ*_{2,β}(k, λ) is an analytic function of the coupling constant λ
 for β < ∞.
 </p>
- The self-energy $\hat{\Sigma}(k, \lambda)$ function is analytic in λ and uniformly C^2 in k for $\beta \to \infty$.

The Honeycomb-Hubbard model at $\mu = 1$, $\lambda = 0$

$$\hat{S}_{2,\beta}(k_0,\mathbf{k},0) = \frac{1}{k_0^2 + |\Omega(\mathbf{k})|^2 - \mu^2 - 2i\mu k_0} \begin{pmatrix} ik_0 + \mu & -\Omega^*(\mathbf{k}) \\ \Omega(\mathbf{k}) & ik_0 + \mu \end{pmatrix}$$

 $k_0 = (2n+1)\pi/\beta$, $\mathbf{k} = (k_1, k_2) \in \mathbb{R}^2/\Lambda$, $\Omega(\mathbf{k}) = 1 + 2e^{-i\frac{3}{2}k_1}\cos(\frac{\sqrt{3}}{2}k_2)$ is called the dispersion relation.

◆□▶ ◆□▶ ◆三▶ ◆三▶ 三三 のへぐ

The Honeycomb-Hubbard model at $\mu = 1$, $\lambda = 0$

$$\hat{S}_{2,eta}(k_0,\mathbf{k},0) = rac{1}{k_0^2 + |\Omega(\mathbf{k})|^2 - \mu^2 - 2i\mu k_0} egin{pmatrix} ik_0 + \mu & -\Omega^*(\mathbf{k}) \ \Omega(\mathbf{k}) & ik_0 + \mu \end{pmatrix}$$

 $k_0 = (2n+1)\pi/\beta$, $\mathbf{k} = (k_1, k_2) \in \mathbb{R}^2/\Lambda$, $\Omega(\mathbf{k}) = 1 + 2e^{-i\frac{3}{2}k_1}\cos(\frac{\sqrt{3}}{2}k_2)$ is called the dispersion relation.

• It is well defined for any $\beta < \infty$.

The Honeycomb-Hubbard model at $\mu = 1$, $\lambda = 0$

$$\hat{S}_{2,eta}(k_0,\mathbf{k},0) = rac{1}{k_0^2 + |\Omega(\mathbf{k})|^2 - \mu^2 - 2i\mu k_0} egin{pmatrix} ik_0 + \mu & -\Omega^*(\mathbf{k}) \ \Omega(\mathbf{k}) & ik_0 + \mu \end{pmatrix}$$

$$k_0 = (2n+1)\pi/\beta$$
, $\mathbf{k} = (k_1, k_2) \in \mathbb{R}^2/\Lambda$,
 $\Omega(\mathbf{k}) = 1 + 2e^{-i\frac{3}{2}k_1}\cos(\frac{\sqrt{3}}{2}k_2)$ is called the dispersion relation.

- It is well defined for any $\beta < \infty$.
- ▶ For $k_0 \rightarrow 0$ ($\beta \rightarrow \infty$), $\mu = 1$, $\hat{S}_{2,\beta}(k_0, \mathbf{k}, 0)$ is singular on the Fermi surface

$$\mathcal{F} = \{\mathbf{k} \in \mathbb{R}^2 / \Lambda, |\Omega(\mathbf{k})| - 1 = 0\}$$

= $\{(k_1, k_2), k_2 = \pm \frac{(2n+1)\pi}{\sqrt{3}}, n \in \mathbb{Z}\}$
 $\cup \{(k_1, k_2), k_2 = \pm \sqrt{3}k_1 \mp \frac{4n+2}{\sqrt{3}}\pi, n \in \mathbb{Z}\}.$

< □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > <

The Fermi surface



<□ > < @ > < E > < E > E のQ @

The Honeycomb Hubbard model at $\mu = 1$, $\lambda \neq 0$.

Theorem (Rivasseau, ZW 2021)

There exists a positive constants β_c = 1/T_c such that for any β ≤ β_c, the "pressure function" log Z_{β,Λ}(λ)/Z_{β,Λ}(0) and the connected two-point function S^c_{2,β}(λ) are analytic functions of the coupling constant λ, in the region

$$|\lambda \log^2 \beta| < 1. \tag{1}$$

The Honeycomb Hubbard model at $\mu = 1$, $\lambda \neq 0$.

Theorem (Rivasseau, ZW 2021)

There exists a positive constants β_c = 1/T_c such that for any β ≤ β_c, the "pressure function" log Z_{β,Λ}(λ)/Z_{β,Λ}(0) and the connected two-point function S^c_{2,β}(λ) are analytic functions of the coupling constant λ, in the region

$$|\lambda \log^2 \beta| < 1. \tag{1}$$

くしゃ (雪) (雪) (雪) (雪) (雪) (

 Fix λ, with |λ| < 1, the transition temperature is
 T_c = C₁e^{-C₂}/_{|λ|^{1/2}}, C₁, C₂ > 0 are two strictly positive constants.
 The Honeycomb Hubbard model at $\mu = 1$, $\lambda \neq 0$.

Theorem (Rivasseau, ZW 2021)

There exists a positive constants β_c = 1/T_c such that for any β ≤ β_c, the "pressure function" log Z_{β,Λ}(λ)/Z_{β,Λ}(0) and the connected two-point function S^c_{2,β}(λ) are analytic functions of the coupling constant λ, in the region

$$|\lambda \log^2 \beta| < 1. \tag{1}$$

- Fix λ, with |λ| < 1, the transition temperature is
 T_c = C₁e^{-C₂/|λ|^{1/2}}, C₁, C₂ > 0 are two strictly positive constants.
- The self-energy function Σ̂(k, λ) is C^{1+ε} differentiable w.r.t. the momentum, so that the ground state is not a Fermi liquid.

• The Hamiltonian H_{Λ} is unbounded. No simple way of prove if $e^{-\beta H_{\Lambda}}$ is trace class. Have to use perturbation theory.

◆□▶ ◆□▶ ◆臣▶ ◆臣▶ 臣 のへぐ

- The Hamiltonian H_{Λ} is unbounded. No simple way of prove if $e^{-\beta H_{\Lambda}}$ is trace class. Have to use perturbation theory.
- Typical term in the perturbation series of $S_{2,\beta}^{c}(\lambda)$ is

$$\int dk \cdots [\hat{S}_{2,\beta}(k,0)]^p.$$

$$\hat{S}_{2,\beta}(k,0) = \frac{1}{k_0^2 + |\Omega(\mathbf{k})|^2 - \mu^2 - 2i\mu k_0} \begin{pmatrix} ik_0 + \mu & -\Omega^*(\mathbf{k}) \\ \Omega(\mathbf{k}) & ik_0 + \mu \end{pmatrix}$$

is singular on the Fermi surfaces \mathcal{F} . Small divisor problem.

(日) (日) (日) (日) (日) (日) (日) (日)

- The Hamiltonian H_Λ is unbounded. No simple way of prove if e^{-βH_Λ} is trace class. Have to use perturbation theory.
- Typical term in the perturbation series of S^c_{2,β}(λ) is

$$\int dk \cdots [\hat{S}_{2,\beta}(k,0)]^p.$$

$$\hat{S}_{2,eta}(k,0) = rac{1}{k_0^2 + |\Omega(\mathbf{k})|^2 - \mu^2 - 2i\mu k_0} egin{pmatrix} ik_0 + \mu & -\Omega^*(\mathbf{k}) \ \Omega(\mathbf{k}) & ik_0 + \mu \end{pmatrix}$$

is singular on the Fermi surfaces \mathcal{F} . Small divisor problem. • Each $\hat{S}_{2,\beta}(k,0)$ is locally L^1 but not L^p for $p \ge 2$.

- The Hamiltonian H_Λ is unbounded. No simple way of prove if e^{-βH_Λ} is trace class. Have to use perturbation theory.
- Typical term in the perturbation series of S^c_{2,β}(λ) is

$$\int dk \cdots [\hat{S}_{2,\beta}(k,0)]^p.$$

$$\hat{\mathcal{S}}_{2,eta}(k,0) = rac{1}{k_0^2 + |\Omega(\mathbf{k})|^2 - \mu^2 - 2i\mu k_0} egin{pmatrix} ik_0 + \mu & -\Omega^*(\mathbf{k}) \ \Omega(\mathbf{k}) & ik_0 + \mu \end{pmatrix}$$

is singular on the Fermi surfaces \mathcal{F} . Small divisor problem.

- Each $\hat{S}_{2,\beta}(k,0)$ is locally L^1 but not L^p for $p \ge 2$.
- $\Omega(\mathbf{k})$ and μ are deformed by interactions, \mathcal{F} is not fixed.

- The Hamiltonian H_Λ is unbounded. No simple way of prove if e^{-βH_Λ} is trace class. Have to use perturbation theory.
- Typical term in the perturbation series of S^c_{2,β}(λ) is

$$\int dk \cdots [\hat{S}_{2,\beta}(k,0)]^p.$$

$$\hat{\mathcal{S}}_{2,eta}(k,0) = rac{1}{k_0^2 + |\Omega(\mathbf{k})|^2 - \mu^2 - 2i\mu k_0} egin{pmatrix} ik_0 + \mu & -\Omega^*(\mathbf{k}) \ \Omega(\mathbf{k}) & ik_0 + \mu \end{pmatrix}$$

is singular on the Fermi surfaces \mathcal{F} . Small divisor problem.

- Each $\hat{S}_{2,\beta}(k,0)$ is locally L^1 but not L^p for $p \ge 2$.
- $\Omega(\mathbf{k})$ and μ are deformed by interactions, \mathcal{F} is not fixed.
- The perturbation series can be unbounded.

Proof of the main theorem -The Grassmann algebra

The Grassmann algebra Gra is an associative, non-commutative, nilpotent algebra generated by the Grassmann variables {ψ^ε_{k,α}}, k = (k₀, k), ε = ±, α = 1, 2 such that ψ^ε_{k,α}ψ^{ε'}_{k',α'} = -ψ^{ε'}_{k',α'}ψ^ε_{k,α} and (ψ^ε_{k,α})² = 0.

Proof of the main theorem -The Grassmann algebra

- ▶ The Grassmann algebra **Gra** is an associative, non-commutative, nilpotent algebra generated by the Grassmann variables $\{\hat{\psi}_{k,\alpha}^{\varepsilon}\}$, $k = (k_0, \mathbf{k})$, $\varepsilon = \pm$, $\alpha = 1, 2$ such that $\hat{\psi}_{k,\alpha}^{\varepsilon}\hat{\psi}_{k',\alpha'}^{\varepsilon'} = -\hat{\psi}_{k',\alpha'}^{\varepsilon'}\hat{\psi}_{k,\alpha}^{\varepsilon}$ and $(\hat{\psi}_{k,\alpha}^{\varepsilon})^2 = 0$.
- ► The Grassmann differentiation and integral: $\partial_{\hat{\psi}_{k,\alpha}^{\varepsilon}}\hat{\psi}_{k',\alpha'}^{\varepsilon'} = \delta_{k,k'}\delta_{\alpha,\alpha'}\delta_{\varepsilon,\varepsilon'}, \int \hat{\psi}_{k,\alpha}^{\varepsilon}d\hat{\psi}_{k',\alpha'}^{\varepsilon'} = \delta_{k,k'}\delta_{\alpha,\alpha'}\delta_{\varepsilon,\varepsilon'}$

Proof of the main theorem -The Grassmann algebra

- ▶ The Grassmann algebra **Gra** is an associative, non-commutative, nilpotent algebra generated by the Grassmann variables $\{\hat{\psi}_{k,\alpha}^{\varepsilon}\}$, $k = (k_0, \mathbf{k})$, $\varepsilon = \pm$, $\alpha = 1, 2$ such that $\hat{\psi}_{k,\alpha}^{\varepsilon} \hat{\psi}_{k',\alpha'}^{\varepsilon'} = -\hat{\psi}_{k',\alpha'}^{\varepsilon'} \hat{\psi}_{k,\alpha}^{\varepsilon}$ and $(\hat{\psi}_{k,\alpha}^{\varepsilon})^2 = 0$.
- ► The Grassmann differentiation and integral: $\partial_{\hat{\psi}_{k,\alpha}^{\varepsilon}}\hat{\psi}_{k',\alpha'}^{\varepsilon'} = \delta_{k,k'}\delta_{\alpha,\alpha'}\delta_{\varepsilon,\varepsilon'}, \int \hat{\psi}_{k,\alpha}^{\varepsilon} d\hat{\psi}_{k',\alpha'}^{\varepsilon'} = \delta_{k,k'}\delta_{\alpha,\alpha'}\delta_{\varepsilon,\varepsilon'}$

• $\int d\psi_{\alpha}^+ d\psi_b^- e^{-\psi_{\alpha}^+ C_{\alpha\beta}\psi_b^-} = C_{\alpha\beta}$ for $C_{\alpha\beta} \in \mathbb{C}$

The Grassmann algebra

The Grassmann Gaussian measure $P(d\psi)$ with covariance $\hat{C}(k) := \hat{S}_{2,\beta}(k,0)$ is defined by :

$$P(d\psi) = N^{-1}D\psi \cdot \exp\left\{-\frac{1}{|\Lambda_L|\beta}\sum_{k\in\mathcal{D}_{\beta,L},\tau=\uparrow\downarrow,\alpha=1,2}\hat{\psi}^+_{k,\tau,\alpha}\hat{C}(k)^{-1}\hat{\psi}^-_{k,\tau,\alpha}\right\}$$

where

$$N = \prod_{\mathbf{k}\in\mathcal{D}_L, \tau=\uparrow\downarrow} \frac{1}{\beta|\Lambda_L|} \begin{pmatrix} -ik_0 - 1 & -\Omega^*(\mathbf{k}) \\ -\Omega(\mathbf{k}) & -ik_0 - 1 \end{pmatrix},$$

 $\mathcal{D}_{\beta,L} = \mathcal{D}_{\beta} \times \mathcal{D}_{L}$, $\mathcal{D}_{\beta} = \{\frac{2\pi}{\beta}(n + \frac{1}{2}), n \in \mathbb{N}\}$, \mathcal{D}_{L} is the dual space of Λ_{L} . We have:

$$\lim_{L \to \infty} \int P(d\psi) \hat{\psi}_{k_1,\tau_1,\alpha_1}^- \hat{\psi}_{k_2,\tau_2,\alpha_2}^+ = \delta_{k_1,k_2} \delta_{\tau_1,\tau_2} [\hat{C}(k_1)]_{\alpha_1,\alpha_2}.$$

・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・

► Define the Grassmann fields $\psi_{x,\tau,\alpha}^{\pm} = \frac{1}{\beta |\Lambda_L|} \sum_{k \in D_{\beta,L}} e^{\pm ikx} \hat{\psi}_{k,\tau,\alpha}^{\pm}, \quad x \in \Lambda_{\beta,L} := [-\beta,\beta) \times \Lambda_L,$

◆□▶ ◆□▶ ◆三▶ ◆三▶ 三三 のへぐ

 Define the Grassmann fields
 ψ[±]_{x,τ,α} = 1/β|Λ_L| Σ_{k∈D_{β,L}} e^{±ikx}ψ[±]_{k,τ,α}, x ∈ Λ_{β,L} := [-β, β) × Λ_L,

 the interacting potential becomes:

$$\mathcal{V}(\psi) = \lambda \sum_{lpha, lpha'=1,2} \int_{\Lambda_{eta,L}} d^3x \, \psi^+_{x,\uparrow,lpha} \psi^-_{x,\uparrow,lpha'} \psi^+_{x,\downarrow,lpha} \psi^-_{x,\downarrow,lpha'} \; ,$$

・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・

 Define the Grassmann fields
 ψ[±]_{x,τ,α} = 1/β|Λ_L| Σ_{k∈D_{β,L}} e^{±ikx} ψ[±]_{k,τ,α}, x ∈ Λ_{β,L} := [-β, β) × Λ_L,

 the interacting potential becomes:

$$\mathcal{V}(\psi) = \lambda \sum_{lpha, lpha'=1,2} \int_{\Lambda_{eta,L}} d^3 x \, \psi^+_{x,\uparrow,lpha} \psi^-_{x,\uparrow,lpha'} \psi^+_{x,\downarrow,lpha} \psi^-_{x,\downarrow,lpha'} \; ,$$

・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・

► The normalized Grassmann measure $\frac{1}{Z}P(d\psi)e^{-\mathcal{V}(\psi)}$, $Z = \int P(d\psi)e^{-\mathcal{V}(\psi)}$ is the partition function.

 Define the Grassmann fields
 ψ[±]_{x,τ,α} = 1/β|Λ_L| Σ_{k∈D_{β,L}} e^{±ikx} ψ[±]_{k,τ,α}, x ∈ Λ_{β,L} := [-β, β) × Λ_L,

 the interacting potential becomes:

$$\mathcal{V}(\psi) = \lambda \sum_{lpha, lpha'=1,2} \int_{\Lambda_{eta,L}} d^3 x \, \psi^+_{x,\uparrow,lpha} \psi^-_{x,\uparrow,lpha'} \psi^+_{x,\downarrow,lpha} \psi^-_{x,\downarrow,lpha'} \; ,$$

► The normalized Grassmann measure $\frac{1}{Z}P(d\psi)e^{-\mathcal{V}(\psi)}$, $Z = \int P(d\psi)e^{-\mathcal{V}(\psi)}$ is the partition function.

► The Schwinger functions:

$$S_{n,\beta}(x_1,\cdots,x_n)=\frac{1}{Z}\int\psi_{x_1,\tau_1,\alpha_1}^{\epsilon_1}\cdots\psi_{x_n,\tau_n,\alpha_n}^{\epsilon_n}e^{-\lambda\mathcal{V}(\psi)}P(d\psi).$$

・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・
 ・

Generating functionals

► Let j^+, j^- be two Grassmann fields. Define $Z(j^+, j^-) = \int e^{-\lambda V(\psi) + \int dx \psi^+(x) j^-(x) + + \int dx j^+(x) \psi^-(x)} P(d\psi)$, the generating functional of the Schwinger's functions.

Generating functionals

- ► Let j^+, j^- be two Grassmann fields. Define $Z(j^+, j^-) = \int e^{-\lambda V(\psi) + \int dx \psi^+(x) j^-(x) + + \int dx j^+(x) \psi^-(x)} P(d\psi)$, the generating functional of the Schwinger's functions.
- W(j[±]) = -¹/_{β|Λ|} log Z(j[±]) is the generating functional of the connected Schwinger's functions.
Generating functionals

- ► Let j^+, j^- be two Grassmann fields. Define $Z(j^+, j^-) = \int e^{-\lambda V(\psi) + \int dx \psi^+(x) j^-(x) + \int dx j^+(x) \psi^-(x)} P(d\psi)$, the generating functional of the Schwinger's functions.
- W(j[±]) = −¹/_{β|Λ|} log Z(j[±]) is the generating functional of the connected Schwinger's functions.
- ► The connected 2*n*-point Schwinger's functions:

$$S_{2n}^c(x_1,\cdots,x_n,y_1,\cdots,y_n) = \frac{\delta^{2n}}{\delta j^+(x_1)\cdots\delta j^+(x_n)\delta j^-(y_1)\cdots\delta j^-(y_n)} W(j^+,j^-)|_{j^\pm=0},$$

Generating functionals

- Let j^+, j^- be two Grassmann fields. Define $Z(j^+, j^-) = \int e^{-\lambda V(\psi) + \int dx \psi^+(x) j^-(x) + + \int dx j^+(x) \psi^-(x)} P(d\psi)$, the generating functional of the Schwinger's functions.
- W(j[±]) = −¹/_{β|Λ|} log Z(j[±]) is the generating functional of the connected Schwinger's functions.
- ▶ The connected 2*n*-point Schwinger's functions:

$$S_{2n}^{c}(x_1,\cdots,x_n,y_1,\cdots,y_n) = \frac{\delta^{2n}}{\delta j^+(x_1)\cdots\delta j^+(x_n)\delta j^-(y_1)\cdots\delta j^-(y_n)} W(j^+,j^-)|_{j^{\pm}=0},$$

► Define $\phi^+(x) = \frac{\delta}{\delta j^-(x)} W(j)$, $\phi^-(x) = \frac{\delta}{\delta j^+(x)} W(j)$. $\Gamma(\phi^+, \phi^-) = W(j^+, j^-) - \int d^3x \left[j^+(x)\phi^-(x) + \phi^+(x)j^-(x)\right]$ is defined as the Legendre transform of W.

Generating functionals

- ► Let j^+, j^- be two Grassmann fields. Define $Z(j^+, j^-) = \int e^{-\lambda V(\psi) + \int dx \psi^+(x) j^-(x) + \int dx j^+(x) \psi^-(x)} P(d\psi)$, the generating functional of the Schwinger's functions.
- W(j[±]) = −¹/_{β|Λ|} log Z(j[±]) is the generating functional of the connected Schwinger's functions.
- ▶ The connected 2*n*-point Schwinger's functions:

$$S_{2n}^c(x_1,\cdots,x_n,y_1,\cdots,y_n) = \frac{\delta^{2n}}{\delta j^+(x_1)\cdots\delta j^+(x_n)\delta j^-(y_1)\cdots\delta j^-(y_n)} W(j^+,j^-)|_{j^{\pm}=0},$$

- ► Define $\phi^+(x) = \frac{\delta}{\delta j^-(x)} W(j)$, $\phi^-(x) = \frac{\delta}{\delta j^+(x)} W(j)$. $\Gamma(\phi^+, \phi^-) = W(j^+, j^-) - \int d^3x \left[j^+(x)\phi^-(x) + \phi^+(x)j^-(x)\right]$ is defined as the Legendre transform of W.
- ► The self-energy $\Sigma(x, y) = \frac{\delta^2}{\delta \phi^+(x) \delta \phi^-(y)} \Gamma(\phi^+, \phi^-)|_{\phi^\pm = 0}$

The partition function

For $|\lambda| < 1$, we do perturbation expansions:

$$Z(\lambda) = S_0(\lambda) = \int P(d\psi) e^{\lambda \int_{\Lambda_{\beta,L}} d^3 x \left[\psi^+_{x,\uparrow} \psi^-_{x,\downarrow} \psi^+_{x,\downarrow} \psi^-_{x,\downarrow}\right]}$$
$$= \sum_{n=0}^{\infty} \frac{\lambda^n}{n!} \int P(d\psi) \left[\int_{\Lambda_{\beta}} d^3 x \left(\psi^+_{x,\uparrow} \psi^-_{x,\uparrow} \psi^+_{x,\downarrow} \psi^-_{x,\downarrow}\right)\right]^n.$$
$$= \sum_n \frac{\lambda^n}{n!} \int_{(\Lambda_{\beta,L})^n} d^3 x_1 \cdots d^3 x_n \begin{cases} x_{1,\varepsilon_1,\tau_1} \cdots x_{n,\varepsilon_n,\tau_n} \\ x_{1,\varepsilon_1,\tau_1} \cdots x_{n,\varepsilon_n,\tau_n} \end{cases},$$

 $\{ \cdot \}$ is a $2n \times 2n$ determinant, Cayley's notation:

$$\begin{cases} x_{i,\tau} \\ x_{j,\tau'} \end{cases} = \det \left[\delta_{\tau\tau'} \left[C(x_i - x_j) \right] \right], C(x - y) = \int_{\Lambda_{\beta,L}} \hat{C}(k) e^{ik(x-y)} d^3x$$

$$\hat{C}(k) = \frac{1}{k_0^2 + |\Omega(\mathbf{k})|^2 - \mu^2 - 2i\mu k_0} \begin{pmatrix} ik_0 + \mu & -\Omega^*(\mathbf{k}) \\ \Omega(\mathbf{k}) & ik_0 + \mu \end{pmatrix}$$

► Q1: Fully expansion of the determinant generates the combinatorial factor (2n)!, which makes the perturbation series divergent.

- Q1: Fully expansion of the determinant generates the combinatorial factor (2n)!, which makes the perturbation series divergent.
- Q2: Ĉ(k) is locally L¹ integrable but not L^p, ∀p ≥ 2;
 C(x − y) decays very slowly when the denominator of Ĉ(k) is close to zero.

- Q1: Fully expansion of the determinant generates the combinatorial factor (2n)!, which makes the perturbation series divergent.
- ► Q2: Ĉ(k) is locally L¹ integrable but not L^p, ∀p ≥ 2; C(x − y) decays very slowly when the denominator of Ĉ(k) is close to zero.
- Solution: partially expand the determinant (fermionic cluster expansions) so that only tree lines appear. Dividing the integral domain of Ĉ(k) into smaller regions (sectors), so that Ĉ(k) and its Fourier transform have optimal decaying property.

くしゃ (雪) (雪) (雪) (雪) (雪) (

• Q3: Due to interactions, $|\Omega(\mathbf{k})|^2 \rightarrow |\Omega(\mathbf{k})|^2 + \Sigma(k_0, \mathbf{k}, \lambda)$ and $\mu \rightarrow \mu + \tilde{\delta}\mu(\lambda)$. The interacting Fermi surface is given by

$$\mathcal{F} = \{ \mathbf{k} | |\Omega(\mathbf{k})|^2 - \mu - \widetilde{\delta}\mu(\lambda) - \Sigma(0, \mathbf{k}, \lambda) = 0 \}.$$

• Q3: Due to interactions, $|\Omega(\mathbf{k})|^2 \rightarrow |\Omega(\mathbf{k})|^2 + \Sigma(k_0, \mathbf{k}, \lambda)$ and $\mu \rightarrow \mu + \tilde{\delta}\mu(\lambda)$. The interacting Fermi surface is given by

$$\mathcal{F} = \{\mathbf{k} \mid |\Omega(\mathbf{k})|^2 - \mu - \tilde{\delta}\mu(\lambda) - \Sigma(0, \mathbf{k}, \lambda) = 0\}.$$

◆□▶ ◆□▶ ◆三▶ ◆三▶ 三三 のへぐ

Solution:

• Q3: Due to interactions,
$$|\Omega(\mathbf{k})|^2 \rightarrow |\Omega(\mathbf{k})|^2 + \Sigma(k_0, \mathbf{k}, \lambda)$$
 and $\mu \rightarrow \mu + \tilde{\delta}\mu(\lambda)$. The interacting Fermi surface is given by

$$\mathcal{F} = \{\mathbf{k} \mid |\Omega(\mathbf{k})|^2 - \mu - \tilde{\delta}\mu(\lambda) - \Sigma(0, \mathbf{k}, \lambda) = 0\}.$$

Solution:

Introducing counter-terms to fix the Fermi surface

$$\delta H_{\Lambda_{L}} = \delta \mu(\lambda) \sum_{k \in \mathcal{D}_{\beta,L}} \sum_{\alpha=1,2} \sum_{\tau=\uparrow,\downarrow} \hat{\psi}^{+}_{k,\tau,\alpha} \hat{\psi}^{-}_{k,\tau,\alpha} + \sum_{k \in \mathcal{D}_{\beta,L}, \tau=\uparrow\downarrow} \sum_{\alpha,\alpha'=1,2} \hat{\nu}(k_{0},\mathbf{k},\lambda) \hat{\psi}^{+}_{k,\tau,\alpha} \hat{\psi}^{-}_{k,\tau,\alpha'}$$
(2)

• Q3: Due to interactions,
$$|\Omega(\mathbf{k})|^2 \rightarrow |\Omega(\mathbf{k})|^2 + \Sigma(k_0, \mathbf{k}, \lambda)$$
 and $\mu \rightarrow \mu + \tilde{\delta}\mu(\lambda)$. The interacting Fermi surface is given by

$$\mathcal{F} = \{\mathbf{k} \mid |\Omega(\mathbf{k})|^2 - \mu - \widetilde{\delta}\mu(\lambda) - \Sigma(0, \mathbf{k}, \lambda) = 0\}.$$

Solution:

Introducing counter-terms to fix the Fermi surface

$$\delta H_{\Lambda_{L}} = \delta \mu(\lambda) \sum_{k \in \mathcal{D}_{\beta,L}} \sum_{\alpha=1,2} \sum_{\tau=\uparrow,\downarrow} \hat{\psi}^{+}_{k,\tau,\alpha} \hat{\psi}^{-}_{k,\tau,\alpha} + \sum_{k \in \mathcal{D}_{\beta,L}, \tau=\uparrow\downarrow} \sum_{\alpha,\alpha'=1,2} \hat{\nu}(k_{0},\mathbf{k},\lambda) \hat{\psi}^{+}_{k,\tau,\alpha} \hat{\psi}^{-}_{k,\tau,\alpha'}$$
(2)

• Choose $\delta\mu(\lambda)$ such that it cancels the term $\tilde{\delta}\mu(\lambda)$.

• Q3: Due to interactions,
$$|\Omega(\mathbf{k})|^2 \rightarrow |\Omega(\mathbf{k})|^2 + \Sigma(k_0, \mathbf{k}, \lambda)$$
 and $\mu \rightarrow \mu + \tilde{\delta}\mu(\lambda)$. The interacting Fermi surface is given by

$$\mathcal{F} = \{\mathbf{k} \mid |\Omega(\mathbf{k})|^2 - \mu - \widetilde{\delta}\mu(\lambda) - \Sigma(0, \mathbf{k}, \lambda) = 0\}.$$

Solution:

Introducing counter-terms to fix the Fermi surface

$$\delta H_{\Lambda_{L}} = \delta \mu(\lambda) \sum_{k \in \mathcal{D}_{\beta,L}} \sum_{\alpha=1,2} \sum_{\tau=\uparrow,\downarrow} \hat{\psi}^{+}_{k,\tau,\alpha} \hat{\psi}^{-}_{k,\tau,\alpha} + \sum_{k \in \mathcal{D}_{\beta,L}, \tau=\uparrow\downarrow} \sum_{\alpha,\alpha'=1,2} \hat{\nu}(k_{0},\mathbf{k},\lambda) \hat{\psi}^{+}_{k,\tau,\alpha} \hat{\psi}^{-}_{k,\tau,\alpha'}$$
(2)

- Choose $\delta\mu(\lambda)$ such that it cancels the term $\tilde{\delta}\mu(\lambda)$.
- Choose $\hat{\nu}(\mathbf{k}_0, \mathbf{k}, \lambda)$ that cancels $\hat{\Sigma}(0, P_F(\mathbf{k}), \lambda)$

• Q3: Due to interactions,
$$|\Omega(\mathbf{k})|^2 \rightarrow |\Omega(\mathbf{k})|^2 + \Sigma(k_0, \mathbf{k}, \lambda)$$
 and $\mu \rightarrow \mu + \tilde{\delta}\mu(\lambda)$. The interacting Fermi surface is given by

$$\mathcal{F} = \{\mathbf{k} \mid |\Omega(\mathbf{k})|^2 - \mu - \widetilde{\delta}\mu(\lambda) - \Sigma(0, \mathbf{k}, \lambda) = 0\}.$$

Solution:

Introducing counter-terms to fix the Fermi surface

$$\delta H_{\Lambda_{L}} = \delta \mu(\lambda) \sum_{k \in \mathcal{D}_{\beta,L}} \sum_{\alpha=1,2} \sum_{\tau=\uparrow,\downarrow} \hat{\psi}^{+}_{k,\tau,\alpha} \hat{\psi}^{-}_{k,\tau,\alpha} + \sum_{k \in \mathcal{D}_{\beta,L}, \tau=\uparrow\downarrow} \sum_{\alpha,\alpha'=1,2} \hat{\nu}(k_{0},\mathbf{k},\lambda) \hat{\psi}^{+}_{k,\tau,\alpha} \hat{\psi}^{-}_{k,\tau,\alpha'}$$
(2)

- Choose $\delta\mu(\lambda)$ such that it cancels the term $\tilde{\delta}\mu(\lambda)$.
- Choose $\hat{\nu}(k_0, \mathbf{k}, \lambda)$ that cancels $\hat{\Sigma}(0, P_F(\mathbf{k}), \lambda)$
- The cancellations are carried in the multi-scale representation using renormalization theory.

The multi-scale analysis

Let G^h₀(ℝ), h > 1, be the Gevrey class of compactly supported functions. Define a cutoff function χ ∈ G^h₀(ℝ) as:

$$\chi(t) = \chi(-t) = \begin{cases} = 0, & \text{for } |t| > 2, \\ \in (0, 1), & \text{for } 1 < |t| \le 2, \\ = 1, & \text{for } |t| \le 1. \end{cases}$$
(3)

◆□▶ ◆□▶ ◆臣▶ ◆臣▶ 臣 の�?

The multi-scale analysis

Let G^h₀(ℝ), h > 1, be the Gevrey class of compactly supported functions. Define a cutoff function χ ∈ G^h₀(ℝ) as:

$$\chi(t) = \chi(-t) = \begin{cases} = 0, & \text{for } |t| > 2, \\ \in (0, 1), & \text{for } 1 < |t| \le 2, \\ = 1, & \text{for } |t| \le 1. \end{cases}$$
(3)

• Given fixed constant $\gamma \geq 10$, construct a partition of unity

$$1 = \sum_{j=0}^{\infty} \chi_{j}(t), \quad \forall t \neq 0;$$

$$\chi_{0}(t) = 1 - \chi(t),$$

$$\chi_{j}(t) = \chi(\gamma^{2j-1}t) - \chi(\gamma^{2j}t) \text{ for } j \ge 1.$$
(4)

The multi-slice expansion

• The free propagator is decomposed as :

$$\hat{C}(k)_{\alpha\alpha'} = \sum_{j=0}^{\infty} \hat{C}_j(k)_{\alpha\alpha'}, \ \alpha, \alpha' = 1, 2,$$

$$\hat{C}_j(k)_{\alpha\alpha'} = \hat{C}(k)_{\alpha\alpha'} \cdot \chi_j[4k_0^2 + e^2(\mathbf{k})],$$
(5)

◆□▶ ◆□▶ ◆臣▶ ◆臣▶ 臣 のへぐ

$$e(\mathbf{k}) = 8[\cos(\sqrt{3}k_2/2)] \cdot [\cos(\frac{1}{4}(3k_1 + \sqrt{3}k_2))] \\ \cdot [\cos(\frac{1}{4}(3k_1 - \sqrt{3}k_2))].$$

The sectors

► Not sufficient to obtain the optimal decaying of propagator.

The sectors

- Not sufficient to obtain the optimal decaying of propagator.
- We introduce a second partition of unity:

$$1 = \sum_{s=0}^{j} v_{s}(t), \quad \begin{cases} v_{0}(t) = 1 - \chi(\gamma^{2}t), \\ v_{s}(t) = \chi_{s+1}(t), \\ v_{j}(t) = \chi(\gamma^{2j}t), \end{cases} \quad \text{for} \quad 1 \le s \le j-1, \ (6)$$

◆□▶ ◆□▶ ◆臣▶ ◆臣▶ 臣 の�?

The sectors

- Not sufficient to obtain the optimal decaying of propagator.
- We introduce a second partition of unity:

$$1 = \sum_{s=0}^{j} v_{s}(t), \quad \begin{cases} v_{0}(t) = 1 - \chi(\gamma^{2}t), \\ v_{s}(t) = \chi_{s+1}(t), \\ v_{j}(t) = \chi(\gamma^{2j}t), \end{cases} \quad \text{for} \quad 1 \le s \le j-1, \ (6)$$

◆□▶ ◆□▶ ◆臣▶ ◆臣▶ 臣 のへぐ

►
$$\hat{C}_{j}(k) = \sum_{\sigma = (s_{a}, s_{b})} \hat{C}_{j,\sigma}(k), \ \hat{C}_{j,\sigma}(k) = \hat{C}_{j}(k) \cdot v_{s_{a}}[t_{a}] \ v_{s_{b}}[t_{b}],$$

 $a, b \in \{1, 2, 3\},$
 $t_{1} = \cos^{2}(\sqrt{3}k_{2}/2), \ t_{2} = \cos^{2}(\frac{1}{4}(3k_{1} + \sqrt{3}k_{2})),$
 $t_{3} = \cos^{2}(\frac{1}{4}(3k_{1} - \sqrt{3}k_{2})).$



Figure: An illustration of the various sectors.

◆□▶ ◆□▶ ◆臣▶ ◆臣▶ 三臣 - のへで

The bounds for the propagators

► Let $[C_{j,\sigma}(x - y)]_{\alpha\alpha'}$ be the Fourier transform of $[\hat{C}_{j,\sigma}(k_0, \mathbf{k})]_{\alpha\alpha'}$,

The bounds for the propagators

► Let $[C_{j,\sigma}(x-y)]_{\alpha\alpha'}$ be the Fourier transform of $[\hat{C}_{j,\sigma}(k_0, \mathbf{k})]_{\alpha\alpha'}$, ► $\|C_{j,\sigma}(x-y)]_{\alpha\alpha'}\|_{L^{\infty}} \leq O(1)\gamma^{-s_a-s_b} e^{-c[d_{j,\sigma}(x,y)]^{\alpha_0}}$, where $s_a, s_b \in [0, j]$, $\alpha_0 = 1/h$, and $d_{j,\sigma}(x, y) = \gamma^{-j}|x_0 - y_0| + \gamma^{-s_a}|x_a - y_a| + \gamma^{-s_b}|x_b - y_b|$

The bounds for the propagators

► Let
$$[C_{j,\sigma}(x-y)]_{\alpha\alpha'}$$
 be the Fourier transform of
 $[\hat{C}_{j,\sigma}(k_0, \mathbf{k})]_{\alpha\alpha'}$,
 $\|C_{j,\sigma}(x-y)]_{\alpha\alpha'}\|_{L^{\infty}} \leq O(1)\gamma^{-s_a-s_b} e^{-c[d_{j,\sigma}(x,y)]^{\alpha_0}}$,
where $s_a, s_b \in [0, j]$, $\alpha_0 = 1/h$, and
 $d_{j,\sigma}(x, y) = \gamma^{-j}|x_0 - y_0| + \gamma^{-s_a}|x_a - y_a| + \gamma^{-s_b}|x_b - y_b|$
 $\|[C_{j,\sigma}(x)]_{\alpha\alpha'}\|_{L^1} \leq O(1)\gamma^j$.

◆□▶ ◆□▶ ◆臣▶ ◆臣▶ 臣 の�?

The fermionic cluster expansions

Recall that

$$Z(\lambda) = \int P(d\psi) e^{\lambda \int_{\Lambda_{\beta,L}} d^3 x \left[\psi^+_{x,\uparrow} \psi^-_{x,\uparrow} \psi^+_{x,\downarrow} \psi^-_{x,\downarrow} \right]}$$

= $\sum_n \frac{\lambda^n}{n!} \int_{(\Lambda_{\beta,L})^n} d^3 x_1 \cdots d^3 x_n \begin{cases} x_{1,\varepsilon_1,\tau_1} \cdots x_{n,\varepsilon_n,\tau_n} \\ x_{1,\varepsilon_1,\tau_1} \cdots x_{n,\varepsilon_n,\tau_n} \end{cases}$,

◆□▶ ◆□▶ ◆臣▶ ◆臣▶ 臣 のへぐ

 $\{\cdot\}$ is a $2n \times 2n$ determinant,

The fermionic cluster expansions

Recall that

$$Z(\lambda) = \int P(d\psi) e^{\lambda \int_{\Lambda_{\beta,L}} d^3 x \left[\psi^+_{x,\uparrow} \psi^-_{x,\uparrow} \psi^+_{x,\downarrow} \psi^-_{x,\downarrow} \right]}$$

= $\sum_n \frac{\lambda^n}{n!} \int_{(\Lambda_{\beta,L})^n} d^3 x_1 \cdots d^3 x_n \begin{cases} x_{1,\varepsilon_1,\tau_1} \cdots x_{n,\varepsilon_n,\tau_n} \\ x_{1,\varepsilon_1,\tau_1} \cdots x_{n,\varepsilon_n,\tau_n} \end{cases}$,

◆□▶ ◆□▶ ◆三▶ ◆三▶ 三三 のへで

 $\{\,\cdot\,\}$ is a 2n imes 2n determinant,

 The expanded terms have can be labeled by graphs, called the Feymann graphs.

The fermionic cluster expansions

Recall that

$$Z(\lambda) = \int P(d\psi) e^{\lambda \int_{\Lambda_{\beta,L}} d^3 x \left[\psi^+_{x,\uparrow} \psi^-_{x,\uparrow} \psi^+_{x,\downarrow} \psi^-_{x,\downarrow} \right]}$$

= $\sum_n \frac{\lambda^n}{n!} \int_{(\Lambda_{\beta,L})^n} d^3 x_1 \cdots d^3 x_n \begin{cases} x_{1,\varepsilon_1,\tau_1} \cdots x_{n,\varepsilon_n,\tau_n} \\ x_{1,\varepsilon_1,\tau_1} \cdots x_{n,\varepsilon_n,\tau_n} \end{cases}$,

 $\left\{ \;\cdot\;
ight\}$ is a 2n imes 2n determinant,

- The expanded terms have can be labeled by graphs, called the Feymann graphs.
- Partially expand the determinant such that only tree graphs appear.

< □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > <

Theorem (The BKAR jungle Formula. Brydges, Kennedy 87, Abdesselam Rivasseau 95)

Let $I_n = \{1, \dots, n\}$, $\mathcal{P}_n = \{\ell = (i, j), i, j \in I_n, i \neq j\}$, \mathcal{S} a set of smooth functions from $\mathbb{R}^{\mathcal{P}_n}$ to some Banach space, $\mathbf{1} \in \mathbb{R}^{\mathcal{P}_n}$ be the vector with every entry equals 1. Then for any $\mathbf{x} = (x_\ell)_{\ell \in \mathcal{P}_n} \in \mathbb{R}^{\mathcal{P}_n}$ and $f \in \mathcal{S}$:

$$f(\mathbf{1}) = \sum_{\mathcal{J}} \left(\int_0^1 \prod_{\ell \in \mathcal{F}} dw_\ell \right) \left(\prod_{k=1}^m \left(\prod_{\ell \in \mathcal{F}_k \setminus \mathcal{F}_{k-1}} \frac{\partial}{\partial x_\ell} \right) \right) f[X^{\mathcal{F}}(w_\ell)],$$

- J = (F₀ ⊂ F₁ · · · ⊂ F_{rmax} = F) is any partially ordered set of forests F_i with n vertices.
- $X^{\mathcal{F}}(w_{\ell})$ is a vector with elements $x_{\ell} = x_{ij}^{\mathcal{F}}(w_{\ell})$:
 - $x_{ij}^{\mathcal{F}} = 1$ if i = j, or if i and j are connected by \mathcal{F}_{k-1} .
 - $x_{ii}^{\mathcal{F}} = 0$ if i and j are not connected by \mathcal{F}_k ,
 - $x_{ij}^{\mathcal{F}} = \inf_{\ell \in P_{ij}^{\mathcal{F}}} w_{\ell}$, if *i* and *j* are connected by the forest \mathcal{F}_k but not \mathcal{F}_{k-1} , where $P_{ij}^{\mathcal{F}_k}$ is the unique path in the forest that connects *i* and *j*,

$$\begin{split} \bullet \ S_{2p}^c &= \sum_n S_{2p,n}^c \lambda^n, \\ S_{2p,n}^c &= \ \frac{1}{n!} \sum_{\{\underline{\tau}\}, \mathcal{G}, \mathcal{T}} \ \sum_{\mathcal{J}}' \epsilon(\mathcal{J}) \prod_{j=1}^n \int d^3 x_j \delta(x_1) \\ &\prod_{\ell \in \mathcal{T}} \int_0^1 dw_\ell C_{\tau_\ell, \sigma_\ell}(x_\ell, \bar{x}_\ell) \prod_{i=1}^n \chi_i(\sigma) \det_{\mathrm{left}}(C_j(w)) . \end{split}$$

◆□ ▶ < 圖 ▶ < 圖 ▶ < 圖 ▶ < 圖 • 의 Q @</p>

$$S_{2p}^{c} = \sum_{n} S_{2p,n}^{c} \lambda^{n},$$

$$S_{2p,n}^{c} = \frac{1}{n!} \sum_{\{\underline{\tau}\}, \mathcal{G}, \mathcal{T}} \sum_{\mathcal{J}}' \epsilon(\mathcal{J}) \prod_{j=1}^{n} \int d^{3}x_{j} \delta(x_{1})$$

$$\prod_{\ell \in \mathcal{T}} \int_{0}^{1} dw_{\ell} C_{\tau_{\ell}, \sigma_{\ell}}(x_{\ell}, \bar{x}_{\ell}) \prod_{i=1}^{n} \chi_{i}(\sigma) \det_{\mathrm{left}}(C_{j}(w)) .$$

◆□ ▶ < 圖 ▶ < 圖 ▶ < 圖 ▶ < 圖 • 의 Q @</p>

$$S_{2p}^{c} = \sum_{n} S_{2p,n}^{c} \lambda^{n},$$

$$S_{2p,n}^{c} = \frac{1}{n!} \sum_{\{\underline{\tau}\},\mathcal{G},\mathcal{T}} \sum_{\mathcal{J}}' \epsilon(\mathcal{J}) \prod_{j=1}^{n} \int d^{3}x_{j} \delta(x_{1})$$

$$\prod_{\ell \in \mathcal{T}} \int_{0}^{1} dw_{\ell} C_{\tau_{\ell},\sigma_{\ell}}(x_{\ell}, \bar{x}_{\ell}) \prod_{i=1}^{n} \chi_{i}(\sigma) \det_{\mathrm{left}}(C_{j}(w)) .$$

◆□▶ ◆□▶ ◆臣▶ ◆臣▶ 臣 のへぐ

► Theorem (Power counting Theorem, Rivasseau, ZW 2021)

$$S_{2\rho}^{c} = \sum_{n} S_{2\rho,n}^{c} \lambda^{n},$$

$$S_{2\rho,n}^{c} = \frac{1}{n!} \sum_{\{\underline{\tau}\},\mathcal{G},\mathcal{T}} \sum_{\mathcal{J}}' \epsilon(\mathcal{J}) \prod_{j=1}^{n} \int d^{3}x_{j} \delta(x_{1})$$

$$\prod_{\ell \in \mathcal{T}} \int_{0}^{1} dw_{\ell} C_{\tau_{\ell},\sigma_{\ell}}(x_{\ell}, \bar{x}_{\ell}) \prod_{j=1}^{n} \chi_{j}(\sigma) \det_{\mathrm{left}}(C_{j}(w)) .$$

► Theorem (Power counting Theorem, Rivasseau, ZW 2021)

• There exists two positive constants C_1 and C_2 , independent of T, such that $||S_{2p,n}||_{L^{\infty}} \leq C_1$, $\forall p \geq 2$, $n \geq 1$, and $\lambda < O(1)/|\log T|^2$.

$$S_{2p}^{c} = \sum_{n} S_{2p,n}^{c} \lambda^{n},$$

$$S_{2p,n}^{c} = \frac{1}{n!} \sum_{\{\underline{\tau}\},\mathcal{G},\mathcal{T}} \sum_{\mathcal{J}}' \epsilon(\mathcal{J}) \prod_{j=1}^{n} \int d^{3}x_{j} \delta(x_{1})$$

$$\prod_{\ell \in \mathcal{T}} \int_{0}^{1} dw_{\ell} C_{\tau_{\ell},\sigma_{\ell}}(x_{\ell}, \bar{x}_{\ell}) \prod_{j=1}^{n} \chi_{j}(\sigma) \det_{\mathrm{left}}(C_{j}(w)) .$$

Theorem (Power counting Theorem, Rivasseau, ZW 2021)

▶ There exists two positive constants C_1 and C_2 , independent of T, such that $||S_{2p,n}||_{L^{\infty}} \leq C_1$, $\forall p \geq 2$, $n \geq 1$, and $\lambda < O(1)/|\log T|^2$.

• $\|S_{2,n}\|_{L^{\infty}} \leq C_2 \log \frac{1}{T}$, for $n \geq 1$ and $\lambda < O(1)/|\log T|^2$.

$$S_{2\rho}^{c} = \sum_{n} S_{2\rho,n}^{c} \lambda^{n},$$

$$S_{2\rho,n}^{c} = \frac{1}{n!} \sum_{\{\underline{\tau}\},\mathcal{G},\mathcal{T}} \sum_{\mathcal{J}}' \epsilon(\mathcal{J}) \prod_{j=1}^{n} \int d^{3}x_{j} \delta(x_{1})$$

$$\prod_{\ell \in \mathcal{T}} \int_{0}^{1} dw_{\ell} C_{\tau_{\ell},\sigma_{\ell}}(x_{\ell}, \bar{x}_{\ell}) \prod_{j=1}^{n} \chi_{j}(\sigma) \det_{\mathrm{left}}(C_{j}(w)) .$$

Theorem (Power counting Theorem, Rivasseau, ZW 2021)

▶ There exists two positive constants C_1 and C_2 , independent of T, such that $||S_{2p,n}||_{L^{\infty}} \leq C_1$, $\forall p \geq 2$, $n \geq 1$, and $\lambda < O(1)/|\log T|^2$.

• $||S_{2,n}||_{L^{\infty}} \leq C_2 \log \frac{1}{T}$, for $n \geq 1$ and $\lambda < O(1)/|\log T|^2$.

• We need renormalizations for the p = 1 case.

The Multi-arch expansions for the self-energy

Theorem (Rivasseau, ZW 2021)

The amplitude of the self-energy is given by:

$$\begin{split} \Sigma(y,z) &= \sum_{n=0}^{\infty} \frac{\lambda^{n+2}}{n!} \int_{\Lambda^n} d^3 x_1 \dots d^3 x_n \sum_{\{\underline{\tau}\}} \sum_{\mathcal{G}_{\mathcal{B}}} \sum_{\mathcal{E} \mathcal{B}} \sum_{\mathcal{T}} \sum_{\{\sigma\}} \\ &\sum_{\substack{m-\text{arch systems} \\ \left(\prod_{(f_1,g_1),\dots,(f_m,g_m)\right) \\ \text{with } m \leq p}} \left(\prod_{\ell \in \mathcal{T}} \int_0^1 dw_\ell \right) \left(\prod_{r=1}^m \int_0^1 ds_r \right) \left(\prod_{\ell \in \mathcal{T}} C_{\sigma(\ell)}(f_\ell,g_\ell) \right) \\ &\left(\prod_{r=1}^m C(f_r,g_r)(s_1,\dots,s_{r-1}) \right) \frac{\partial^m \det_{left,\mathcal{T}}}{\prod_{r=1}^m \partial C(f_r,g_r)} (\{w_\ell\},\{s_r\}) \;. \end{split}$$

 $\blacktriangleright \exists C, K > 0 \text{ s.t. } \|\Sigma(y, z)\|_{L^{\infty}} \leq K \log \frac{1}{T}, \forall n \geq 1, \lambda < \frac{C}{|\log T|^2}.$

Need renormalization.

Renormalization theory

► A systematic way of removing divergences in QMB and QFT.

Renormalization theory

- A systematic way of removing divergences in QMB and QFT.
 - The divergent terms are expressed in the multi-scale representation and organized according to the descending order of the scaling indices.

< □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > <
Renormalization theory

- A systematic way of removing divergences in QMB and QFT.
 - The divergent terms are expressed in the multi-scale representation and organized according to the descending order of the scaling indices.

 Cancel any divergent term with the corresponding counter-term.

Renormalization theory

- A systematic way of removing divergences in QMB and QFT.
 - The divergent terms are expressed in the multi-scale representation and organized according to the descending order of the scaling indices.

- Cancel any divergent term with the corresponding counter-term.
- The divergent terms are organized into a tree structure, called the Gallavotti-Nicolo tree.

Renormalization theory

- A systematic way of removing divergences in QMB and QFT.
 - The divergent terms are expressed in the multi-scale representation and organized according to the descending order of the scaling indices.
 - Cancel any divergent term with the corresponding counter-term.
- The divergent terms are organized into a tree structure, called the Gallavotti-Nicolo tree.



The renormalization of the two-point function

▶ Let ϕ_{e_1} , ϕ_{e_2} some smooth, compactly supported functions.

$$\int dy' dz \ \phi_{e_1}(y') S_r^c(y, z) \phi_{e_2}(z)$$
(7)
= $\int dy \ [\int dz \ S_r^c(y, z)] \phi_{e_1}(y) \phi_{e_2}(y)$
+ $\int dy dz \ S_r^c(y, z) \phi_{e_1}(y) [\phi_{e_2}(z) - \phi_{e_1}(y)].$
:= $\int dy dz \ \phi_{e_1}(y) [\tau + (1 - \tau)] S_r^c(y, z) \phi_{e_2}(z),$

The renormalization of the two-point function

• Let ϕ_{e_1} , ϕ_{e_2} some smooth, compactly supported functions.

$$\int dy' dz \ \phi_{e_1}(y') S_r^c(y, z) \phi_{e_2}(z)$$
(7)
= $\int dy \ [\int dz \ S_r^c(y, z)] \phi_{e_1}(y) \phi_{e_2}(y)$
+ $\int dy dz \ S_r^c(y, z) \phi_{e_1}(y) [\phi_{e_2}(z) - \phi_{e_1}(y)].$
:= $\int dy dz \ \phi_{e_1}(y) [\tau + (1 - \tau)] S_r^c(y, z) \phi_{e_2}(z),$

► The local term: $\delta \mu^r(y) = -[\int dz \ S_r^c(y, z)]$ will be canceled by the counter-term at scale r: $\delta \mu^r + \tilde{\delta} \mu^r = 0$,

The renormalization of the two-point function

• Let ϕ_{e_1} , ϕ_{e_2} some smooth, compactly supported functions.

$$\int dy' dz \ \phi_{e_1}(y') S_r^c(y, z) \phi_{e_2}(z)$$
(7)
= $\int dy \ [\int dz \ S_r^c(y, z)] \phi_{e_1}(y) \phi_{e_2}(y)$
+ $\int dy dz \ S_r^c(y, z) \phi_{e_1}(y) [\phi_{e_2}(z) - \phi_{e_1}(y)].$
:= $\int dy dz \ \phi_{e_1}(y) [\tau + (1 - \tau)] S_r^c(y, z) \phi_{e_2}(z),$

- ► The local term: $\delta \mu^r(y) = -[\int dz \ S_r^c(y, z)]$ will be canceled by the counter-term at scale r: $\delta \mu^r + \tilde{\delta} \mu^r = 0$,
- Similarly, cancellation of the self-energy with the counter-term

$$\hat{\Sigma}_{s_+,s_-}^r \big[(2\pi T, P_F(\mathbf{k}))_{s_+,s_-}, \hat{\nu}^{\leq (r-1)}, \lambda \big] + \hat{\nu}_{s_+,s_-}^r (P_F(\mathbf{k})_{s_+,s_-}, \lambda) = 0.$$

The renormalizations are performed from the high scale terms to the low scale terms, following the Gallavotti-Nicolò tree.



Figure: A Gallavotti-Nicolò tree with 16 nodes and 8 bare vertices. The round dots represent the nodes and bare vertices, and the big square represents the root, which has the scaling index $r \leq -1$. The dash lines are the inclusion relations between these nodes and the thin lines are the external fields of the nodes.

• We have the desired L^1 bounds for the free propagators.



- ▶ We have the desired *L*¹ bounds for the free propagators.
- ▶ Using Gram-Hadammard inequality to bound the determinant: If *M* is a square matrix with elements $M_{ij} = \langle A_i, B_j \rangle$, with $A_i, B_j \in L^2$, then $\|\det M\|_{L^{\infty}} \leq \prod_i \|A_i\|_{L^{\infty}} \cdot \|B_i\|_{L^{\infty}}$.

- ▶ We have the desired *L*¹ bounds for the free propagators.
- ▶ Using Gram-Hadammard inequality to bound the determinant: If *M* is a square matrix with elements $M_{ij} = \langle A_i, B_j \rangle$, with $A_i, B_j \in L^2$, then $\|\det M\|_{L^{\infty}} \leq \prod_i \|A_i\|_{L^{\infty}} \cdot \|B_i\|_{L^{\infty}}$.
- Summation over the sector indices using the Sector counting lemma (ZW 2021);

- ▶ We have the desired *L*¹ bounds for the free propagators.
- Using Gram-Hadammard inequality to bound the determinant: If M is a square matrix with elements $M_{ij} = \langle A_i, B_j \rangle$, with $A_i, B_j \in L^2$, then $\|\det M\|_{L^{\infty}} \leq \prod_i \|A_i\|_{L^{\infty}} \cdot \|B_i\|_{L^{\infty}}$.
- Summation over the sector indices using the Sector counting lemma (ZW 2021);

• Bounds over spanning trees with *n* vertices $\sum_T \leq n!$;

◆□ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ < □ ▶ <

► Theorem (Benffatto, Giuliani, Mastropietro, 2007) For $0 < \mu \le 1$, the ground state is a Fermi liquid for $T \ge T_c$, with $T_c = K_1 \exp(-\frac{C_1}{|\lambda|})$.

► Theorem (Benffatto, Giuliani, Mastropietro, 2007)

For $0 < \mu \le 1$, the ground state is a Fermi liquid for $T \ge T_c$, with $T_c = K_1 \exp(-\frac{C_1}{|\lambda|})$.

◆□▶ ◆□▶ ◆三▶ ◆三▶ 三三 のへぐ

- ► Theorem (Benffatto, Giuliani, Mastropietro, 2007) For $0 < \mu \le 1$, the ground state is a Fermi liquid for $T \ge T_c$, with $T_c = K_1 \exp(-\frac{C_1}{|\lambda|})$.
- ► Theorem (Afchain, Magnen, Rivasseau, 2004) For $\mu = 2$, the ground state is a not a Fermi liquid for $T \ge T_c$, with $T_c = K_2 \exp(-\frac{C_2}{|\lambda|^{1/2}})$.

- ► Theorem (Benffatto, Giuliani, Mastropietro, 2007) For $0 < \mu \le 1$, the ground state is a Fermi liquid for $T \ge T_c$, with $T_c = K_1 \exp(-\frac{C_1}{|\lambda|})$.
- ► Theorem (Afchain, Magnen, Rivasseau, 2004) For $\mu = 2$, the ground state is a not a Fermi liquid for $T \ge T_c$, with $T_c = K_2 \exp(-\frac{C_2}{|\lambda|^{1/2}})$.

- ► Theorem (Benffatto, Giuliani, Mastropietro, 2007) For $0 < \mu \le 1$, the ground state is a Fermi liquid for $T \ge T_c$, with $T_c = K_1 \exp(-\frac{C_1}{|\lambda|})$.
- Theorem (Afchain, Magnen, Rivasseau, 2004)
 For µ = 2, the ground state is a not a Fermi liquid for T ≥ T_c, with T_c = K₂ exp(-^{C₂}/_{|λ|^{1/2}}).
- ▶ Theorem (ZW, 2022)

For $\mu = 2 - \mu_0$, $\mu_0 \ll 1$, the ground state is a not a Fermi liquid for $T \ge T_c$, with $T_c = \frac{K_3}{\mu_0} \exp(-\frac{C_3}{|\lambda|^{1/2}})$.

Conclusions and perspectives

► We provide rigorous proof that the ground state of the Honeycomb Hubbard model at µ = 1 is not a Fermi liquid.

◆□▶ ◆□▶ ◆三▶ ◆三▶ 三三 のへぐ

Conclusions and perspectives

► We provide rigorous proof that the ground state of the Honeycomb Hubbard model at µ = 1 is not a Fermi liquid.

• Case of $0 < \mu < 1$, crossover of $\mu = 0$ and $\mu = 1$?

Conclusions and perspectives

- ▶ We provide rigorous proof that the ground state of the Honeycomb Hubbard model at µ = 1 is not a Fermi liquid.
- Case of $0 < \mu < 1$, crossover of $\mu = 0$ and $\mu = 1$?
- Metal-Insulator transitions and many-body localization in Hubbard model.

< □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > <

References

- V. Rivasseau, Z. Wang,
 - Honeycomb Hubbard Model at van Hove Filling Part I: Construction of the Schwinger Functions, arXiv:2108.10852,
 - Honeycomb Hubbard Model at van Hove Filling II: Lower Bounds of the Self-Energy, arXiv:2108.10415.

< □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > <

Thanks for your attention!