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I. STARGRAPH AND ITS PROPERTIES

A. Hamiltonian and wavefunctions

As shown in the previous section, at the heart of multi channel Kondo there is a stargraph problem which is the zero mode of the lowenergy fixed point Hamiltonian of the Multi channel Kondo. In this section our focus will be on the stargraph problem itself. As shown in the Fig.1 above one central spin (\vec{S}_d) is connected with K spins forming a K channel stargraph problem corresponding to the K -channel Kondo problem. The Hamiltonian of this above model is given as

$$H = \alpha \vec{S}_d \cdot \sum_{i=1}^K \vec{S}_i = \alpha \vec{S}_d \cdot \vec{S}$$

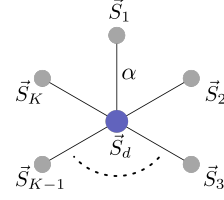


FIG. 1. This is a stargraph model with one central spin \vec{S}_d and K outer spin $1/2$ s. The central spin is coupled with all the outer spins with Heisenberg coupling with coupling strength α .

$$= \frac{\alpha}{2} (\vec{J}^2 - \vec{S}^2 - \vec{S}_d^2), \quad \alpha > 0. \quad (1)$$

where $\vec{J} = \vec{S} + \vec{S}_d$ and $\vec{S} = \sum_{i=1}^K \vec{S}_i$. One can see that the large spin S can take many possible values as this is made out of K spin $1/2$ s. We can see $[H, J] = 0 = [H, J^i]$ for $i = x, y, z$, $[J^i, J^j] \neq 0$ for $i \neq j$. Also the operators $\hat{Z} = 2^{K+1} S_d^z \prod_{i=1}^K S_i^z$ and $\hat{X} = 2^{K+1} S_d^x \prod_{i=1}^K S_i^x$ commutes with the Hamiltonian, $[H, \hat{Z}] = 0 = [H, \hat{X}]$ though $[\hat{Z}, \hat{X}] \neq 0$, in general. Thus we can form the CSCO with H, J, J^z, S, S_d . For a particular value S , J can take two values $S \pm S_d$ with the energies $-\alpha S_d(S_d + 1 \mp J)$ respectively. S_d is fixed thus the energy values only depends on the corresponding J value of the state. As the energy does not depend on the J^z , all $2J + 1$ J^z states labeled by $|S_d, S; J, J^z\rangle$ are degenerate. It is easy to see that the ground state has energy $E = -\alpha S_d(S_d + 1 + J)$ with $J = S - S_d$, thus $E_g = -\alpha S_d(S_{max} + 1)$ with S taking the maximum possible value $S_{max} = K/2$.

1. Ground state wavefunction

Above calculations show stargraph with K channels has K -fold ground state degeneracy, states labeled as $|S_d, S; J, J^z\rangle$ where J^z takes $K = 2J + 1$ distinct values, $J^z = \{-J, \dots, +J\}$. Here our goal is to find the wavefunction in the fundamental spin basis $|S_d^z, S_1^z, \dots, S_K^z\rangle$. To achieve that we use the Clebsch-gordon coefficients, we know for two spins j_1 and j_2 with total spin J and $J^z = M$ one can expand the state as

$$|j_1, j_2; J, M\rangle = \sum_{\substack{m_1 = \{-j_1, \dots, j_1\} \\ m_2 = \{-j_2, \dots, j_2\}}} \mathcal{C}_{j_1, j_2; J, M}^{j_1, m_1; j_2, m_2} |j_1, m_1; j_2, m_2\rangle, \quad (2)$$

where $j_1^z = m_1$ and $j_2^z = m_2$ and $\mathcal{C}_{j_1, j_2; J, M}^{j_1, m_1; j_2, m_2}$ is the Clebsch-Gordon coefficient. In our problem two spins S_d and S are forming one large spin J represented by the state $|S_d, S; J, J^z\rangle$, in the ground state $J = S - 1/2$. Thus the ground state $|S_d, S; (S - \frac{1}{2}), M\rangle \equiv |\alpha\rangle$ can be expanded as

$$|\alpha\rangle = \sum_{\substack{m_1 = \{-S_d, \dots, S_d\} \\ m_2 = \{-S, \dots, S\}}} \mathcal{C}_{S_d, S; J, M}^{S_d, m_1; S, m_2} |S_d, m_1\rangle \otimes |S, m_2\rangle, \quad (3)$$

Because in the ground state S is maximum, the state $|S, S^z\rangle$ can be decomposed in terms of a two spin problem of spin $1/2$ and $S - 1/2$ forming the state $|1/2, S - 1/2; S, S^z\rangle$ which can be further expanded in the basis $|1/2, m'_1\rangle \otimes |S - 1/2, m'_2\rangle$ where m'_1, m'_2 are the z -components of the spin $1/2$ and $S - 1/2$. Next we decompose the state $|S - 1/2, m'_2\rangle$ in further smaller spin problem and so on. This way we finally arrive at the ground state wavefunction in terms of the fundamental basis $|S_d^z, S_1^z, \dots, S_K^z\rangle$, give as

$$|S_d, S; J, M\rangle = \sum_{S_d^z, \{S_i^z\}} \mathcal{D}_{S_d^z, \{S_i^z\}} |S_d^z, S_1^z, \dots, S_K^z\rangle \quad (4)$$

where $\{S_i^z\} \equiv (S_1^z, \dots, S_K^z)$ and the coefficient $\mathcal{D}_{S_d^z, \{S_i^z\}}$ is given as

$$\mathcal{D}_{S_d^z, \{S_i^z\}} = \sum_{\substack{m_2=[-S, S] \\ m_4=[-(S-1/2), (S-1/2)] \\ m_{2N-4}=[-1, 1]}} \mathcal{M}_{m_2, m_4, \dots, m_{2N-2}}^{S_d^z, S_1^z, \dots, S_K^z}, \quad (5)$$

and further the coefficients $\mathcal{M}_{m_2, m_4, \dots, m_{2N-2}}^{S_d^z, S_1^z, \dots, S_K^z} \equiv \Sigma$, are made out of the products of Clebsh-Gordon Coefficients

$$\Sigma = \mathcal{C}_{S_d, S; J, M}^{S_d, S_d^z, S, m_2} \mathcal{C}_{S_1, (S-1/2), m_4}^{S_1, S_1^z, (S-1/2), m_4} \dots \mathcal{C}_{S_{K-1}, S_K^z, 1, m_{2N-4}}^{S_{K-1}, S_{K-1}^z, S_K, S_K^z}, \quad (6)$$

Using these wavefunctions we can calculate various entanglement features of this stargraph problem.

B. Entanglement properties

Using the above exact wavefunctions we can numerically compute various entanglement properties of this stargraph model.

1. Impurity entanglement entropy

Here we are interested in finding the entanglement of the central impurity spin with the rest of the multi-channel bath zero mode. We know that for multi channel case the ground state is degenerate, thus for each degenerate state we can calculate the entanglement entropy. As we have already discussed for N_{ch} number of channels there are N_{ch} number of degenerate ground states, labeled by the J_z quantum number.

For single channel case the the ground state is unique and is

$$|J = 0, J_z = 0\rangle = \frac{1}{\sqrt{2}}(|\uparrow_d \downarrow_0\rangle - |\downarrow_d \uparrow_0\rangle) \quad (7)$$

Thus, one can easily calculate the reduced density matrix of the impurity by tracing out the 0^{th} state. From that reduced density matrix the entanglement entropy is $\log 2$, which is maximum possible for a spin-1/2.

Next, we are interested in finding this entanglement entropy for the multi-channel case $N_{ch} > 1$. We have generated the degenerate ground states analytically using the Clebsch-Gordon coefficients. On this wavefunctions we do various entanglement and correlation studies. We start with the ground state wavefunction, $|\psi_g\rangle_{J_z} = |1/2_d, S; J, J_z\rangle$, starting with one of those degenerate ground states labeled by J_z we calculate the density matrix.

$$\rho_{J_z} = |\psi_g\rangle_{J_z} \langle \psi_g|_{J_z} \quad (8)$$

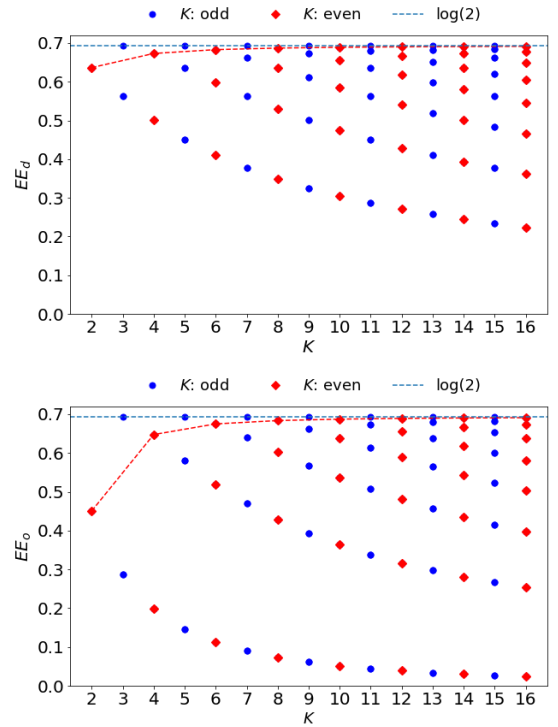


FIG. 2. Y-axis shows the impurity entanglement entropy with the rest and X-axis shows the number of channels N_{ch} . For one value of $N_{ch} > 1$, there are N_{ch} degenerate ground states and their corresponding impurity entanglement entropy.

Using the wavefunction we can do the partial tracing on the impurity spin to get the reduced density matrix of the rest. From that we get entanglement entropy for different number of channels. For even number of channel cases J_z and $-J_z$ sector has same entanglement entropy. For the case of odd N_{ch} there also $\pm J_z$ sectors shares same entanglement entropy, but there is a state $J_z = 0$ which has entanglement entropy $\log 2$. Thus from the above we can see that $J_z = 0$ state corresponding to odd number of channels shows perfect screening in terms of the maximum entanglement entropy and $J_z = 0$ magnetization.

2. Mutual Information

Mutual information between two sub-systems A, B is defined as,

$$I^2(A : B) = S_A + S_B - S_{A \cup B} \quad (9)$$

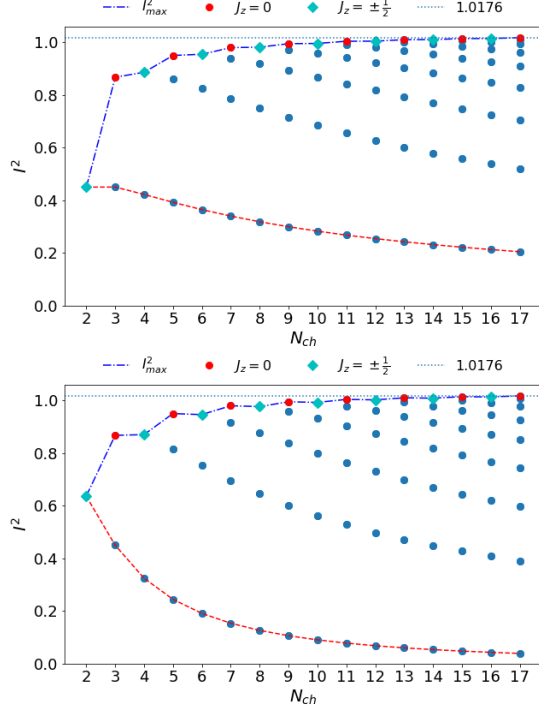


FIG. 3. *Left:* Y-axis is the mutual information between the impurity-spin and one outer spin and x-axis is the number of channel N_{ch} . *Right:* This shows the variation of mutual information between two outer spins.

We first study the mutual information between the impurity-spin and one outer spin $I^2(d : o)$. Because all the outer spins are symmetric, the mutual information between impurity-spin with any one outer spin is same. Next we calculate the mutual information between two outer spin $I^2(o : o)$. One can see that the maximum mutual information for both the cases are from the $(J_z)_{min}$ state. For odd number of channels $(J_z)_{min} = 0$ and for even number of channels $\pm(J_z)_{min} = \pm 1/2$. And the minimum mutual information corresponds to the $\pm(J_z)_{max} = \pm J$ state.

As can be seen from the above Figure.3 that both the maximum mutual information of $I^2(d : o)$ and $I^2(o : o)$ try to saturate near the same value. Thus in the lowest J_z sector the entanglement distribution between impurity and outer vs outer and outer is same. Which suggests that even though there were no coupling among the outer spins explicitly, there are entanglement among them. The minimum of the mutual information coming from the largest J_z sector, the $I^2(o : o)$ corresponding to this sector asymptotically approach zero for large

N_{ch} limit. Which shows that in the large channel limit $N_{ch} \gg 1$ though the $\min I^2(o : o)$ approach zero the $\min I^2(d : o)$

3. Tri-partite information

Apart from Mutual information we can calculate the tri-partite information also. This tri-partite information is defined among three sub-systems A, B, C , as

$$I^3_{A,B,C} = (S_A + S_B + S_C) - (S_{AB} + S_{BC} + S_{CA}) + S_{ABC} \quad (10)$$

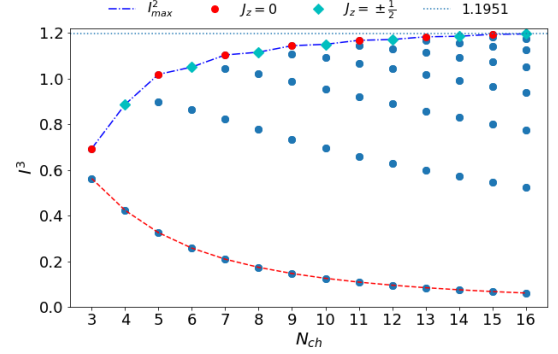


FIG. 4. This plot shows the variation of the tri-partite information with the number of channels N_{ch} .

Our tri-partite information measure (Fig.4) among the outer spins shows that the just like the mutual information the minimum tri-partite information asymptotically approaches zero in the $N_{ch} \gg 1$ limit. Same is the case for 4-partite information also.

4. I_N vs N

Here we are interested in understanding the entanglement distribution among different number of outer spins. To find that out we take a multi-channel stargraph model with N_{ch} number of outer spins. We calculate different multi partite informations $I^m, m \in \{2, \dots, N_{ch}\}$ on it's N_{ch} degenerate ground states.

Our above study (Fig.5) shows that the mutual information has the maximum value and higher order multi-partite informations is smaller and smaller approaching zero.

C. Correlation Studies

1. Quantum energy

Here we are interested in the energy expectation value arising from the quantum fluctuation part of the Hamiltonian (??). We define using energy E_{ising} and quantum

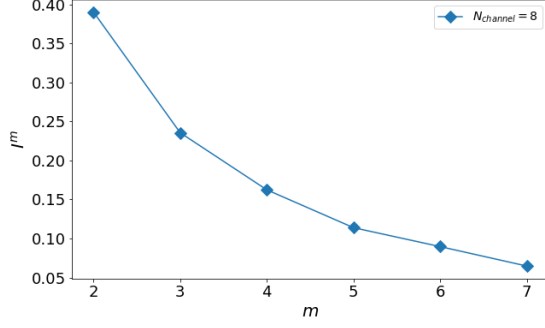


FIG. 5. This figure shows how the m -partite information varies with m showing the entanglement distribution among different number of outer-spins.

energy E_Q in the ground state $|\psi_g\rangle$ as,

$$E_{ising} = |\langle \psi_g | \mathcal{H}_0^c | \psi_g \rangle|, \quad E_Q = |\langle \psi_g | \mathcal{H}_0^Q | \psi_g \rangle| = |E_g| - |E_{ising}|, \\ = \frac{\alpha(S+1)}{2} - |E_{ising}|. \quad (11)$$

This quantum energy E_Q is generated by the spin-flips between the impurity spin and the outer spin. Thus it is interesting to find out how this quantum energy changes with the increase of the number of channels (N_{ch}). Another important parameter is the Quantum energy per channel, $e_Q = |E_Q|/N_{ch}$, and its variation towards the infinity channel limit. We know in the ground state $N_{ch} = 2S$ thus,

$$e_Q = \frac{\alpha(S+1)}{4S} - \frac{|E_{ising}|}{2S}, \\ = \frac{\alpha}{4} - \frac{1}{N_{ch}} \left(|E_{ising}| - \frac{\alpha}{2} \right), \quad (12)$$

Thus from the above equation of quantum energy per channel one can see that in the infinity-channel limit $N_{ch} \rightarrow \infty$,

$$e_Q = \frac{\alpha}{4} - \frac{|E_{ising}|}{N_{ch}} < \frac{\alpha}{4}, \quad (13)$$

2. Correlation study

Here we want to calculate various correlation functions in the ground state. We start with the expectation value of J_z^2 in the ground state $|\psi_g\rangle$,

$$j_z^2 = \frac{1}{N_{ch}} \langle \psi_g | J_z^2 | \psi_g \rangle = \frac{1}{N_{ch}} \left[J^2 - (J_x^2 + J_y^2) \right], \\ = \frac{1}{N_{ch}} \left[\frac{(N_{ch}^2 - 1)}{4} - (J_x^2 + J_y^2) \right] \quad (14)$$

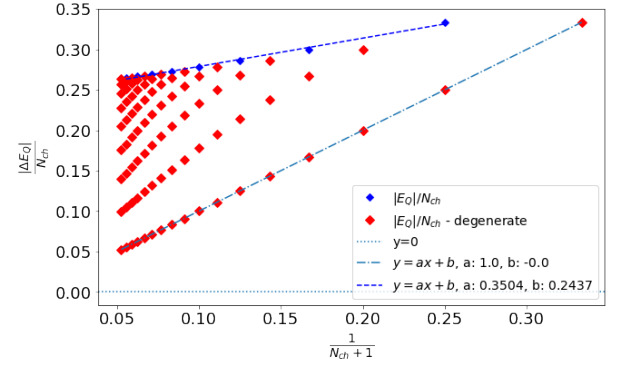


FIG. 6. This shows the variation of quantum energy per channel with $1/N$, where $N = N_{ch} + 1$ is the total number of spins in the systems including the impurity spins.

$$\langle \psi_g | J_z^2 | \psi_g \rangle_{max} = \left(\frac{N_{ch} - 1}{2} \right)^2, \\ \langle \psi_g | J_z^2 | \psi_g \rangle_{min} = \left(\frac{N_{ch} - (N_{ch} - 1)}{2} \right)^2 = \frac{1}{4}, \text{ for } N_{ch} \text{ even.} \\ = 0, \text{ for } N_{ch} \text{ odd.} \quad (15)$$

Thus,

$$\langle (J_x)^2 + (J_y)^2 \rangle_{max} = \langle J^2 \rangle - \langle J_z^2 \rangle_{min}, \\ = \frac{N_{ch}^2 - 1}{4} - \frac{1}{4} = \frac{N_{ch}^2 - 2}{4}, \text{ for } N_{ch} \text{ even} \\ = \frac{N_{ch}^2 - 1}{4}, \text{ for } N_{ch} \text{ odd}$$

Which shows, for single channel case $N_{ch} = 1$, this quantum fluctuation expectation value vanishes, while for multi channel this takes non-zero values. Thus

$$\frac{1}{N_{ch}} \langle (J_x)^2 + (J_y)^2 \rangle_{max} = \frac{N_{ch}}{4} - \frac{1}{2N_{ch}}, \text{ for } N_{ch} \text{ even} \\ = \frac{N_{ch}}{4} - \frac{1}{4N_{ch}}, \text{ for } N_{ch} \text{ odd}$$

Thus in the large channel number limit $N_{ch} \gg 1$, one can see that

$$\frac{1}{N_{ch}} \langle (J_x)^2 + (J_y)^2 \rangle_{max} \rightarrow \frac{N_{ch}}{4}, \text{ for both even and odd } N_{ch} \quad (16)$$

J_z can take values $\{-J, -J+1, \dots, J-1, J\}$ in the ground state, where $J = S - 1/2 = (N_{ch} - 1)/2$.

3. Staggered magnetization

One can analytically check the single channel Kondo results. $N_{ch} = 1$, thus here the ground state is unique.

$$|\psi_g\rangle = \frac{1}{\sqrt{2}} \left(|\uparrow\downarrow\rangle - |\downarrow\uparrow\rangle \right) = |J=0, J_z=0\rangle \quad (17)$$

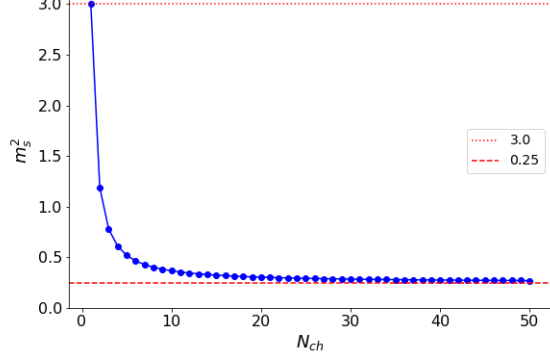


FIG. 7. This shows how the staggered magnetization changes with the number of channels N_{ch} .

Thus in this ground state $\langle J_z \rangle = 0$ and the entanglement entropy is $\log 2$ the maximum value. Now we define a staggered magnetization (7) operator M_s measure,

1. For the single channel case both the \vec{S}_d and \vec{S} are spin-1/2 object and the total spin is defined as $\vec{J} = \vec{S}_d + \vec{S}$. Here \vec{S} is the total outer spin.

$$\begin{aligned} M_s^2 &= \langle \psi_g | (\vec{S}_d - \vec{S})^2 | \psi_g \rangle = \langle \psi_g | 2(\vec{S}_d^2 + \vec{S}^2) - \vec{J}^2 | \psi_g \rangle \\ &= 2 \langle \psi_g | \vec{S}_d^2 + \vec{S}^2 | \psi_g \rangle \\ &= 3 \end{aligned} \quad (18)$$

2.

$$\begin{aligned} M_s^2 &= \langle \psi_g | (\vec{S}_d - \vec{S})^2 | \psi_g \rangle = \langle \psi_g | 2(\vec{S}_d^2 + \vec{S}^2) - \vec{J}^2 | \psi_g \rangle \\ &= \left\langle \psi_g \left| 2 \left(\frac{3}{4} + \frac{N_{ch}(N_{ch}+2)}{4} \right) - \frac{(N_{ch}+1)(N_{ch}-1)}{4} \right| \psi_g \right\rangle \\ &= \frac{N_{ch}^2}{4} + N_{ch} + \frac{7}{4} \end{aligned} \quad (19)$$

Thus, staggered magnetization squared per channel,

$$\begin{aligned} m_s^2 &= \frac{M_s^2}{N_{ch}^2} = \frac{1}{4} + \frac{1}{N_{ch}} + \frac{7}{4N_{ch}^2} \xrightarrow{N_{ch} \gg 1} \frac{1}{4} + \frac{1}{N_{ch}} \quad (20) \\ \frac{\vec{J}^2}{N_{ch}^2} &= \frac{(\vec{S}_d + \vec{S})^2}{N_{ch}^2} = \frac{N_{ch}^2 - 1}{4N_{ch}^2} \xrightarrow{N_{ch} \gg 1} \frac{1}{4}. \end{aligned}$$

Using the $SU(2)$ property of our problem, we can define m_s for each spatial direction, in 3D there are three independent spatial direction.

$$\langle (m_s^x)^2 \rangle = \langle (m_s^y)^2 \rangle = \langle (m_s^z)^2 \rangle = \frac{1}{3} m_s^2, \quad (21)$$

We define $m_Q^2 = m_s^2 - 1/4$ which in the the large N_{ch} limit becomes $1/N_{ch}$ and eventually vanishes in the $N_{ch} \rightarrow \infty$ limit.

Thus to summarize,

1. In the single channel problem, $\langle (m_s^z)^2 \rangle = 1$.
2. In the multi-channel case, $\langle (m_s^z)^2 \rangle = \frac{1}{3N_{ch}} + \frac{1}{12} + \frac{7}{12N_{ch}^2} \xrightarrow{N_{ch} \gg 1} \frac{1}{12} + \frac{1}{3N_{ch}} \rightarrow \frac{1}{12}$. This show partial screening.

Thus it is important to find the scaling of

1. M_s^2 vs N_{ch} .
2. Uncompensated Quantum Fluctuation vs N_{ch} .

D. Calculating thermodynamic quantities

We take the stargraph Hamiltonian and do the exact diagonalization and find the entire eigenspectrum. Using the full eigenspectrum we calculate the partition function $Z = \sum_i e^{-\beta E_i}$, where E_i is the energy of the i^{th} state. One can rewrite this partition function interms of the state degeneracies as $Z = \sum_{\epsilon} d(\epsilon) e^{-\beta \epsilon}$, where $d(\epsilon)$ is the degeneracy of the state with energy ϵ . Thus the free energy is given as $\mathcal{F} = -k_B T \log Z$.

1. Thermal entropy

Thermal entropy is defined as $S = -(\frac{\partial \mathcal{F}}{\partial T})_H$, where H represents a constant magnetic field. In ourcase we will be interested in the zero field case.

$$\begin{aligned} \mathcal{F} &= -k_B T \log Z \\ S &= -(\frac{\partial \mathcal{F}}{\partial T}) = -k_B \log Z - k_B T \frac{1}{Z} \frac{dZ}{dT} \end{aligned} \quad (22)$$

Now

$$\begin{aligned} Z &= \sum_{\epsilon} d(\epsilon) e^{-\beta \epsilon} \Rightarrow \frac{dZ}{dT} = \sum_{\epsilon} d(\epsilon) e^{-\beta \epsilon} (-\epsilon) \frac{d\beta}{dT} \\ &= \sum_{\epsilon} d(\epsilon) e^{-\beta \epsilon} \frac{\epsilon}{k_B T^2} = k_B \sum_{\epsilon} \epsilon d(\epsilon) e^{-\beta \epsilon} \end{aligned} \quad (23)$$

Thus we get

$$\begin{aligned} S &= -k_B \log \sum_{\epsilon} d(\epsilon) e^{-\beta \epsilon} - \frac{1}{\beta} \frac{k_B \sum_{\epsilon} \epsilon d(\epsilon) e^{-\beta \epsilon}}{\sum_{\epsilon} d(\epsilon) e^{-\beta \epsilon}} \\ \lim_{\beta \rightarrow \infty} S &= -k_B \log_2 d(\epsilon_G), \text{ and } \lim_{\beta \rightarrow 0} S = -k_B \log \sum_{\epsilon} d(\epsilon) \end{aligned} \quad (24)$$

Thus at the extreme temperature it is easy to calculate the thermal entropy, but it is difficult to visualize for any intermediate temperatures. Thus we plot the thermal entropy (unit $\log 2$) for different temperatures and for different channels in Fig.8. This shows at the extreme temperature the thermal entropy saturates and at the intermediate temperature it changes from one to the other like a soliton like solution. At low temperature the thermal entropy in the unit of $\log 2$ is not always quantized but at high temperature it is. To understand this let's say

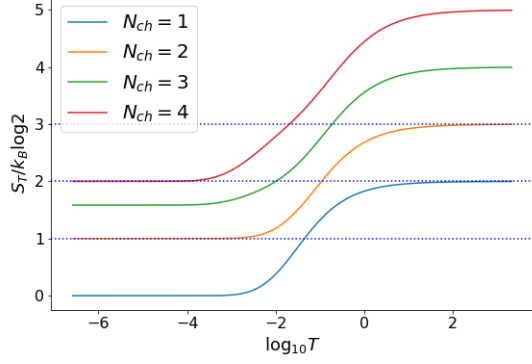


FIG. 8. This shows the variation of thermal entropy with the temperature.

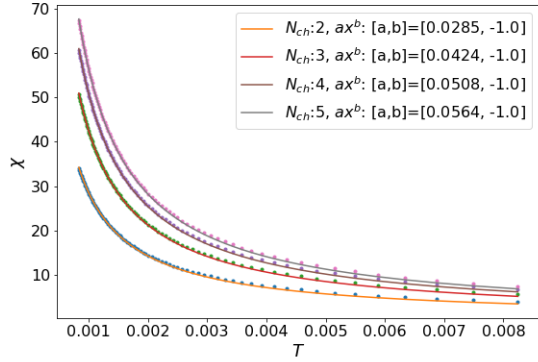


FIG. 9. Impurity susceptibility vs temperature.

at low temperature $S_T/k_B \log T = \Omega = \log_2 d(\epsilon_G)$. Here $d(\epsilon_G)$ is always integer and equal to the channel number K . Thus

$$d(\epsilon_G) = 2^\Omega = K \quad (25)$$

From our study we can see that for some channel number we get $S_T/k_B \log T$ to be integer. One can easily see that those $K = 2^n$ channel cases, n is an integer and will have integer thermal entropy. Thus between m and $m+1$ integer thermal entropy plateau there will be $(2^m - 1)$ number of fractional thermal entropy plateau, which is always odd.

2. Suseptibility

1. Impurity-field susceptibility : Now we add small magnetic field to the impurity spin and calculate the suseptibility as a function of temperatur for different channel cases.
2. Outer-field suseptibility: Now we add a unifor magnetic field on the outer spins and calculate the suseptibility for different chanel cases.

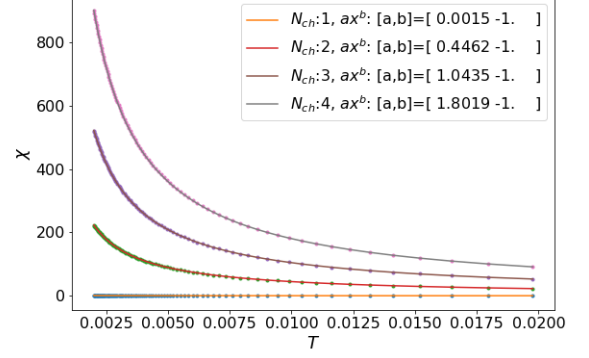


FIG. 10. Outer suseptibility vs temperature.

As can be seen from the above results that the both impurity and outer field suseptibility follows a power law in temperature for $K > 1$ case.

$$\chi_{imp}(T) \sim T^{-1}, \quad \chi_{outer} \sim T^{-1} \quad (26)$$

This shows an universality among the greater than one channels.

II. EXCITATION PROPERTIES

Here we start with the low energy zero mode fixed point Hamiltonian for the two-channel Kondo obtained under URG treatment. Our goal here is to add excitations on top of this ground state and using the perturbation theory. In doing so we consider the leading order scatter appereing in the lowest order in the perturbative expansion. Note, in the fixed point zero mode Hamiltonian impurity spin is coupled with the bath electrons via spin exchange coupling, there is no charge exchange coupling present.

$$H^{(2)} = \alpha \vec{S}_{imp} \cdot (\vec{S}_1 + \vec{S}_2), \quad \vec{S}_i = \frac{1}{2} c_{i,\alpha}^\dagger \vec{\sigma}_{\alpha,\beta} c_{i,\beta} \quad (27)$$

Here $\vec{S}_i = 1, 2$ represents the spin degree of freedom present in the origin of the i^{th} channel. Thus one can rewrite the Eq.(27) interms of the electronic degree of freedom as,

$$\begin{aligned} \mathcal{H}_0 = H_{0_1} + H_{0_2} = & \alpha S_{imp}^z \left(\frac{n_{1,\uparrow} - n_{1,\downarrow}}{2} \right) + \frac{1}{2} \left(S_{imp}^+ c_{1,\downarrow}^\dagger c_{1,\uparrow} - S_{imp}^- c_{1,\uparrow}^\dagger c_{1,\downarrow} \right) \\ & + \alpha S_{imp}^z \left(\frac{n_{2,\uparrow} - n_{2,\downarrow}}{2} \right) + \frac{1}{2} \left(S_{imp}^+ c_{2,\downarrow}^\dagger c_{2,\uparrow} - S_{imp}^- c_{2,\uparrow}^\dagger c_{2,\downarrow} \right) \end{aligned}$$

Here the basis we are interested in is $\{\mathcal{B}\} = \{|S_{imp}^z, n_{1,\uparrow}, n_{1,\downarrow}, n_{2,\uparrow}, n_{2,\downarrow}\rangle\}$. In this basis we write down all $2^3 = 8$ states of the spin Hamiltonian eq.(27)

$$|\alpha_0\rangle = \frac{1}{\sqrt{6}} \left[2|\downarrow, 1, 0, 1, 0\rangle - |\uparrow, 0, 1, 1, 0\rangle - |\uparrow, 1, 0, 0, 1\rangle \right]$$

$$|\alpha_1\rangle = \frac{1}{\sqrt{6}} \left[2|\uparrow, 0, 1, 0, 1\rangle - |\downarrow, 1, 0, 0, 1\rangle - |\downarrow, 0, 1, 1, 0\rangle \right] \\ E_{\alpha_0} = E_{\alpha_1} = -\alpha \quad (29)$$

$$|\beta_0\rangle = \frac{1}{\sqrt{2}} \left[|\downarrow, 1, 0, 0, 1\rangle - |\downarrow, 0, 1, 1, 0\rangle \right] \\ |\beta_1\rangle = \frac{1}{\sqrt{2}} \left[|\uparrow, 1, 0, 0, 1\rangle - |\uparrow, 0, 1, 1, 0\rangle \right] \\ E_{\beta_1} = E_{\beta_2} = 0 \quad (30)$$

$$|\gamma_0\rangle = \frac{1}{\sqrt{3}} \left[|\uparrow, 0, 1, 0, 1\rangle + |\downarrow, 1, 0, 0, 1\rangle + |\downarrow, 0, 1, 1, 0\rangle \right] \\ |\gamma_1\rangle = \frac{1}{\sqrt{3}} \left[|\uparrow, 1, 0, 0, 1\rangle + |\uparrow, 0, 1, 1, 0\rangle + |\downarrow, 1, 0, 1, 0\rangle \right] \\ |\gamma_2\rangle = |\uparrow, 1, 0, 1, 0\rangle \\ |\gamma_3\rangle = |\downarrow, 0, 1, 0, 1\rangle \quad (31) \\ E_{\gamma_0} = E_{\gamma_1} = E_{\gamma_2} = E_{\gamma_3} = \frac{\alpha}{2} \quad (32)$$

Note this spin Hamiltonian has total 8 states, where ground state is double degenerate with energy $-\alpha$ and the lowest excited state with energy 0 is also double degenerate and the second excited state with energy $\alpha/2$ is 4-fold degenerate.

On top of the zero mode spin-channel interactions we want to consider the effect of the momentum space kinetic energy term $\sum_k \epsilon_k (n_{1,k} + n_{2,k})$. We assume identical dispersion for both the channels. Our zero mode spin Hamiltonian is written in the real space, thus it will be easier to do the perturbation theory on the real space. The Fourier transformation of the momentum space kinetic energy term leads to real space hopping term. For simplicity we will be considering the nearest neighbor hoppings between sites within each channel. The real space hopping contributes to the Hamiltonian H_X with the hopping strength is

$$H_X = -t \sum_{\substack{\langle 1, l_1 \rangle \\ \langle 2, l_2 \rangle}} (c_{1,\sigma}^\dagger c_{l_1,\sigma} + c_{2,\sigma}^\dagger c_{l_2,\sigma} + \text{h.c.}) \quad (33)$$

Here l_i represents the nearest site to the origin of the i^{th} channel. Thus here we are interested in studying the Hamiltonian $\mathcal{H}_0 + H_X$. The perturbing term H_X contains scattering which takes the states of the zero mode Hamiltonian out of spin-channel scattering space. Thus before we start the perturbation theory we must expand the basis of the zero mode Hamiltonian itself. Thus the basis which contains the nearest neighbor sites of both the channels is given as

$$\{\mathcal{B}_{\text{extnd}}\} \equiv \{\mathcal{B}\} \otimes \{ |n_{1,\uparrow}, n_{1,\downarrow}, n_{2,\uparrow}, n_{2,\downarrow}\rangle \} \\ = \{ |S_{\text{imp}}^z, n_{1,\uparrow}, n_{1,\downarrow}, n_{2,\uparrow}, n_{2,\downarrow}\rangle \otimes |n_{1,\uparrow}, n_{1,\downarrow}, n_{2,\uparrow}, n_{2,\downarrow}\rangle \} \quad (34)$$

Note there are $2^4 = 16$ elements in the subspace $\{ |n_{1,\uparrow}, n_{1,\downarrow}, n_{2,\uparrow}, n_{2,\downarrow}\rangle \}$. In the spin-channel basis $\{\mathcal{B}\}$

there are 2 degenerate ground states. Because the unperturbed Hamiltonian \mathcal{H}_0 has no scattering terms outside the spin sector, in the extended basis $\{\mathcal{B}_{\text{extnd}}\}$ the total ground state degeneracy becomes simply $2 \times 2^4 = 32$ fold. Thus the total Hamiltonian we are interested in is

$$\mathcal{H} = \frac{\alpha \hbar}{2} \vec{S}_d \cdot \sum_{i=\{1,2\}} \sum_{\alpha, \beta \in \{\uparrow, \downarrow\}} c_{i\alpha}^\dagger \vec{\sigma}_{\alpha\beta} c_{i\alpha} - t \sum_{i=\{1,2\}} \sum_{\langle i, l_i \rangle} (c_{i,\sigma}^\dagger c_{l_i,\sigma} + \text{h.c.}) \quad (35)$$

There are two degenerate states $|\alpha_1\rangle$ and $|\alpha_2\rangle$. Thus the two degenerate ground states in the above basis is given as,

$$|\tilde{\alpha}_0\rangle = |\alpha_0\rangle \otimes |n_{1,\uparrow}, n_{1,\downarrow}, n_{2,\uparrow}, n_{2,\downarrow}\rangle \quad (36)$$

$$|\tilde{\alpha}_1\rangle = |\alpha_1\rangle \otimes |n_{1,\uparrow}, n_{1,\downarrow}, n_{2,\uparrow}, n_{2,\downarrow}\rangle \quad (37)$$

Using degenerate perturbation theory we calculate the first and the second order corrections to the Hamiltonian. The first order and the second order low energy effective Hamiltonian is given as

$$H^{(1)} = \sum_{ij} |\alpha_i\rangle \langle \alpha_i | V | \alpha_j \rangle \langle \alpha_j | \\ H^{(2)} = \sum_{ij} \sum_l |\alpha_i\rangle \frac{\langle \alpha_i | V | \mu_l \rangle \langle \mu_l | V | \alpha_j \rangle}{E_0 - E_l} \langle \alpha_j | \quad (38)$$

where $|\alpha_i\rangle$ represents the ground states with energy E_0 and μ_l represents the excited states with energy E_l . One can easily check the diagonal term in effective Hamiltonian coming from the 1st or any odd order correction is zero as the state never returns to the original starting state via odd number of scatterings. For the case of first order it is easy to see that the off-diagonal correction to the effective Hamiltonian is also zero. Thus the lowest order where we expect the non-zero corrections is the second order. We first calculate the diagonal correction in the first order

$$H_{\text{diag}}^{(2)} = \sum_i \sum_l |\alpha_i\rangle \langle \alpha_i | \frac{|\langle \alpha_i | V | \mu_l \rangle|^2}{E_0 - E_l} \quad (39)$$

As single scattering always takes the state outside of the spin-channel, the state $V|\alpha_i\rangle$ is an excited state with energy 0 ($= E_l$). This α_i representing the ground state can take total 32 configurations as discussed. Out of this 32 states $|\alpha_0\rangle \otimes |n_{1,\uparrow}, n_{1,\downarrow}, n_{2,\uparrow}, n_{2,\downarrow}\rangle$ can take 16 configurations and $|\alpha_1\rangle \otimes |n_{1,\uparrow}, n_{1,\downarrow}, n_{2,\uparrow}, n_{2,\downarrow}\rangle$ can take the rest 16 configurations. Note the two degenerate states $|\tilde{\alpha}_0\rangle$ and $|\tilde{\alpha}_1\rangle$ is labeled by the J^z eigenvalue $\pm 1/2$ respectively, where J^z is the total z-component of the impurity and zero modes of the baths. Thus these 32 ground states can be separated in two 16 state groups labeled by the $J^z = \pm 1/2$ eigenvalues. For the $J^z = 1/2$ sector we get the effective Hamiltonian

$$H_{\text{diag}}^{(2)}(1/2) = \frac{2t^2}{E_0} \hat{I} - \frac{2t^2}{3E_0} \left[(n_{1\uparrow} - n_{1\downarrow}) + (n_{2\uparrow} - n_{2\downarrow}) \right] \quad (40)$$

Similarly for $J^z = -1/2$ we get

$$H_{\text{diag}}^{(2)}(1/2) = \frac{2t^2}{E_0} \hat{I} + \frac{2t^2}{3E_0} \left[(n_{1\uparrow} - n_{1\downarrow}) + (n_{2\uparrow} - n_{2\downarrow}) \right] \quad (41)$$

Thus the diagonal part of the effective Hamiltonian in the second order is

$$\begin{aligned} H_{diag}^{(2)} &= \sum_i \sum_l |\alpha_i\rangle \langle \alpha_i| \frac{|\langle \alpha_i | V | \mu_l \rangle|^2}{E_0 - E_l} \\ &= -4t^2 \hat{I} \end{aligned} \quad (42)$$

where \hat{I} is the identity operator made out of all $2^4 = 16$ possible number operator combinations as shown below

$$\hat{I} = \sum_{\hat{o}=\hat{n},(1-\hat{n})} \hat{o}_{l_1,\uparrow} \hat{o}_{l_1,\downarrow} \hat{o}_{l_2,\uparrow} \hat{o}_{l_2,\downarrow} \quad (43)$$

Now we are interested in calculating the offdiagonal term present in the second order low energy effective Hamiltonian which is given as

$$\begin{aligned} H_{off}^{(2)} &= \sum_{i \neq j} \sum_l |\alpha_i\rangle \frac{\langle \alpha_i | V | \mu_l \rangle \langle \mu_l | V | \alpha_j \rangle}{E_0 - E_l} \langle \alpha_j | \\ &= -\frac{8t^2}{3} \left[(S_1^z)^2 c_{2\uparrow}^\dagger c_{2\downarrow} \left(c_{l_1\uparrow} c_{l_1\downarrow}^\dagger + c_{l_2\uparrow} c_{l_2\downarrow}^\dagger \right) \right. \\ &\quad \left. + (S_2^z)^2 c_{1\uparrow}^\dagger c_{1\downarrow} \left(c_{l_1\uparrow} c_{l_1\downarrow}^\dagger + c_{l_2\uparrow} c_{l_2\downarrow}^\dagger \right) \right] + \text{h.c.} \end{aligned} \quad (44)$$

A. Studying the LEH

1. Self energy and Specific heat

Here we study this LEH to understand various features like change in the self-energy, and the nature of the momentum space Hamiltonian. To calculate the self energy we use Hartree-Fock to get the shift in the kinetic energy. In the real space one can get a diagonal piece from the above equation by using the Fermionic anticommutation relations.

$$H_{eff}^{off,(2)}|_{diag} = -\frac{16t^2}{3} [(S_1^z)^2 + (S_2^z)^2] \quad (45)$$

We can find the corresponding momentum space Hamiltonian by doing the Fourier transformation to the Hamiltonian $H_{eff}^{off,(2)}|_{diag}$.

$$\begin{aligned} H_{eff}^{off,(2)}|_{diag} &= -\frac{4t^2}{3} \frac{1}{N} \left[\sum_{k,\sigma} n_{k\sigma} \left(1 - \frac{1}{N} \sum_{k_2} n_{k_2,-\sigma} \right) \right. \\ &\quad \left. + \sum_{k,\sigma} \tilde{n}_{k\sigma} \left(1 - \frac{1}{N} \sum_{k_2} \tilde{n}_{k_2,-\sigma} \right) \right] \end{aligned} \quad (46)$$

in momentum space we can do again the Hartree-Fock to measure the self-energy

$$\bar{\epsilon}_k - \epsilon_k = \Sigma_k = -\frac{4t^2}{3N} \left(1 - \frac{1}{N} \sum_{k_2} \langle n_{k_2,-\sigma} \rangle \right)$$

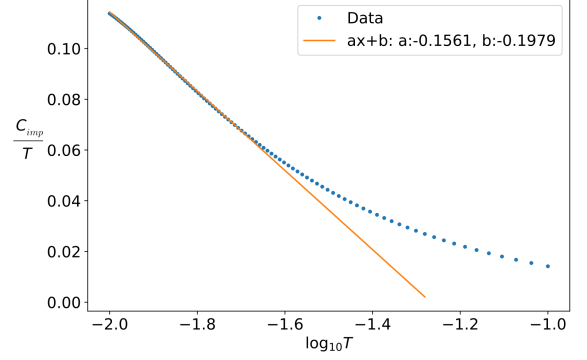


FIG. 11. This figure shows the variation of the impurity specific heat with the temperature.

$$= -\frac{4t^2}{3N^2} \left(1 - \frac{N}{e^{(\epsilon_k - \mu)/k_B T} + 1} \right) \quad (47)$$

Using the self-energy we calculate the impurity specific heat defined as $C_{imp} = C(J^*) - C(0)$.

$$C_{imp} = \sum_{\Lambda, \sigma} \beta \left[\frac{(\bar{\epsilon}_\Lambda)^2 e^{\beta \bar{\epsilon}_\Lambda}}{(e^{\beta \bar{\epsilon}_\Lambda} + 1)^2} - \frac{(\epsilon_\Lambda)^2 e^{\beta \epsilon_\Lambda}}{(e^{\beta \epsilon_\Lambda} + 1)^2} \right] \quad (48)$$

Here we are interested in the low temperature behavior of this impurity specific-heat. Using the self-energy obtained above we calculate the impurity susceptibility for $t = 0.1$ and $\alpha = 1$. The above Fig.11 shows that at low temperature follows logarithmic behavior for the single channel case which is in agreement with the results known in the literature [CITE].

$$\frac{C_{imp}}{T} \propto \log T \quad (49)$$

2. Exact diagonalization

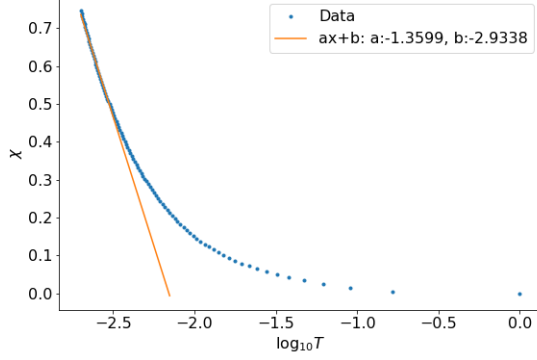
Here we do the exact diagonalization of the NFL Hamiltonian eq.(44) and find the eigenspectrum. Using this eigen spectrum we compute various quantities like χ . We can compute other quantities from this exact diagonalization.

$$\chi = \beta \left[\frac{\sum e^{-\beta \bar{\epsilon}_\Lambda} \langle \bar{S}^z 2 \rangle}{\sum e^{-\beta \bar{\epsilon}_\Lambda}} - \frac{\sum e^{-\beta \epsilon_\Lambda} \langle S^z 2 \rangle}{\sum e^{-\beta \epsilon_\Lambda}} \right] \quad (50)$$

We numerically compute the χ from the definition above and study its dependence on temperature. The above figure shows that at low temperature

$$\chi(T) \propto \log T \quad (51)$$

This matches with the already known result of the literature. The ratio $\frac{\chi(T)}{C_{imp}/T} \sim 8.3$ greater than the Wilson ratio. One can check this quantity for larger system size and at lower temperature for a better agreement.



3. Momentum space structure of the NFL

Direct fourier transformation on the $H_{eff}^{(2)}$ leads to the LEH in momentum space showing higher order scatter-

Thus we get the diagonal correction of effective Hamiltonian in the momoentum spcae,

$$\begin{aligned} H_{eff}^{(2)} \Big|_{diag} &= -\frac{8t^2}{3} \left[\frac{1}{4N} \sum_{k,\sigma} n_{k\sigma} - \frac{1}{2N^2} \sum_{k_1,k_2} n_{k_1\uparrow} n_{k_2\downarrow} \right] \times \left[\frac{1}{N^2} \sum_{k_1,k_2} \tilde{n}_{k_1\uparrow} (1 - \tilde{n}_{k_2\downarrow}) + (\uparrow \rightarrow \downarrow) \right] \times 2 \\ &= -\frac{4t^2}{3} \left[\frac{1}{N} \sum_{k,\sigma} n_{k\sigma} - \frac{1}{N^2} \sum_{k_1,k_2,\sigma} n_{k_1,\sigma} n_{k_2,-\sigma} \right] \times \left[\frac{1}{N} \sum_{k_1,\sigma} \tilde{n}_{k_1,\sigma} - \frac{1}{N^2} \sum_{k_1,k_2,\sigma} \tilde{n}_{k_1,\sigma} \tilde{n}_{k_2,-\sigma} \right], \end{aligned} \quad (53)$$

Where $n_{k,\sigma}$ and $\tilde{n}_{k,\sigma}$ are the occupation of the states k, σ corresponding to two different channels. One can define the charge density of i^{th} channel with the spin- σ as

$$q_{i,\sigma} = \frac{Q_{i,\sigma}}{N} = \frac{1}{N} \sum_k n_{k,\sigma} \quad (54)$$

Thus one can re-write the above equation as,

$$H_{eff}^{(2)} \Big|_{diag} = -\frac{4t^2}{3} \prod_{i=\{1,2\}} \left[q_{i,\uparrow} + q_{i,\downarrow} - q_{i,\uparrow} q_{i,\downarrow} \right] \quad (55)$$

This shows the inter-channel charge coupling present in the excitation spectrum.

B. Third order

Thus the off-diagonal part of the effective low energy Hamiltonian in the 2nd order is

$$\begin{aligned} H_{eff,off}^{(2)} &= \frac{J^2}{\alpha} \left[c_{1\uparrow} c_{1\downarrow}^\dagger \left(-\alpha\beta \left\{ \Sigma_{3,2} + c_{3\uparrow}^\dagger c_{3\downarrow} c_{2\uparrow}^\dagger c_{2\downarrow}^\dagger \right\} - \alpha\gamma \Omega_{3,2} \right) \right. \\ &\quad + c_{2\uparrow} c_{2\downarrow}^\dagger \left(-\alpha\beta \left\{ \Sigma_{1,3} + c_{1\uparrow}^\dagger c_{1\downarrow} c_{3\uparrow}^\dagger c_{3\downarrow}^\dagger \right\} - \alpha\gamma \Omega_{1,3} \right) \\ &\quad \left. + c_{3\uparrow} c_{3\downarrow}^\dagger \left(-\alpha\beta \left\{ \Sigma_{2,1} + c_{2\uparrow}^\dagger c_{2\downarrow} c_{1\uparrow}^\dagger c_{1\downarrow}^\dagger \right\} - \alpha\gamma \Omega_{2,1} \right) + \text{h.c.} \right] \end{aligned}$$

ing involving electronic degree of freedoms. The Fourier transformation of the electronic creation operator gives

Thus one gets,

$$\begin{aligned} (S_1^z)^2 \Big|_{diag} &= \frac{1}{4} (n_{1\uparrow} - n_{1\downarrow})^2 \Big|_{diag} = (n_{1\uparrow} + n_{1\downarrow} - 2n_{1\uparrow} n_{1\downarrow}) / 4 \Big|_{diag} \\ &= \frac{1}{4N} \sum_{k,\sigma} n_{k\sigma} - \frac{1}{4N^2} \sum_{k_1,k_2,\sigma} n_{k_1,\sigma} n_{k_2,-\sigma} \end{aligned} \quad (52)$$

We know that the diagoanl peice of the following terms is given as

Thus the diagonal piece coming from the term,

$$c_{2\uparrow}^\dagger c_{2\downarrow} \left(c_{l_1\uparrow} c_{l_1\downarrow}^\dagger + c_{l_2\uparrow} c_{l_2\downarrow}^\dagger \right) \Rightarrow \frac{1}{N} \sum_{k,k'} e^{-i(k-k')r} c_{k,\sigma}^\dagger c_{k',-\sigma}$$

$$\otimes \left[c_{l_1\uparrow}^\dagger c_{l_1\downarrow} + c_{l_2\uparrow}^\dagger c_{l_2\downarrow} + c_{l_3\uparrow}^\dagger c_{l_3\downarrow} + \text{h.c.} \right]$$

$$\Sigma_{i,j} = -8S_i^z S_j^z [(C_i^z - C_j^z)^2 + (S_i^z - S_j^z)^2] \quad (56)$$

$$\Omega_{i,j} = 4S_i^z S_j^z (C_i^z + C_j^z)(C_j^z + C_i^z) \quad (57)$$

Now we are interested calculating the diagonal part of the effective Hamiltonian. In this case we get non-zero contribution from all the three ground states $|J^z = -1\rangle$, $|J^z = 0\rangle$ and $|J^z = 1\rangle$. We get the diagonal contribution to the LEH in the second order is

$$H_{eff,diag}^{(2)} = -\frac{7.2J^2}{\alpha} \hat{I} \quad (58)$$

The contribution associater to different ground states $|J^z = -1\rangle$, $|J^z = 0\rangle$ and $|J^z = 1\rangle$ is given respectively as $-\frac{2.4J^2}{\alpha} \hat{I} + \hat{\mathcal{F}}$, $-\frac{2.4J^2}{\alpha} \hat{I}$, $-\frac{2.4J^2}{\alpha} \hat{I} - \hat{\mathcal{F}}$, where $\hat{\mathcal{F}}$ is a function of diagonal number operators of the nearest-neighbor site degree of freedom l_1, l_2, l_3 etc.

III. CHANNEL ANISOTROPY AND BERRY PHASE

noindent Here we start with the multi-channel zero mode model, any put anisotropy in the coumping α . Let's

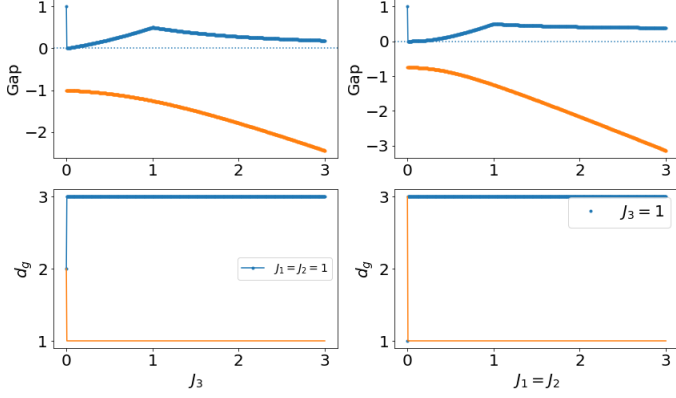


FIG. 12. Add proper figures.

start with a simple possibility where there are two coupling α_1 and α_2 , where both the α_1, α_2 are positive and negative. We write down the Hamiltonian corresponding to these cases,

$$H_{an} = \alpha_1 \vec{S}_d \cdot \vec{S}_1 + \alpha_2 \vec{S}_d \cdot \vec{S}_2 \quad (59)$$

There are K_1 and K_2 numbers of outer spins respectively which are connected to the central spin with the coupling with α_1 and α_2 . Thus the total number of channels is $K = K_1 + K_2$. When $\alpha_1 = 0, \alpha_2 \neq 0$ then we end up with a multi-channel problem with K_1 number of channels thus the degeneracy becomes K_1 and for $\alpha_2 = 0, \alpha_1 \neq 0$ the degeneracy becomes K_2 . Though this can be trivially verified from the above Hamiltonian, the important property of this problem is the robustness of the ground state degeneracy. As we vary α_1 and α_2 , to study the ground state degeneracy. We find that for any values α_1 and α_2 such that $\alpha_1/\alpha_2 \neq 0$ and $\alpha_2/\alpha_1 \neq 0$ the ground state degeneracy remains constant at $K = K_1 + K_2$. It is known in the literature [CITE someone] that under above mentioned an-isotropy the low energy effective Hamiltonian becomes a multi-channel Kondo with smaller number of channels.

A. Berry phase

We start with the general multi-channel problem where the coupling strength with the central impurity spin of the i^{th} channel is α_i . Here our goal is to use the robustness of the ground state degeneracy of the star-graph model to build the gauge theoretic structure. We define the hamiltonian of star-graph as

$$H_{st}(\vec{\alpha}) = \sum_{i=1}^K \alpha_i \vec{S}_d \cdot \vec{S}_i \quad (60)$$

We define the K -dimensional coupling space where one possible configuration is denoted by the vector $\vec{\alpha} = \mathcal{N}(\alpha_1, \alpha_2, \dots, \alpha_K)$, where \mathcal{N} is the normalization factor,

where $\alpha_i \neq 0$ and $\alpha_i \neq \infty, \forall i \in [K]$. The Schrodinger equation for the Hamiltonian reads as

$$H_{st}(\vec{\alpha}) \Psi_m(\vec{\alpha}) = E_m(\vec{\alpha}) \Psi_m(\vec{\alpha}) \quad (61)$$

Here, we are only interested in the ground state manifold, where the energy of the state is E_g . Due to the degeneracy (K -fold) there are K states in that manifold with the same energy labeled by say k . Thus the equation for the ground state manifold reads

$$H_{st}(\vec{\alpha}) \Psi_g^k(\vec{\alpha}) = E_g(\vec{\alpha}) \Psi_g^k(\vec{\alpha}) \quad (62)$$

One can do a simple transformation to make the ground state energy zero. Thus

$$H_{st}(\vec{\alpha}) \Psi_g^k(\vec{\alpha}) = 0$$

$$i\hbar \frac{\partial \Psi(\vec{\alpha}(t))}{\partial t} = H_{st}(\vec{\alpha}(t)) \Psi(\vec{\alpha}(t)) = 0 \quad (63)$$

As the ground state is degenerate one can take any linear combinations of the states to create arbitrary basis state (orthogonal and normalized). Let's say $\psi_k(\vec{\alpha}(t))$ represents the arbitrary basis state, which should be smooth in t locally and follow the Schrodinger equation,

$$H_{st}(\vec{\alpha}(t)) \psi_k(\vec{\alpha}(t)) = 0 \quad (64)$$

We here try to find the solution of the Eq.(63), we start with the trial solution

$$\eta_a(t) = U_{ab}(t) \psi_b(t) \quad (65)$$

Finish this section.

IV. TOPOLOGY AND GAUGE THEORY

A. In our case of stargraph model

In our model we find two string operators $\hat{\mathbb{Z}} = \sigma_d^z \prod_{i=1}^K \sigma_i^z$ and $\hat{\mathbb{X}} = \sigma_d^x \prod_{i=1}^K \sigma_i^x$, where K is the number of channels. These two operators commute with the Hamiltonian H of this stargraph.

$$[H, \hat{\mathbb{Z}}] = 0 = [H, \hat{\mathbb{X}}] \quad (66)$$

We can rewrite the operators in the form of a twist operator as,

$$\mathbb{Z} = \sigma_d^z \prod_{i=1}^K \sigma_i^z = \frac{1}{i^{K+1}} e^{i\frac{\pi}{2} \sum_i \sigma_i^z} = e^{i\frac{\pi}{2} \sum_i (\sigma_i^z - 1)} = e^{i\pi \sum_i (S_i^z - \frac{1}{2})} \quad (67)$$

Similarly one can rewrite

$$\mathbb{X} = \sigma_d^x \prod_{i=1}^K \sigma_i^x = e^{i\pi \sum_i (S_i^x - 1/2)} \quad (68)$$

The commutation between these two operators are given as

$$[\mathbb{Z}, \mathbb{X}] = \prod_{i=d,1}^K \sigma_i^z \prod_{i=d,1}^K \sigma_i^x \left[1 - (-1)^{K+1} \right] \quad (69)$$

Thus we can see that for case where number of channels is odd that means $K+1$ is even, these two string operators commutes with each other forming a CSCO (H, \hat{Z}, \hat{X}) . On the other hand in the case where number of channels K is even, these two operators does not commute with each other. Let's discuss these two cases separately.

1. $K = \text{odd}$

In the case where the number of channels K is odd, we get $[\mathbb{Z}, \mathbb{X}] = 0$. Thus forming a CSCO with elements H, \hat{Z}, \hat{X} . Thus one can label the states using the eigenvalues of these operators. For the case of $K = 2n+1$ channels, the ground state is K fold degenerate and labeled by the eigenvalues of the operator J^z which can take values $J^z = [-J, -J+1, \dots, J-1, J]$, where $J = (K-1)/2$. For the odd channel case J^z can take zero value. Thus

$$|J^z\rangle \equiv |Z, X\rangle \quad (70)$$

where the eigenvalue Z can take $(K+1)/2$ values related to the absolute $|J^z|$ but cannot distinguish between $\pm J^z$ and X can take two values ± 1 which breaks the degeneracy between $\pm J^z$. Thus these two operators together can uniquely label all the degenerate states.

B. Main

We can define two operators, for convenience let's call translation and twist operator respectively, \hat{T} and \hat{O} defined as

$$\hat{T} = e^{i\frac{2\pi}{K}\hat{\Sigma}}, \quad \hat{O} = e^{i\hat{\phi}}, \quad \hat{\Sigma} = [\hat{J}^z - (K-1)/2] \quad (71)$$

One can see that the generators of these above operators commutes with the Hamiltonian, $[H, J^x] = [H, J^z] = [H, J^y] = 0$. In the large channel limit we can do semiclassical approximation. As J^x and J^y both commutes with the Hamiltonian H then $[H, J^y(J^x)^{-1}] = 0$, and any non-singular function of these operators must also commute with the Hamiltonian, thus we can say that

$$\hat{\phi} = \tan^{-1}(\hat{J}^y(\hat{J}^x)^{-1}) \quad (72)$$

Then we get $[\hat{\phi}, \hat{H}] = 0$. We can label the ground states by the eigen values of the translation operator \hat{T} . We can also use the ground states labeled by the eigenvalues of J^z (M , say). Then the ground states are

$$|M_1\rangle, |M_2\rangle, |M_2\rangle, \dots, |M_K\rangle \quad (73)$$

Then the operations of the translation operators on this states are give below.

$$\hat{T}|M_i\rangle = e^{i\frac{2\pi}{K}[M_i - (K-1)/2]}|M_i\rangle = e^{i2\pi\frac{p_i}{K}}|M_i\rangle, \quad p_i \in [1, \dots, K] \quad (74)$$

Now we check the braiding rule between the twist and the translation operators, we find

$$\hat{T}\hat{O}\hat{T}^\dagger\hat{O}^\dagger = e^{\frac{2\pi}{K}[i\hat{\Sigma}, i\hat{\phi}]} = e^{i\frac{2\pi}{K}}$$

$$\hat{T}\hat{O}^m\hat{T}^\dagger\hat{O}^{\dagger m} = e^{\frac{2\pi}{K}[i\hat{\Sigma}, im\hat{\phi}]} = e^{i2\pi\frac{m}{K}} \quad (75)$$

Next we shown that the states $\hat{O}^m|M_i\rangle$ are orthogonal to each other and with the state $|M_i\rangle$.

$$\begin{aligned} \hat{T}\hat{O}^m|M_i\rangle &= \hat{O}^m\hat{T}e^{i2\pi\frac{m}{K}}|M_i\rangle = \hat{O}^me^{i\frac{2\pi(m+p_i)}{K}}|M_i\rangle \\ \hat{T}\left(\hat{O}^m|M_i\rangle\right) &= e^{i\frac{2\pi(m+p_i)}{K}}\left(\hat{O}^m|M_i\rangle\right) \end{aligned} \quad (76)$$

Thus different states $\hat{O}^m|M_i\rangle$ for different m are labeled by different eigenvalues of the translation operations, thus are orthogonal to each other. Now as the twist operator \hat{O} commutes with the Hamiltonian, thus we can see that

$$\langle M_i|\hat{O}^{m\dagger}H\hat{O}^m|M_i\rangle = \langle M_i|H|M_i\rangle \quad (77)$$

Thus the energy eigen values of all the orthogonal states are equal. Thus all the K mutually orthogonal states are degenerate.

C. Fractional statistics

V. LOCAL MOTT LIQUID

Using the unitary renormalization group method we decouple the impurity spin from the zero-modes of the K channels. We start with the Hamiltonian. The zero mode Hamiltonian is a stargraph model

$$\begin{aligned} H &= \alpha \vec{S}_d \cdot \sum_i \vec{S}_i = \alpha S_d^z S^z + \frac{\alpha}{2}(S_d^+ S^- + S_d^- S^+) \\ H_D &= \alpha S_d^z S^z, \quad H_X = \frac{\alpha}{2}(S_d^+ S^- + S_d^- S^+) \end{aligned} \quad (78)$$

Here we want to remove the quantum fluctuations between the impurity spin and the rest, for that we do one step URG.

$$\Delta H = H_X \frac{1}{\hat{\omega} - H_D} H_X \quad (79)$$

In the zero mode IR fixed point ground state (stargraph) of multi-channel ground state J^z is a good quantum number but S^z is not. There is no net S^z field thus in the ground state manifold the average S^z is vanishing, $\langle S^z \rangle = 0$. We use this expectation value to replace the denominator of the above RG equation.

Thus we get the effective Hamiltonian is

$$\begin{aligned} H_{eff} &= \frac{\alpha^2 \Gamma_\uparrow}{2} (S^2 - S^{z2}) (\tau_d^2 - \tau_d^z) \\ &= \beta_\uparrow(\alpha, \omega_\uparrow) (S^2 - S^{z2}) (\tau_d^2 - \tau_d^z), \quad \frac{\alpha^2 \Gamma_\uparrow}{2} = \beta_\uparrow(\alpha, \omega_\uparrow) \end{aligned}$$

$$= \frac{\beta_{\uparrow}(\alpha, \omega_{\uparrow})}{2}(S^2 - S^z) = \frac{\beta_{\uparrow}(\alpha, \omega_{\uparrow})}{4}(S^+ S^- + S^- S^+) \quad (80)$$

. In terms of the electronics degree of freedom this looks

$$H_{eff}(\omega, \alpha) = \frac{\beta_{\uparrow}(\alpha, \omega_{\uparrow})}{4} \sum_{i \neq j} \sum_{\substack{\alpha_i, \beta_i \in \{\uparrow, \downarrow\} \\ \alpha_j, \beta_j \in \{\uparrow, \downarrow\}}} \vec{\sigma}_{\alpha_i \beta_i} \vec{\sigma}_{\alpha_j \beta_j} c_{0\alpha_i}^{(i)\dagger} c_{0\beta_i}^{(i)} c_{0\alpha_j}^{(j)\dagger} c_{0\beta_j}^{(j)} + \text{h.c.}$$

We will study the two cases of the above effective Hamiltonian depending on the sign of the prefactor. The case $\frac{\beta_{\uparrow}(\alpha, \omega_{\uparrow})}{2} > 0$ is the ferromagnetic case where in the ground state S takes the minimum value and S^z takes the possible maximum value. For an example, in the case of K channels the minimum value of S will be 0 or $1/2$ for K being *even* or *odd* respectively. We will be interested in the other case where $\frac{\beta_{\uparrow}(\alpha, \omega_{\uparrow})}{2} < 0$. In this case the ground state corresponds to S being maximum and S^z being minimum case. The effective Hamiltonian can be re-written in this case as

$$H_{eff} = -\frac{|\beta_{\uparrow}(\alpha, \omega_{\uparrow})|}{4}(S^+ S^- + S^- S^+) \quad (82)$$

The CSCO of this Hamiltonian contains H, S, S^z . In the ground state S is maximum thus $S = K/2$ and S^z can take $2S + 1 = K + 1$ values. Thus in the largest S sector there are $K + 1$ states. One can explore different states via a flux tuning mechanism. In the presence of the flux the Hamiltonian looks like

$$H = -\frac{|\beta_{\uparrow}(\alpha, \omega_{\uparrow})|}{2} S^2 + \frac{|\beta_{\uparrow}(\alpha, \omega_{\uparrow})|}{2} \left(S^z - \frac{\Phi}{\Phi_0} \right)^2 \quad (83)$$

$$\begin{aligned} S_d^z = +1/2, \{J^z\} &= \{-(K-1)/2, -(K-3)/2, \dots, (K-1)/2, (K+1)/2\} \\ S_d^z = -1/2, \{J^z\} &= \{-(K+1)/2, -(K-1)/2, \dots, (K-3)/2, (K-1)/2\} \end{aligned} \quad (85)$$

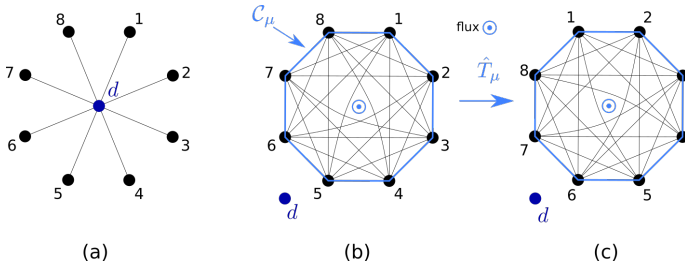


FIG. 13. This figure shows the comparison and mapping between the stargraph degenerate ground states and the states of the quantum fluctuation resolved all to all model.

Thus one can see that the ground state degeneracy of the stargraph model with different J^z values gets lifted in the quantum fluctuation resolved all-to-all model. Depending on the value of the flux there is an unique ground

A. Mapping with the degenerate ground state of stargraph

Now we are going to discuss about the mapping of K degenerate ground states of the parent K -channel stargraph model with the states of this flux-inserted all-to-all model obtained by removing impurity-bath quantum fluctuations. In the stargraph model the K ground states is labeled by the J^z eigenvalues. In the stargraph ground state $J = S - 1/2$, where S is maximum with value $S = K/2$. Thus there are $2J + 1 = K$ unique J^z eigenvalues corresponding to the J sector labeling different eigenstates. The ground state Hilbert space \mathcal{H}_{gr} contains the J^z states

$$\mathcal{H}_{gr} = \{-(K-1)/2, -(K-3)/2, \dots, (K-3)/2, (K-1)/2\} \quad (84)$$

Now we shift our focus to the low energy Hamiltonian with flux shown in the eq.(83). We can see that there is a unique ground state labeled by the S^z eigenvalues. By tuning the flux (Φ/Φ_0) one can go from one ground state to another ground state or change the ground state. We can see that ground state energy is independent of the configuration of the impurity spin ($S_d^z = \pm 1/2$). Thus for each of those two possibilities we get two set of ground states labeled by the S^z and the $J^z = S^z + S_d^z$. Similar to the stargraph case here also the ground state corresponds to the largest S sector, $S = K/2$. Thus total number of S^z eigenvalues are $2S + 1 = K + 1$ taking values $\{-K/2, -(K-2)/2, \dots, (K-2)/2, K/2\}$.

state labeled by the J^z value. Tuning the flux one can go from one ground state to the other. This shows how one can explore different degenerate ground states of the parent stargraph model via the insertion of flux in the corresponding quantum fluctuation resolved all-to-all effective Hamiltonian.

B. Local mott-liquid

Impurity-bath quantum fluctuation resolved all-to-all effective Hamiltonian is written in terms of the spinor spin operators which are individually made out of two electronic degree of freedom. This is defined as $\vec{S}_i = \frac{\hbar}{2} \sum_{\alpha, \beta \in \{\uparrow, \downarrow\}} c_{0\alpha}^{(i)\dagger} \vec{\sigma}_{\alpha\beta} c_{0\beta}^{(i)}$. In the above eq.(81) we see the $U(1)$ symmetry of the effective Hamiltonian. Using the

spinor representation one can see the spin creation operation in terms of the electronic degree of freedom is

$$S_i^+ = \frac{\hbar}{2} \sum_{\alpha, \beta \in \{\uparrow, \downarrow\}} c_{0\alpha}^{(i)\dagger} \sigma_{\alpha\beta}^+ c_{0\beta}^{(i)} = \frac{\hbar}{2} c_{0\uparrow}^{(i)\dagger} c_{0\downarrow}^{(i)} \quad (86)$$

$$S_i^z = \frac{\hbar}{2} \sum_{\alpha, \beta \in \{\uparrow, \downarrow\}} c_{0\alpha}^{(i)\dagger} \sigma_{\alpha\beta}^z c_{0\beta}^{(i)} = \frac{\hbar}{2} (c_{0\uparrow}^{(i)\dagger} c_{0\uparrow}^{(i)} - c_{0\downarrow}^{(i)\dagger} c_{0\downarrow}^{(i)}) \quad (87)$$

Thus this spinor spins are nothing but the Anderson pseudospin formulation in the spin channel. The spin creation operator (S_i^+) shows simultaneous creation of an electron-hole pair at the realspace origin of the i^{th} channel. The condensation of such electron-hole pairs has already been shown in [CITE: mott]. Thus for this effective Hamiltonian one can define twist-translation operations to construct the gauge theory and unveiling any hidden degeneracy. Let's recall the effective Hamiltonian

$$H_{eff} = \frac{\beta_{\uparrow}(\alpha, \omega_{\uparrow})}{4} \left[\sum_{ij} S_i^+ S_j^- + \text{h.c.} \right], \quad (88)$$

where i, j are the channel indices. Due to the all-to-all nature of the connectivity one can draw total $K!$ possible unique closed paths (\mathcal{C}_{μ}) where each nodes (channels) is being touched only once. One can thus define total $K!$ number of translation operators (\hat{T}_{μ}) which keeps the Hamiltonian invariant. Let's define one of such translation operator \hat{T}_{μ} which gives a periodic shifts along the closed path \mathcal{C}_{μ} . Let's define twist operator along the path \mathcal{C}_{μ}

$$\hat{\mathcal{O}}_{\mu} = \exp(i \frac{2\pi}{K} \sum_{j=1}^K j S_j^z), \quad \hat{T}_{\mu} = e^{i \hat{P}_{\mu}^{cm}} \quad (89)$$

The operation of the translation operator \hat{T}_{μ} is defined as $\hat{T}_{\mu} S_j^z = S_{j+1}^z$ where $j+1$ and j are the nearest neighbor on the closed path \mathcal{C}_{μ} . Then the braiding rule between the twist and translation operators are give as

$$\begin{aligned} \hat{T}_{\mu} \hat{\mathcal{O}}_{\mu} \hat{T}_{\mu}^{\dagger} \hat{\mathcal{O}}_{\mu}^{\dagger} &= \exp\{i[2\pi S_1^z - \frac{2\pi}{K} S^z]\} \\ &= \exp\{i[\pm\pi - \frac{2\pi}{K} S^z]\} = \exp(i \frac{2\pi p}{q}) \quad (90) \end{aligned}$$

The availability of the non-trivial braiding statistics between these twist and translation operators is possible if $p \neq 0$ and $q \neq \infty$. Further simplification leads to the condition

$$\begin{aligned} \pm\pi - \frac{2\pi}{K} S^z = \frac{2\pi p}{q} &\Rightarrow \pm\frac{1}{2} - \frac{S^z}{K} = \frac{p}{q} \\ \frac{(\pm K - 2S^z)}{2K} &= \frac{p}{q}, \quad (91) \end{aligned}$$

p, q are mutual prime, We know that the S^z can take values $(\mp K/2 \pm m)$ where m is a integer $0 \leq m \leq K$. Putting this value in the above equation leads to two possible solutions.

$$\frac{(K-m)}{K} = \frac{p}{q}, \quad \frac{-m}{K} = \frac{p}{q} \quad (92)$$

For the first case $K-m = p$ we can see that $m = K$ makes p trivial has the allowed values $m = 0, \dots, K-1$, which represents the corresponding S^z eigen values

$$-K/2, -K/2+1, \dots, K/2-2, K/2-1 \quad (93)$$

Similarly the second case implies, $-m = p$, but $p = 0$ is not allowed as this makes the braiding statistics trivial, thus the possible S^z values are

$$-K/2+1, -K/2+2, \dots, K/2-1, K/2 \quad (94)$$

Thus we get the general braiding statistics between the twist and the translation

$$\hat{T}_{\mu} \hat{\mathcal{O}}_{\mu} \hat{T}_{\mu}^{\dagger} \hat{\mathcal{O}}_{\mu}^{\dagger} = e^{i \frac{2\pi p}{K}} \quad (95)$$

where p corresponds to different S^z states, related as $p = \pm K/2 - S^z$. Thus we can see that there are K possible S^z plateau states in the all-to-all model where each plateaux is K fold degenerate. This p/K is similar to the filling factor of the fractional quantum Hall effect. There are possibly $K!$ pairs of twist and translation operator corresponding to different closed paths \mathcal{C}_{μ} which can probe this degeneracy.

1. Action on the Hamiltonian

Now we find out the action of these twist operators on the Hamiltonian. The Hamiltonian is written as

$$H_{eff} = \frac{\beta_{\uparrow}(\alpha, \omega_{\uparrow})}{2} (S^{x2} + S^{y2}) \quad (96)$$

Due the all-to-all nature of the effective Hamiltonian one can find $K!$ possible relative arrangement of those K channels which keeps the Hamiltonian invariant. Here we briefly discuss the choice of the closed loop \mathcal{C}_{μ} and the insertion of the flux. As shown in the Fig.13(b) we have chosen a particular closed path which crosses all the outer spin only once. We embed that closed loop on a plane and put the flux perpendicular to the plane through the closed loop. One can find a different closed loop where the ordering of the outer spins will be different. The action of the translation operator shifts the outer spins along this closed curve by one step.

The action of the twist operator on the S^x is determined in the following calculation

Thus in the large channel limit

$$\begin{aligned} &\lim_{K \rightarrow \infty} \hat{\mathcal{O}}_{\mu} H_{eff} \hat{\mathcal{O}}_{\mu}^{\dagger} \\ &= \frac{\beta_{\uparrow}(\alpha, \omega_{\uparrow})}{2} \left[\hat{\mathcal{O}}_{\mu} S^x \hat{\mathcal{O}}_{\mu}^{\dagger} \hat{\mathcal{O}}_{\mu} S^x \hat{\mathcal{O}}_{\mu}^{\dagger} + \hat{\mathcal{O}}_{\mu} S^y \hat{\mathcal{O}}_{\mu}^{\dagger} \hat{\mathcal{O}}_{\mu} S^y \hat{\mathcal{O}}_{\mu}^{\dagger} \right] \\ &= \frac{\beta_{\uparrow}(\alpha, \omega_{\uparrow})}{2} [S^{x2} + S^{y2}] = H_{eff} \\ &\lim_{K \rightarrow \infty} [\hat{\mathcal{O}}, H_{eff}] = 0 \quad (97) \end{aligned}$$

The translation operator $\hat{T}_{\mu} = e^{i \hat{P}_{\mu}}$ commutes with the Hamiltonian thus the generator \hat{P}_{μ} also commutes with

the Hamiltonian. We can label the j^{th} state with the eigenvalues of this operator $\hat{\mathcal{P}}_\mu$ as $|p_\mu^j\rangle$. We here discuss the action of these twist and translation operators on these states.

$$\begin{aligned}\hat{T}_\mu |p_\mu^j\rangle &= e^{i\hat{\mathcal{P}}_\mu} |p_\mu^j\rangle = e^{ip_\mu^j} |p_\mu^j\rangle \\ \hat{T}_\mu \hat{\mathcal{O}}_\mu |p_\mu^j\rangle &= \hat{\mathcal{O}}_\mu \hat{T}_\mu e^{i\frac{2\pi m}{K}} |p_\mu^j\rangle, \\ &\text{here } m \text{ signifies different } S^z \text{ plateaux.} \\ &= \hat{\mathcal{O}}_\mu e^{i(\frac{2\pi m}{K} + p_\mu^j)} |p_\mu^j\rangle, \quad \frac{2\pi m}{K} \equiv p_\mu^m\end{aligned}$$

$$\hat{T}_\mu \left(\hat{\mathcal{O}}_\mu |p_\mu^j\rangle \right) = e^{i(p_\mu^m + p_\mu^j)} \left(\hat{\mathcal{O}}_\mu |p_\mu^j\rangle \right) \quad (98)$$

Thus we can see that $\hat{\mathcal{O}}_\mu |p_\mu^j\rangle$ is a state of the traslation operator with different eiven value than $|p_\mu^j\rangle$, thus orthogonal to each other. Simiarly one can generally show that $\langle p_\mu^j | \hat{\mathcal{O}}_\mu^q | p_\mu^j \rangle = 0$, where q is any integer. Also in the large K limit we can see from the eq.(97) that these differnt twisted states has same energy. Which shows that at each plateau state labeled by the S^z eigenvalue has K fold degenerate eigenstates labeled by the eigenvalue of the translation operators. Thus the CSCO is formed by H, S^z, \hat{T} . Thus states can be lebeled as $|E, S_j^z, P_\mu^j\rangle$.