

Self Energy effect in frequency dependent Vertex flow equation

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I. INTRODUCTION

The formulation of the Wetterich equation¹ laid the basis for a class of powerful new methods that generally go under the name of functional Renormalization Group (fRG).^{2,3} Given the generality of the Wetterich equation, different conjugations of the fRG are more or less suited to treat different classes of systems. For example, while the derivative expansion is most suited for bosonic systems, for fermionic systems on a lattice an expansion into the fields is likely the most sensible choice. *Is it correct to write this? Are derivative and field expansion mutually exclusive?*

In the field expansion implementation of the fRG, the two-particle vertex is the core object to study in order to understand the physical properties of the system. However due to its dependence on three frequency and three momentum arguments the two-particle vertex is an extremely difficult object to deal with, unless drastic approximations are made. For fermions on a lattice the momentum-structure of the vertex has been understood while mostly neglecting the vertex frequency-dependence. In particular, the first results have been obtained using the N -patch scheme,⁴⁻⁷ in which all momenta are treated on the same footing. Once an understanding was obtained this way, it was possible to tailor more refined treatments of the momentum dependence, like the form factor decomposition, which focuses on the most relevant physics.^{8,9} *is there some older reference?*

A comparable understanding is lacking when considering both momentum and frequency dependencies of the vertex. The very high computational cost and the conceptual difficulty of interpreting data on the imaginary frequency axis did not allow for a frequency treatment comparable to the momentum one. Indeed in fRG (for lattice models), frequency has been either completely neglected, by assuming a static vertex, or treated in approximate schemes¹⁰ which capture part of the physics only.

In the present paper we start filling in this gap, by presenting fully-fledged fRG results with frequency and momentum dependent vertex. The vertex frequency dependence plays an increasingly important role when the strong coupling regime is approached, as it has been confirmed mostly for impurity models,^{11,12} or in the dynamical mean field theory¹³⁻¹⁵ (and its diagrammatic extensions)¹⁶ framework.¹⁷ Hence a proper treatment of the vertex frequency-dependence is mandatory for meth-

ods that aim at the study of the interplay between fluctuations in all the channels in the presence of strong coupling, like DMF²RG.¹⁸

Our results show the feasibility, and in some respects, the necessity for an improved frequency treatment of the vertex, also at moderate coupling. A simpler treatment of the frequency dependence was already pursued by Husemann *et al.* in Ref. 10, where they have found the emergence of regions in the parameters space characterized by an unusual divergence in the charge channel for finite frequency transfer, which cannot be connected with a phase transition. An analysis of the vertex structure in this region will allow us to explain this feature in terms of simple diagrammatic elements.

Another drawback of a static vertex approximation is that the difficulty in obtaining a frequency dependent self-energy as well its feedback in the flow, unless some further approximation is employed.⁷ On the contrary, in this work, we have evaluated also the flow equation of the self-energy. We will show that the self-energy feedback affects the results also qualitatively: by preventing the charge channel from diverging. This is in contrast with the common belief that the self-energy feedback plays a minor role in the flow.

The paper is structured as follows. In Sec. II we will briefly introduce the model under consideration and the basic equations of the fRG in its expansion in the fields. After discussing in some more detail our treatment for the two-particle vertex in Sec. III, we will move on to the discussion of the main results in Sec. IV. Here we start by a stability analysis, then we show the flow evolution and frequency dependence of the vertex. We then consider the charge divergence with a diagrammatic explanation, before turning to the analysis of the obtained self-energy. We draw our conclusions in Sec. V. In the Appendix VIA we report all the final expression for the vertex flow equations, while in the Appendix VIB we show the frequency dependence also in the pairing channel.

II. FORMALISM

A. Model

The Hubbard model¹⁹ describes spin- $\frac{1}{2}$ fermions with a local interaction:

$$\mathcal{H} = \sum_{i,j,\sigma} t_{ij} c_{i,\sigma}^\dagger c_{j,\sigma} + U \sum_i n_{i,\uparrow} n_{i,\downarrow}, \quad (1)$$

where $c_{i,\sigma}^\dagger$ and $c_{i,\sigma}$ are, respectively, creation and annihilation operators for fermions on site i with spin orientation σ (\uparrow or \downarrow). We consider the two-dimensional case on a square lattice and repulsive interaction $U > 0$ at finite temperature T . The hopping amplitude is restricted to $t_{ij} = -t$ for nearest neighbors, $t_{ij} = -t'$ for next-to-nearest neighbors. Fourier transforming the hopping matrix yields the bare dispersion relation

$$\varepsilon_{\mathbf{k}} = -2t(\cos k_x + \cos k_y) - 4t' \cos k_x \cos k_y. \quad (2)$$

B. Flow equations

In the following paragraph we will provide some details about the functional renormalization group,^{2,20} defining in particular the notation used for the vertex.

Generally speaking, the fRG implements the renormalization group idea in the functional integral formalism. This is done by endowing the bare action with an additional dependence on a scale-parameter Λ ,^{2,20}

$$\mathcal{S}^\Lambda[\bar{\psi}, \psi] = -(\bar{\psi}, G_0^{\Lambda^{-1}} \psi) + \mathcal{S}_{\text{int}}, \quad (3)$$

where \mathcal{S}_{int} is the interaction part, and $(\bar{\psi}, \psi)$ summarizes the summation over all the quantum numbers of the fermionic fields $\bar{\psi}$ and ψ . The scale dependence, acquired through the non-interacting propagator G_0^Λ , generates flow equations (with known initial conditions) for generating functionals defined via functional integrals with the action \mathcal{S}^Λ , such as the generating functional for the connected Green's function and its Legendre transform, the so-called average effective action.¹ The final result is recovered for some final Λ -value restoring the original bare propagator, $G_0^{\Lambda_f} = G_0$, so that the physical action of interest is recovered.

We will apply this approach to the effective action, whose expansions in the fields generates the one-particle irreducible (1PI) vertex functions. By expanding the functional flow equation,¹ one obtains a hierarchy of flow equations for the 1PI functions, involving vertices of arbitrarily high orders. We will restrict ourselves to the two-particle level truncation by retaining only the two lowest nonvanishing orders in the expansion, that is, we consider the flow of the self-energy Σ^Λ and of the two-particle vertex V^Λ , neglecting the effects of higher order vertices.

This truncation restricts the applicability of the approach to the weak-to-moderate coupling regime.²¹ It can be further shown that, at the two-particle level truncation, the

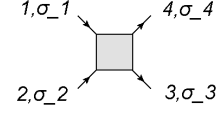


Figure 1: Notation of the two-particle vertex. **placeholder**

fRG sums up efficiently, although approximately, the so-called parquet-diagrams.²⁴

Due to SU(2) symmetry, the self-energy is diagonal in spin-space:

$$\Sigma_{\sigma\sigma'}^\Lambda(k) = \Sigma(k) \delta_{\sigma,\sigma'}, \quad (4)$$

where $k = (\nu, \mathbf{k})$, ν is a fermionic Matsubara frequency and \mathbf{k} a momentum in the first Brillouin zone.

For the notation of the two-particle vertex function $V_{\sigma_1\sigma_2\sigma_3\sigma_4}(k_1, k_2, k_3)$ we refer to Fig. 1, where $k_i = (\mathbf{k}_i, \nu_i)$. The momentum $k_4 = k_1 + k_2 - k_3$ is fixed by momentum conservation. The SU(2) spin-rotation symmetry guarantees that the vertex does not vanish only for six spin combinations: $V_{\uparrow\uparrow\uparrow\uparrow}^\Lambda = V_{\downarrow\downarrow\downarrow\downarrow}^\Lambda$, $V_{\uparrow\downarrow\uparrow\downarrow}^\Lambda = V_{\downarrow\uparrow\downarrow\uparrow}^\Lambda$, and $V_{\uparrow\downarrow\downarrow\uparrow}^\Lambda = V_{\downarrow\uparrow\uparrow\downarrow}^\Lambda$. Finally, due to the SU(2) symmetry and crossing relation one has¹⁷

$$V_{\uparrow\uparrow\uparrow\uparrow}^\Lambda(k_1, k_2, k_3) = V_{\uparrow\downarrow\uparrow\downarrow}^\Lambda(k_1, k_2, k_3) - V_{\uparrow\downarrow\downarrow\uparrow}^\Lambda(k_1, k_2, k_1 + k_2 - k_3), \quad (5)$$

$$V_{\uparrow\downarrow\downarrow\uparrow}^\Lambda(k_1, k_2, k_3) = -V_{\uparrow\downarrow\uparrow\downarrow}^\Lambda(k_1, k_2, k_1 + k_2 - k_3). \quad (6)$$

This allows us to express the vertex by only one function of three frequency-momentum arguments: $V^\Lambda(k_1, k_2, k_3) \equiv V_{\uparrow\downarrow\uparrow\downarrow}^\Lambda(k_1, k_2, k_3)$, all the others spin components being obtained by Eqs. (5-6).⁸

The flow equation for the self energy can then be written as²

$$\frac{d}{d\Lambda} \Sigma^\Lambda(k) = - \int_p S^\Lambda(p) [2V^\Lambda(k, p, p) - V^\Lambda(k, p, k)], \quad (7)$$

with $p = (\mathbf{p}, \omega)$ and $k = (\mathbf{k}, \nu)$. We use the notation $\int_p = T \sum_\omega \int_{\mathbf{p}}$, where \sum_ω is the Matsubara frequency sum, and $\int_{\mathbf{p}} = \int \frac{d\mathbf{p}}{(2\pi)^2}$ is the normalized integration over the first Brillouin zone.

$$S^\Lambda = \left. \frac{dG^\Lambda}{d\Lambda} \right|_{\Sigma^\Lambda = \text{const}} \quad (8)$$

is the so-called single-scale propagator, and G^Λ is the full propagator, which is related to the bare propagator and the self-energy by the Dyson equation $(G^\Lambda)^{-1} = (G_0^\Lambda)^{-1} - \Sigma^\Lambda$.

The flow equation for the vertex can be written as^{2,8}

$$\frac{d}{d\Lambda} V^\Lambda(k_1, k_2, k_3) = \mathcal{T}_{\text{pp}}^\Lambda(k_1, k_2, k_3) + \mathcal{T}_{\text{ph}}^\Lambda(k_1, k_2, k_3) + \mathcal{T}_{\text{phc}}^\Lambda(k_1, k_2, k_3), \quad (9)$$

where³²

$$\begin{aligned} \mathcal{T}_{\text{pp}}^\Lambda(k_1, k_2, k_3) = & -\frac{1}{2} \int_p \mathcal{P}^\Lambda(p, k_1 + k_2 - p) \left\{ V^\Lambda(k_1, k_2, k_1 + k_2 - p) V^\Lambda(k_1 + k_2 - p, p, k_3) \right. \\ & \left. + V^\Lambda(k_1, k_2, p) V^\Lambda(p, k_1 + k_2 - p, k_3) \right\}; \end{aligned} \quad (10)$$

$$\begin{aligned} \mathcal{T}_{\text{ph}}^\Lambda(k_1, k_2, k_3) = & -\int_p \mathcal{P}^\Lambda(p, k_3 - k_1 + p) \left\{ 2V^\Lambda(k_1, k_3 - k_1 + p, k_3) V^\Lambda(p, k_2, k_3 - k_1 + p) \right. \\ & \left. - V^\Lambda(k_1, k_3 - k_1 + p, p) V^\Lambda(p, k_2, k_3 - k_1 + p) - V^\Lambda(k_1, k_3 - k_1 + p, k_3) V^\Lambda(k_2, p, k_3 - k_1 + p) \right\}; \end{aligned} \quad (11)$$

$$\mathcal{T}_{\text{phc}}^\Lambda(k_1, k_2, k_3) = \int_p \mathcal{P}^\Lambda(p, k_2 - k_3 + p) V^\Lambda(k_1, k_2 - k_3 + p, p) V^\Lambda(p, k_2, k_3). \quad (12)$$

The subscripts pp, ph and phc stand respectively for *particle-particle*, *particle-hole* and *particle-hole crossed*. Here we have defined the quantity

$$\mathcal{P}^\Lambda(p, p') = G^\Lambda(p) S^\Lambda(p') + S^\Lambda(p) G^\Lambda(p'), \quad (13)$$

which is the scale-derivative, at fixed self energy, of the product of two Green's functions.

III. VERTEX APPROXIMATION

In order to deal with the frequency and momentum dependence of the vertex, we start by decomposing it as follows:¹⁰

$$V^\Lambda(k_1, k_2, k_3) = U - \phi_p^\Lambda(k_1 + k_2; k_1, k_3) + \phi_m^\Lambda(k_3 - k_1; k_1, k_2) + \frac{1}{2} \phi_m^\Lambda(k_2 - k_3; k_1, k_2) - \frac{1}{2} \phi_c^\Lambda(k_2 - k_3; k_1, k_2), \quad (14)$$

where we introduced the pairing channel ϕ_p , the magnetic channel ϕ_m and the charge channel ϕ_c . Substituting Eq. (14) in Eq. (9) we obtain:

$$\begin{aligned} -\dot{\phi}_p^\Lambda(k_1 + k_2; k_1, k_3) + \dot{\phi}_m^\Lambda(k_3 - k_1; k_1, k_2) + \frac{1}{2} \dot{\phi}_m^\Lambda(k_2 - k_3; k_1, k_2) - \frac{1}{2} \dot{\phi}_c^\Lambda(k_2 - k_3; k_1, k_2) = \\ \mathcal{T}_{\text{pp}}^\Lambda(k_1, k_2, k_3) + \mathcal{T}_{\text{ph}}^\Lambda(k_1, k_2, k_3) + \mathcal{T}_{\text{phc}}^\Lambda(k_1, k_2, k_3). \end{aligned} \quad (15)$$

We associate the momentum transfer argument of \mathcal{P}^Λ in Eqns. (10-12) to the momentum transfer argument of the ϕ_x on the right hand side of Eq. 14. This way, it is easy to attribute $\mathcal{T}_{\text{pp}}^\Lambda$ to the flow equation of the only function in Eq. (15) that depends explicitly on $k_1 + k_2$: $-\dot{\phi}_p^\Lambda = \mathcal{T}_{\text{pp}}^\Lambda$. The same is true for the particle hole crossed channel: $\mathcal{T}_{\text{phc}}^\Lambda = \dot{\phi}_m^\Lambda$. We associate to the particle-hole diagram the third and fourth term on the left hand side of Eq. (15): $\mathcal{T}_{\text{ph}}^\Lambda(k_1, k_2, k_3) = \frac{1}{2} \dot{\phi}_m^\Lambda(k_2 - k_3; k_1, k_2) - \frac{1}{2} \dot{\phi}_c^\Lambda(k_2 -$

$k_3; k_1, k_2)$. Hence the flow equations for ϕ then read:

$$\dot{\phi}_p^\Lambda(Q; k_1, k_3) = -\mathcal{T}_{\text{pp}}^\Lambda(k_1, Q - k_1, k_3), \quad (16)$$

$$\begin{aligned} \dot{\phi}_c^\Lambda(Q; k_1, k_2) = & -2\mathcal{T}_{\text{ph}}^\Lambda(k_1, k_2, k_2 - Q) + \\ & \mathcal{T}_{\text{phc}}^\Lambda(k_1, k_2, Q + k_1), \end{aligned} \quad (17)$$

$$\dot{\phi}_m^\Lambda(Q; k_1, k_2) = \mathcal{T}_{\text{phc}}^\Lambda(k_1, k_2, Q + k_1), \quad (18)$$

where $Q = (\Omega, \mathbf{Q})$ is a frequency and momentum transfer.

Following Refs. 8,10, we address first the momentum dependence. To this end, we introduce a decomposition of unity by means of a set of orthonormal form factors for the two fermionic momenta $\{f_l(\mathbf{k})\}$ obeying the completeness relation²⁵:

$$\int_{\mathbf{k}} f_l(\mathbf{k}) f_m(\mathbf{k}) = \delta_{l,m}. \quad (19)$$

The procedure outlined here is described in detail, e.g., in Ref. 25.

We can then project each channel on a subset of form factors, whose choice is physically motivated⁸. If one could keep all the form factors the expansion would be exact.

For the pairing channel we keep only $f_s(\mathbf{k}) = 1$ and $f_d(\mathbf{k}) = \cos k_x - \cos k_y$:

$$\phi_p^\Lambda(Q; k_1, k_3) = \mathcal{S}_{\mathbf{Q}, \Omega}^\Lambda(\nu_1, \nu_3) + f_d\left(\frac{\mathbf{Q}}{2} - \mathbf{k}_1\right) f_d\left(\frac{\mathbf{Q}}{2} - \mathbf{k}_3\right) \mathcal{D}_{\mathbf{Q}, \Omega}^\Lambda(\nu_1, \nu_3). \quad (20)$$

The divergence in the channel \mathcal{S} is associated to the emergence of s -wave superconductivity, while \mathcal{D} to d -wave superconductivity.^{2,20}

For the charge and magnetic channels we restrict ourselves to $f_s(\mathbf{k}) = 1$ only:

$$\phi_c^\Lambda(Q; k_1, k_2) = \mathcal{C}_{\mathbf{Q}, \Omega}^\Lambda(\nu_1, \nu_2), \quad (21)$$

$$\phi_m^\Lambda(Q; k_1, k_2) = \mathcal{M}_{\mathbf{Q}, \Omega}^\Lambda(\nu_1, \nu_2), \quad (22)$$

corresponding to instabilities in the charge and magnetic channels, respectively.

For each channel in Eq. (14) we have defined *its own* frequency notation, consisting of one transfer frequency in the specific channel and two remaining independent fermionic frequencies. At finite temperature the frequency transfer is a bosonic Matsubara frequency.

The choice of the mixed notation is the most natural since the transferred momentum and frequency play a special role in the diagrammatics. Indeed, it is the

only dependence generated in second order perturbation theory and the main dependence in finite order perturbation theory. This notation is also convenient to express the Bethe-Salpeter equations, which are deeply related to parquet-approximations and fRG. Although one expects a leading dependence in the bosonic frequency, in particular in the weak coupling regime, we will see that in some cases the dependence on fermionic frequencies can become strong and not negligible. In Refs. 8, with a simplified frequency dependence, the channel functions above are interpreted as bosonic exchange propagators. Such an interpretation is missing with full frequency-dependence.

The flow equations for the channels \mathcal{S} , \mathcal{D} , \mathcal{C} and \mathcal{M} can be derived from the projection onto form factors of Eq. (10)-(12):

$$\dot{\mathcal{S}}_{\mathbf{Q}, \Omega}^\Lambda(\nu_1, \nu_3) = - \int_{\mathbf{k}_1, \mathbf{k}_3} \mathcal{T}_{\text{pp}}^\Lambda(k_1, Q - k_1, k_3); \quad (23)$$

$$\dot{\mathcal{D}}_{\mathbf{Q}, \Omega}^\Lambda(\nu_1, \nu_3) = - \int_{\mathbf{k}_1, \mathbf{k}_3} f_d\left(\frac{\mathbf{Q}}{2} - \mathbf{k}_1\right) f_d\left(\frac{\mathbf{Q}}{2} - \mathbf{k}_3\right) \mathcal{T}_{\text{pp}}^\Lambda(k_1, Q - k_1, k_3); \quad (24)$$

$$\dot{\mathcal{C}}_{\mathbf{Q}, \Omega}^\Lambda(\nu_1, \nu_2) = \int_{\mathbf{k}_1, \mathbf{k}_2} \mathcal{T}_{\text{phc}}^\Lambda(k_1, k_2, Q + k_1) - 2\mathcal{T}_{\text{ph}}(k_1, k_2, k_2 - Q); \quad (25)$$

$$\dot{\mathcal{M}}_{\mathbf{Q}, \Omega}^\Lambda(\nu_1, \nu_2) = \int_{\mathbf{k}_1, \mathbf{k}_2} \mathcal{T}_{\text{phc}}^\Lambda(k_1, k_2, Q + k_1). \quad (26)$$

The final equations are then obtained by substituting the decomposition (14) into the equations above, and using trigonometric equalities.

As an example we report here the equation for the magnetic channel, while the expression for the other channels are reported in the Appendix VIA:

$$\dot{\mathcal{M}}_{\mathbf{Q}, \Omega}^\Lambda(\nu_1, \nu_2) = \sum_{\nu} L_{\mathbf{m}, \mathbf{Q}, \Omega}^\Lambda(\nu_1, \nu) P_{\mathbf{Q}, \Omega}^\Lambda(\nu) L_{\mathbf{m}, \mathbf{Q}, \Omega}^\Lambda(\nu, \nu_2 - \Omega), \quad (27)$$

with:

$$P_{\mathbf{Q}, \Omega}^\Lambda(\omega) = \int_{\mathbf{p}} G^\Lambda(\mathbf{p}, \omega) S^\Lambda(\mathbf{Q} + \mathbf{p}, \Omega + \omega) + G^\Lambda(\mathbf{Q} + \mathbf{p}, \Omega + \omega) S^\Lambda(\mathbf{p}, \omega), \quad (28)$$

and:

$$\begin{aligned} L_{\mathbf{m}, \mathbf{Q}, \Omega}^\Lambda(\nu_1, \nu_2) &= U + \mathcal{M}_{\mathbf{Q}, \Omega}^\Lambda(\nu_1, \nu_2) + \int_{\mathbf{p}} \left\{ -\mathcal{S}_{\mathbf{p}, \nu_1 + \nu_2}^\Lambda(\nu_1, \nu_1 + \Omega) - \frac{1}{2} \mathcal{D}_{\mathbf{p}, \nu_1 + \nu_2}^\Lambda(\nu_1, \nu_1 + \Omega) [\cos(Q_x) + \cos(Q_y)] \right. \\ &\quad \left. + \frac{1}{2} [\mathcal{M}_{\mathbf{p}, \nu_2 - \nu_1 - \Omega}^\Lambda(\nu_1, \nu_2) - \mathcal{C}_{\mathbf{p}, \nu_2 - \nu_1 - \Omega}^\Lambda(\nu_1, \nu_2)] \right\} \end{aligned} \quad (29)$$

Note that after the momentum integrals in P and L are performed, the right hand side can be expressed as a matrix-matrix multiplication in frequency space, where Ω and \mathbf{Q} appear as parameters.

After this decomposition, the evaluation of the vertex-flow equation, depending on six arguments, is reduced to

the flow of the four functions \mathcal{S} , \mathcal{D} , \mathcal{C} , \mathcal{M} each of them depending on three frequencies and one momentum only.

In order to compute these equations numerically we discretize the momentum dependence on patches covering the Brillouin zone and truncate the frequency dependence to some maximal frequency value.

A. Interaction flow

To use the flow equations defined above we need to specify the Λ -dependence of the non interacting propagator G_0^Λ . We used the *interaction cutoff*, introduced in Ref. 26:

$$G_0^\Lambda(k) = \Lambda G_0(k) = \frac{\Lambda}{i\nu + \mu - \varepsilon_{\mathbf{k}}}, \quad (30)$$

Where the scale-parameter Λ flows from 0 to 1. μ is the chemical potential needed to fix the occupation at the desired value n . Correspondingly the interacting Green's function reads:

$$G^\Lambda(k) = \frac{\Lambda}{i\nu - \varepsilon_{\mathbf{k}} + \mu^\Lambda - \Lambda \Sigma^\Lambda(k)} \quad (31)$$

We have introduced a Λ -dependent chemical potential to maintain the occupation fixed during the flow. The chemical potential becomes a functional of the flowing self-energy, $\mu = \mu[\Sigma^\Lambda]$, whose value is found by solving the equation:

$$n = n^\Lambda(\mu) \equiv \int_{\mathbf{k}} \frac{e^{i\nu 0^+}}{i\nu - \varepsilon_{\mathbf{k}} + \mu^\Lambda - \Lambda \Sigma^\Lambda(k)}. \quad (32)$$

for μ .

The main advantage of the interaction cutoff is that the Λ -dependent action can be interpreted²⁶ as the physical action of the system with rescaled interaction $\tilde{U}^\Lambda = \Lambda^2 U$.

Since T acts as an infrared cutoff, for our purposes we do not need to worry about the fact that this cutoff is not scale-selective, and hence does not regularize possible divergences in the bubbles.

IV. RESULTS

In this section we present our results obtained by the full frequency dependent fRG. All the results are presented in units of the nearest-neighbor hopping $t = 1$. Unless specified otherwise, the next-nearest-neighbor hopping is $t' = -0.32t$, the interaction strength $U = 4t$, and the temperature $T = 0.08t$.

We have implemented numerically the flow equations reported in the appendix. To take into account the distinct momentum dependences of the self-energy and the vertex, we have defined two different patching schemes of the respective Brillouin zones. Similarly to what is done in Ref. 8, the vertex patching describes more accurately the corners around $(0,0)$ and (π, π) , where we expect the instability vectors. For the self-energy, the most relevant physics happens in the vicinity of the Fermi surface.

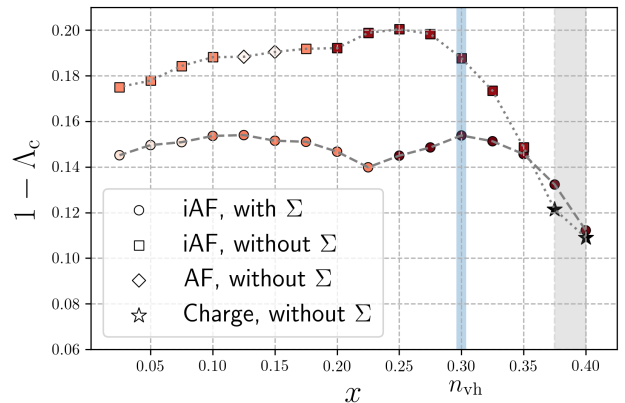


Figure 2: Critical scale $1 - \Lambda_c$ as a function of doping $x = 1 - n$, for $T = 0.08t$, $t' = -0.32t$ and $U = 4t$. Square symbols and circles refer to incommensurate antiferromagnetism (iAF) without and with self-energy feedback, respectively. The black stars refer to a divergence in the charge channel at $\mathbf{Q} = (0,0)$. The color of squares and circles encodes the distance of the incommensurate magnetic \mathbf{Q} -vector from (π, π) : darker color corresponds to a larger distance. The darkest color corresponds to $\delta = 1.13$. The vertical light blue line marks van Hove filling.

Therefore we concentrate the patches along the Fermi surface and in its immediate vicinity (see Figs. 9 and 10), with more points close to the *antinodal* region near $(\pi, 0)$, relevant for antiferromagnetism. In the calculations presented in the following we have used 29 patches for the vertex and 44 for the self-energy.

For the implementation of the frequency dependence we found it convenient to rewrite \mathcal{S} , \mathcal{D} , \mathcal{C} and \mathcal{M} as functions of three bosonic frequencies. For each frequency argument we restricted ourselves to at least 40 positive and 40 negative Matsubara frequencies. Beyond these frequencies we have used the asymptotic values.

A. Analysis of instabilities

By means of the fRG one can perform an instability analysis of the system: for some value of the flow parameter Λ one of the channels shows a divergence. The value Λ_c for which this happens is called *critical scale*, and from the diverging channel one can infer the leading instability of the system. In Fig. 2 we show the critical scale $1 - \Lambda_c$ as a function of the doping $x = 1 - n$ with and without self-energy feedback. For the physical interpretation of Λ_c in the interaction flow, we refer to the rescaled interaction²⁶ \tilde{U}^Λ discussed in Sec. ??.

We defined the critical scale as the flow parameter for which the value of the largest channel exceeds $200t$. We checked that these results are also consistent with a stability analysis based on the susceptibilities.

A divergence of the vertex at finite temperature is as-

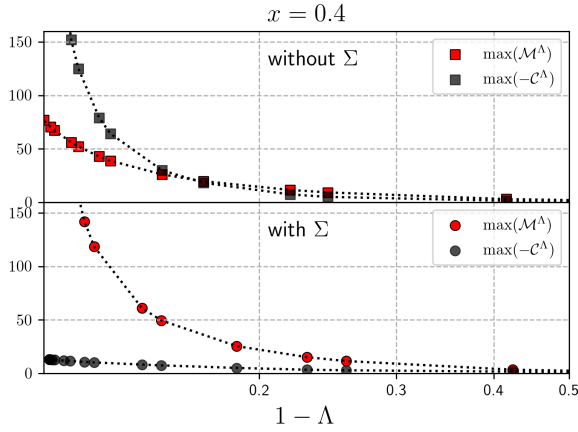


Figure 3: Flow of the maximal values of the charge (\mathcal{C}) and magnetic (\mathcal{M}) channels as functions of $1 - \Lambda$, for $x = 0.4$, $t' = -0.32$, $U = 4t$ and $T = 0.08t$. Top: without self-energy feedback; bottom: with self-energy feedback.

sociated with spontaneous symmetry breaking, in violation of the Mermin-Wagner theorem.²⁷ This is a consequence of the truncation of the flow equations. Instead, we should interpret the finite temperature vertex divergence as the signal of the appearance of strong bosonic fluctuations that cannot be treated within the approximation-scheme we are using.²¹ Even though in our framework the flow cannot be continued beyond the critical scale, from the analysis of vertex and self-energy we can identify the relevant effective interactions of the system.

For the parameter sets shown in Fig. 2, and without self-energy feedback, there are two possible instabilities. For doping smaller than 0.35 the leading fluctuations of the system are either commensurate or incommensurate antiferromagnetic. The incommensurability δ is defined through $\mathbf{Q} = (\pi, \pi - \delta)$, the momentum where the magnetic channel \mathcal{M}^Λ has its maximum. The region of commensurate antiferromagnetism for $0.125 \leq x \leq 0.150$ has to be attributed to the presence of a large plateau around (π, π) in the bare bubble. Correspondingly, the commensurate antiferromagnetic instability is almost degenerate with an incommensurate one.

The most striking feature in Fig. 2 is the presence of a divergence in the charge channel \mathcal{C}^Λ at $\mathbf{Q} = (0, 0)$ for the largest values of doping, marked by black stars. This feature was already observed in a fRG calculation with a simplified frequency parametrization by Husemann *et al.* in Ref. 8, and named *scattering instability*. The charge channel \mathcal{C}^Λ diverges for finite frequency transfer $\Omega = 2\pi T$, which does not allow for a natural interpretation in terms of a physical instability. The frequency structure of the charge channel \mathcal{C}^Λ together with its origin will be discussed further in paragraph ??.

The self-energy feedback has three effects. First, it decreases $1 - \Lambda_c$. Second, the incommensurability vector is affected, the region of commensurate antiferromagnetism

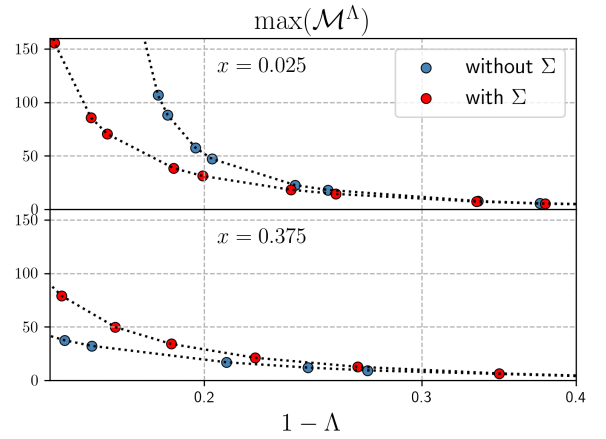


Figure 4: Flow of the maximal values of the magnetic (\mathcal{M}) channel as functions of $1 - \Lambda$, for $x = 0.025$ (top) and $x = 0.375$ (bottom). The other parameters are $t' = -0.32$, $U = 4t$ and $T = 0.08t$. Red symbols: with self-energy feedback; blue symbols: without self-energy feedback.

disappears, and one can observe a more regular trend of increasing δ with x . Third, the divergence in the charge channel is completely suppressed, and the leading instability in the doping region $0.375 \leq x \leq 0.4$ remains incommensurate antiferromagnetic. This can be also seen from Fig. 3, where we compare the flow of the maximum (of the absolute value) of magnetic and charge channels with and without the self-energy feedback for doping $x = 0.4$. Without self-energy feedback, the charge channel reaches large and negative values. The presence of such a large (and negative) charge channel inhibits the magnetic channel. The effect of the self-energy in the flow is evident: the charge channel is strongly damped. At the same time the magnetic channel is enhanced.

This is confirmed by Fig. 4, where we show the maximum of \mathcal{M} with and without self-energy feedback for $x = 0.025$ (top) and $x = 0.375$ (bottom). One can see that the enhancement of \mathcal{M} due to the self-energy is specific of the large doping region, while, in the small doping region the self energy decreases \mathcal{M} . The self-energy affects the magnetic channel directly by reducing the particle-hole bubble, and indirectly through the feedback of other channels, that is, reducing the charge channel. The former effect dominates for small doping, the latter at large doping.

To better understand these effects we looked for possible changes in the Fermi surface structure, analyzing the momentum distribution

$$n^\Lambda(\mathbf{k}) = 2T \sum_{\nu} \frac{e^{i\nu 0^+}}{i\nu - \varepsilon_{\mathbf{k}} + \mu^\Lambda - \Lambda \Sigma^\Lambda(\mathbf{k}, \nu)}. \quad (33)$$

The factor 2 accounts for the spin degree of freedom. In Fig. 5 we show the non-interacting (top left) and interacting (top right) occupation in the first quadrant of the Brillouin zone for doping $x = 0.025$. The latter is computed at the critical scale Λ_c .

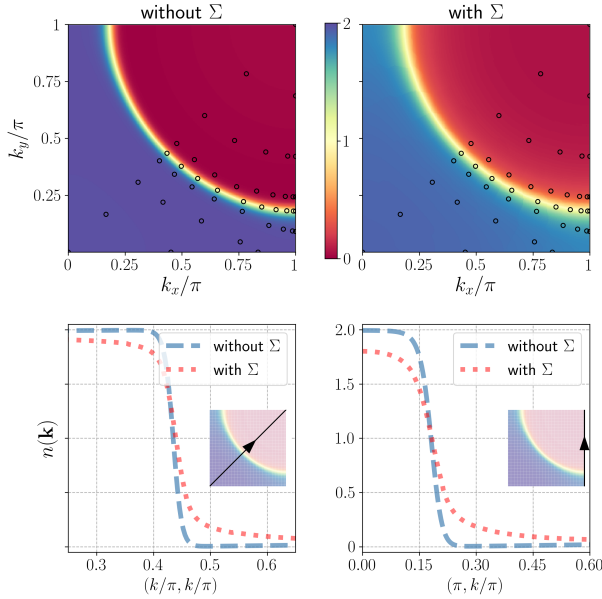


Figure 5: Top row: momentum distribution for $t' = -0.32t$, $T = 0.08t$ and doping $x = 0.025$. Left panel: non-interacting case. Right panel: interacting case for $U = 4t$. The black circles mark the points used to patch the self-energy. Bottom row: cut of the occupation along the Brillouin zone paths reported as arrows in the insets. Blue dashed curves are results for the non-interacting system, while red dotted curves are for $U = 4t$.

Comparing the two panels, one does not observe any relevant shift of the Fermi surface position, but the Fermi surface broadening is appreciably larger in the interacting case, due to the self-energy. Similar results apply for doping $x = 0.4$, as one can see from Fig. 6, where the broadening is more evident.

B. Frequency dependence of vertex

We now discuss the remarkable frequency dependence of the vertex. In particular, we will look at the channels that show a divergence, that is, the charge and the magnetic instabilities observed in Fig. 2, while we refer to the Appendix for the pairing channel.

As mentioned in the previous section the divergences of the charge and magnetic channels are quite different. The charge channel diverges for a finite frequency transfer, and only when we neglect the self-energy feedback. Since the dependence on the transfer momentum and frequency (\mathbf{Q}, Ω) has already been discussed in Ref 10, we focus on the dependence on the fermionic frequencies. Therefore we present various color plots for fixed (\mathbf{Q}, Ω) , showing the dependence on ν_1 and ν_2 .

In the top left panel of Fig. 7 we show the magnetic channel $\mathcal{M}_{\mathbf{Q},\Omega}^{\Lambda_c}(\nu_1, \nu_2)$ in the small doping region, where

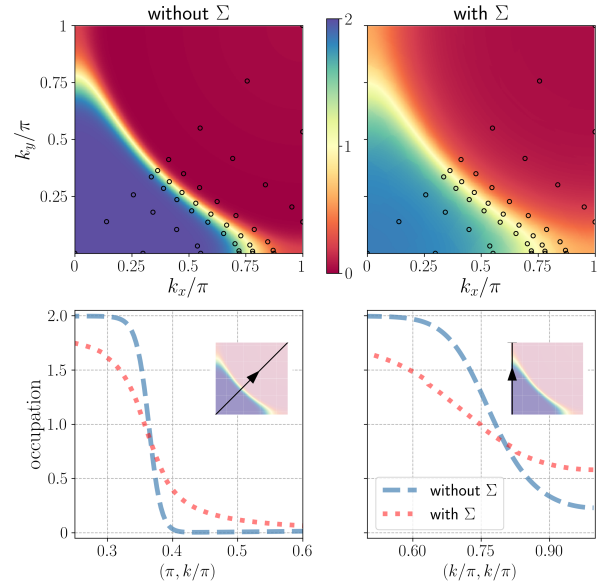


Figure 6: Top row: momentum distribution for $t' = -0.32t$, $T = 0.08t$ and doping $x = 0.4$. Left panel: non-interacting case. Right panel: interacting case for $U = 4t$. The black circles mark the points used to patch the self-energy. Bottom row: cut of the occupation along the Brillouin zone paths reported as arrows in the insets. Blue dashed curves are results for the non-interacting system, while red dotted curves are for $U = 4t$.

antiferromagnetism is the leading instability. The results shown have been calculated with self-energy feedback, but the frequency structures we discuss do not depend strongly on the presence of the self-energy. For clarity we restrict the plots to the first 20 positive and negative Matsubara frequencies, larger frequencies can be deduced by the asymptotic behavior.²⁸ When only one channel in Eq. (9) is taken into account, the fRG equations are equivalent to the RPA. The magnetic channel calculated with RPA would depend only on the frequency and momentum transfer, not on ν_1 and ν_2 . Hence any variation in the frequency structure has to be ascribed to the presence of the other channels in the fRG. The channel competition suppresses the magnetic channel: the largest value of \mathcal{M} is reduced compared to the RPA, and the frequency dependent structure at the center is further reduced compared to the asymptotic values.

In the bottom left panel of Fig. 7 we show the frequency dependence of the charge channel $\mathcal{C}_{\mathbf{Q},\Omega}^{\Lambda_c}(\nu_1, \nu_2)$ for a finite frequency transfer $\Omega = 2\pi/\beta$, related to the charge instability discussed in Ref. 8 and above. The frequency structure is completely different from the magnetic channel. The charge channel assumes negative values, and the maximum is for frequencies $\nu_1 = \pi T$ and $\nu_2 = -\pi T$. This structure cannot be explained in terms of standard ladder diagrams. It might be related to the behavior of

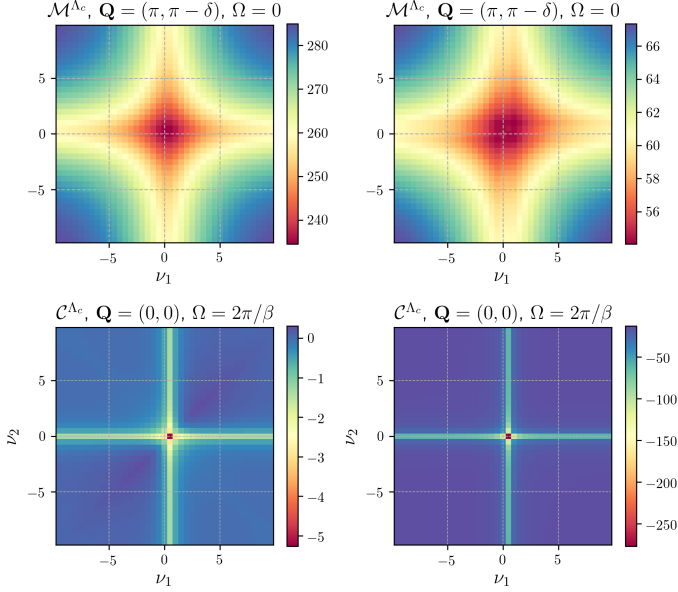


Figure 7: Frequency dependence of the magnetic (top) and charge (bottom) channel for $t' = -0.32$, $U = 4t$ and $T = 0.08t$. *Top left*: Magnetic channel $\mathcal{M}_{\mathbf{Q},\Omega}^{\Lambda_c}(\nu_1, \nu_2)$ with self-energy feedback at the instability vector and for vanishing frequency transfer, for doping $x = 0.025$. *Top right*: Magnetic channel $\mathcal{M}_{\mathbf{Q},\Omega}^{\Lambda_c}(\nu_1, \nu_2)$ without self-energy feedback at $\mathbf{Q} = (\pi, \pi - \delta)$, $\delta = 1.13$ (corresponding to the largest magnetic coupling) and for vanishing frequency transfer, for doping $x = 0.4$. *Bottom left*: Frequency dependence of the charge channel $\mathcal{C}_{\mathbf{Q},\Omega}^{\Lambda_c}(\nu_1, \nu_2)$ with self-energy feedback at $\mathbf{Q} = (0, 0)$ and frequency transfer $\Omega = 2\pi T$, for doping $x = 0.025$. *Bottom right*: Frequency dependence of the charge channel $\mathcal{C}_{\mathbf{Q},\Omega}^{\Lambda_c}(\nu_1, \nu_2)$ without self-energy feedback at $\mathbf{Q} = (0, 0)$ and frequency transfer $\Omega = 2\pi T$, for doping $x = 0.4$.

the retarded interaction described in Ref. 29.

In the two right panels of Fig. 7 we show the same quantities but for $x = 0.4$ and without self-energy feedback. In this case, the localized peak in the charge channel is the leading interaction. The position and shape of the frequency structures are similar to the one described above.

C. Origin of charge singularity

To gain insight in the origin of the singular frequency structures observed in the charge channel, we identify a simple set of diagrams reproducing the same features. The main idea is that the magnetic channel, which is generated first, is responsible for the singular structure in the charge channel.

To check this qualitatively, we first compute an effective interaction by means of an RPA in the magnetic

channel, and then insert this effective magnetic interaction into a subsequent RPA equation for the charge channel. Of course one does not expect quantitative agreement with the fRG, since we overestimate both interactions, but the approximation is sufficient to reproduce and explain the qualitative features we are interested in.

We start by introducing an effective interaction that includes the magnetic fluctuations as computed by RPA¹⁷ in the particle-hole crossed channel:

$$U_{\mathbf{Q},\Omega}^{\text{eff}} = \frac{U}{1 - U\Pi_{\mathbf{Q},\Omega}}. \quad (34)$$

Since the bare interaction U is local, U^{eff} depends only on the transfer momentum \mathbf{Q} and frequency Ω of the particle-hole bubble

$$\Pi_{\mathbf{Q},\Omega} = -T \sum_{\nu} \int_{\mathbf{p}} G_0(\mathbf{p}, \nu) G_0(\mathbf{p} + \mathbf{Q}, \nu + \Omega). \quad (35)$$

The magnetic effective interaction in Eq. (34) will be now used to compute the RPA equation for the charge channel. Adopting the simplified momentum dependences of the effective interactions used in the fRG calculation, only the momentum integrated, that is, local part of the magnetic interaction $U_{\Omega}^{\text{eff}} = \int_{\mathbf{Q}} U_{\mathbf{Q},\Omega}^{\text{eff}}$ contributes to the charge channel. We thus obtain

$$\mathcal{C}_{\mathbf{Q},\Omega}(\nu_1, \nu_3) = U_{\text{eff},\nu_1-\nu_3} [\delta_{\nu_1,\nu_3} + U_{\nu_1-\nu_3}^{\text{eff}} \Pi_{\mathbf{Q},\Omega}(\nu_1)]^{-1}, \quad (36)$$

where

$$\Pi_{\mathbf{Q},\Omega}(\nu) = - \int_{\mathbf{p}} G_0(\mathbf{p}, \nu) G_0(\mathbf{p} + \mathbf{Q}, \nu + \Omega). \quad (37)$$

Here the charge channel is expressed in terms of ν_1 and $\nu_3 = \nu_2 - \Omega$ rather than in terms of ν_1 and ν_2 . Note that the fermion frequencies ν are not summed in $\Pi_{\mathbf{Q},\Omega}(\nu)$. Eq. (36) is nothing more than an RPA equation with a frequency dependent interaction in the particle-hole channel.¹⁷ U^{eff} depends on $\nu_1 - \nu_3$ due to the frequency exchange from particle-hole crossed to particle-hole notation. We note that, in the case of a frequency independent effective interaction U_{eff} , Eq. (36) becomes ν_1 and ν_3 independent and only the summed bubble $\Pi_{\mathbf{Q},\Omega}$ appears. The frequency dependence of U_{eff} qualitatively affects the results.

In Fig. 8, we show the charge channel as computed from Eq. (36) for $\mathbf{Q} = (0, 0)$ and different Ω as a function of ν_1 and ν_2 , for $T = t$ and $x = 0.375$. We have to choose such a high temperature to stay in a stable paramagnetic phase, due to the above-mentioned overestimation of the fluctuations within the RPA. In the more accurate fRG calculation the magnetic instability occurs at lower temperatures. The frequency structure in Fig. 8 for $\Omega = 2\pi T$ is very similar to the one shown in Fig. 7. The simple diagrams considered here reproduce the position of the main structures, as well as the correct sign of the charge channel. This is true also for the other

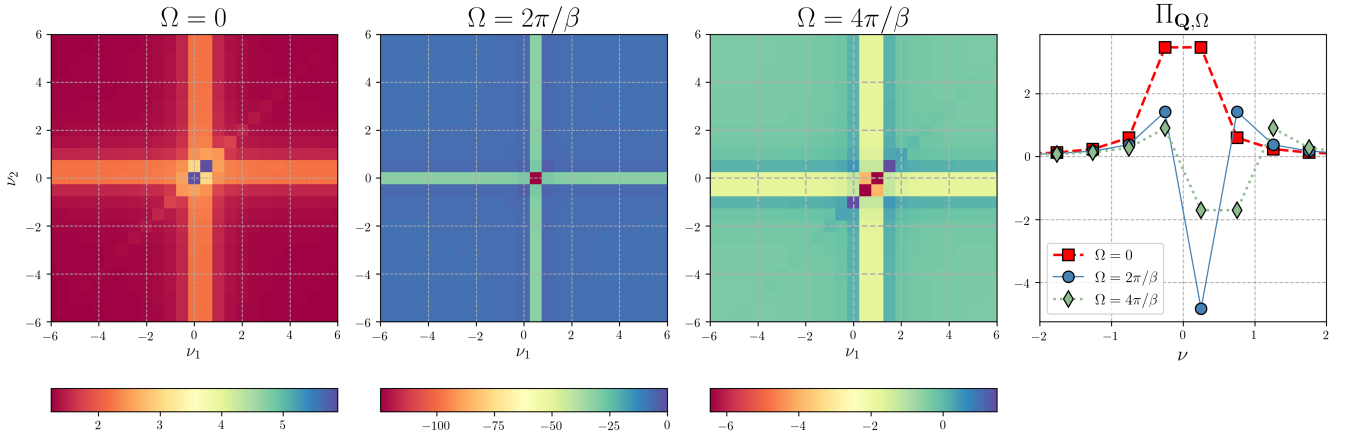


Figure 8: In the first three panels from the left, the charge channel $C_{\mathbf{Q}=(0,0),\Omega}(\nu_1, \nu_2 - \Omega)$ computed from Eq. (36) is shown as a function of ν_1 and ν_2 for transfer frequencies $\Omega = 0$, $\Omega = 2\pi T$ and $\Omega = 4\pi T$, respectively. In the right panel, the bubble $\Pi_{\mathbf{Q}=(0,0),\Omega}(\nu)$ is shown as a function of ν for $\Omega = 0$, $\Omega = 2\pi T$ and $\Omega = 4\pi T$. The model parameters are $t' = -0.32$ and $U = 4$, the doping $x = 0.375$, and the temperature $T = t$.

bosonic Matsubara frequency shown here, for which we do not report the fRG results. Furthermore, upon lowering the temperature the charge channel diverges also for other finite bosonic Matsubara frequencies, while it does not diverge for $\Omega = 0$. From this we conclude that the diagrams described here are responsible for the frequency structure of the charge channel observed in the fRG.

To understand why the divergence appears for a finite frequency Ω , we notice that in Eq.(36) the Ω dependence appears only through the bubble $\Pi_{\mathbf{Q},\Omega}(\nu)$. The frequency summed particle-hole bubble obeys the following relation:

$$\Pi_{\mathbf{Q} \rightarrow (0,0),\Omega} = \frac{1}{\beta} \sum_{\nu} \Pi_{\mathbf{Q} \rightarrow (0,0),\Omega}(\nu) = C \delta_{\Omega,0}, \quad (38)$$

where C is a constant that, at low temperature, approaches the density of states at the Fermi level. In the rightmost panel of Fig. 8, we show the bubble $\Pi_{\mathbf{Q}=(0,0),\Omega}(\nu)$ as a function of ν for different values of Ω . We note that it has a large negative peak for $\Omega = 2\pi T$. This is due to the property (38): the summed bubble must vanish for $\Omega \neq 0$, hence a large negative value is needed to cancel the positive contributions at large frequency. We have thus identified the origin of the frequency structure observed in the charge channel, which seems to be quite general and arises from simple diagrams.

Including the self-energy in the calculation of the bubble, Eq. (38) does not evaluate to a δ -function anymore, and the difference between the summed bubble at vanishing frequency and for frequency $2\pi T$ is diminished. This is probably the reason why the inclusion of the self-energy feedback prevents the unphysical divergence of the charge channel.

D. Self energy

We now discuss the frequency and momentum dependence of the self energy. In Fig. 9 we show the frequency dependence of the imaginary part of the self-energy at $T = 0.08t$ and low doping $x = 0.025$. The spread between the maximal and minimal self-energy at each frequency is rather small, indicating that the self-energy did not develop a large momentum dependence even when the flow parameter reached the critical scale. For all momenta, $|\text{Im}\Sigma(\mathbf{k}, \nu)|$ decreases as a function of decreasing frequency, as in a Fermi liquid. One would generally expect the antinodal region to be more affected by correlation effects. However, there is only a slight increase of $|\text{Im}\Sigma(\mathbf{k}, \nu)|$ in this region. At the temperature we are considering, we do not observe a tendency towards the opening of a momentum selective gap.

In Fig. 10 we show the imaginary part of the self-energy for a larger doping $x = 0.4$. As in the previous case, we do not see much momentum differentiation.

The self-energy enters directly in the calculation of the momentum distribution through the Green's function, already discussed above, and shown in Figs. 5 and 6. In the bottom panels of these figures, we show how the momentum distribution evolves along two different cuts in the Brillouin zone, crossing the *nodal* and *antinodal* regions, respectively. The drop in the momentum distribution is sharper along the diagonal, and the self-energy effects are stronger along the antinodal cut. For doping $x = 0.4$ the broadening of the Fermi surface, already larger at the non interacting level, is further enhanced by the self-energy.

To study further the difference between nodal and antinodal regions in the iAF case, we studied the quasi-particle weight^{2,30} $Z_{\mathbf{k}}$, and the decay rate $\gamma_{\mathbf{k}}$. Instead of relying on analytical continuation, we have extracted the parameters directly from the imaginary axis data. To do

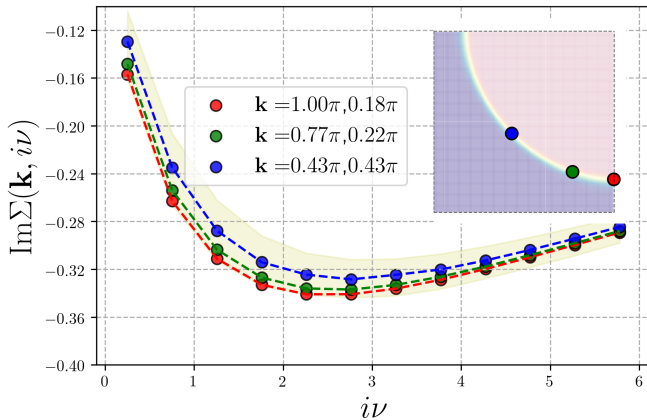


Figure 9: Self-energy as a function of frequency for $U = 4t$, $t' = -0.32t$ and low doping $x = 0.025$ at a temperature $T = 0.08t$. The location of the \mathbf{k} -point in the Brillouin zone is color coded in the inset. The position of all the patching points taken into account for the self-energy is shown as black circles in the top row of Figs. 5 and 6, and does not change during the flow. The shaded area highlights the region between the maximal and minimal value of the self-energy for each frequency.

so we have fitted the first few frequencies of the imaginary part of the self-energy with a polynomial of degree l : $\text{Im}\Sigma(\mathbf{k}, i\nu) \approx a_0(\mathbf{k}) + a_1(\mathbf{k})\nu + \dots + a_l(\mathbf{k})\nu^l$ and we identified $\gamma_{\mathbf{k}} = a_0(\mathbf{k})$ and $Z_{\mathbf{k}} = \frac{1}{1-a_1(\mathbf{k})}$. The procedure only works if the temperature is small enough, and if the frequencies used for the fitting are not too high. We checked that the results were stable by changing the number of frequencies and the order of the polynomial used for the fit. In Fig. 11 we plot $Z_{\mathbf{k}}$ and $\gamma_{\mathbf{k}}$ against the angle θ along the Fermi surface, $\theta = 0$ corresponding to the antinodal point and $\theta = \pi/4$ to the nodal one. The variation of the quasiparticle weight along the Fermi surface is extremely small with Z assuming values between 0.754 and 0.760. On the other hand, the relative variation of the decay rate γ along the Fermi surface is sizable, varying from $\gamma \approx 0.056t$ and $\gamma \approx 0.082t$. These values are comparable with the temperature $T = 0.08t$. We conclude that at the critical scale, the system still has coherent quasiparticles along the Fermi surface, with a higher decay rate in the antinodal region. This is consistent with the observations of Ref.³¹, where non-Fermi liquid behavior of the self-energy was observed only very close to the critical temperature and in the immediate vicinity of the hot spots. *Honerkamp*

V. CONCLUSIONS

- order of the paragraph? shall we discuss first self energy or vertex?

• do we want to show more vertex structure in the

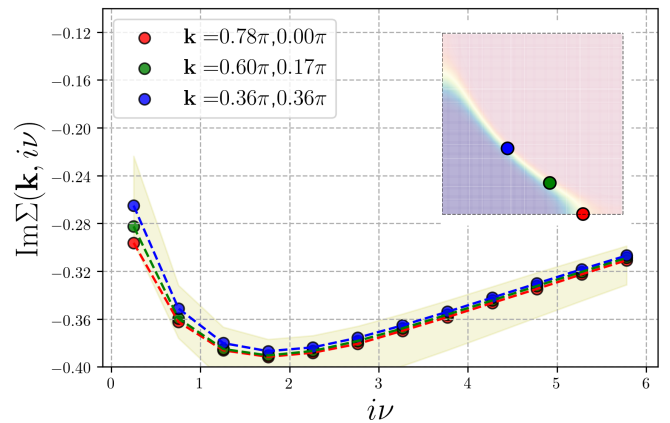


Figure 10: Self-energy as a function of frequency for $U = 4t$, $t' = -0.32t$ and high doping $x = 0.4$ at a temperature $T = 0.08$. The location of the \mathbf{k} -point in the Brillouin zone is color coded in the inset. The position of all the patching points taken into account for the self-energy is shown as black circles in the top row of Figs. 5 and 6, and does not change during the flow. The shaded area is the region between the maximal and minimal value of the self-energy for each frequency.

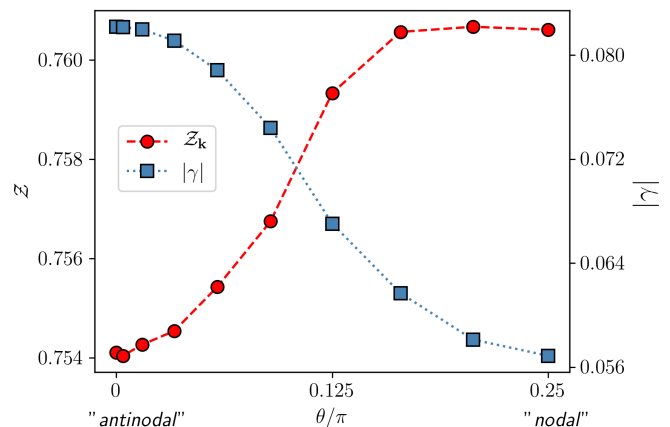


Figure 11: Quasiparticle weight $Z_{\mathbf{k}}$ and decay rate $\gamma_{\mathbf{k}}$ as function of the angle θ for the same parameters as in Fig. 9. The values on the left axis refer to the quasiparticle weight, the values on the right axis refer to the decay rate.

appendix?

Acknowledgments

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VI. APPENDICES

A. Flow equations

We here report the final expression for the flow equations of each channel. To simplify the notation from now on we omit the Λ dependencies. The flow equation for the s -wave superconductivity channel \mathcal{S} reads:

$$\dot{\mathcal{S}}_{\mathbf{Q}}^{\Omega;\nu_1,\nu_3} = \frac{1}{2} \sum_{\nu} L_{s\mathbf{Q}}^{\Omega;\nu_1,\nu} P_{s,\mathbf{Q}}^{\Omega,\nu} L_{s\mathbf{Q}}^{\Omega;\nu,\Omega-\nu_3} + \frac{1}{2} \sum_{\nu} L_{s\mathbf{Q}}^{\Omega;\Omega-\nu_1,\nu} P_{s,\mathbf{Q}}^{\Omega,\nu} L_{s\mathbf{Q}}^{\Omega;\nu,\nu_3}, \quad (39)$$

with:

$$P_{s,\mathbf{Q}}^{\Omega;\omega} = \int_{\mathbf{p}} G_{\mathbf{p},\omega} S_{\mathbf{Q}-\mathbf{p},\Omega-\omega} + G_{\mathbf{Q}-\mathbf{p},\Omega-\omega} S_{\mathbf{p},\omega}, \quad (40)$$

and:

$$L_{s\mathbf{Q}}^{\Omega;\nu_1,\nu_3} = U - \mathcal{S}_{\mathbf{Q}}^{\Omega;\nu_1,\nu_3} + \int_{\mathbf{p}} \left[\mathcal{M}_{\mathbf{p}}^{\nu_3-\nu_1;\nu_1,\Omega-\nu_1} + \frac{1}{2} \mathcal{M}_{\mathbf{p}}^{\Omega-\nu_1-\nu_3;\nu_1,\Omega-\nu_1} - \frac{1}{2} \mathcal{C}_{\mathbf{p}}^{\Omega-\nu_1-\nu_3;\nu_1,\Omega-\nu_1} \right]. \quad (41)$$

The flow equation for the d -wave superconductivity channel \mathcal{D} reads:

$$\dot{\mathcal{D}}_{\mathbf{Q}}^{\Omega;\nu_1,\nu_3} = \frac{1}{2} \sum_{\nu} L_{d\mathbf{Q}}^{\Omega;\nu_1,\nu} P_{d,\mathbf{Q}}^{\Omega,\nu} L_{d\mathbf{Q}}^{\Omega;\nu,\Omega-\nu_3} + \frac{1}{2} \sum_{\nu} L_{d\mathbf{Q}}^{\Omega;\Omega-\nu_1,\nu} P_{d,\mathbf{Q}}^{\Omega,\nu} L_{d\mathbf{Q}}^{\Omega;\nu,\nu_3}, \quad (42)$$

with:

$$P_{d,\mathbf{Q}}^{\Omega;\omega} = \int_{\mathbf{p}} f_d \left(\frac{\mathbf{Q}}{2} - \mathbf{p} \right)^2 [G_{\mathbf{p},\omega} S_{\mathbf{Q}-\mathbf{p},\Omega-\omega} + G_{\mathbf{Q}-\mathbf{p},\Omega-\omega} S_{\mathbf{p},\omega}], \quad (43)$$

and:

$$L_{d\mathbf{Q}}^{\Omega;\nu_1,\nu_3} = -\mathcal{D}_{\mathbf{Q}}^{\Omega;\nu_1,\nu_3} + \frac{1}{2} \int_{\mathbf{p}} (\cos p_x + \cos p_y) \left[\mathcal{M}_{\mathbf{p}}^{\nu_3-\nu_1;\nu_1,\Omega-\nu_1} + \frac{1}{2} \mathcal{M}_{\mathbf{p}}^{\Omega-\nu_1-\nu_3;\nu_1,\Omega-\nu_1} - \frac{1}{2} \mathcal{C}_{\mathbf{p}}^{\Omega-\nu_1-\nu_3;\nu_1,\Omega-\nu_1} \right]. \quad (44)$$

Since \mathcal{D} is generated during the flow by the other channels only, see Eq. (44), it is the most sensitive channel to the frequency approximation made. Neglecting the vertex frequency one will likely overestimate L_d , as already mentioned in Ref. 10.

The flow equation for the charge channel \mathcal{C} reads:

$$\dot{\mathcal{C}}_{\mathbf{Q}}^{\Omega;\nu_1,\nu_2} = \sum_{\nu} L_{c\mathbf{Q}}^{\Omega;\nu_1,\nu} P_{\mathbf{Q}}^{\Omega,\nu} L_{c\mathbf{Q}}^{\Omega;\nu,\nu_2-\Omega}, \quad (45)$$

with:

$$L_{c\mathbf{Q}}^{\Omega;\nu_1,\nu_2} = U - \mathcal{C}_{\mathbf{Q}}^{\Omega;\nu_1,\nu_2} + \int_{\mathbf{p}} \left[-2\mathcal{S}_{\mathbf{p}}^{\nu_1+\nu_2;\nu_1,\nu_2-\Omega} + \mathcal{S}_{\mathbf{p}}^{\nu_1+\nu_2;\nu_1,\Omega+\nu_1} \right. \quad (46)$$

$$\left. + \left(\mathcal{D}_{\mathbf{p}}^{\nu_1+\nu_2;\nu_1,\nu_2-\Omega} - \frac{1}{2} \mathcal{D}_{\mathbf{p}}^{\nu_1+\nu_2;\nu_1,\Omega+\nu_1} \right) [\cos(Q_x) + \cos(Q_y)] \right. \quad (47)$$

$$\left. + \frac{3}{2} \mathcal{M}_{\mathbf{p}}^{\nu_2-\nu_1-\Omega;\nu_1,\nu_2} + \frac{1}{2} \mathcal{C}_{\mathbf{p}}^{\nu_2-\nu_1-\Omega;\nu_1,\nu_2} \right], \quad (48)$$

while $P_{\mathbf{Q}}^{\Omega;\omega}$ is given in Eq. (28). The equation for the magnetic channel is reported in Eq. (27). The form factor decomposition allows to decouple the momentum integrals, in the calculation of the L 's, Eqns. (29), (41), (44) and (46), from the frequency summations in the flow equations, hence reducing the numeric effort.

B. Pairing channel

In Figs. 12 and 13 we display the frequency dependence of the pairing functions \mathcal{S} and \mathcal{D} . As a consequence of Eq. (44) the frequency structure of \mathcal{D} is decaying in all directions.¹² The small numerical values are due to three main

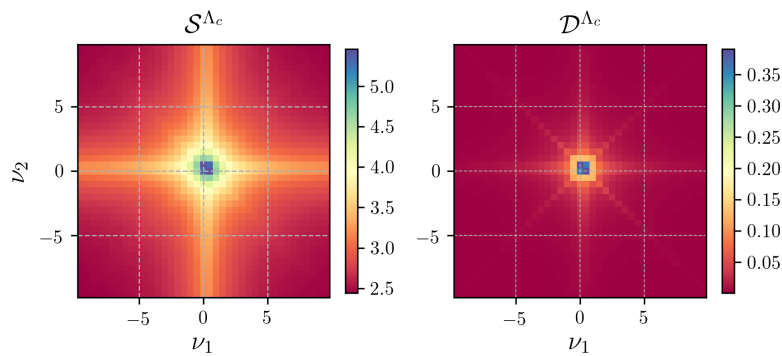


Figure 12: Frequency dependence of \mathcal{S} (left panel) and \mathcal{D} (right panel). Here $T = 0.08t$, $t' = -0.32t$, $U = 4t$ and $x = 0.025$. The transfer momentum and frequency are $\mathbf{Q} = (0, 0)$ and $\Omega = 0$.

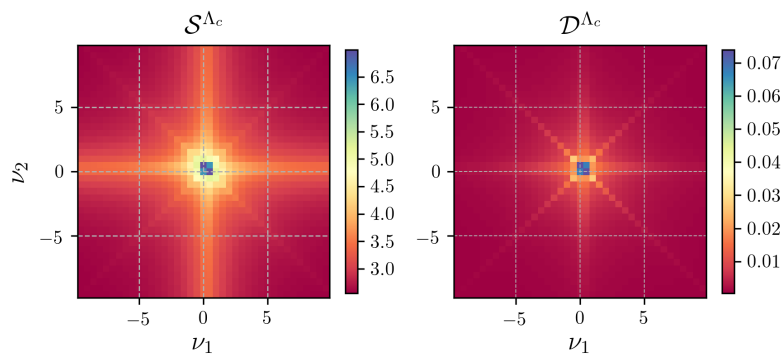


Figure 13: Frequency dependence of \mathcal{S} (left panel) and \mathcal{D} (right panel). Here $T = 0.08t$, $t' = -0.32t$, $U = 4t$ and $x = 0.4$. The transfer momentum and frequency are $\mathbf{Q} = (0, 0)$ and $\Omega = 0$.

reasons: first, the d -wave pairing is expected to increase suddenly for temperatures immediately above its critical temperature. Second, as argued in Ref. 10, previous fRG calculations with a static vertex are likely to overestimate the d -wave channel by neglecting the frequency dependence in Eq. (44). Finally, the interaction scheme itself has a tendency to suppress the d -wave pairing, since in the terms of Ref. 12, its diagrammatic contributions can be classified as *rest function*.

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