Charge and current orders in the spin-fermion model with overlapping hot spots.

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Experiments carried over the last years on the underdoped cuprates have revealed a variety of symmetry-breaking phenomena in the pseudogap state. Charge-density waves, breaking of $C_4$ rotational symmetry as well as time-reversal symmetry breaking have all been observed in several cuprate families. In this regard, theoretical models where multiple non-superconducting orders emerge are of particular interest. We consider the recently introduced (Phys. Rev. B 93, 085131 (2016)) spin-fermion model with overlapping 'hot spots' on the Fermi surface. Focusing on the particle-hole instabilities we obtain a rich phase diagram with the chemical potential relative to the dispersion at $(0, \pi)$; $(\pi, 0)$ and the Fermi surface curvature in the antinodal regions being the control parameters. We find evidence for d-wave Pomeranchuk instability, d-form factor charge density waves as well as commensurate and incommensurate staggered bond current phases similar to the d-density wave state. The current orders are found to be promoted by the curvature. Considering the appropriate parameter range for the hole-doped cuprates, we discuss the relation of our results to the pseudogap state and incommensurate magnetic phases of the cuprates.

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I. INTRODUCTION

Origin of the pseudogap state remains one of the main puzzles in the physics of the high-$T_c$ cuprate superconductors. First observed in NMR measurements, it is characterized by the loss of the density of states due to the opening of a partial gap at the Fermi level below the pseudogap temperature $T^* > T_c$. Studies of the pseudogap by means of ARPES and Raman scattering have revealed that it opens around $(0, \pi)$ and $(\pi, 0)$ points of the 2D Brillouin zone, the so-called antinodal regions. With increasing hole doping both $T^*$ and the gap magnitude decrease monotonously and eventually disappear. However, modern experiments add many more unconventional details to this picture, showing that a crucial role in the pseudogap state is played by the various ordering tendencies.

To begin with, the point-group symmetry of the CuO$_2$ planes appears to be broken. Namely, scanning tunneling microscopy (STM) and transport studies show the absence of $C_4$ rotational symmetry in the pseudogap state. More recently, magnetic torque measurements of the bulk magnetic susceptibility confirmed $C_4$ breaking occurring at $T^*$. Additionally, an inversion symmetry breaking associated with pseudogap has been discovered by means of second harmonic optical anisotropy measurement.

Other experiments suggest that an unconventional time-reversal symmetry breaking is inherent to the pseudogap. Polarized neutron diffraction studies of different cuprate families reveal a magnetic signal commensurate with the lattice appearing below $T^*$ and interpreted as being due to a $Q = 0$ intra-unit cell magnetic order. The signal has been observed to develop above $T^*$ with a finite correlation length and breaks the $C_4$ symmetry (note, however, that a recent report does not bear evidence of such a signal). Additionally, at a temperature $T_K$ that is below $T^*$ but shares a similar doping dependence polar Kerr effect has been observed, which implies that time-reversal symmetry is broken. Additional signatures of a temporally fluctuating magnetism below $T^*$ are also available from the recent studies.

While the signatures described above indicate that the pseudogap state is a distinct phase with a lower symmetry, there exist only few experiments that yield a thermodynamic evidence for a corresponding phase transition. On the other hand, transport measurements suggest the existence of quantum critical points (QCPs) of the pseudogap phase, accompanied by strong mass enhancement in line with the existence of a QCP.

Additionally, in the recent years the presence of charge density waves (CDW) has been discovered in a similar doping range. The CDW onset temperature can be rather close to $T^*$ but has been shown to have a distinct dome-shaped doping dependence in YBCO and Hg-1201. Diverse probes such as resonant X-ray and hard X-ray scattering, STM and NMR have observed CDW with similar properties in most of the hole-doped cuprate compounds with the exception of La-based ones (in which the spin and charge modulations are intertwined). Generally, the CDWs have the following common properties. The modulation wavevectors are confined along the Brillouin zone axes (axial CDW) and increase with doping. While the modulations along both directions are usually observed, there is an experimental evidence that the CDW is unidirectional locally. The intra-unit cell structure of the CDW is characterized by a d-form factor with the charge being modulated at two oxygen sites of the unit cell in antiphase with each...
other.

From the theoretical perspective, one of the initial interpretations was that the pseudogap was a manifestation of fluctuating superconductivity, either in a form of preformed Bose pairs\cite{46,47} or strong phase fluctuations\cite{48}. However, the onset temperatures of superconducting fluctuations observed in the experiments\cite{49-51} are considerably below $T^*$ and have a distinct doping dependence. Another scenario dating back to the seminal paper\cite{52} attributes the pseudogap to the short-range correlations due to strong on-site repulsions\cite{53,54}. Numerical quantum Monte Carlo simulations\cite{55-58} of the Hubbard model support this idea. However, this scenario ‘as is’ does not explain the broken symmetries of the pseudogap state. More recently, these results have been interpreted as being due to topological order\cite{59,60}, that can also coexist with the breaking of discrete symmetries\cite{61}.

A different class of proposals for explaining the pseudogap behavior involves a competing symmetry-breaking order. One of the possible candidates discussed in the literature is the $Q = 0$ orbital loop current order\cite{62}. Presence of circulating currents explicitly breaks the time reversal symmetry, allowing one, with appropriate modifications, to describe the phenomena observed in polarized neutron scattering\cite{63} and polar Kerr effect\cite{64} experiments. However, it does not lead to a gap on the Fermi surface at the mean-field level. Numerical studies of the three-band Hubbard model give arguments both for\cite{65,66} and against\cite{67-69} this type of order. Other proposals for $Q = 0$ magnetic order include spin-nematic\cite{15,70}, oxygen orbital moment\cite{71} or magnetoelastic multipole\cite{72,73} order.

Charge nematic order\cite{74} and the related d-wave Pomern anchuk instability\cite{75,76} of the Fermi surface have also been considered in the context of the pseudogap state. It breaks the $C_4$ rotational symmetry of the $CuO_2$ planes in agreement with numerous experiments\cite{9-13}. While not opening a gap, fluctuating distortion of the Fermi surface can result in an arc-like momentum distribution of the spectral weight and non-Fermi liquid behaviour\cite{77-79}. Evidence for this order comes from numerical studies of the Hubbard model with functional renormalization group\cite{75}, dynamical mean-field theory\cite{80,81} and other\cite{82,83} methods as well as analytical studies of forward-scattering\cite{84} and spin-fermion\cite{85,86} models.

Another possibility is the CDW\cite{87-89}. More recent studies focus on the important role of the interplay between CDW and superconducting fluctuations\cite{90}, preemptive orders and time reversal symmetry breaking\cite{91} (that can result in the polar Kerr effect\cite{92,93}), vertex corrections for the interactions\cite{94,99}, CDW phase fluctuations\cite{96} and possible SU(2) symmetry\cite{97-99}. Additionally, pair density wave - a state with modulated Cooper pairing amplitude has been proposed to explain the pseudogap and CDW\cite{100,102}, which can be also understood with the concept of 'intertwined' SC and CDW orders\cite{103}.

An interesting alternative is the d-density wave\cite{104} (DDW) state (also known as flux phase\cite{105,106}) which is characterized by a pattern of bond currents modulated with the wavevector $Q = (\pi, \pi)$ that is not generally accompanied by a charge modulation. This order leads to a reconstructed Fermi surface consistent with the transport\cite{27,107} and ARPES\cite{108} signatures of the pseudogap. Moreover, the time-reversal symmetry is also broken and a modified version of DDW can explain the polar Kerr effect\cite{109} observation. Additionally, model calculations show\cite{110,111} that the system in the DDW state can be unstable to the formation of axial CDWs. Studies aimed at a direct detection of magnetic moments created by the DDW have yielded results both supporting\cite{112,113} and against\cite{114} their existence (or with the conclusion that the signal is due to impurity phases\cite{115}). However, theoretical estimates of the resulting moments are model-dependent\cite{116,117}. There also exists indirect evidence from superfluid density measurements\cite{118}. Theoretical support for the DDW comes from renormalization group\cite{119} and variational Monte Carlo\cite{120} studies of the Hubbard model, DMRG studies of $t - J$ ladders\cite{121}, and mean-field studies of $t - J$\cite{122} as well as single\cite{123} and three-orbital\cite{124} models. However, the regions of DDW stability found in these studies vary significantly and depend on the value of particular interactions\cite{123,124} or details of the Fermi surface\cite{120}.

Overall, the question of possible competing orders in the cuprates has turned out to be a rather complicated one. Interestingly, state-of-art numerical calculations comparing different methods show that the energy difference between distinct ground states can be miniscule\cite{125} explaining some of the difficulties. Thus, analytical approaches which allow one to study the influence of different parameters in detail can be of interest.

In this paper, we deduce leading non-superconducting orders using a low-energy effective theory for fermions interacting with antiferromagnetic (AF) fluctuations. While such theories can be in principle derived from the microscopic Hubbard or $t$-$J$ Hamiltonians\cite{126}, we employ here a semiphenomenological approach in the spirit of the widely used spin-fermion (SF) model\cite{90,91,97,127,128}. Our take on this problem differs in that we relax the usual assumption that the interaction, being peaked at $(\pi, \pi)$, singles out eight isolated ‘hot spots’ on the Fermi surface. In contrast, we consider that neighboring hot spots may strongly overlap and form antinodal ‘hot regions’. This assumption agrees well with the ARPES results\cite{3} demonstrating the pseudogap covers the full antinodal region without pronounced maxima at the ‘hot spots’ of the standard SF model. Moreover, the electron spectrum in the antinodal regions has been found\cite{108,129} to be shallow with respect to the pseudogap energy scale for the hole-underdoped samples, i.e. the pseudogap opens also at points that are not in immediate vicinity of the Fermi surface. From the spin fluctuation perspective this can be anticipated if the AF fluctuations correlation length is small enough such that the resulting interaction between fermions is uniformly smeared covering the full antinodal regions. Indeed, the neutron scatter-
ing experiments\textsuperscript{130,131} show that the correlation lengths at the temperatures and dopings relevant for the pseudogap amount to several unit cell lengths.

SF model with overlapping hot spots has been introduced in our recent publications\textsuperscript{85,86}, where we have considered normal state properties as well as charge orders corresponding to intra-region particle-hole pairing. For the case of a small Fermi surface curvature, it has been shown that the d-wave Pomeranchuk instability is the leading one for sufficiently shallow electron spectrum in the hot regions. This is in contrast to the diagonal d-form factor CDW usually being the leading particle-hole instability in the standard SF model\textsuperscript{90,97,126}. As a result of Pomeranchuk transition, the $C_4$ symmetry gets broken by a deformation of the Fermi surface and an intra-unit-cell charge redistribution. Additionally, as the Pomeranchuk order leaves the Fermi surface ungapped, we have shown that at lower temperatures an axial CDW with dominant d-form factor and d-wave superconductivity may appear. These results are in line with the simultaneous observation of the commensurate $C_4$ breaking\textsuperscript{9-13} and axial d-form factor CDWs\textsuperscript{39,45}. At the same time, these order parameters, although being in agreement with the experimental observations, do not readily explain the time-reversal symmetry breaking phenomena as well as the possible Fermi surface reconstruction into hole pockets\textsuperscript{37}.

In this paper, we consider a possibility of an inter-region particle-hole pairing, akin to the excitonic insulator proposed long ago\textsuperscript{132}. The resulting state is similar in properties to the d-density wave\textsuperscript{104}, having staggered bond currents. In addition, we find also evidence for an incommensurate version thereof. It turns out (Sec. III,IV) that the Fermi surface curvature in the antinodal regions (assumed to be small in Refs.\textsuperscript{85 and 86}) is the most important ingredient that stabilizes this state against the charge orders, thus leading to a rich phase diagram. We further discuss the relation of our findings to the pseudogap state in Sec.V.

The paper is organized as follows. In Sec. II we present the model and assumptions that we use. In Sec. III we analyze a simplified version of the model ignoring retardation effects and identify the emerging orders. In Sec. IV we present the results for the full model and discuss the approximations used. In Sec. V we discuss the relation of our results to the physics of underdoped cuprates and in Sec.VI we summarize our findings.

II. MODEL.

The spin-fermion model describes the low-energy physics of the cuprates in terms of low-energy fermions interacting via the antiferromagnetic paramagnons. The latter are assumed to be remnants of the parent insulating AF state destroyed by hole doping\textsuperscript{127}. The resulting interaction is strongly peaked at the wavevector $Q_0 = (\pi, \pi)$ corresponding to the antiferromagnetic order periodicity and is described by a propagator

$$\frac{1}{(q - Q_0)^2 + \frac{1}{\xi_{AF}^2} - \left(\frac{\omega}{v_\nu}\right)^2}.$$ Then, one can identify eight 'hot spots' on the Fermi Surface mutually connected by $Q_0$ where the interaction is expected to be strongest (see the left part of Fig.1). A conventional approximation motivated by the proximity to AF quantum critical point (QCP) where $\xi_{AF} \to \infty$ is to consider only small $\delta p \sim 1/\xi_{AF}$ vicinities of the 'hot spots' to be strongly affected by the interaction. However, at temperatures relevant for the pseudogap state this argument does not have to hold — the experimentally reported correlation lengths\textsuperscript{130,131} are indeed rather small. Moreover, ARPES experiments\textsuperscript{8} show that the effects of the pseudogap extend well beyond the 'hot spots' to the Brillouin zone edges $(\pi, 0), (0, \pi)$ without being significantly weakened.

![Fig. 1. Overlapping hot spots on the Fermi surface typical for the hole-doped cuprates.](image)

A different approach has been introduced in\textsuperscript{85,86}. As $\xi_{AF}$ becomes smaller, the 'hot spots' expand and and can eventually overlap and merge forming two 'hot regions' (see the right part of Fig.1). For the latter to occur the fermionic dispersion in the antinodal region should be shallow (see also the formal definition below), which is supported by the experimental data\textsuperscript{108,129}. To describe this situation we consider the following Lagrangian:

$$L_{\text{SF}} = \sum_{p,\nu=1,2} c^\dagger_p [\partial_\tau + \varepsilon_\nu(p)] c^\nu_p + \frac{1}{2} \sum_q \varepsilon_q (-v_s^2 \partial_\tau^2 + \mathbf{q}^2 + 1/\xi^2) \varphi_q + \lambda \sum_{p,q} \left[ c^\dagger_{p+\mathbf{q}} \varphi_q c^\nu_p + c^\dagger_{p+\mathbf{q}} \varphi_q c^\nu_p \right],$$

where $c^\nu_p$ and $\varphi_q$ are the fermionic and bosonic (paramagnon) fields, respectively and $\varepsilon_\nu(p)$ is the fermionic dispersion where the index $\nu$ enumerates the two 'hot regions'. Additionally, as the quantity $|\varepsilon((0, \pi)/(\pi, 0) - \mu|$ has been observed\textsuperscript{108,129} to be of the order of the pseudogap energy or smaller, one expects the fermions in the whole
region to participate in the interaction. Consequently one has to consider the dispersion relation not linearized near the Fermi surface. Due to the saddle points present at $(\pi,0)$ and $(0,\pi)$ the minimal model for the dispersion is:

\[
\varepsilon_1(p) = \alpha p_x^2 - \beta p_y^2 - \mu, \quad \varepsilon_2(p) = \alpha p_y^2 - \beta p_x^2 - \mu, \quad (2.2)
\]

where $\alpha$ has the meaning of the inverse fermion mass and $\beta/\alpha$ controls the Fermi surface curvature. Note that the chemical potential of the system is determined by the full Fermi surface. As the Fermi energies (measured from the $\Gamma$-point) in hole-doped cuprates are quite large we neglect the temperature dependence of the chemical potential, and, consequently, $\mu$. On the other hand, as the relevant temperatures for the pseudogap onset are still sizable, we shall not consider the effects of simultaneous development of multiple instability channels due to the van Hove singularities.\(^\text{133}\)

In Refs. 85 and 86 the limit $\beta \to 0$ has been considered. For this case the condition of ‘shallowness’ leading to the merging of the hot spots reads $1/\xi_{AF} \gg \sqrt{\mu/\alpha}$. However, in order to keep the simple quadratic form of the paramagnon dispersion we also assume $\xi_{AF} \gg a_0$, where $a_0$ is the lattice spacing. Here we will consider the consequences of finite $\beta$ for the model 2.1 under the same assumption of the strong hot spot overlap. As in Refs. 85 and 86 we will concentrate on the particle-hole (non-superconducting) orders. Note that for interaction being via the antiferromagnetic paramagnons only, d-wave superconductivity is expected to overcome the particle-hole orders.\(^\text{86,128}\) However, additional interactions present in real systems, such as nearest-neighbor\(^\text{134}\) or remnant low-energy Coulomb repulsion\(^\text{85}\) should suppress it with respect to the particle-hole orders.

### III. PHASE DIAGRAM FOR A SIMPLIFIED MODEL.

To address qualitative features of the emerging orders we can use a simplified version of the model (2.1) also introduced previously by us.\(^\text{85}\) It amounts to substitution of the paramagnon part of the Lagrangian with a constant interregion interaction between the fermions. This is also equivalent to taking the $\xi \to 0$ limit for the paramagnon propagator. Additionally, we neglect the self-energy effects for this case.

We start with the following Lagrangian

\[
\sum_{p,\nu=1,2,\sigma} c_p^{\nu\dagger} \partial_\tau \varepsilon_\nu(p) c_p^{\nu}\nonumber
\]

\[
-\lambda_0 \sum_{p,p',q} (c_{p+q}^{1\dagger} \bar{\sigma} c_{p}^{2} + c_{p+q}^{2\dagger} \bar{\sigma} c_{p}^{1}) \nonumber
\]

\[
x(c_{p+q}^{1\dagger} \bar{\sigma} c_{p}^{2} + c_{p}^{1\dagger} \bar{\sigma} c_{p+q}^{2}) \nonumber
\]

\[
\times (c_{p+q}^{1\dagger} \bar{\sigma} c_{p}^{2} + c_{p}^{2\dagger} \bar{\sigma} c_{p+q}^{1}). \quad (3.1)
\]

The spin structure of the interaction is taken here in full analogy to the original spin-fermion model. Additionally, the integrals that appear below are cut off at momenta $p_x, p_y \sim 1/\xi$ and we assume that the inequality $\alpha/\xi^2 \gg \mu, T$ (strong overlap of hot spots) is fulfilled.

In addition to the d-wave superconductivity, one can identify two attractive singlet channels (the triplet channels, as in the SF model, are subleading with the effective coupling being three times weaker). Corresponding order parameters can be written in terms of the averages (spin indices are suppressed):

\[
W = \langle (c_{p+Q/2}^{1\dagger} c_{p}^{2} - c_{p}^{1\dagger} c_{p+Q/2}^{2}) \rangle = -\langle (c_{p+Q/2}^{2\dagger} c_{p}^{1} - c_{p}^{2\dagger} c_{p+Q/2}^{1}) \rangle, \quad (3.2)
\]

\[
D = \langle (c_{p+Q/2}^{1\dagger} c_{p}^{2} - c_{p}^{1\dagger} c_{p+Q/2}^{2}) \rangle = -\langle (c_{p+Q/2}^{2\dagger} c_{p}^{1} - c_{p}^{2\dagger} c_{p+Q/2}^{1}) \rangle. \quad (3.3)
\]

The interaction for the orders without the sign change between the regions (corresponding to s-form factor charge order) is repulsive and consequently such orders are not expected to appear on the mean-field level. As is shown in detail below, the order parameter $W$, Eq.(3.2), corresponds to a charge order with a d-form factor. Due to the change of the sign between the regions the charge is modulated only at the oxygen orbitals of the unit cell corresponding to the bonds in a single-band model. In the case $Q \neq 0$ (Fig. 2a) this order represents the d-form factor charge density wave while for $Q = 0$ (Fig. 2b) it leads to an intra-unit cell redistribution of charge accompanied by a d-wave Pomeranchuk deformation of the Fermi surface.

![FIG. 2. Illustration of charge modulations for a) d-form factor CDW with Q along the diagonal b) d-wave Pomeranchuk phase](image-url)
we present the resulting phase diagram where as well as a detailed description or DDW 3.2 ax-

\[ \chi_{DDW} \] contrast to the previous studies of the SF model \[ \beta/\alpha \] at moderate charge and current phases, respectively.  

To stimulate the Pomeranchuk and IDDW phases among the currents between nearest neighbors are depicted.  

\[ \mu/T = \chi_{DDW} + \chi_{IDDW} \]

where \( \mu/T \) is the stability temperature here. To map out the phase diagram we fix the leading instabilities here: to the d-wave Pomeranchuk or DDW phases at a lower temperature (see text).  

\[ \beta = \alpha \text{ one has } \varepsilon_1 = -\varepsilon_2 \text{ for } \alpha p^2 \gg \mu. \]  

As \( \alpha/\xi^2 \gg \mu \), the Fermi surfaces of the two regions are nearly nested in the large part of the regions, in contrast to a CDW, for which nesting is restricted to a vicinity of a single point in \( \mathbf{p} \)-space. It is, however, surprising that the current phases start dominating at \( \beta/\alpha \) considerably smaller than 1. Below we provide analytical results leading to the phase diagram of Fig. 4 as well as a detailed description of the emerging orders.  

- **Charge Density Wave** is represented by Eq. (3.2) with finite value of \( Q \). Due to the sign change between the two FS regions the amplitude of the on-site charge modulation proportional to \( \sum_{\mathbf{p}} W(\mathbf{p}) \sim W_1 + W_2 \) vanishes.  

It has been shown\(^{85,90} \), however, that the charge modulation on oxygen sites is related to bond operators in the single-band model as:

\[ \delta n_{CO}(\mathbf{r}) \sim e^{iQ\mathbf{r}}(W_1 - W_2) + c.c. \]

where \( i \) and \( i + 1 \) are two neighboring copper sites. The proportionality coefficient depends on additional assumptions: in Ref. 85 \( p/8 \) has been obtained using Zhang-Rice singlet doping picture. Transforming the expression to the momentum space one obtains

\[ \delta n_{CO}(\mathbf{r}) \sim e^{iQ\mathbf{r}}(W_1 - W_2) + c.c. \]

Let us compute now the corresponding susceptibility, Eq. (3.6). First of all, one can conclude from Eq. (3.4)
that only \( Q_x = Q_y \equiv Q \) satisfies \( \chi_{\text{LDW}} = \chi_{\text{CDW}} \), i.e. the wavevector is directed along the diagonal. This is in line with the previous results on the spin-fermion model\(^{90,128}\). The momentum integrals for the present model with overlapping hotspots can be evaluated explicitly for two limiting cases (for the details of calculation see Appendix A). In the limit \( \beta \to 0 \) one obtains:

\[
\chi_{\text{CDW}}(T, \mathbf{Q}) = \frac{T}{\alpha} \sum_{\omega_n} \frac{i \pi \text{sgn}[\omega_n] \sqrt{\alpha/\xi^2}}{(i \omega_n + \mu - \alpha Q^2/4) \sqrt{i \omega_n + \mu}}.
\]

In the opposite limit \( \beta/\xi^2 \gg \mu, T \) one gets

\[
\chi_{\text{CDW}}(\mathbf{Q}) \approx \frac{T}{\sqrt{\alpha \beta}} \sum_{|\omega_n| < \beta/\xi^2} -i \pi \text{sgn}[\omega_n] \arctanh \frac{\gamma - i \omega_n - \mu}{\sqrt{\gamma^2 - i \omega_n - \mu}} \frac{\sqrt{\gamma}}{\sqrt{\gamma - i \omega_n - \mu}}.
\]

where \( \gamma = (\alpha - \beta) Q^2/4 \). The expression in the sum in Eq. (3.9) is obtained for \( \omega_n \ll \beta/\xi^2 \). However, as the resulting sum is logarithmically large for \( \beta/\xi^2 \gg \mu, T \), we can simply disregard the contribution from higher Matsubara frequencies.

- **Pomeranchuk instability** corresponds to the anomalous average 3.2 with \( \mathbf{Q} = 0 \). It leads to a d-wave-like deformation of the Fermi surface breaking \( C_4 \) symmetry without opening a gap. Additionally, the Pomeranchuk order should be accompanied by an intra-unit cell redistribution of the charge on the two oxygen orbitals, which can be readily seen from (3.7).

The expressions for the susceptibilities can be obtained from Eqs. (3.8) and (3.9) by taking the \( \mathbf{Q} \to 0 \) limit. Moreover, the sign of \( \frac{\partial \chi_{\text{CDW}}}{\partial Q^2} \bigg|_{\mathbf{Q}=0} \) allows one to check the stability of the Pomeranchuk phase with respect to the CDW. For \( \beta \to 0 \) we get

\[
\chi_{\text{Pom}} = \frac{T}{\alpha} \sum_{\omega_n} \frac{i \pi \text{sgn}[\omega_n] \sqrt{\alpha/\xi^2}}{(i \omega_n + \mu)^{3/2}}.
\]

and

\[
\frac{\partial \chi_{\text{CDW}}}{\partial Q^2} \bigg|_{\mathbf{Q}=0} = \frac{T}{4} \sum_{\omega_n} \frac{i \pi \text{sgn}[\omega_n] \sqrt{\alpha/\xi^2}}{(i \omega_n + \mu)^{5/2}}.
\]

Numerical calculation shows that the expression (3.11) changes sign from positive to negative for \( \mu/T < 1.1 \). Therefore, the Pomeranchuk phase is stable for \( \mu/T_{\text{ins}} < 1.1 \) for \( \beta \to 0 \) (in agreement with Ref. 85). In the opposite limit \( \beta/\xi^2 \to \infty \), however, the expansion of (3.9) yields

\[
\frac{\partial \chi_{\text{CDW}}}{\partial Q^2} \approx \frac{\alpha - \beta}{4} \frac{2T}{\alpha^2} \sum_{|\omega_n| < \beta/\xi^2} \frac{i \pi \text{sgn}[\omega_n]}{(i \omega_n + \mu)}
\]

\[
= \frac{\alpha - \beta}{4} \frac{2T}{\alpha^2} \sum_{0 < \omega_n < \beta/\xi^2} \frac{4 \pi \omega_n \mu}{(\omega_n^2 + \mu^2)}
\]

which is always positive. Thus, to obtain the phase boundary between CDW and Pomeranchuk phases at finite \( \beta \) one needs to perform the momentum integration assuming finite values of \( \beta/\xi^2 \). The general result is rather cumbersome and is presented in Appendix A (Eq. A1). For \( \mu/T \ll 1 \) a simple expression is found

\[
\left( \frac{\mu}{T} \right)_{\text{CDW}} \approx \frac{\alpha}{\beta} \frac{T}{\alpha - \beta/\xi^2} \approx 1.17 \frac{\alpha}{\beta} \frac{T}{\alpha - \beta/\xi^2}.
\]

One can see that \( (\mu/T)_{\text{CDW}} \) decreases for \( \beta < 0.5 \) and then starts to increase. However, as we shall see, this upturn is located in the region where DDW is the leading instability. Note that in Fig. 4 the Pomeranchuk/CDW boundary is found from the full expression (A1) numerically for \( (\alpha/\xi^2)/T = 50 \).

- **Current Phases (DDW/IDDW)** are represented by the anomalous average (3.3). This order parameter does not result in any charge modulations on both the copper and the oxygen orbitals. In the former case this is guaranteed by the d-wave symmetry, while for the oxygen orbitals (e.g., \( O_x \)):

\[
\delta n_{O_x}(r) \sim \text{Re} \left[ e^{i(Q+(\pi,\pi))r} \sum_p \cos(p_x(c^\dagger_p + c^\dagger_{-p}) + [p_x-(\pi,\pi)-Q]/2) \right]
\]

\[
= 2\text{Re} \left[ e^{i(Q+(\pi,\pi))r} \sum_p \sin(p_x)D \right] = 0.
\]

However, it can be shown that the order parameter \( D \) induces a staggered pattern of currents flowing through the lattice. The current between the lattice sites \( i \) and \( j \) is given by\(^{123}\):

\[
I_{ij} = -\frac{t_{ij}}{\hbar} (\hat{c}_{i}^\dagger \hat{c}_{j} - \hat{c}_{j}^\dagger \hat{c}_{i}),
\]

where \( t_{ij} \) is the hopping parameter. For example, the current \( I_{i,i+x} \) between nearest neighbor sites along \( x \) can be estimated as

\[
I_{i,i+x} \approx 4\text{Re} \left[ \frac{it}{\hbar} e^{i(Q+(\pi,\pi))R_i - iQ_x/2} D \right].
\]

In Fig. 3 an illustration of the current patterns is presented. Note that, in general, currents between non-nearest neighbors are induced, too. However, as this effect depends on the structure of the DDW amplitude in the entire Brillouin zone, we will not consider it in this work.

To calculate the resulting magnetic fields one should however calculate the current density \( \mathbf{j}_q \) rather then the current. For a square lattice with nearest-neighbor hopping \( t \) one can obtain the following result neglecting the smearing of atomic wavefunctions with respect to the cur-
the result is presented for

\[ j_{x,y}(q) = -8\pi iD_{\alpha}(\pi,0)e_{x,y}e_{x,y} \]

\[ \times (q e_{x,y})^{-1} \sum_{K_n} \delta (q - Q_{AF} + K_n), \]  

(3.16)

where \( e_{x,y} \) is a unit vector along \( x(y) \) axis and \( K_n \) is a reciprocal lattice vector. Additionally, one can calculate the magnetic field along the \( z \)-direction produced by the DDW \( B_z \) assuming that DDWs are aligned in-phase along \( z \) axis

\[ B_z(q) = \frac{4\pi i}{c} \frac{q_y j_y - q_y j_x}{q^2}. \]  

(3.17)

Note that in the full model (Sec. IV) the order parameter is frequency-dependent due to the retardation effects and \( D \) has to be calculated using a corresponding anomalous Green’s function. Thus, \( D \) is not, in general, simply related to the magnitude of the pseudogap or \( T^* \) as would be the case for the constant interaction.

Let us now present the results for the thermodynamic susceptibilities. For the commensurate (\( Q = 0 \)) state, one obtains in the limit \( \beta \to 0 \)

\[ \chi_D(T) = \frac{\beta^2}{2\alpha} \tanh \mu/2T. \]  

(3.18)

In the opposite limit \( \beta/\xi^2 \to \infty \), we have instead

\[ \chi_D(T) \approx \frac{4\arctanh(\sqrt{\beta/\alpha})}{\alpha + \beta} \sum_{|\omega_n| < \beta/\xi^2} \frac{i\pi \text{sgn}[\omega_n]}{\omega_n + \mu}. \]  

(3.19)

For the case \( Q \neq 0 \), let us first consider the stability of the DDW with respect to an infinitesimal commensuration vector \( Q \). The general expression obtained from

\[ \frac{\partial^2 \chi_{DDW}}{\partial Q_1^2} \bigg|_{Q=0} = \frac{\partial^2 \chi_D}{\partial Q_1^2} \bigg|_{Q=0} = 0 \]

is presented in Appendix A (Eq. A4). Numerical solution of the resulting equation is represented by the DDW/IDDW critical line in the phase diagram of Fig. 4. Qualitatively, low values of \( \mu \) favor IDDW. For \( \mu/T \ll 1 \), \( \alpha - \beta \ll \alpha \) the result can also be expressed analytically

\[ \left( \frac{\mu}{T} \right)^{DDW}_{cr} \approx 0.7(1 - \beta/\alpha)^2. \]  

(3.20)

Furthermore, we have studied the dependence of the direction and magnitude of \( Q \) maximizing \( \chi_{DDW}(Q) \) numerically. For this purpose, expressions (A5,A6,A7) have been used. In Fig. 5 the result is presented for \( \mu/T = 0.1 \). Interestingly, the orientation of \( Q \) in the IDDW phase is almost always along the axes. While there seems to be a transition to a diagonal phase at low curvatures, the charge order is dominant in that region, as is shown below.

**Competition between charge and current orders**

We are now in position to compare the tendencies to form charge (CDW/Pomeranchuk) and current (DDW/IDDW) order. For \( \beta \ll \alpha \), comparing (3.8) and (3.10) to (3.18) one finds an additional factor \( \sqrt{\alpha/\xi^2} \) in the former. After the summation this translates into a large parameter of the order of \( \sim \sqrt{\alpha/\xi^2}/[T,\mu] \) present in the susceptibilities for charge instabilities. Incommensurability of the DDW does not change this conclusion (see Eq. A3).

In the case \( \beta \sim \alpha \), \( \beta/\xi^2 \gg \mu, T \) one gets large logarithmic contributions in Matsubara sums for both charge and current orders. Therefore, one can estimate the transition line by equating the prefactors of the sums (3.9) (3.19), which leads to the following equation

\[ \frac{1}{\sqrt{\alpha/\beta}} = \frac{4\arctanh(\sqrt{\beta/\alpha})}{\alpha + \beta}. \]

Numerical solution of this equation yields \( \langle \beta/\alpha \rangle_{cr} \approx 0.29 \), this value being independent of \( \mu/T \). The finite slope of the charge/current boundary in Fig. 4 results from finite values of \( \langle \alpha/\xi^2 \rangle/T \) taken in the numerical calculations. One can however show, that the slope is strongly suppressed being proportional to \( \log^{-1}(\beta/\xi^2)/T \).

**IV. CHARGE AND CURRENT ORDERS IN THE FULL MODEL**

We turn now to the analysis of the full SF model (2.1). Let us first consider the effects of interactions in the normal state. The interactions renormalize the Green’s function \( G \) of the fermions and \( D \) of the paramagnons leading to:
$$G_{\alpha\beta}(\varepsilon_n, p) = \frac{\delta_{\alpha\beta}}{if(\varepsilon_n, p) - \varepsilon_\nu(p)},$$  \hspace{1cm} (4.1)

$$\mathcal{D}_{mn'}(\omega_n, q) = -\frac{\delta_{mm'}}{\Omega(\omega_n, q)/v_s^2 + q^2 + 1/\xi^2},$$

where $\alpha, \beta$ are fermion spin indices, $\nu = 1, 2$ is the 'hot region' index and $m, m' = (x, y, z)$ enumerate the components of the paramagnon field $\varphi$. Additionally, $f(\varepsilon_n, p) = \varepsilon_n + i\Sigma(\varepsilon_n, p)$ and $\Omega(\omega_n, q) = \omega_n^2 + v_s^2\Pi(\omega_n, q)$, $\Sigma(\varepsilon_n, p)$ and $\Pi(\omega_n, q)$ being the fermionic self-energy and polarization operator for paramagnons, respectively. In this section $\varepsilon_n = (2n + 1)\pi T$ and $\omega_n = 2n\pi T$ stand for fermionic and bosonic Matsubara frequencies, respectively.

To calculate the self-energies we use the approximations illustrated diagrammatically in Fig. 6. This is justified by a small parameter $\sqrt{(T, \mu, v_s^2/\alpha)/(\alpha/\xi^2)}$ for $\beta/\xi^2 \ll \mu$ (see Appendix C1). To study the formation of the DDW we will, however, use these approximations for all values of $\beta$ assuming that the results to be nevertheless correct at least qualitatively. The resulting momentum-dependent self-consistency equations have the same form as the ones presented in Ref. 85. Assuming the strong overlap of hot spots expressed by the inequality $\alpha/\xi^2 \gg \mu$, one can perform the momentum integration in the self-energies. In the limit $\beta \ll \alpha$ the latter can be shown to be momentum-independent. For larger $\beta$ we approximate the self energies by their values at zero incoming momentum. Additionally, we have evaluated the momentum integral in the polarization operator for the paramagnons without any cutoff. It turns out that to reproduce our previous results one needs to introduce a cutoff $\Lambda$ such that $a\Lambda^2 \rightarrow \infty$ while $b\Lambda^2 \rightarrow 0$. Physically, $\Lambda$ is related to the deviation of the fermionic spectrum from the form (2.2) outside the 'hot regions'. Here we will assume for simplicity that $\beta\Lambda^2 \gg \mu, T$ which should be valid for not too small $\beta$. Further details of the calculations can be found in Appendix C2. Introducing an energy scale $\Gamma = \sqrt{\lambda^2 v_s^2/\alpha}$ (note that in our previous work 85 a different scale ($\lambda^2 v_s/\sqrt{\alpha}^{2/3}$ has been used) the resulting equations can be cast in a dimensionless form where all quantities are assumed to be normalized by $\Gamma$ to the appropriate power

$$f(\varepsilon) - \varepsilon = \frac{1.5T}{\pi} \sum_{\varepsilon'} \text{sgn}[\text{Re}f_{\varepsilon'}] \arctanh \left\{ \frac{\Omega_{\varepsilon', \varepsilon} + \alpha + (i\varepsilon')^2/\beta}{\Omega_{\varepsilon', \varepsilon} + \alpha} \right\}$$

$$\Omega(\omega) - \omega^2 = \frac{T}{1 + \beta/\alpha} \sum_{\varepsilon} \text{sgn}[\text{Re}f_{\varepsilon}^+] \text{sgn}[\text{Re}f_{\varepsilon}^-] \left\{ \frac{4i\text{arctanh}(\sqrt{\beta/\alpha})/\pi\text{sgn}[\text{Re}f_{\varepsilon}^-] + 1}{if_{\varepsilon}^- + \mu} \right\}$$

$$- \frac{2i}{\pi} \text{sgn}[\text{Re}f_{\varepsilon}^+] \text{arctanh} \left\{ \frac{\sqrt{\frac{i}{f_{\varepsilon}^+ + \mu}}}{\sqrt{\frac{i}{f_{\varepsilon}^+ + \mu}}} \right\} + \frac{\text{sgn}[\text{Re}f_{\varepsilon}^-] \text{arctanh} \left\{ \sqrt{\frac{i}{f_{\varepsilon}^- + \mu}} \right\}}{\sqrt{\frac{i}{f_{\varepsilon}^- + \mu}}},$$

where $a = v_s^2/\xi^2$, $\Omega_\omega = \omega^2 + v_s^2\Pi(\omega)$, $f_+ = (\alpha f_{\varepsilon} + \beta f_{\varepsilon})/(\alpha + \beta)$ and $f_- = (\alpha f_{\varepsilon} + \beta f_{\varepsilon})/(\alpha + \beta)$. The value $\Omega(0)$ has been absorbed into redefinition of $1/\xi^2$. Now we proceed to the analysis of the emergence of the particle-hole orders. The general mean-field equation for the Pomeranchuk order has been derived in Ref. 85. For the charge density wave order parameter $W_Q(\varepsilon_n, p)$ one obtains

$$W_Q(\varepsilon_n, p) = 3\lambda^2 T \sum_{\varepsilon',p'} \mathcal{D}(\varepsilon - \varepsilon', p - p') W_Q(\varepsilon'_n, p'_n).$$

$$A_W = G_a^{-1}(\varepsilon', p' + q/2) G_a^{-1}(\varepsilon', p' - q/2) - |W_Q(\varepsilon', p')|^2,$$

while the equations for the DDW and IDDW can be written as

$$D(\varepsilon, p) = 3\lambda^2 T \sum_{\varepsilon',p'} \mathcal{D}(\varepsilon - \varepsilon', p - p') D(\varepsilon'_n, p'_n),$$

$$A_D = G_a^{-1}(\varepsilon', p' + q/2) G_b^{-1}(\varepsilon', p' - q/2) - |D_Q(\varepsilon', p')|^2,$$
These equations can be used to study the full temperature, momentum, and Matsubara frequency dependence of the order parameters, provided the expressions for the normal self-energies are suitably modified (see Ref.\textsuperscript{85} for the case of Pomeranchuk order at small $\beta$). Here we will concentrate on the critical temperatures of the emerging orders in order to check if the general trends observed in Sec.\textsuperscript{III} still hold. Therefore, we simply use Eqs. (4.2) for the normal state self-energies in our study.

Assuming the order parameters to be momentum independent we integrate over momenta (for details see Appendix C\textsuperscript{2}) in the equations for charge and current orders. The resulting equations for the Pomeranchuk $P(\varepsilon)$, CDW $C(\varepsilon)$, DDW $D(\varepsilon)$ and IDDW $D_I(\varepsilon)$ order parameters near the critical temperature are rather cumbersome and we present them in Appendix C\textsuperscript{2} (Eq. C\textsuperscript{6}). All quantities in (C\textsuperscript{6}) are normalized by $\Gamma$. Additionally, to obtain a closed-form answer for IDDW, we have assumed $Q$ along diagonal. While this assumption does not allow one to study the orientation of $Q$, for small $Q$ it is supposed to yield a correct critical temperature, allowing to draw conclusions about the commensurate DDW stability. From Eq. C\textsuperscript{6} it is already evident that for $\beta/\xi^2 \ll 1$ r.h.s. of the equation for the charge orders contains a large factor $1/\sqrt{v_x^2/\alpha}$, ultimately meaning that current orders do not appear in this case.

In Fig.\textsuperscript{7} results of numerical solution of the equations (C\textsuperscript{6}) are presented for two sets of parameters. The value of parameters characterizing the incommensurability for CDW and IDDW have been chosen to maximize the critical temperatures. Considering the qualitative character of our approximations for finite $\beta$ below we analyze only the general features of the obtained results. One can see that the Pomeranchuk instability is considerably more robust to increasing $\beta$ than is expected from the simplified model (3.13). Additionally, the IDDW seems to play a more important role. Actually, both the results can be qualitatively understood as being a consequence of the renormalization of the fermionic self-energy resulting in the replacement $\varepsilon \to f(\varepsilon)$. At low frequencies one obtains $Re f(\varepsilon) \sim \varepsilon/\mathcal{Z}$, where $\mathcal{Z} < 1$. Therefore, the parameter $\mu$ is renormalized to a smaller value $\mathcal{Z} \mu$. As smaller values of $\mu$ qualitatively favor Pomeranchuk and IDDW phases, this explains the observed tendency.

As for charge/current order competition one can draw a conclusion that the boundary between charge and current phases $(\beta/\alpha)_c$ appears to be remarkably close to the one obtained in the simplified model. The dip in critical temperatures at intermediate $\beta/\alpha$ is actually also qualitatively present in the simplified model, however there we concentrated on competition between phases at a given $T$. One could expect that in this region superconductivity will re-emerge as a leading instability even if the remnant Coulomb interaction acts against it. One should also keep in mind that closeness of the different phases in energy may induce strong fluctuations that can modify the results obtained here in the mean field approximation. These effects, however, should be important only close to phase boundaries. Thus, we suggest that the fluctuations will not change the results qualitatively, but leave detailed investigations for future studies.

V. DISCUSSION.

Considering the obtained non-superconducting phases two tentative scenarios can be anticipated for the pseudogap state. In the first one, Pomeranchuk instability is the leading one and is expected to occur at $T_{Pom} \gtrsim T^*$. Then, at $T^*$ the pseudogap would open due to the formation of an axial CDW\textsuperscript{85}. However, the time reversal symmetry breaking does not appear naturally in this scenario unless more complicated form-factors for the CDW are considered\textsuperscript{91,93}. Moreover, while in some compounds\textsuperscript{29} $T_{CDW}$ has been observed to be close to $T^*$ this does not seem to be the general case\textsuperscript{30,31}. Additionally, more recent transport data suggests a Fermi surface reconstruction taking place at $T_{CDW} < T < T^*$\textsuperscript{27} to be distinct from the one caused by the CDW\textsuperscript{135}.

In the other scenario the leading instability is the DDW that has its onset at $T^*$. This is consistent with transport\textsuperscript{27} and ARPES\textsuperscript{108} signatures of the pseudogap. It has also been shown that an axial d-form factor CDW can emerge on the Fermi surface reconstructed by the DDW\textsuperscript{110,111}. DDW breaks time reversal symmetry but, as it breaks also the translational symmetry, additional Bragg peaks at $(\pi + Q_x, \pi + Q_y)$ are expected to appear in the DDW phase. While definitive experimental evidence for these peaks seems to be lacking\textsuperscript{112-115}, we note that the magnitude of the signal predicted by the BCS-like theories\textsuperscript{104} should change in the case of a strongly frequency-dependent order parameter, such as the one in the SF model. Thus it is possible that the magnitude of the additional peaks predicted using a frequency-independent order parameter could be overestimated. Moreover, the $Q = 0$ signal, observed experimentally\textsuperscript{15,16} might originate from higher-order processes but this possibility has not been investigated theoretically so far.

Additionally, at low dopings there is evidence for a $(\pi + Q_x, \pi + Q_y)$ order not accompanied by the CDW from neutron scattering experiments\textsuperscript{130}. We suggest that
this observation can be explained in terms of the incommensurate IDDW order studied in the present work. Unlike previous works, where a similar state has been suggested, in our case IDDW is not accompanied by a charge modulation.

Let us discuss now the values of parameters of the model (2.1) that might be most suitable for the cuprates. Relevant values of $\mu$ have been identified in our previous work and are usually of the order of $T^*$ for the underdoped case. The Fermi surface curvature $\beta/\alpha$ appears now to be another crucial parameter controlling the phase diagram. One can relate $\beta/\alpha$ to the tight-binding parametrization of the dispersion in the full Brilloin zone $\varepsilon(p) = -2t(\cos p_x + \cos p_y) - 4t' \cos p_x \cos p_y - 2t''(\cos 2p_x + \cos 2p_y)$.

$$\frac{\beta}{\alpha} = \frac{t + 2t' - 4t''}{t - 2t' + 4t''}$$

Taking result from literature one obtains $\beta/\alpha \approx 0.11$ for Bi-2201 and $\beta/\alpha \approx 0.15$ for BSCCO. Tight binding fits for YBCO (Ref.138 and references therein) yield negative values for $\beta/\alpha$ not considered here. At the same time, the electronic structure around the antinodes in YBCO has also interpreted in terms of an extended Van Hove singularity corresponding to $\beta = 0$ case. Moreover, there are theoretical arguments that such a behavior should be stabilized by interactions. Anyway, the curvatures are rather low and seemingly constrain us to the regime where the Pomeranchuk instability is the leading one. However, drawing quantitative conclusions about the appearance of the current orders demands taking into account renormalization of the Fermi surface curvature by the low-energy interactions, which is beyond the scope of the present paper.

**VI. CONCLUSION**

We have studied particle-hole instabilities in the spin-fermion model in the regime where shallowness of the antinodal dispersion combined with finite AF correlation length leads to a strong overlap of the 'hot spots' on the Fermi surface. A rich phase diagram has been obtained as a function of the chemical potential (doping) relative to the dispersion saddle-points and Fermi surface curvature in the antinodal regions. The phases obtained include Pomeranchuk and current-ordered phases previously not encountered in the SF model. We have shown that for small curvatures $\beta/\xi^2 \ll \mu, T$ an Eliashberg-like approximation is justified by a small parameter $\sqrt{T, \mu, v_s^2/\alpha}/(\alpha/\xi^2)$. The self-energy effects have been found to promote Pomeranchuk and incommensurate current orders. The current orders possess attractive features for an explanation of the pseudogap phase, namely, the particle-hole asymmetric gap in the antinodal regions, time reversal symmetry breaking and Fermi surface reconstruction into hole pockets. Moreover, the incommensurate current order obtained in this work can potentially explain the incommensurate magnetism observed at low dopings. Finally, we expect our results to be also of relevance to other itinerant systems with strong antiferromagnetic fluctuations.

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**Appendix A: Matsubara Susceptibilities for Simplified Model**

1. **Charge Orders**

In the limit $\beta/\xi^2 \rightarrow 0$ the integral over one of the momenta in (3.6) yields a factor of $2/\xi$, while for the second one limits can be taken to $\pm \infty$ for $\alpha/\xi^2 \gg \mu, \omega_n$. As the resulting sum over $\omega_n$ converges we can neglect the contribution from $\omega_n \gtrsim \alpha/\xi^2$ resulting in (3.8).

In the opposite case $\beta/\xi^2 \gg \omega_n, \mu$ one can extend the integration limits for both momenta to $\pm \infty$. The resulting integrals can be evaluated in this case using:

$$\frac{1}{a_1a_2} = \int_{-1/2}^{1/2} dx \frac{1}{[(a_1 - a_2)x + (a_1 + a_2)/2]^2}.$$ 

One has:

$$(a_1 - a_2)x + (a_1 + a_2)/2 \equiv (2\alpha p_1 Q_1 - 2\beta p_2 Q_2)x + \alpha(p_1^2 + Q_1^2/4) - [\beta(p_2^2 + Q_2^2/4) + i\omega + \mu] = \alpha(p_1 + Q_1x)^2 - \beta(p_2 + Q_2x)^2 - (i\omega + \mu + (\alpha Q_1^2 - \beta Q_2^2)(x^2 - 1/4)).$$
The sum over Matsubara frequencies appears to diverge at large \( \omega_n \), but only logarithmically. This allows one to obtain the leading contribution to \( \chi_{CDW}(Q) \) by introducing a cutoff at \( \omega_n \sim \beta/\xi^2 \) in the sum and neglecting the region \( \omega_n \gtrsim \beta/\xi^2 \) provided that \( \beta/\xi^2 \gg \mu, T \).

To study the stability of the \( Q = 0 \) phase at finite \( \beta \) we expand the CDW susceptibility in Eq. (3.6) in powers of \( Q \). One obtains that \( \frac{\partial \chi}{\partial \xi_1^2} \bigg|_{Q=0} \) as well as \( \frac{\partial^2 \chi}{\partial \xi_1 \xi_2} \bigg|_{Q=0} \) vanish. Since \( Q \) is along the diagonal the condition for the critical value of \( \mu \) is \( \frac{\partial^2 \chi}{\partial \xi_1^2} + \frac{\partial^2 \chi}{\partial \xi_2^2} \bigg|_{Q=0} = 0 \). Performing the expansion and integrating over momenta \( (\alpha \xi^2 \gg \mu, \omega_n \) is assumed) one obtains:

\[
\begin{align*}
\frac{\alpha Q^2}{4} \nu_0 T \sum_{\omega_n} i \pi \text{sgn}[\omega_n] \frac{3(\omega_n + \mu) + 2\beta/\xi^2}{3\sqrt{\alpha} \beta} (\omega_n + \mu)^2 \sqrt{\beta/\xi^2} \bigg( (\beta/\xi^2 + \omega_n + \mu)^3 
\end{align*}
\]

\[
\frac{-\beta Q^2}{4 \nu_0 T \sum_{\omega_n} i \pi \text{sgn}[\omega_n]} \frac{\sqrt{\beta/\xi^2}}{6 \sqrt{\alpha} \beta} \bigg( (\omega_n + \mu)^2 (\omega_n + \mu + \beta/\xi^2)^3/2 \bigg) = 0.
\]

Expanding this result for \( \mu \ll T \) one obtains (3.13) after summation.

2. Current Orders

For \( Q = 0 \) we rewrite the expression in (3.6)

\[
\begin{align*}
\frac{1}{\alpha^2 - (\omega_n + \mu_+)^2} & = \\
\frac{1}{\alpha + \beta (\omega_n - \alpha p_1^2 + \beta p_2^2 - \mu_1)} & = \\
\frac{\alpha}{\alpha + \beta (\omega_n - \alpha p_1^2 + \beta p_2^2 - \mu_1)} & - \frac{\alpha - \beta}{\alpha + \beta (\omega_n - \alpha p_1^2 + \beta p_2^2 - \mu_1)} \\
+ & \frac{1}{\alpha + \beta (\omega_n - \alpha p_1^2 + \beta p_2^2 - \mu_1)}. 
\end{align*}
\]

Assuming only \( \alpha/\xi^2 \gg \mu, \omega_n \) one can evaluate the integral over momenta analytically. This yields

\[
\chi_D(T) = T \sum_{\omega_n} \frac{4 \pi \text{sgn}[\omega_n] \text{arctanh} \left\{ \frac{\alpha/\xi^2}{\sqrt{\omega_n - \beta \omega_n + \mu}} \right\}}{\alpha + \beta (\omega_n + \mu)} + \frac{4}{\alpha + \beta} \text{arctanh} \left\{ \frac{\alpha - \beta}{\omega_n + \mu} \right\} \text{arctanh} \left\{ \frac{\omega_n + \mu}{\sqrt{\omega_n + \mu}} \right\}. \quad (A2)
\]

From this equation one can obtain (3.18) and (3.19). For finite \( Q \) a closed form for \( \chi_D(Q) \) can be obtained for \( \beta/\xi^2 \gg T, \mu \) using:

\[
\frac{1}{a_1 a_2} = \int_{-1/2}^{1/2} dx \frac{1}{(a_1 - a_2) x + (a_1 + a_2) / 2}. 
\]
For the momentum integral in (3.6) we obtain:

\[
- \int_{-\infty}^{\infty} dp_1 \int_{-\infty}^{\infty} dp_2 \int_{-1/2}^{1/2} dx \frac{1}{iQ^2},
\]

\[
[...] = \left[ (\alpha + \beta)x + \frac{\alpha - \beta}{2} \right] (p_1^2 + Q_1^2/4) + \left[ (\alpha - \beta)x + \frac{\alpha + \beta}{2} \right] p_1 Q_1
\]

\[
+ \left[ -(\alpha + \beta)x + \frac{\alpha - \beta}{2} \right] (p_2^2 + Q_2^2/4) + \left[ (\alpha - \beta)x - \frac{\alpha + \beta}{2} \right] p_2 Q_2 - (i\omega + \mu).
\]

To change the integration order we need to assume \(Q_1, Q_2 \neq 0\) as \(\pm(\alpha + \beta)x + (\alpha - \beta)/2\) can vanish inside the \(x\) integration region while \((\alpha - \beta)x \pm (\alpha + \beta)/2\) does not cross zero for all \(x\). To integrate over \(p_1\) we rewrite \([...]\):

\[
[...] = \left[ (\alpha + \beta)x + \frac{\alpha - \beta}{2} \right] \left( p_1 + \frac{(\alpha - \beta)x + (\alpha + \beta)/2 Q_1}{(\alpha + \beta)x + (\alpha - \beta)/2} \right)^2 + \frac{Q_1^2}{4} \frac{4\alpha\beta(x^2 - 1/4)}{(\alpha + \beta)x + (\alpha - \beta)/2}
\]

\[- \left[ (\alpha + \beta)x - \frac{\alpha - \beta}{2} \right] \left( p_2 + \frac{-(\alpha - \beta)x + (\alpha + \beta)/2 Q_2}{(\alpha + \beta)x - (\alpha - \beta)/2} \right)^2 - \frac{Q_2^2}{4} \frac{4\alpha\beta(x^2 - 1/4)}{(\alpha + \beta)x - (\alpha - \beta)/2} - (i\omega + \mu).
\]

Now one can simplify the calculation by shifting the integration variables. First let us integrate over \(p_1\). The answer depends on the sign of \((\alpha + \beta)x + \frac{\alpha - \beta}{2}\):

\[
-1 \frac{2}{(\alpha + \beta)x - \frac{\alpha - \beta}{2}} \left( p_2^2 + \frac{Q_2^2\alpha\beta(x^2 - 1/4)}{(\alpha + \beta)x - (\alpha - \beta)/2} - \frac{Q_1^2\alpha\beta(x^2 - 1/4)}{(\alpha + \beta)x + (\alpha - \beta)/2} + i\omega + \mu \right)^{(3/2)}.
\]

The remaining integral over \(p_2\) yields:

\[
\int_{-\infty}^{\infty} dp_2 \frac{1}{(\alpha + \beta)x - \frac{\alpha - \beta}{2}} \left( p_2^2 + \frac{Q_2^2\alpha\beta(x^2 - 1/4)}{(\alpha + \beta)x - (\alpha - \beta)/2} - \frac{Q_1^2\alpha\beta(x^2 - 1/4)}{(\alpha + \beta)x + (\alpha - \beta)/2} + i\omega + \mu \right)^{(3/2)} =
\]

\[
- \frac{1}{i\omega + \mu + \frac{Q_2^2\alpha\beta(x^2 - 1/4)}{(\alpha + \beta)x - (\alpha - \beta)/2} - \frac{Q_1^2\alpha\beta(x^2 - 1/4)}{(\alpha + \beta)x + (\alpha - \beta)/2}} \left\{ \begin{array}{ll}
\frac{2}{\sqrt{(\alpha + \beta)x - (\alpha - \beta)/2}} & \text{if } x > \frac{\alpha - \beta}{2(\alpha + \beta)} \\
-\frac{2\text{sgn}(\omega_0)}{\sqrt{(\alpha + \beta)x + (\alpha - \beta)/2}} & \text{if } x < \frac{\alpha - \beta}{2(\alpha + \beta)}. 
\end{array} \right.
\]

Combining the results above one obtains two contributions. The first one is:

\[
I_1 = \int_{-1/2}^{1/2} dx \frac{\pi}{(\alpha + \beta)^2 x^2 - (\alpha - \beta)^2/4}
\]

\[
(i\omega + \mu)[(\alpha + \beta)^2 x^2 - (\alpha - \beta)^2/4] + \alpha\beta(x^2 - 1/4)\{(Q_2^2 - Q_1^2)(\alpha + \beta)x + (Q_2^2 + Q_1^2)(\alpha - \beta)/2\}.
\]

Or, after a change of variables \(x \rightarrow (\alpha - \beta)x/(\alpha + \beta)/2\) and some algebra:

\[
\int_{-1}^{1} dx \frac{\pi}{\alpha + \beta} \frac{\sqrt{1 - x^2}}{(i\omega + \mu)[1 - x^2] + \frac{\alpha\beta}{\alpha - \beta} \left(1 - x^2 \left(\frac{\alpha - \beta}{\alpha + \beta}\right)^2\right)} \{\delta Q^2 x + Q_2^2\}
\]

where \(Q^2 = (Q_1^2 + Q_2^2)/2\) and \(\delta Q^2 = (Q_2^2 - Q_1^2)/2\). For \(\delta Q^2 = 0\) one can evaluate the integral analytically to obtain

\[
I_1(Q, Q) = \frac{\pi^2}{\alpha + \beta} \frac{1}{i\omega + \mu + \frac{\alpha\beta(\alpha - \beta)Q_1^2}{(\alpha + \beta)^2}} \left[ 1 - \sqrt{\frac{4\alpha^2\beta^2 Q_1^2}{(\alpha + \beta)^2(\alpha - \beta)^2} \frac{i\omega + \mu + \frac{\alpha\beta Q_1^2}{\alpha - \beta}}{i\omega + \mu + \frac{\alpha\beta Q_1^2}{\alpha - \beta}} \right].
\]
Matsubara sum diverges logarithmically, however one can subtract the divergent part. The divergent part is then evaluated with a cutoff at 

\[ M = \frac{1}{i\omega + \mu + \frac{\alpha(\alpha - \beta)Q^2}{(\alpha + \beta)^2}} \left( \arctanh \sqrt{\frac{\beta}{\alpha}} - \sqrt{\frac{\alpha}{\alpha + \beta} \frac{\alpha^2 Q^2}{\alpha - \beta}} \arctanh \sqrt{\frac{\beta}{\alpha + \beta}} \right). \]

For numerical calculations of \( \delta Q = 0 \), the integral in \( I_2 \) can be evaluated to obtain

\[ \chi_{\text{DDW}}(Q) = \frac{\pi^2}{\alpha + \beta} \frac{1}{i\omega + \mu + \frac{\alpha(\alpha - \beta)Q^2}{(\alpha + \beta)^2}} \left( \arctanh \sqrt{\frac{\beta}{\alpha}} - \sqrt{\frac{\alpha}{\alpha + \beta} \frac{\alpha^2 Q^2}{\alpha - \beta}} \arctanh \sqrt{\frac{\beta}{\alpha + \beta}} \right). \]

Combining this with the first contribution we get:

The expression (A3) can be used to calculate the IDDW/DDW phase boundary because for small \( Q \) one can show that \( \chi_D(Q \to 0) \approx a + b(Q_0^2 + Q_0^2) \). From the condition \( \frac{\partial \chi_D}{\partial Q^2} = 0 \) we get after evaluation of the Matsubara sum:

\[ -\frac{i}{2\pi(\alpha + \beta)^2} \left[ \frac{(\alpha \beta)^{3/2}}{\alpha - \beta} + \frac{\alpha \beta(\alpha - \beta)}{\alpha + \beta} \arctanh \sqrt{\frac{\beta}{\alpha}} \right] \left( \psi \left[ \frac{1}{2} + i\mu \frac{1}{2\pi T} \right] - \psi \left[ \frac{1}{2} - i\mu \frac{1}{2\pi T} \right] \right) \]

\[ + \frac{(\alpha - \beta)\alpha \beta \pi^2}{8(\alpha + \beta)^3 \cosh^2 \left( \mu/2T \right)} = 0. \]

This expression has been used to calculate the IDDW/DDW boundary numerically. As is evident from Fig. 4 \( \mu/T \) becomes small for \( (\alpha - \beta) \ll \alpha \). Expanding (A4) for \( \mu/T \ll 1, (\alpha - \beta) \ll \alpha \) one obtains Eq. 3.20.

For numerical calculations of \( \chi_{\text{DDW}} \) and orientation of \( Q \) we have evaluated the Matsubara sum before the integrals in \( I_1, I_2 \). Using \( T \sum \frac{1}{i\omega + \alpha} = \tanh(\alpha/2T)/2 \) one obtains for \( I_1 \):

\[ \chi_{\text{DDW}} = \frac{\pi}{2(\alpha + \beta)} \int_{-1}^{1} \frac{dx}{\sqrt{1 - x^2}} \tanh \left( \frac{\mu}{2T} \right) + \frac{\alpha}{1 - \frac{(\alpha - \beta)Q^2}{2T}} \left[ \psi \left( \frac{1}{2} + i\kappa \frac{1}{2\pi T} \right) - \psi \left( \frac{1}{2} - i\kappa \frac{1}{2\pi T} \right) \right] \]
Orientation and magnitude of $\mathbf{Q}$ are found maximizing $\chi_{DDW}(\mathbf{Q})$. One can get some analytical insight on the possible orientation of $\mathbf{Q}$ in the case $\beta \ll \alpha$. $I_2$ can be approximately evaluated taking $x \approx \pm \frac{\alpha - \beta}{2(\alpha + \beta)}$ in the integral:

$$I_2 \approx \frac{2\pi \text{sgn}[\omega_n]\sqrt{\frac{\beta}{\alpha}}}{\alpha [\omega + \mu + \beta Q_2^2]} \left(1 - \sqrt{\frac{\beta Q_2^2}{\omega + \mu + \beta Q_2^2}} \frac{\text{arctanh} \sqrt{\frac{\omega + \mu + \beta Q_1^2}{\beta Q_2^2}}}{\beta Q_2^2} \right)$$

$$+ \frac{2\pi \text{sgn}[\omega_n]\sqrt{\frac{\beta}{\alpha}}}{\alpha [\omega + \mu + \beta Q_2^2]} \left(1 - \sqrt{\frac{\beta Q_2^2}{\omega + \mu + \beta Q_2^2}} \frac{\text{arctanh} \sqrt{\frac{\omega + \mu + \beta Q_1^2}{\beta Q_2^2}}}{\beta Q_2^2} \right)$$

At $T, \mu \ll \beta Q^2$ one can go to integration over $\omega$. The second terms in the brackets yield after integration over $\omega$:

$$\chi^2_{DDW} \approx \frac{\beta}{\alpha^3} \log \left(\frac{(\beta/\epsilon)^2}{(\beta Q^2 + \mu)^2 - (\beta Q^2)^2} \frac{\beta}{\alpha} \right) + \frac{\beta}{\alpha^3} \text{const} + O(\mu/\beta Q^2) \sim$$

$$2\pi \frac{\beta}{\alpha^3} \log \left(\frac{\beta/\epsilon^2}{(\beta Q^2 + \mu)^2} \right) + O(\mu/\beta Q^2) + \text{const.}$$

$\chi_{DDW}^1(\mathbf{Q})$ can be shown to be bounded from above by $\frac{\pi^2}{2(\alpha + \beta)}$. Consequently, for $\beta/\epsilon^2 \gg \beta Q^2 \gg \mu, T$ contribution from $\chi^2_{DDW}$ is dominant and maximizing $\delta Q^2$ is favorable. The maximal absolute value of $\delta Q^2$ is $Q^2$ that is reached is reached for $\mathbf{Q}$ along one of the BZ axes. However as $\beta$ decreases $\chi^2_{DDW}$ eventually becomes less important due to the factor $\sqrt{\beta/\alpha}$. As $\chi_{DDW}$ has been shown above to be maximal for $\delta Q^2 = 0$ one expects a transition to diagonal $\mathbf{Q}$ at low $\beta/\alpha$. This is in line with the results of numerical calculation in Fig. 5.

Appendix B: Expression for the current density

The definition of the current density directly follows from the standard expression for the magnetic part of the action $S_{\text{mag}}$ in the presence of a vector potential $\mathbf{A}(\mathbf{r})$. Here we derive the bare part $S_0[\mathbf{A}]$ of the action $S$ in the presence of an external potential $\mathbf{A}(\mathbf{r})$ and derive the expression for the current using a standard formula from electrodynamics for the magnetic part $S_{\text{mag}}$ of the action

$$S_{\text{mag}} = -\frac{1}{c} \int (J_\mathbf{r} \mathbf{A}(\mathbf{r}, \xi) d\tau d\mathbf{r}, \quad (B1)$$

For simplicity we consider the energy operator to be of the form

$$\mathcal{E}(-i\nabla) = J(2 - \cos(-i\nabla \mathbf{a}_x) - \cos(-i\nabla \mathbf{a}_y)). \quad (B2)$$

In Eq. (B2) $\mathbf{a}_x$ and $\mathbf{a}_y$ are lattice vectors directed along $x$ and $y$ bonds of the $CuO$ lattice, respectively, and $|\mathbf{a}_x| = |\mathbf{a}_y| = a_0$.

The energy operator for the system with the vector potential can easily be written using the minimal coupling equivalent to the Peierls’ substitution in Eq. (B2)

$$-i\nabla \rightarrow -i\nabla - \frac{e}{c} \mathbf{A} \quad (B3)$$

Then, the energy operator takes the form

$$\mathcal{E}\left(-i\nabla - \frac{e}{c} \mathbf{A}(\mathbf{r})\right) = J\left[\cos\left(-i\nabla - \frac{e}{c} \mathbf{A}(\mathbf{r})\right) \mathbf{a}_x\right]$$

$$\cos\left(-i\nabla - \frac{e}{c} \mathbf{A}(\mathbf{r})\right) \mathbf{a}_y\right) - 2] \quad (B4)$$

The current operator should be defined calculating the linear term of the expansion of the action $S_{\text{mag}}$ in $\mathbf{A}(\mathbf{r})$.

As the vector potential $\mathbf{A}(\mathbf{r})$ does not commute with the gradient, the expansion in $\mathbf{A}(\mathbf{r})$ is not trivial and we use time ordering products. For any non-commuting operators $A$ and $B$ one has

$$\exp(A + B) = \exp AT_\alpha \exp \left(\int_0^1 \tilde{B}(\alpha) d\alpha\right)$$

$$= T_\alpha \exp \left(\int_0^1 \tilde{B}^A(\alpha) d\alpha\right) \exp A, \quad (B5)$$
\[ \hat{B} (\alpha) = e^{-\mathcal{A}_t} B e^{\mathcal{A}_t}, \quad \hat{B}^A (\alpha) = e^{\mathcal{A}_t} B e^{-\mathcal{A}_t}, \] (B6)

and \( T^t \) and \( T^A \) are time ordering and anti time ordering operators, respectively.

Introducing operators
\[ \tilde{A} (\mathbf{r}, \alpha) = e^{-\mathcal{A}_t} \mathbf{A} (\mathbf{r}) e^{\mathcal{A}_t} = \mathbf{A} (\mathbf{r} - \alpha \mathbf{a}_x), \]
\[ \tilde{A}^A (\mathbf{r}, \alpha) = e^{\mathcal{A}_t} \mathbf{A} (\mathbf{r}) e^{-\mathcal{A}_t} = \mathbf{A} (\mathbf{r} + \alpha \mathbf{a}_x) \] (B7)

one comes to the following expressions
\[ \exp \left[ \pm \left( \nabla - \frac{ie}{c} \mathbf{A} (\mathbf{r}) \right) \alpha \right] \]
\[ = \exp [\pm \alpha \nabla] \exp \left[ \frac{ie}{c} \int_0^1 \alpha_x \mathbf{A} (\mathbf{r} \pm \alpha \mathbf{a}_x) \, d\alpha \right] \]
\[ = \exp \left[ \frac{ie}{c} \int_0^1 \alpha_x \mathbf{A} (\mathbf{r} \pm \alpha \mathbf{a}_x) \, d\alpha \right] \exp [\pm \alpha \nabla] \] (B8)

Using Eq. (B8) we rewrite the energy operator \( \hat{\varepsilon} \left( -i \nabla - \frac{e}{c} \mathbf{A} (\mathbf{r}) \right) \). This leads to the following expression for the action
\[ S_0 [\mathbf{A}] = -t \int c^\dagger (\tau, \mathbf{r}) \left[ \exp \left[ \mp \frac{ie}{c} \int_0^1 \alpha_x \mathbf{A} (\mathbf{r} + \alpha \mathbf{a}_x) \, d\alpha \right] \right. \]
\[ \left. \times \int_0^1 \left[ \langle c^\dagger (\tau, \mathbf{r} + \alpha \mathbf{a}_x, \mathbf{a}_y) c (\tau, \mathbf{r} + (1 - \alpha) \mathbf{a}_x, \mathbf{a}_y) \rangle \right] \right. \]
\[ - \langle c^\dagger (\tau, \mathbf{r} + \alpha \mathbf{a}_x, \mathbf{a}_y) c (\tau, \mathbf{r} - (1 - \alpha) \mathbf{a}_x, \mathbf{a}_y) \rangle \right] d\tau \] (B9)

Now, expanding the exponentials in the vector potential \( \mathbf{A} (\mathbf{r}) \), and comparing the linear in \( \mathbf{A} (\mathbf{r}) \) term with \( S_{\text{mag}} \), Eq. (B1), we bring the correlation function for the current density to the form
\[ j_{x,y} (\tau, \mathbf{r}) = -\frac{i}{2} eJ \mathbf{a}_{x,y} \]
\[ \times \int_0^1 \left[ \langle c^\dagger (\tau, \mathbf{r} - \alpha \mathbf{a}_x, \mathbf{a}_y) c (\tau, \mathbf{r} + (1 - \alpha) \mathbf{a}_x, \mathbf{a}_y) \rangle \right] \frac{d\alpha}{2}, \] (B10)

where \( j_{x,y} \) are \( x \)- and \( y \)- components of the current density. The angular brackets in Eq. (B10) stand for averaging with the action \( S \) of the system. We will use for this averaging the action in the mean field approximation.

We emphasize that the current density \( j_{x,y} (\tau, \mathbf{r}) \) is a function of the continuous coordinate \( \mathbf{r} \) and Eq. (B10) is valid not only on the sites of the lattice. This is very important because in some cases non-zero circulating currents turn to zero at these points.

In order to calculate physical quantities can expand the fields \( c (\mathbf{r}) \) in the Bloch functions \( \psi_\mathbf{P} \)
\[ c (\mathbf{r}) = \int c_\mathbf{P} \psi_\mathbf{P} (\mathbf{r}) \frac{d\mathbf{P}}{(2\pi)^2}, \] (B11)

where \( \psi_\mathbf{P} (\mathbf{r}) \) has the standard form
\[ \psi_\mathbf{P} (\mathbf{r}) = e^{i\mathbf{P} \cdot \mathbf{r}} u_\mathbf{P} (\mathbf{r}), \] (B12)

\( u_\mathbf{P} (\mathbf{r}) \) is periodic function with the period \( \mathbf{a}_{x,y} \) and \( \mathbf{P} \) is a quasimomentum in the first Brillouin zone. Note that in the main text the quasimomenta are defined in units of the inverse lattice spacing.

However, as we use the spectrum, Eq. (B4), corresponding to a tight binding limit, the eigenfunctions of the Hamiltonian are localized near the lattice sites and it is convenient to expand the Bloch functions in Wannier functions \( w_{\mathbf{R}_n} (\mathbf{r}) \) representing the functions \( \psi_\mathbf{P} (\mathbf{r}) \) as
\[ \psi_\mathbf{P} (\mathbf{r}) = \sum_{\mathbf{R}_n} e^{i\mathbf{P} \cdot \mathbf{R}_n} w_{\mathbf{R}_n} (\mathbf{r}), \] (B13)
where \( \mathbf{R}_n = a_x n_x + a_y n_y, \mathbf{n} = (n_x, n_y), \ n_x, n_y = 0, \pm 1, \pm 2, \pm 3, \ldots \) and \( N \) is the total number of the sites. Then,

\[
c(\mathbf{r}) = \sum_{\mathbf{R}_n} c_\mathbf{p} e^{i \mathbf{p} \cdot \mathbf{R}_n} w_{\mathbf{R}_n}(\mathbf{r}) \tag{B14}
\]

The functions \( w_{\mathbf{R}_n}(\mathbf{r}) \) are localized near the sites with the coordinates \( \mathbf{r} - \mathbf{R}_n \). The function \( w_0(\mathbf{r}) \) is localized near \( \mathbf{r} = 0 \), and \( w_{\mathbf{R}_n}(\mathbf{r}) = w_0(\mathbf{r} - \mathbf{R}_n) \). The Wannier functions are normalized as follows

\[
\int |w_{\mathbf{R}_n}(\mathbf{r})|^2 d\mathbf{r} = 1 \tag{B15}
\]

Taking the Fourier transform of the current

\[
\mathbf{j}_{x,y}(\tau, \mathbf{q}) = \int \mathbf{j}_{x,y}(\tau, \mathbf{r}) e^{-i \mathbf{q} \cdot \mathbf{r}} d\mathbf{r} \tag{B16}
\]

we substitute Eq. (B14) into Eq. (B10) and the latter into Eq. (B16). The we shift \( \mathbf{r} \to \mathbf{r} - \mathbf{a}_{x,y} \) in the first term in \( \mathbf{j}_{x,y}(\mathbf{r}) \) and \( \mathbf{r} \to \mathbf{r} + \mathbf{a}_{x,y} \) in the second one and use the fact that the product \( w_{\mathbf{R}_n}(\mathbf{r})w_{\mathbf{R}_n}(\mathbf{r} \pm \mathbf{a}_{x,y}) \) is essentially different from zero only for \( \mathbf{R}_n = \mathbf{R}_{n_2} \pm \mathbf{a}_{x,y} \). The integral over \( \mathbf{r} \) reduces to the following expression

\[
\int e^{-i q r} w_0^2(\mathbf{r}) d\mathbf{r} \simeq 1 \tag{B17}
\]

for \( |\mathbf{q}| \ll l_c^{-1} \), where \( l_c \) is the localization radius of the function \( w_0(\mathbf{r}) \).

The calculation of the sum over \( \mathbf{R}_n \) is performed using the Poisson formula

\[
\sum_{\mathbf{R}_n} e^{i \mathbf{R}_n (\mathbf{P}_2 - \mathbf{P}_1 - \mathbf{q})} = \left( \frac{2\pi}{a_0} \right)^2 \sum_{\mathbf{K}_n} \delta(\mathbf{P}_2 - \mathbf{P}_1 - \mathbf{q} - \mathbf{K}_n), \tag{B18}
\]

where \( \mathbf{K}_n \) is the vector of the reciprocal lattice, \( \mathbf{K}_n = \frac{2\pi}{a_0} (n_x, n_y) \). The summation over the vectors of the reciprocal lattice is important because \( \mathbf{q} \) is not necessarily located in the first Brillouin zone.

As a result, we come to the following expression for the current density (as it does not depend on time, we omit from now on the variable \( \tau \))

\[
\mathbf{j}_{x,y}(\mathbf{q}) = e \mathbf{e}_{x,y} \sum_{\mathbf{K}_n} \int \delta(\mathbf{P}' - \mathbf{P} - \mathbf{q} - \mathbf{K}_n)
\]

\[
\times \left\langle \mathbf{c}_\mathbf{p}(\tau) \mathbf{v}_{x,y}(\mathbf{P}) \mathbf{c}_\mathbf{p}'(\tau) \right\rangle \frac{d\mathbf{P}d\mathbf{P}'}{(2\pi)^2}, \tag{B19}
\]

where \( \mathbf{e}_{x,y} = \mathbf{a}_{x,y}/a_0 \) is the unit vector along \( x \) or \( y \) bond. In Eq. (B19) integration is performed over all \( \mathbf{P} \) and \( \mathbf{P}' \) inside the first Brillouin zone. The effective velocity \( \mathbf{v}_{x,y}(\mathbf{P}) \) equals

\[
\mathbf{v}_{x,y}(\mathbf{P}) = J a_0 \int_0^1 \sin \left[ (\mathbf{P} + \alpha \mathbf{a}_{x,y}) \mathbf{a}_{x,y} \right] d\alpha. \tag{B20}
\]

In the limit \( \mathbf{q} \to 0 \), the function \( \mathbf{v}_{x,y}(\mathbf{P}) \) is just the conventional velocity

\[
\mathbf{v}_{0,y}(\mathbf{P}) = J a_0 \sin(\mathbf{P}a_{x,y}). \tag{B21}
\]

Averaging in Eq. (B19) we reduce the latter to the form

\[
\mathbf{j}_{x,y}(\tau, \mathbf{q}) = 2 e \mathbf{e}_{x,y} \sum_{\mathbf{K}_n} \int \delta(\mathbf{P}' - \mathbf{P} - \mathbf{q} - \mathbf{K}_n)
\]

\[
\times \mathbf{v}_{x,y}(\mathbf{P}) g_{\mathbf{p}' \mathbf{p}}(0) \frac{d\mathbf{P}d\mathbf{P}'}{(2\pi)^2}, \tag{B22}
\]

where

\[
g_{\mathbf{p}' \mathbf{p}}(0) = - \left\langle \mathbf{c}_\mathbf{p}'(\tau) \mathbf{c}_\mathbf{p}^*(\tau) \right\rangle. \tag{B23}
\]
The factor 2 in Eq. (B22) is due to spin.

The main contribution in the integral over \( P \) in Eq. (B19) comes from the hot regions and it is again convenient to change to the variables \( c_{1,2} \), Eq. (2.1) and the momenta \( p \) counted from the middle of the edges of the reciprocal lattice. Then, using the symmetry relation

\[
g_{p}^{12}(0) = -g_{p}^{21}(0) \tag{B24}
\]

and the fact that

\[
e^{iQ_{x}a_{x,y}} = -e^{iQ_{y}a_{x,y}} \tag{B25}
\]

Where \( Q_{x} = (\pi/a_{0}, 0) \), \( Q_{y} = (0, \pi/a_{0}) \) Eq. (B22) can be written in the form

\[
\mathbf{j}_{x,y}(\mathbf{q}) = 2\epsilon e_{x,y} \sum_{K_{n}} \frac{d\mathbf{p}}{(2\pi)^{2}} \bar{v}_{x,y}(\mathbf{p}) g_{p}^{12}(0) \times \left[ \delta(\mathbf{q} + K_{n} + Q_{AF}) + \delta(\mathbf{q} + K_{n} - Q_{AF}) \right], \tag{B26}
\]

where

\[
\bar{v}_{x,y}(\mathbf{p}) = ta_{0} \int_{0}^{1} \sin \left[ (\mathbf{p} + \alpha \mathbf{q}) a_{x,y} \right] d\alpha = \frac{t \cos \mathbf{p} a_{x,y} - \cos \left( (\mathbf{p} + \mathbf{q}) a_{x,y} \right)}{\mathbf{q} e_{x,y}}. \tag{B27}
\]

Integrating in Eq. (B26) over \( p \) we reduce this equation to the form

\[
\mathbf{j}_{x,y}(\mathbf{q}) = -8et \bar{D} e^{iQ_{x}a_{x,y}} e_{x,y} \times \sin^{2} \left( \frac{\mathbf{q} a_{x,y}}{2} \right) \sum_{K_{n}} \delta(\mathbf{q} - Q_{AF} + K_{n}), \tag{B28}
\]

where

\[
\bar{D} = i \int g_{p}^{12}(0) \cos \left( \mathbf{p} a_{x,y} \right) \frac{d\mathbf{p}}{(2\pi)^{2}} \approx \frac{i}{(2\pi)^{2}} \left( \langle \hat{c}_{1} \hat{c}_{2}^{\dagger} \rangle - \langle \hat{c}_{2} \hat{c}_{1}^{\dagger} \rangle \right). \tag{B29}
\]

As the order parameter, Eq. (3.3), is imaginary, the coefficient \( \bar{D} \) is real. The current has peaks at points \( Q_{AF} + K_{n_{x}} \) and \( Q_{AF} + K_{n_{y}} \). The function \( \mathbf{j}_{x,y}(\mathbf{q}) \) can be further simplified at the peak values and represented in the form

\[
\mathbf{j}_{x,y}(\mathbf{q}) = -8et \bar{D} e^{iQ_{x}a_{x,y}} e_{x,y} \times \left( \mathbf{q} e_{x,y} \right)^{-1} \sum_{K_{n}} \delta(\mathbf{q} - Q_{AF} + K_{n}), \tag{B30}
\]

Nevertheless, the sine function can be relevant if the peaks are smeared.

The particle conservation reads

\[
\mathbf{q} \mathbf{j}(\mathcal{T}, \mathbf{q}) = \mathbf{q} \mathbf{j}_{x}(\mathcal{T}, \mathbf{q}) + \mathbf{q} \mathbf{j}_{y}(\mathcal{T}, \mathbf{q}) = 8et \bar{D} \sum_{K_{n}} \left( e^{iQ_{x}a_{x}} + e^{iQ_{y}a_{y}} \right) \delta(\mathbf{q} - Q_{AF} + K_{n}) = 0. \tag{B31}
\]

The current density can also be written in the real space. It is important to emphasize that \( \mathbf{q} \) is a momentum (not a quasimomentum) and therefore the current density in the real space \( \mathbf{j}(\mathbf{r}) \) can be written for the continuous coordinate \( \mathbf{r} \) using the standard Fourier transform, Eq. (B16). A simple calculation leads to the following expression

\[
\mathbf{j}_{x,y}(\mathbf{r}) = -\frac{et \bar{D} a_{0}^{3} e^{iQ_{x}a_{x,y}}}{\pi^{2}} e_{x,y} \times \sum_{R_{n}} \left[ \cos(Q_{AF} R_{n}) \int_{0}^{1} \delta(\mathbf{r} - R_{n} + \alpha a_{x,y}) d\alpha \right]. \tag{B32}
\]
Eq. (B32) describes currents circulating around the elementary cells. The currents oscillate with the period $Q_{AF}$ and form the bond current antiferromagnet. This picture corresponds to the one proposed in Ref. 104.

In order to avoid a confusion we would like to note that $j_{x,y}(\tau, \mathbf{r})$ is the two dimensional current density. The three dimensional current density $J_{x,y}(\mathbf{r})$ can be written as

$$J_{x,y}(\mathbf{r}) = \sum_{m=-\infty}^{\infty} j_{x,y}(\mathbf{r}, m) \delta(z - mc_0), \quad (B33)$$

where $z$ is the coordinate perpendicular to the planes and $c_0$ is the distance between the layers. Provided the currents in different layers are in phase (the function $j_{x,y}(\mathbf{r}, m)$ does not depend on $m$) one can approximate for rough estimates the 3D current density as follows

$$J_{x,y}(\mathbf{r}) \approx j_{x,y}(\mathbf{r})/c_0 \quad (B34)$$

In the next subsection we use this approximation in order to visualize roughly the structure of the magnetic field.

The circulating currents produce magnetic fields that can be measured by various techniques. As the explicit expression for the spontaneous currents has been obtained, the magnetic field can be determined without difficulties. The Fourier transform of the magnetic field $B(\tau, \mathbf{q})$ can easily be written using the Maxwell equation

$$B(\tau, \mathbf{q}) = \frac{4\pi i}{c} \mathbf{q} \times J(\tau, \mathbf{q}) \quad (B35)$$

With the approximation (B34), only the $z$-component $B_z$ of the field,

$$B_z(\tau, \mathbf{q}) = \frac{4\pi i}{c} q_x j_y - q_y j_x \quad (B36)$$

is not equal to zero.

**Appendix C: Details of calculations for the SF model**

1. Estimate of vertex corrections

Here we show that the vertex corrections contain a small parameter $\sqrt{T, \mu, v_s^2/\alpha}/(\alpha/\xi^2)$ if $\beta \ll \alpha$. As an example we compare the two self-energy diagrams presented in Fig. 8 for $\beta = 0$. The bare propagators for fermions $G_a = (i\varepsilon + \mu - \alpha p_a^2)^{-1}$ are assumed to be much 'sharper' in momentum space than the bosonic ones $\mathcal{D} = (\omega^2/v_s^2 + p^2 + 1/\xi^2)^{-1}$ due to $\alpha/\xi^2 \gg \mu$. The incoming momenta $\mathbf{p}$ are taken to be $\sim \sqrt{\mu/\alpha} \ll 1/\xi$ in magnitude. This allows one to simplify the resulting integrals neglecting the dispersion in the bosonic propagators along $a, b$ or both. We get

$$\lambda^4 T^2 \sum_{\varepsilon', \varepsilon''} \int dp' dp'' \mathcal{D}(\varepsilon - \varepsilon', \mathbf{p} - \mathbf{p}') G_a^2(\varepsilon', \mathbf{p}') \mathcal{D}(\varepsilon' - \varepsilon'', \mathbf{p}' - \mathbf{p}'') G_b(\varepsilon'', \mathbf{p}'') =$$

$$\lambda^4 T^2 \sum_{\varepsilon', \varepsilon''} \int dp' dp'' \frac{1}{(\varepsilon - \varepsilon')^2/v_s^2 + (\mathbf{p} - \mathbf{p}')^2 + \frac{1}{\xi^2} (i\varepsilon' + \mu - \alpha (p_a')^2)^2} \frac{1}{(\varepsilon' - \varepsilon'')^2/v_s^2 + (\mathbf{p}' - \mathbf{p}'')^2 + \frac{1}{\xi^2} i\varepsilon'' + \mu - \alpha (p_a'')^2}$$

![FIG. 8. Self-energy diagrams of the same order a) without b) with vertex corrections](image-url)
Using the strong overlap $1/\xi \gg \sqrt{|\mu, T|/\alpha}$ we can neglect $\mathbf{p}$ and $\mathbf{p}_{a}'$ in the first bosonic propagator and $\mathbf{p}_{a}'$ and $\mathbf{p}_{a}''$ in the second. We proceed to obtain

$$\lambda^4 T^2 \sum_{\epsilon', \epsilon''} \int d\mathbf{p}' d\mathbf{p}'' \frac{1}{(\epsilon - \epsilon')^2/v_s^2 + (\mathbf{p}')^2 + \frac{1}{\xi} (i\epsilon' + \mu - \alpha(\mathbf{p}_{a}')^2) (\epsilon' - \epsilon'')^2/v_s^2 + (\mathbf{p}_{a}')^2 + \frac{1}{\xi} i\epsilon'' + \mu - \alpha(\mathbf{p}_{a}'')^2} \sim \frac{\lambda^4 T^2}{L^{5/2}} \sum_{\epsilon', \epsilon''} \sqrt{(\epsilon - \epsilon')^2 + v_s^2} (i\epsilon' + \mu)^{3/2} \sqrt{(\epsilon' - \epsilon'')^2 + v_s^2} \sqrt{\epsilon'' + \mu},$$

where all quantities having dimensions of energy after $\sum$ are normalized to $\Gamma$. For diagram b) we get

$$b) \sim \lambda^4 T^2 \sum_{\epsilon', \epsilon''} \int d\mathbf{p}' d\mathbf{p}'' \mathcal{D}(\epsilon - \epsilon', \mathbf{p} - \mathbf{p}') G_b(\epsilon', \mathbf{p}') G_b(\epsilon', \mathbf{p}'') \mathcal{D}(\epsilon' - \epsilon'', \mathbf{p}' - \mathbf{p}'') G_b(\epsilon + \epsilon'' - \epsilon', \mathbf{p} + \mathbf{p}'' - \mathbf{p}') \approx$$

$$\sim \frac{\lambda^4}{\Gamma^{5/2}} \sum_{\epsilon', \epsilon''} \sqrt{(\epsilon - \epsilon')^2 + v_s^2} (i\epsilon' + \mu)^{3/2} \sqrt{(\epsilon' - \epsilon'')^2 + v_s^2} \sqrt{\epsilon'' + \mu}.$$
Integral over \( p'_n \) has relevant momenta \( \sim 1/\xi \). We can neglect the dependence of \( \Omega \) on \( p'_n \) provided that \( |\Omega(\varepsilon_n - \varepsilon'_n, 1/\xi, 0) - \Omega(\varepsilon_n - \varepsilon'_n, 0, 0)| / \nu_s^2 \ll 1/\xi^2 \) (see calculation for \( \Omega \) below). For \( f \) the condition is \( |f(\varepsilon'_n, 1/\xi, 0) - f(\varepsilon'_n, 0, 0)| \ll \beta/\xi^2 + \mu \). Using

\[
\int dx \frac{1}{\sqrt{x^2 + a(x^2 + b)}} = \frac{\text{arctanh}(\frac{\sqrt{b-a}}{\sqrt{b}})}{\sqrt{b} \sqrt{b-a}}
\]

We then obtain:

\[
f^a(\varepsilon_n) - \varepsilon_n = \frac{3\Lambda^2 T}{2\pi \sqrt{\alpha \beta}} \frac{\text{sgn}[\text{Re}f^b(\varepsilon'_n)]}{\sqrt{\Omega(\varepsilon_n - \varepsilon'_n)/\nu_s^2 + \frac{1}{\xi^2}}} \frac{\text{arctanh} \left\{ \frac{\sqrt{\Omega(\varepsilon_n - \varepsilon'_n)/\nu_s^2 + \frac{1}{\xi^2}}}{\Omega(\varepsilon_n - \varepsilon'_n)/\nu_s^2 + \frac{1}{\xi^2} - (if^b(\varepsilon'_n) + \mu)/\beta} \right\}}{\Omega(\varepsilon_n - \varepsilon'_n)/\nu_s^2 + \frac{1}{\xi^2}}
\]

Now we turn to calculation of \( \Pi(\omega_n, p) = (\Omega(\omega_n, p) - \omega_n^2)/\nu_s^2 \):

\[
\Pi(\omega_n, p) = 2\Lambda^2 T \sum_{\varepsilon_n + p} G^1(\varepsilon_n + \omega_n, p + q) G^2(\varepsilon_n, p) + G^2(\varepsilon_n + \omega_n, p + q) G^1(\varepsilon_n, p).
\]

Let us consider the momentum integral in the first term:

\[
\int \frac{d^2p}{(2\pi)^2} \frac{1}{if^1(\varepsilon_n + \omega_n) - \alpha(p_1 + q_1)^2 + \beta(p_2 + q_2)^2 + \mu if^2(\varepsilon_n) - \alpha p_1^2 + \beta p_2^2 + \mu^2} = \frac{1}{\alpha + \beta \left[ (\alpha - \beta)p_1^2 - \frac{\alpha + \beta}{\alpha - \beta} \right] [\alpha z_+ + \beta z + (\beta^2 - \alpha^2)p_1^2] + \frac{\alpha}{\alpha + \beta} \left[ (\alpha - \beta)p_2^2 - \frac{\alpha + \beta}{\alpha - \beta} \right] [\alpha z_+ + \beta z + (\beta^2 - \alpha^2)p_2^2]} = \frac{\alpha}{\alpha + \beta} \left[ (\alpha - \beta)p_1^2 - \frac{\alpha + \beta}{\alpha - \beta} \right] [\alpha z_+ + \beta z + (\beta^2 - \alpha^2)p_1^2] + \frac{\alpha}{\alpha + \beta} \left[ (\alpha - \beta)p_2^2 - \frac{\alpha + \beta}{\alpha - \beta} \right] [\alpha z_+ + \beta z + (\beta^2 - \alpha^2)p_2^2]
\]

where \( z_+ = if_{\varepsilon + \omega} + \mu \), \( z = if_{\varepsilon + \omega} + \mu \). The result of the integral is:

\[
\lambda^2 T \sum_{\varepsilon} \frac{4}{\pi^2} \text{arctanh} \left\{ \frac{\sqrt{\Omega(\varepsilon + \omega)/\nu_s^2 + \frac{1}{\xi^2}}}{\sqrt{\Omega(\varepsilon + \omega)/\nu_s^2 + \frac{1}{\xi^2} - \beta(\alpha + \beta)/(\alpha + \beta)} \sqrt{\Omega(\varepsilon + \omega)/\nu_s^2 + \frac{1}{\xi^2} - \beta(\alpha + \beta)/(\alpha + \beta)} \sqrt{\Omega(\varepsilon + \omega)/\nu_s^2 + \frac{1}{\xi^2} - \beta(\alpha + \beta)/(\alpha + \beta)}} \right\}
\]

It can be seen that for \( \beta \to 0 \), \( \beta \Lambda^2 \to 0 \), \( \alpha \Lambda^2 \to \infty \) one recovers our previous result\(^{85}\). Note that taking the limit \( \Lambda \to \infty \) before \( \beta \to 0 \) leads to a different answer. However, assuming \( \Lambda \gg 1/\xi \) in the region \( \beta \sim \alpha \) we can simplify
For evaluating the momentum integral under the same assumptions as for the fermion self-energy we obtain:

\[ \Pi(\omega_n \neq 0, p) = \lambda^2 T \sum_{\epsilon} -\frac{2i}{\pi} \frac{\text{sgn}[\text{Re}(\epsilon f_{\epsilon +} + \beta f_{\epsilon}]) \arctanh \left\{ \sqrt{\frac{2}{3} \sqrt{\frac{i(\epsilon f_{\epsilon +} + \beta f_{\epsilon}) + (\alpha + \beta)\mu}{i(\epsilon f_{\epsilon +} + \beta f_{\epsilon}) + (\alpha + \beta)\mu}} \right\}}{\sqrt{i(\epsilon f_{\epsilon +} + \beta f_{\epsilon}) + (\alpha + \beta)\mu}} \]

\[ -\frac{2i}{\pi} \frac{\text{sgn}[\text{Re}(\beta f_{\epsilon +} + \alpha f_{\epsilon})] \arctanh \left\{ \sqrt{\frac{2}{3} \sqrt{\frac{i(\beta f_{\epsilon +} + \alpha f_{\epsilon}) + (\alpha + \beta)\mu}{i(\beta f_{\epsilon +} + \alpha f_{\epsilon}) + (\alpha + \beta)\mu}} \right\}}{\sqrt{i(\beta f_{\epsilon +} + \alpha f_{\epsilon}) + (\alpha + \beta)\mu}} \]

\[ + \frac{\text{sgn}[\text{Re}(\beta f_{\epsilon +} + \alpha f_{\epsilon})] \text{sgn}[\text{Re}(\beta f_{\epsilon +} + \alpha f_{\epsilon})]}{\sqrt{i(\alpha f_{\epsilon +} + \beta f_{\epsilon}) + (\alpha + \beta)\mu}} \]

\[ \Pi(0, p) = \lambda^2 T \sum_{\epsilon} -\frac{4i\arctanh(\sqrt{\beta/\alpha}/\pi\text{sgn}[\text{Re}(f_{\epsilon})] - 1)}{(\alpha + \beta)(i f_{\epsilon} + \mu)}. \]

As the main objective of the current work is to study the effects of finite \( \beta \) this expression has been used for numerical calculations. One notes however that \( \Pi(0, p) \) is logarithmically divergent. In what follows we absorb this divergence into the value of \( 1/\xi^2 \) by subtracting \( \Pi(0, p) \) from \( \Pi(\omega_n \neq 0, p) \).

Let us now derive the self-consistency equations for the competing order parameters. For the Pomeranchuk instability order parameter \( P(\varepsilon_n) \) one has \( (P(\varepsilon_n) \equiv P(\varepsilon_n) = -P_2(\varepsilon_n)) \):

\[ P(\varepsilon_n) = -3\lambda^2 T \sum_{\varepsilon_n} \int \frac{dp'_d dp'_b}{(2\pi)^2} \frac{1}{\Omega(\varepsilon_n - \varepsilon_n' + \mu) + (p - p')^2 + \frac{i}{\varepsilon}} \left\{ \frac{\text{sgn}[\text{Re}(f_{\epsilon'})] \arctanh \left\{ \sqrt{\frac{i f_{\epsilon'} + \mu - \beta(\Omega_{\varepsilon - \epsilon'}/v_z^2 + 1/\xi^2)}{\beta(\Omega_{\varepsilon - \epsilon'}/v_z^2 + 1/\xi^2)} \right\}}{\sqrt{\Omega_{\varepsilon - \epsilon'} + 1/\xi^2(i f_{\epsilon'} + \mu - \beta(\Omega_{\varepsilon - \epsilon'}/v_z^2 + 1/\xi^2))}} \right\}. \]

Evaluating the momentum integral under the same assumptions as for the fermion self-energy we obtain:

\[ P(\varepsilon_n) = 3i\lambda^2 T \sum_{\varepsilon_n} \frac{P(\varepsilon_n') \text{sgn}[\text{Re}(f_{\epsilon'})]}{4\pi \sqrt{\alpha}} \left\{ \arctanh \left\{ \sqrt{\frac{i f_{\epsilon'} + \mu - \beta(\Omega_{\varepsilon - \epsilon'}/v_z^2 + 1/\xi^2)}{\beta(\Omega_{\varepsilon - \epsilon'}/v_z^2 + 1/\xi^2)} \right\} \right\}. \]

where we used:

\[ \int_{-\infty}^{\infty} dx \frac{1}{(x^2 + a^2)^{3/2}} = -\frac{2\arctan\left\{ \sqrt{\frac{b^2 - x^2}{x}} \right\}}{b(a^2 - b^2)^{3/2}}. \]

For \( \beta \ll \alpha \) we can also obtain the equation for CDW. In this case it is clear that \( Q_{CDW} = (Q, Q) \) and the equation for the CDW order parameter \( C(\varepsilon_n) \) is:

\[ C(\varepsilon_n) = -3\lambda^2 T \sum_{\varepsilon_n} \int \frac{dp'_d dp'_b}{(2\pi)^2} \frac{1}{\Omega(\varepsilon_n - \varepsilon_n' + \mu) + (p - p')^2 + \frac{i}{\varepsilon}} \left\{ \frac{\text{sgn}[\text{Re}(f_{\epsilon'})] \arctanh \left\{ \sqrt{\frac{i f_{\epsilon'} + \mu - \beta(\Omega_{\varepsilon - \epsilon'}/v_z^2 + 1/\xi^2)}{\beta(\Omega_{\varepsilon - \epsilon'}/v_z^2 + 1/\xi^2)} \right\}}{\sqrt{\Omega_{\varepsilon - \epsilon'} + 1/\xi^2(i f_{\epsilon'} + \mu - \beta(\Omega_{\varepsilon - \epsilon'}/v_z^2 + 1/\xi^2))}} \right\}. \]

Evaluating the momentum integral under the same assumptions as for the fermion self-energy we obtain:

\[ C(\varepsilon_n) = 3i\lambda^2 T \sum_{\varepsilon_n} \frac{C(\varepsilon_n') \text{sgn}[\text{Re}(f_{\epsilon'})]}{4\pi \sqrt{\alpha}} \left\{ \arctanh \left\{ \sqrt{\frac{i f_{\epsilon'} + \mu - \beta(\Omega_{\varepsilon - \epsilon'}/v_z^2 + 1/\xi^2)}{\beta(\Omega_{\varepsilon - \epsilon'}/v_z^2 + 1/\xi^2)} \right\} \right\}. \]
where we have used:

\[
\int dx \frac{1}{x^2 + c^2 x^2 - a^2 \sqrt{x^2 + b^2}} = \\
\arctan \left( \frac{\sqrt{a^2 + c^2}}{c(a^2 + c^2)\sqrt{b^2 - c^2}} \right) - \frac{1}{c(a^2 + c^2)\sqrt{a^2 + b^2}} \arctanh \left( \frac{\sqrt{a^2 + b^2}}{a(a^2 + c^2)\sqrt{a^2 + b^2}} \right)
\]

For DDW the equations take form (we use \(D(\varepsilon_n) \equiv D^1(\varepsilon_n) = -D^2(\varepsilon_n)\)):

\[
D(\varepsilon_n) = -3\lambda^2 T \sum_{\varepsilon_n'} \int \frac{dp'dp}{(2\pi)^2} \frac{1}{\Omega(\varepsilon_n - \varepsilon_n')/v_s^2 + (p - p')^2 + \frac{1}{\xi^2}} \Omega(\varepsilon_n - \varepsilon_n')/v_s^2 + (p - p')^2 + \frac{1}{\xi^2}
\]

(C3)

First we rewrite the integrand in a similar way to the polarization operator \(\Pi(q, \omega_n)\):

\[
\frac{1}{(i\varepsilon - \alpha p_1^2 + \beta p_2^2 + \mu)(i\varepsilon - \alpha p_1^2 + \beta p_2^2 + \mu)} = \\
\frac{1}{(\alpha + \beta)(i\varepsilon - \alpha p_1^2 + \beta p_2^2 + \mu)} \left[ \frac{\beta}{\alpha + \beta} \frac{\beta p_2^2 - \alpha p_1^2 + i\varepsilon - \alpha p_1^2 + \beta p_2^2 + \mu}{\alpha p_1^2 - \beta p_2^2 + i\varepsilon - \alpha p_1^2 + \beta p_2^2 + \mu} \right]
\]

(C4)

If \(\alpha - \beta \approx \alpha\) we can neglect the dependence of the bosonic propagator on \(p_1(p_2)\) for the first (second) term in the integral and we can neglect the momenta in the bosonic propagator altogether for the third term. We obtain as an intermediate result:

\[
D(\varepsilon_n) = \frac{3\lambda^2 T}{\alpha + \beta} \sum_{\varepsilon_n'} iD(\varepsilon_n') \frac{1}{\pi} \frac{\arctanh(\sqrt{\beta}/\alpha) - i\pi \sgn[\text{Ref}e\varepsilon']/2}{i\varepsilon - \alpha p_1^2 + \beta p_2^2 + \mu} \left\{ \frac{\Omega_{\varepsilon - \varepsilon'}/v_s^2 + 1/\xi^2 + (i\varepsilon - \alpha p_1^2 + \beta p_2^2 + \mu)}{(\alpha - \beta)[(\alpha - \beta)p_2^2 - (i\varepsilon - \alpha p_1^2 + \beta p_2^2 + \mu)]} \right\}
\]

(C5)

We use our assumptions \(\alpha - \beta \approx \alpha\) and \(|\sqrt{(i\mu + \mu)}/\alpha| \ll |\Omega + 1/\xi^2|\) to further simplify the answer:

\[
D(\varepsilon_n) = \frac{3\lambda^2 T}{\alpha + \beta} \sum_{\varepsilon_n'} \frac{D(\varepsilon_n') \frac{\sgn[\text{Ref}e\varepsilon']\arctanh(\sqrt{\beta}/\alpha)/\pi + 1/4}{i\varepsilon - \alpha p_1^2 + \beta p_2^2 + \mu}}{\Omega_{\varepsilon - \varepsilon'}/v_s^2 + 1/\xi^2}
\]

The result is reminiscent of the expression for \(\chi_{DDW}\) in the toy model. To study qualitatively the presence of IDDDW within spin-fermion model we use the following equation that can be easily derived assuming \(\beta/\xi^2 \gg |i\mu + \mu|\) and incommensurability \((Q, Q)\) along the diagonal using the result (A3):

\[
D_I(\varepsilon_n) = \frac{3\lambda^2 T}{\alpha + \beta} \sum_{\varepsilon_n'} \frac{D_I(\varepsilon_n') \frac{\sgn[\text{Ref}e\varepsilon']\arctanh(\sqrt{\beta}/\alpha)/\pi + 1/4}{i\varepsilon - \alpha p_1^2 + \beta p_2^2 + \mu}}{\Omega_{\varepsilon - \varepsilon'}/v_s^2 + 1/\xi^2}
\]

\[
- \frac{\sgn[\text{Ref}e\varepsilon']/\pi}{i\varepsilon - \alpha p_1^2 + \beta p_2^2 + \mu} \left[ \frac{\alpha \beta}{\alpha + \beta} \frac{\alpha \beta Q^2}{\alpha + \beta} \arctanh \left( \frac{\alpha \beta Q^2}{\alpha + \beta} \right) \right].
\]
Introducing an energy scale $\Gamma = \sqrt{\lambda^2 v^2 / \alpha}$ (note that in our previous work a larger scale $(\lambda^2 v / \sqrt{\alpha})^{2/3}$ has been used) we can bring the equations to a dimensionless form:

$$P(\varepsilon) = 3iT \sum_{\varepsilon''} \frac{P(\varepsilon') \text{sgn}[\text{Re} \varepsilon']}{4\pi} \begin{cases} v^2_\varepsilon / \alpha \cdot \text{arctanh} \left( \frac{\{(if' - \mu)v^2_\varepsilon/\alpha - \beta/\alpha(\Omega_{\varepsilon''} + a)\}}{\sqrt{\Omega_{\varepsilon''} + a}[(if' - \mu)v^2_\varepsilon/\alpha - \beta/\alpha(\Omega_{\varepsilon''} + a)]} \right) & \text{for } \text{sgn} \left( \text{Re} \varepsilon' \right) = 1, \\ v^2_\varepsilon / \alpha \cdot \text{arctanh} \left( \frac{\{(if' - \mu)v^2_\varepsilon/\alpha - \beta/\alpha(\Omega_{\varepsilon''} + a)\}}{\sqrt{\Omega_{\varepsilon''} + a}[(if' - \mu)v^2_\varepsilon/\alpha - \beta/\alpha(\Omega_{\varepsilon''} + a)]} \right) & \text{for } \text{sgn} \left( \text{Re} \varepsilon' \right) = -1 \end{cases},$$

$$C(\varepsilon) = \frac{3iT}{4\pi} \sum_{\varepsilon''} C(\varepsilon') \text{sgn}[\text{Re} \varepsilon'] \begin{cases} \sqrt{\alpha Q^2 / 4 - if' - \mu} / \sqrt{\alpha Q^2 / 4 - (if' - \mu)v^2_\varepsilon / \alpha - (\Omega_{\varepsilon''} + a)\beta / \alpha} & \text{for } \text{sgn} \left( \text{Re} \varepsilon' \right) = 1, \\ \sqrt{\alpha Q^2 / 4 - if' - \mu} / \sqrt{\alpha Q^2 / 4 - if' - \mu} & \text{for } \text{sgn} \left( \text{Re} \varepsilon' \right) = -1 \end{cases},$$

$$D(\varepsilon) = \frac{0.75T}{1 + \beta / \alpha} \sum_{\varepsilon''} D(\varepsilon') \frac{4i\text{sgn}[\text{Re} \varepsilon'] \text{arctanh}(\sqrt{\beta / \alpha}) / \pi + 1}{if' + \mu},$$

$$D_1(\varepsilon) = \frac{0.75T}{1 + \beta / \alpha} \sum_{\varepsilon''} D_1(\varepsilon') \frac{4i\text{sgn}[\text{Re} \varepsilon'] \text{arctanh}(\sqrt{\beta / \alpha}) / \pi + 1}{if' + \mu + \frac{\beta Q^2}{(\alpha + \beta)^2}},$$

and

$$D(\varepsilon) = \frac{0.75T}{1 + \beta / \alpha} \sum_{\varepsilon''} D_1(\varepsilon') \frac{4i\text{sgn}[\text{Re} \varepsilon'] \text{arctanh}(\sqrt{\beta / \alpha}) / \pi + 1}{if' + \mu + \frac{\beta Q^2}{(\alpha + \beta)^2}}.$$

(C6)

The equations have been solved numerically by an iteration method with nonlinearities $1 / 10 |D(\varepsilon')|^2 + 1, 1 / (10 |P(\varepsilon')|^2 + 1)$ introduced to r.h.s. of the order parameter equations to enforce convergence below critical temperature. The values of critical temperatures are not affected by this procedure. The number of Matsubara frequencies taken has been $300 + 1 / (\pi T)$, but not larger then 800.

While solving equations numerically two obstacles were encountered. First, the equations contain nonanalytic functions of a complex argument. To exclude ambiguity, we exclude the half-axis $\text{Re} \varepsilon < 0, \text{Im} \varepsilon = 0$ and check that arguments never cross it. In practice this means that the square roots should be evaluated from combinations like $if + \mu$ which always do have an imaginary part and never cross $\text{Re} \varepsilon < 0, \text{Im} \varepsilon = 0$ as functions of $\varepsilon$ but not of $(if_1 + \mu)$ or $(if + \mu)$ as these can cross the negative axis as functions of $\varepsilon$.

The second obstacle is that at low $\mu$ spurious solutions for $D$ appear. They don’t converge even for large numbers of iterations. However the convergence can be greatly improved by the following trick, which is a simplified version of Newton’s method. For $T > T_{Pom,DDW}$ we have an equation:

$$\vec{X} = A_{\mu} \vec{X},$$

and the Newton’s method looks:

$$\vec{X}_{n+1} = \vec{X}_n + (1 - A)^{-1}(A\vec{X}_n - \vec{X}_n).$$

Evaluating the matrix $(1 - A)^{-1}$ is a rather slow operation. However in our case $A_{\mu} \sim (\Omega_{\mu} + a)$ and thus diagonal elements dominate. We can approximately use then:

$$\vec{X}_{n+1} = \vec{X}_n + \text{diag}(1 - A_{\mu})^{-1}(A\vec{X}_n - \vec{X}_n).$$

This allows us to get rid of the spurious solutions and improve convergence to obtain consistent $T_{Pom/DDW}$ values.

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