

Coulomb interaction driven instabilities of sliding Luttinger liquids

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We study systems made of periodic arrays of one-dimensional quantum wires coupled by Coulomb interaction. Using bosonization an interacting metallic fixed point is obtained, which is shown to be a higher-dimensional analog of the Tomonaga-Luttinger liquid, or a sliding Luttinger liquid. This non-Fermi liquid metallic state, however, is unstable in the presence of weak interwire backscatterings, which favor charge density wave states and suppress pairing. Depending on the effective strength of the Coulomb repulsion and the size of interwire spacing, various charge density wave states are stabilized, including Wigner crystal states. Our method allows for the determination of the specific ordering patterns and corresponding energy and temperature scales.

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I. INTRODUCTION

Tomonaga-Luttinger liquid (or simply Luttinger liquid, abbreviated as LL) is the generic metallic state realized in interacting one-dimensional systems [1–4]. Phase space constraints and nesting enable weak short-range interactions to destroy quasiparticle coherence while preserving metallicity, making LL the earliest example of a metal that is not described by Fermi liquid theory. Thus LLs have long motivated theoretical modeling of non-Fermi liquid states in higher dimensions [5–8]. Indeed, within a coupled-wire construction, it was shown that a higher-dimensional analog of LL, the sliding Luttinger liquid (SLL), can be stabilized above one dimension in the presence of short-range repulsive interactions and/or vanishing interwire hoppings [9–12]. Despite the coupling between wires, the SLL possesses an emergent “sliding” symmetry corresponding to independent translation invariance on each wire, whose nature will be made precise in Sec. II. The SLL is an anisotropic metal, which behaves like an LL along the wires, while transport is suppressed in the transverse direction(s).

In parallel to these, coupled LLs have also been used as a paradigm to study competing orders above one dimension [13,14]. In one dimension, strong quantum fluctuations prevent spontaneous breaking of continuous symmetries [15]. Instead, tendencies toward ordering in different channels manifest themselves in the power-law correlations of various local order parameters, with the power law exponents (or scaling dimensions) indicating the strength of the ordering tendency. Indeed, the *unstable* coupled LL fixed point can be used as the starting point of systematic analyses of the physics of ordering in coupled LLs, based on renormalization group (RG) arguments. In recent examples, the paradigm of wires coupled by short- [16] and long-ranged [17] interactions was applied to understand the physics of magnetic field driven catalysis in metals with low carrier density [18–20]. Quenching of the kinetic energy on the plane perpendicular to the applied field makes the metal susceptible to density wave ordering. Owing to a small or vanishing Fermi energy, the quantum limit is reached at moderate magnetic field strength, and the lowest Landau level dominates the low-energy physics. The degeneracy of the lowest Landau level is utilized to map the problem to that of coupled wires, where the number of wires is controlled by the degeneracy.

In this work, we revisit the problem of quantum wires coupled by Coulomb interaction in the regime where single-electron interwire hoppings are suppressed. The purpose of our work is the following. In reality, the interwire couplings that lead to SLL physics and those that lead to charge density wave (CDW) ordering come from the same Coulomb interaction. They should, therefore, be treated on equal-footing. We will demonstrate that such a treatment leads to specific predictions of the leading CDW instability and resultant ordering pattern, as well as corresponding energy scales.

The paper is organized as follows. In Sec. II, we introduce the model, and derive the bosonized action for a system of infinite number of quantum wires in d dimensions coupled by Coulomb interaction in the forward scattering channel. In Sec. III, we show that the action describes an SLL fixed point in d dimensions, and deduce the exponents that characterize the fixed point. In Sec. IV, we analyze the stability of the SLL fixed point against various symmetry breaking perturbations. Within a tree-level RG analysis we show that multiple charge density wave (CDW) states compete for dominance in the absence of a dimensional crossover, and interwire pairing instabilities are suppressed. We establish the zero (finite) temperature phase diagram as a function of Coulomb interaction strength (temperature) and interwire spacing. Finally, we close with a discussion of our results in Sec. V.

II. MODEL

We consider a $(d - 1)$ -dimensional lattice of identical wires of spinless fermions in d dimensions with the wires lying along the \hat{x} axis [21]. The wires are labeled by a $(d - 1)$ -dimensional vector, \mathbf{n} , such that $\mathbf{n} \cdot \hat{x} = 0$. The fermion field on the \mathbf{n} th wire is expressed in terms of the hydrodynamic modes as [4]

$$\psi_{\mathbf{n}}(\tau, x) \approx e^{-i\pi\rho_0 x} \psi_{L,\mathbf{n}}(\tau, x) + e^{i\pi\rho_0 x} \psi_{R,\mathbf{n}}(\tau, x), \quad (1)$$

where modes carrying momenta of magnitude larger than $\pi\rho_0$ are ignored. The left and right moving fermions are expressed as

$$\begin{aligned} \psi_{L,\mathbf{n}}(\tau, x) &= e^{i(\vartheta_{\mathbf{n}}(\tau, x) - \varphi_{\mathbf{n}}(\tau, x))} \sqrt{\rho_0 + \frac{1}{\pi} \partial_x \varphi_{\mathbf{n}}(\tau, x)}, \\ \psi_{R,\mathbf{n}}(\tau, x) &= e^{i(\vartheta_{\mathbf{n}}(\tau, x) + \varphi_{\mathbf{n}}(\tau, x))} \sqrt{\rho_0 + \frac{1}{\pi} \partial_x \varphi_{\mathbf{n}}(\tau, x)}. \end{aligned} \quad (2)$$

Here, ρ_0 is the mean density, $\partial_x \varphi_{\mathbf{n}}$ is the local density modulation of fermions on each wire, and $\vartheta_{\mathbf{n}}$ is the phase of the fermion field. The action for free fermions is given by

$$S_0 = \frac{1}{2} \sum_{\mathbf{n}} \int_{(\tau,x)} \left[\frac{2i}{\pi} (\partial_x \varphi_{\mathbf{n}}) (\partial_\tau \vartheta_{\mathbf{n}}) + \frac{v_F}{\pi} (\partial_x \varphi_{\mathbf{n}})^2 + \frac{v_F}{\pi} (\partial_x \vartheta_{\mathbf{n}})^2 \right], \quad (3)$$

where $\int_x \equiv \int dx$ and $v_F = k_F/m$ with $k_F = \pi \rho_0$ being the Fermi wave vector, and m being the mass of the noninteracting fermions.

We introduce an instantaneous interaction among the fermions,

$$S_I = \frac{1}{2} \sum_{\mathbf{n}, \mathbf{m}} \int_{(\tau,x,x')} V(x-x', \mathbf{n}-\mathbf{m}) \rho_{\mathbf{n}}(\tau,x) \rho_{\mathbf{m}}(\tau,x'), \quad (4)$$

where $\rho_{\mathbf{n}}(\tau,x) = \rho_0 + \frac{1}{\pi} \partial_x \varphi_{\mathbf{n}}(\tau,x)$ is the density on the \mathbf{n} th wire. We assume the wires are uniformly spaced with lattice spacing a , and define the Fourier components through

$$\varphi_{\mathbf{n}}(\tau,x) = l_{\text{BZ}}^{d-1} \int_{-\infty}^{\infty} \frac{dk_0 dk_x}{(2\pi)^2} \times \int_{\text{BZ}} \frac{d^{d-1} \mathbf{K}}{(2\pi)^{d-1}} e^{i\tau k_0 + i x k_x + i \mathbf{a} \cdot \mathbf{K}} \varphi(k), \quad (5)$$

where BZ indicates the first Brillouin zone, l_{BZ}^{-1} is a measure of linear dimension of BZ such that its volume is $(\frac{2\pi}{l_{\text{BZ}}})^{d-1}$, and \mathbf{K} represents points inside BZ. Thus the bosonized action for the interacting theory is given by

$$S = \frac{l_{\text{BZ}}^{d-1}}{2\pi} \left(\frac{l_{\text{BZ}}}{a} \right)^{d-1} \int dk [2i k_0 k_x \varphi(-k) \vartheta(k) + V_\varphi(\vec{k}) k_x^2 \varphi(-k) \varphi(k) + v_F k_x^2 \vartheta(-k) \vartheta(k)], \quad (6)$$

where $dk = \frac{dk_0 dk_x d^{d-1} \mathbf{K}}{(2\pi)^{d+1}}$, $\vec{k} \equiv (k_x, \mathbf{K})$, and $V_\varphi(\vec{k}) = v_F + \pi^{-1} (l_{\text{BZ}}/a)^{d-1} V(\vec{k})$ with $V(\vec{k})$ being the Fourier conjugate of $V(x-x', \mathbf{n}-\mathbf{m})$. In $d=2$, the lattice geometry dependent ratio l_{BZ}/a equals 1, while in $d=3$ it equals 1 and $(\sqrt{3}/2)^{1/2}$ for the square and triangular lattices, respectively. From the coordinate space representation of Eq. (6), we deduce that the action is invariant under wire dependent shifts $\varphi_{\mathbf{n}}(\tau,x) \mapsto \varphi_{\mathbf{n}}(\tau,x) + \varphi_{\mathbf{n}}^{(0)}$ and $\vartheta_{\mathbf{n}}(\tau,x) \mapsto \vartheta_{\mathbf{n}}(\tau,x) + \vartheta_{\mathbf{n}}^{(0)}$, where $\varphi_{\mathbf{n}}^{(0)}$ and $\vartheta_{\mathbf{n}}^{(0)}$ are constants. The former invariance is the sliding symmetry advertised in Sec. I, and it corresponds to translation invariance along each wire. It allows the fermions on distinct wires to “slide” with respect to each other. The latter invariance corresponds to particle number conservation on each wire which prevents single-particle interwire hoppings. This results in two flat patches of Fermi surface that are nested by the wave vector $2k_F \hat{x}$. Generally, such extensive nesting makes the metallic state exceedingly susceptible to weak coupling instabilities. However, with a suitable choice of short-range interwire interactions, it is possible to stabilize this metallic state in $d > 1$ [9–12]. Although Eqs. (3) and (6) are both Gaussian actions, they describe a noninteracting and an interacting fixed point, respectively. We will elucidate this point further through the computation of scaling exponents in subsequent sections.

III. COULOMBIC SLIDING LUTTINGER LIQUID

In this section, we characterize the fixed point described by the action in Eq. (6). We consider both the interwire and intrawire interactions that arise from unscreened Coulomb interaction among fermions, $V(x-x', \mathbf{n}-\mathbf{m}) = \frac{e^2}{4\pi\epsilon \sqrt{(x-x')^2 + a^2 |\mathbf{n}-\mathbf{m}|^2}}$ with e being the electric charge of the fermions and ϵ being the permittivity, such that

$$V(\vec{k}) = \frac{2e^2}{4\pi\epsilon} \sum_{\mathbf{n}} e^{i \mathbf{a} \cdot \mathbf{n} \cdot \mathbf{K}} \mathcal{K}_0(a |\mathbf{n}| k_x), \quad (7)$$

where $\mathcal{K}_0(x)$ is the modified Bessel function of the second kind, and $\mathcal{K}_0(x) \sim -\ln x$ ($\mathcal{K}_0(x) \sim e^{-x}$) for $x \ll 1$ ($x \gg 1$) [22]. In the limit $a |\vec{k}| \ll 1$, the summand varies slowly as a function of \mathbf{n} . Therefore, in this limit, we replace the sum by an integral to obtain

$$V_\varphi(\vec{k}) \approx v_F + \frac{A'_d e^2}{4\pi\epsilon a^{d-1}} \frac{(l_{\text{BZ}}/a)^{d-1}}{(k_x^2 + |\mathbf{K}|^2)^{(d-1)/2}}, \quad (8)$$

where $A'_d > 0$, and $(A'_2, A'_3) = (2, 4)$.

In order to obtain the analog of Luttinger parameters, we project $V_\varphi(\vec{k})$ in the forward scattering channel along the wires by setting $k_x = 0$. Thus we consider

$$V_\varphi(\mathbf{K}) = v_F + \frac{A_d e^2}{4\pi\epsilon (a |\mathbf{K}|)^{(d-1)}}, \quad (9)$$

where $A_d = A'_d (l_{\text{BZ}}/a)^{d-1}$, as the effective Coulomb interaction for scatterings in the forward scattering channel. By utilizing the two component basis, $(\varphi, \vartheta)^T$, the propagators for $\vartheta(k)$ and $\varphi(k)$ are easily deduced from Eq. (6),

$$G_\vartheta(k) = \frac{\pi (a/l_{\text{BZ}})^{d-1} V_\varphi(\vec{k})}{k_0^2 + v_F V_\varphi(\mathbf{K}) k_x^2}, \quad G_\varphi(k) = \frac{\pi (a/l_{\text{BZ}})^{d-1} v_F}{k_0^2 + v_F V_\varphi(\mathbf{K}) k_x^2}. \quad (10)$$

Furthermore, they are correlated as

$$\langle \varphi(-k) \vartheta(k) \rangle = \frac{\pi (a/l_{\text{BZ}})^{d-1} k_0}{i k_x (k_0^2 + v_F V_\varphi(\mathbf{K}) k_x^2)}. \quad (11)$$

The equal-time correlation between the simplest vertex operators,

$$\langle e^{i \vartheta_{\mathbf{n}}(\tau,x)} e^{-i \vartheta_{\mathbf{m}}(\tau,x+\Delta x)} \rangle = \frac{\delta_{\mathbf{n},\mathbf{m}}}{|\lambda_0 \Delta x|^{2\eta_\vartheta(\mathfrak{h}_d)}}, \quad (12)$$

$$\langle e^{i \varphi_{\mathbf{n}}(\tau,x)} e^{-i \varphi_{\mathbf{m}}(\tau,x+\Delta x)} \rangle = \frac{\delta_{\mathbf{n},\mathbf{m}}}{|\lambda_0 \Delta x|^{2\eta_\varphi(\mathfrak{h}_d)}}, \quad (13)$$

where $\lambda_0^{-1} > k_F^{-1}$ is the short-distance cutoff along the wires, and

$$\mathfrak{h}_d = \frac{A_d e^2}{4\pi^d \epsilon v_F} \quad (14)$$

is the effective fine structure constant. We derive the exponents $\eta_\varphi(\mathfrak{h}_d)$ and $\eta_\vartheta(\mathfrak{h}_d)$ for various wire-stacking geometries in Appendix A. To illustrate the general features of these

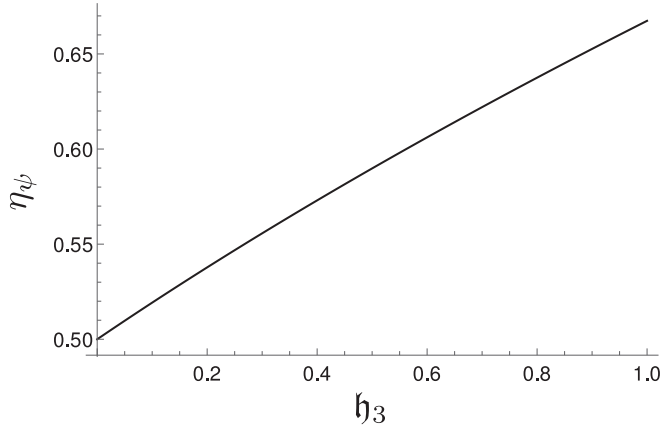


FIG. 1. The dependence of η_ψ on h_3 for the square lattice geometry in $d = 3$.

exponents here we quote the results for the square lattice,

$$\eta_\vartheta(h_d) = \frac{1}{4} \int_0^1 d^{d-1} w \left(1 + \frac{h_d}{|\mathbf{w}|^{d-1}} \right)^{\frac{1}{2}}, \quad (15)$$

$$\eta_\varphi(h_d) = \frac{1}{4} \int_0^1 d^{d-1} w \left(1 + \frac{h_d}{|\mathbf{w}|^{d-1}} \right)^{-\frac{1}{2}}, \quad (16)$$

where the integrations are over the unit $(d-1)$ -dimensional cube, and \mathbf{w} is a vector inside the cube. Since the integrand in Eq. (15) [Eq. (16)] is larger (smaller) than 1, $\eta_\vartheta(h_d) > \eta_\vartheta(0)$ [$\eta_\varphi(h_d) < \eta_\varphi(0)$] for any $h_d > 0$. Therefore intrawire phase (density) fluctuations are suppressed (enhanced) compared to that at the noninteracting fixed point as the effective strength of the Coulomb interaction increases.

The fermion propagator on the \mathbf{n} th wire is given by

$$\langle \psi_{\mathbf{n}}(0, x) \psi_{\mathbf{n}}^\dagger(0, 0) \rangle \sim \frac{\rho_0}{|\lambda_0 x|^{2\eta_\psi(h_d)}}, \quad (17)$$

where $\eta_\psi(h_d) = \eta_\vartheta(h_d) + \eta_\varphi(h_d)$, and we have ignored an overall phase factor arising from the correlation between φ and ϑ . In $d = 2$, we obtain $\eta_\psi(h_2) = \frac{1}{2} \sqrt{1 + h_2}$, while it is numerically computed in $d = 3$ and its behavior as a function of h_3 is shown in Fig. 1. In both cases, $\eta_\psi > \frac{1}{2}$ for $h_d > 0$. The faster decay of the fermion-fermion correlation compared to the noninteracting limit ($h_d = 0$), implies that Eq. (6) describes a Luttinger-liquid-like metallic state in $d > 1$. Indeed, this is an example of a sliding Luttinger liquid state. Due to the central role played by interfermion Coulomb repulsion, we refer to it as *Coulombic sliding Luttinger liquid* (CSLL). Unlike SLLs arising in systems with short-ranged interactions, the CSLL is controlled by a single parameter, h_d .

Due to the absence of interwire hoppings, the single-particle correlation functions are diagonal in the wire index. If this were true for all correlation functions, then the CSLL would be equivalent to a collection of noninteracting LLs, albeit with renormalized exponents. The distinction is easily demonstrated with the aid of the density-density correlation between two distinct wires,

$$\langle \rho_{\mathbf{n}}(0, 0) \rho_{\mathbf{m}}(0, 0) \rangle = \rho_0^2 - \left(\frac{\lambda_0}{2\pi} \right)^2 f(h_d, \mathbf{n} - \mathbf{m}), \quad (18)$$

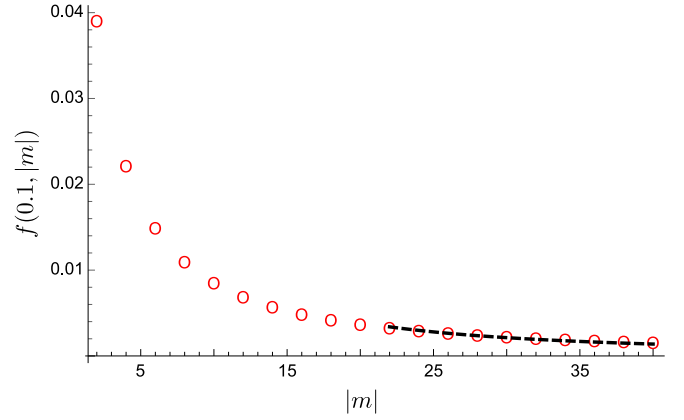


FIG. 2. Deviation of the interwire density-density correlation at the CSLL fixed point from that at the noninteracting fixed point, as a function of the separation between the wires. Here we have chosen $d = 2$ and $h_2 = 0.1$. The empty circles are numerically computed values of $f(0.1, |n - m|)$, and the dashed line is the asymptotic form, $f(h_2, |m|) \sim |m|^{-3/2}$.

where $\mathbf{n} \neq \mathbf{m}$ and the dimensionless function

$$f(h_d, \mathbf{n} - \mathbf{m}) = -\int_{BZ} \frac{d^{d-1} K}{(2\pi)^{d-1}} \frac{\cos(\mathbf{aK} \cdot (\mathbf{n} - \mathbf{m}))}{\sqrt{1 + h_d \left(\frac{\pi}{a|\mathbf{K}|} \right)^{d-1}}}. \quad (19)$$

We note that $f(h_d, \mathbf{m})$ depends on \mathbf{m} only through $|\mathbf{m}|$. Due to the cosine factor, the integration over $|\mathbf{K}|$ obtains dominant contribution from the region where $a|\mathbf{K}| \lesssim |\mathbf{n} - \mathbf{m}|^{-1}$. However, the denominator of the integrand suppresses it at small $|\mathbf{K}|$. Thus $f(h_d, |\mathbf{m}|)$ decreases as $|\mathbf{m}|$ increases, and $f(h_d, |\mathbf{m}|) \sim h_d^{-1/2} |\mathbf{m}|^{-3(d-1)/2}$ for $|\mathbf{m}| \gg 1$. This nontrivial interwire correlation between the densities is demonstrated in $d = 2$ with the aid of Fig. 2. Therefore the CSLL is distinct from a simple collection of LLs in d dimensions. While the low-energy mode disperses as $\omega = \sqrt{v_F V_\varphi(\mathbf{K})} |k_x|$ along the wires, the CSLL is a charge insulator in the transverse direction due to a lack of interwire single-particle hopping.

The physics of spinful electrons confined to one dimension, and interacting through a three-dimensional Coulomb potential was considered by Schulz in Ref. [23]. The unscreened tail of the Coulomb interaction was shown to lead to anomalous logarithmic dependencies of the correlation functions in Eqs. (12) and (13). We recover these anomalous logarithms by setting $d = 1 + \epsilon$, which leads to

$$\eta_\vartheta(h_{1+\epsilon}) = \frac{\sqrt{h_{1+\epsilon} + 1} + h_{1+\epsilon} \text{csch}^{-1}(\sqrt{h_{1+\epsilon}})}{4\epsilon}, \quad (20)$$

$$\eta_\varphi(h_{1+\epsilon}) = \frac{\sqrt{h_{1+\epsilon} + 1} - h_{1+\epsilon} \text{csch}^{-1}(\sqrt{h_{1+\epsilon}})}{4\epsilon}. \quad (21)$$

As $\epsilon \rightarrow 0$ these exponents diverge as ϵ^{-1} , indicating the presence of additional singularities in $d = 1$. In contrast to the qualitative modification of simple LL behavior by long-range Coulomb interaction in $d = 1$, the properties of the CSLL obtained here are qualitatively similar to those obtained in

SLLs with short-range interactions. This difference between $d > 1$ and $d = 1$ is attributed to interwire screening, which removes the singularities arising from the long-range tail of the unscreened Coulomb interaction through the enlarged phase space in the transverse direction [24]. These conclusions are further supported by diagrammatic computations within the parquet approximation [25]. In the earlier works on CSLL in Refs. [24–26], the authors focused on the renormalization of the intrawire interactions due to interwire Coulomb interaction, and showed that density wave (pairing) susceptibilities are enhanced (suppressed). Here we have fully characterized both intra- and interwire correlations. In the following section, we investigate the effects of interwire backscatterings at the CSLL fixed point.

IV. INSTABILITIES OF THE COULOMBIC SLIDING LUTTINGER LIQUID

Weak perturbations that do not break either the sliding symmetry or particle number conservation on each wire will not destabilize the CSLL state. However, even at low

energies, there exists processes, viz. interwire backscatterings and single-particle hoppings, that break either or both the above symmetries. Therefore it is necessary to consider the effect of these symmetry breaking perturbations on the CSLL fixed point in order to establish the true ground state of the system. The purpose of this section is to determine the parameter regime, if any, where the CSLL phase is stable, and identify the potential symmetry broken states where it is unstable.

Although the Coulomb interaction primarily contributes to the forward scattering channel due to the dominance of small-momentum exchange processes, it also mediates weak but nonvanishing $2k_F$ backscatterings. In general, these backscatterings can destabilize the CSLL phase by utilizing the extensive nesting between the two chiral segments of the Fermi surface in $d > 1$. The nesting can alter particle-particle and particle-hole pair hopping amplitudes between different wires. Further, interwire hoppings are always possible due to nonvanishing single-particle tunneling amplitude between wires. In order to investigate the stability of the CSLL against these destabilizing tendencies, we consider the effect of the following operators:

$$O_{\text{CDW}}^{(\mathbf{m})}(\tau, x, \mathbf{n}) = \frac{1}{2}[\psi_{L,\mathbf{n}}^\dagger(\tau, x)\psi_{R,\mathbf{n}}(\tau, x)\psi_{R,\mathbf{n}+\mathbf{m}}^\dagger(\tau, x)\psi_{L,\mathbf{n}+\mathbf{m}}(\tau, x) + \text{H.c.}] = \rho_0^2 \cos[2(\varphi_{\mathbf{n}}(\tau, x) - \varphi_{\mathbf{n}+\mathbf{m}}(\tau, x))], \quad (22)$$

$$O_{\text{SC}}^{(\mathbf{m})}(\tau, x, \mathbf{n}) = \frac{1}{2}[\psi_{L,\mathbf{n}}^\dagger(\tau, x)\psi_{R,\mathbf{n}}^\dagger(\tau, x)\psi_{R,\mathbf{n}+\mathbf{m}}(\tau, x)\psi_{L,\mathbf{n}+\mathbf{m}}(\tau, x) + \text{H.c.}] = \rho_0^2 \cos[2(\vartheta_{\mathbf{n}}(\tau, x) - \vartheta_{\mathbf{n}+\mathbf{m}}(\tau, x))], \quad (23)$$

$$O_{\text{sp}}^{(\mathbf{m})}(\tau, x, \mathbf{n}) = \frac{1}{4}[\psi_{L,\mathbf{n}}^\dagger(\tau, x)\psi_{L,\mathbf{n}+\mathbf{m}}(\tau, x) + \psi_{R,\mathbf{n}}^\dagger(\tau, x)\psi_{R,\mathbf{n}+\mathbf{m}}(\tau, x) + \text{H.c.}] \\ = \rho_0 \cos[\vartheta_{\mathbf{n}}(\tau, x) - \vartheta_{\mathbf{n}+\mathbf{m}}(\tau, x)] \cos[\varphi_{\mathbf{n}}(\tau, x) - \varphi_{\mathbf{n}+\mathbf{m}}(\tau, x)]. \quad (24)$$

An instability driven by $O_{\text{CDW}}^{(\mathbf{m})}$ ($O_{\text{SC}}^{(\mathbf{m})}$) leads to a CDW (superconducting) state which breaks the continuous sliding symmetry (particle number conservation on each wire) to a discrete symmetry. The strengthening of $O_{\text{sp}}^{(\mathbf{m})}$ enhances the energy scale for a crossover from the CSLL state to a d -dimensional Fermi liquid metal. Below the crossover scale both the aforementioned symmetries are broken. We note that the spinlessness of the fermions does not allow intrawire backscatterings, as a result of which $\mathbf{m} \neq 0$. Thus, for a fixed wire-stacking, geometry we obtain a set of operators parameterized by the label \mathbf{m} , which compete with each other.

The leading instability is identified by comparing the scaling dimensions of susceptibilities of various operators. To compute the scaling dimension of the susceptibility of operator $O_X^{(\mathbf{m})}$, we perturb the action, Eq. (6), with the vertex

$$S_X^{(\mathbf{m})} = g_X^{(\mathbf{m})} \sum_{\mathbf{n}} \int d\tau dx O_X^{(\mathbf{m})}(\tau, x, \mathbf{n}), \quad (25)$$

and with the help of the equal time correlation on the \mathbf{n} th wire, $\langle O_X^{(\mathbf{m})}(\tau, x, \mathbf{n}) O_X^{(\mathbf{m})}(\tau, x + \Delta x, \mathbf{n}) \rangle$, we obtain the scaling exponent of $O_X^{(\mathbf{m})}$. The scale invariance of $S_X^{(\mathbf{m})}$ fixes the scaling dimension of the coupling $g_X^{(\mathbf{m})}$, $[g_X^{(\mathbf{m})}] = 2 - [O_X^{(\mathbf{m})}]$. Here the scaling dimension $[\mathcal{V}]$ of an operator \mathcal{V} is defined through the relationship $\mathcal{V}(\lambda) = \mathcal{V}(\lambda_0)e^{[\mathcal{V}]\ell}$, where $\ell = \ln(\lambda_0/\lambda)$ is the RG time/distance, and $\lambda < \lambda_0$ is the running momentum scale. The asymptotic behavior of the equal-time correlation function of

$O_X^{(\mathbf{m})}$ on the \mathbf{n} th wire at the CSLL fixed point is computed in Appendix B, and it takes the form

$$\langle O_X^{(\mathbf{m})}(0, \Delta x, \mathbf{n}) O_X^{(\mathbf{m})}(0, 0, \mathbf{n}) \rangle = \frac{2\rho_0^4}{|\lambda_0 \Delta x|^{2\eta_X(\mathbf{h}_d, \mathbf{m})}}. \quad (26)$$

On coarse-graining, $g_X^{(\mathbf{m})}$ evolves as

$$\partial_\ell \tilde{g}_X^{(\mathbf{m})}(\ell) = (2 - \eta_X(\mathbf{h}_d, \mathbf{m})) \tilde{g}_X^{(\mathbf{m})}(\ell), \quad (27)$$

where $\tilde{g}_X^{(\mathbf{m})}(\ell) \equiv \tilde{g}_X^{(\mathbf{m})}(\lambda) = g_X^{(\mathbf{m})} \lambda^{-(2-\eta_X(\mathbf{h}_d, \mathbf{m}))}$ is the corresponding dimensionless coupling. This implies that the operator $O_X^{(\mathbf{m})}$ is a relevant (irrelevant) perturbation at the CSLL fixed point if $\eta_X(\mathbf{h}_d, \mathbf{m}) < 2$ [$\eta_X(\mathbf{h}_d, \mathbf{m}) > 2$].

We note that, within a weak-coupling framework, generally, symmetry breaking in two and higher-dimensional metals is driven by marginal operators. Consequently, the sign and magnitude of quantum corrections determine the dominant instability. In the present case, however, we are able to fully account for interactions along the wires, which leads to the interwire operators picking up nontrivial scaling dimensions at tree-level as demonstrated in Eq. (27). Thus, approaching from weak-coupling side, the dominant instability is determined by the coupling which reaches a order of 1 value quickest in terms of the RG time, ℓ . This also leads to a non-BCS form of various energy scales associated with symmetry breaking transitions in the CSLL metal.

A. Charge density wave instabilities

As noted in Sec. III, the repulsive nature of the Coulomb interaction suppresses (enhances) fluctuations in the particle-particle (particle-hole) channel. This renders the Josephson (SC) couplings irrelevant at the CSLL fixed point, while the CDW couplings become relevant. In this subsection we consider the effect of the CDW couplings.

As a representative example, let us consider the wires stacked into a $(d-1)$ -dimensional square lattice. The scaling dimension of $O_{\text{CDW}}^{(\mathbf{m})}$ is

$$\eta_{\text{CDW}}(h_d, \mathbf{m}) = 2^{2-d} \int_{-1}^1 d^{d-1} w \frac{1 - \cos(\pi \mathbf{m} \cdot \mathbf{w})}{\sqrt{1 + \frac{h_d}{|\mathbf{w}|^{d-1}}}}. \quad (28)$$

With the help of Eq. (19), we obtain $\eta_{\text{CDW}}(h_d, \mathbf{m}) = 2 \int_0^1 d^{d-1} w (1 + h_d |\mathbf{w}|^{d-1})^{-1/2} + 2f(h_d, \mathbf{m})$. Since the first term is independent of \mathbf{m} , η_{CDW} follows the trend of $f(h_d, \mathbf{m})$ as $|\mathbf{m}|$ is tuned. Thus η_{CDW} decreases as $|\mathbf{m}|$ increases. When $h_d > 0$, in the region where $|\mathbf{w}|^{d-1} \ll h_d$ the integrand is suppressed by a factor of $\sqrt{|\mathbf{w}|^{d-1}}$ compared to the noninteracting fixed point. Thus, at the CSLL fixed point $\eta_{\text{CDW}} < 2$ for any $h_d, |\mathbf{m}| > 0$, and the fixed point is *always* unstable to the formation of a CDW in any $d \geq 2$. We note that, while in $d = 2$, η_{CDW} is completely determined by h_d and \mathbf{m} , in $d > 2$, η_{CDW} also depends on the wire-stacking geometry through the cosine term in the numerator of Eq. (28). In Appendix B, we demonstrate the stacking geometry dependence by comparing the results for square and triangular lattices.

In order to identify the leading CDW instability as a function of effective fine structure constant, h_d , and lattice geometry, we assume that the coupling $g_X^{(\mathbf{m})}$ results from $2k_F$ backscatterings mediated by Coulomb repulsion, which implies

$$\begin{aligned} \tilde{g}_{\text{CDW}}^{(\mathbf{m})}(h_d, \Upsilon; \lambda_0) &= \frac{e^2}{4\pi\epsilon v_F} \int dx \frac{e^{i2k_F x}}{\sqrt{x^2 + a^2|\mathbf{m}|^2}} \\ &\approx \frac{2\pi^{d-1}}{A_d} h_d e^{-\Upsilon|\mathbf{m}|}, \end{aligned} \quad (29)$$

where $\Upsilon = 2k_F a$. We have also assumed that the density of fermions on a given wire is larger than the density of wires, which implies $\Upsilon > 1$. Equations (27) and (28) imply that at fixed h_d the scaling dimension of $\tilde{g}_{\text{CDW}}^{(\mathbf{m})}$ increases with $|\mathbf{m}|$, which suggests that the CDW operator with the largest allowed $|\mathbf{m}|$ drives the dominant instability. However, the CDW gap, whose magnitude determines the depth of the free energy minimum, depends on both the scaling dimension and the strength of interwire backscattering. In particular, the gap is proportional to $\tilde{g}_{\text{CDW}}^{(\mathbf{m})}(h_d, \Upsilon; \lambda_0)$, which appears to favor the CDW with $|\mathbf{m}| = 1$. These opposing tendencies generically lead to a CDW state with a wave vector whose magnitude lies in between the largest and smallest allowed transverse momenta. The leading CDW instability at fixed h_d , Υ , and wire-stacking geometry is the one that *minimizes* the ratio

$$\frac{\lambda_0}{\lambda_{\text{CDW}}^{(\mathbf{m})}(h_d, \Upsilon)} = \left(\frac{1}{\tilde{g}_{\text{CDW}}^{(\mathbf{m})}(h_d, \Upsilon; \lambda_0)} \right)^{1/(2-\eta_{\text{CDW}}(h_d, \mathbf{m}))}, \quad (30)$$

where $\lambda_{\text{CDW}}^{(\mathbf{m})}$ is such that $\tilde{g}_{\text{CDW}}^{(\mathbf{m})}(h_d, \Upsilon; \lambda_{\text{CDW}}^{(\mathbf{m})}) \sim 1$. The algebraic dependence of $\lambda_{\text{CDW}}^{(\mathbf{m})}$ on the interwire coupling, $\tilde{g}_{\text{CDW}}^{(\mathbf{m})}$, is a nonperturbative effect that results from the inclusion of *all* intrawire interactions. This relationship, however, is subject to the bare couplings $\tilde{g}_{\text{CDW}}^{(\mathbf{m})}(h_d, \Upsilon; \lambda_0)$ being small. The resultant $T = 0$ phase diagrams in $d = 2$ and 3 are shown in Fig. 3. We note that for lattices with C_n point group symmetry, CDW states with wave vectors related by the C_n symmetry are degenerate.

From the phase diagrams, we deduce that CDW states with larger wave vectors are favored at stronger interaction strengths and larger interwire spacings. MacDonald and Fisher identified the CDW state formed between adjacent wires in $d = 2$ as a Wigner crystal [27]. However, as shown in Fig. 3, the possible symmetry broken states extend beyond Wigner crystals, as CDW states with smaller wave vectors and multiple sites per unit cell (in the transverse lattice) are stabilized through the competition between bare coupling strengths and scaling dimensions of interwire CDW operators. While the leading instability fixes the ordering wave vector

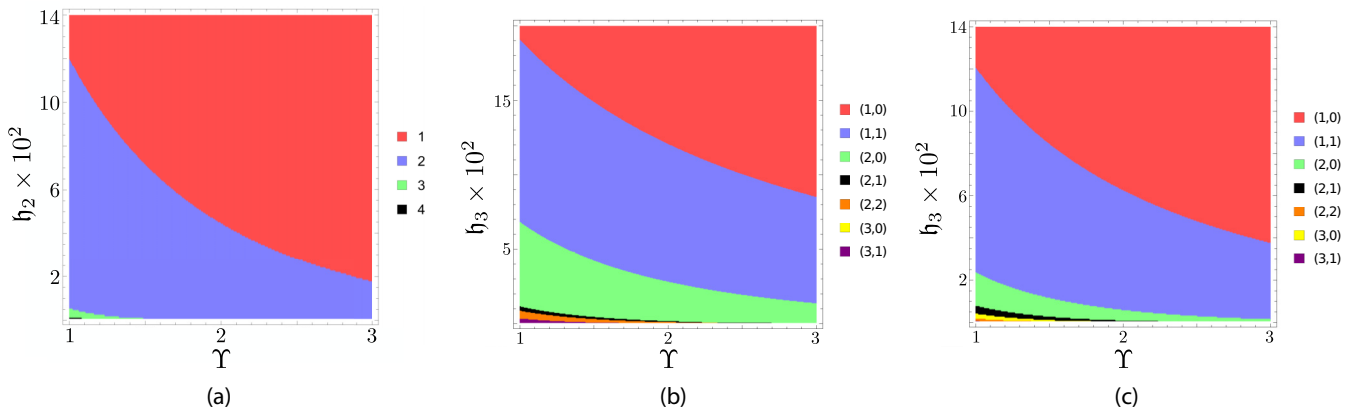


FIG. 3. Zero-temperature phase diagrams in (a) $d = 2$, and $d = 3$ with wires stacked in (b) square lattice and (c) triangular lattice geometry. Here, h_d is the effective fine structure constant, and $\Upsilon = 2k_F a$. The colors represent distinct CDW states which become dominant as (h_d, Υ) are tuned. The legends on the right indicate the direction and periodicity of the ordering vector in the dominant CDW state. We note that directions related by the C_n point group symmetry of the lattice in $d = 3$ are degenerate.

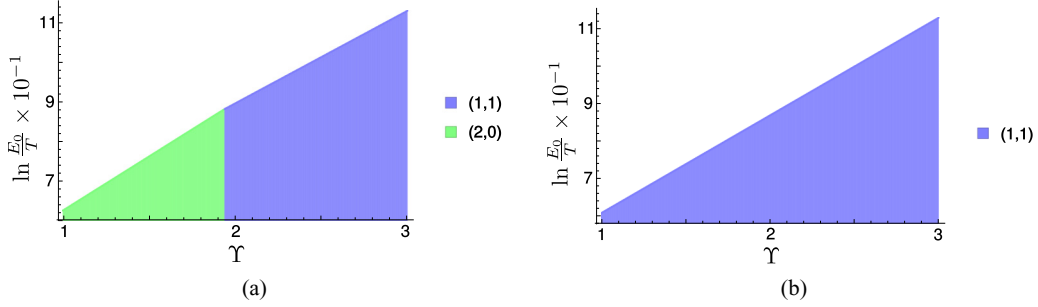


FIG. 4. Finite temperature phase diagram in $d = 3$ showing the transition temperature T_c (top boundary of the colored regions) as a function of $\Upsilon = 2k_F a$ for the (a) square and (b) triangular lattice geometries, respectively. The colors represent distinct CDW states which become dominant as Υ is tuned. The legends on the right of each diagram indicate the direction and periodicity of the ordering vector in the dominant CDW state. Here $E_0 = v_F \lambda_0$, and $\hbar_3 = \frac{A_3}{137} \frac{(c/v_F)}{\pi^2 \epsilon_r}$ with $c/v_F = 100$ and relative permittivity $\epsilon_r = 10$. We note that directions related by the C_n point group symmetry of the lattice are degenerate.

and, consequently, the number of sites per unit cell, it cannot determine the intra-unit-cell ordering pattern, which is fixed by sub-leading operators. We note that the staggered pattern of CDW modulation is most readily realized in CDW states with two sites per unit cell.

Although the CSLL state is unstable at $T = 0$, it exists above a critical temperature,

$$T_c^{(\mathbf{m})}(\hbar_d, \Upsilon) \sim v_F \lambda_0 (\tilde{g}_{\text{CDW}}^{(\mathbf{m})}(\hbar_d, \Upsilon; \lambda_0))^{-\frac{1}{2-\eta_{\text{CDW}}(\hbar_d, \mathbf{m})}}. \quad (31)$$

The corresponding finite- T phase diagrams in $d = 3$ are shown in Figs. 4(a) and 4(b) for the square and triangular lattice geometries, respectively. We note that at the noninteracting fixed point described by Eq. (3) $g_{\text{CDW}}^{(\mathbf{m})}$ is marginal, which implies

$$T_c^{(\mathbf{m})}(\hbar_d, \Upsilon) \sim v_F \lambda_0 \exp \left\{ -\frac{1}{\alpha \tilde{g}_{\text{CDW}}^{(\mathbf{m})}(\hbar_d, \Upsilon; \lambda_0)} \right\}, \quad (32)$$

has the BCS form with α being a nonuniversal numerical factor. Here, the $T_c^{(\mathbf{m})}$ is solely determined by the strength of interwire Coulomb repulsion. Since the interaction between nearest-neighbor wires is strongest, Eq. (32) implies that CDW states that modulate over a lattice spacing is the dominant instability for any \hbar_d and Υ . This is in sharp contrast to the result obtained in Eq. (31) by perturbing at the CSLL fixed point with the same operator, where more general CDW states are possible.

B. Dimensional crossover

While discussing the CDW instabilities, we implicitly assumed that the energy scale below which interwire single particle hoppings become important is small compared to $\lambda_{\text{CDW}}^{(\mathbf{m})}$. The effects of interwire hoppings in quasi-one-dimensional metals have undergone extensive investigations [22,28–31]. Here we will estimate the crossover scale that is accessible within our approach, below which the interwire hoppings cannot be ignored, and the coupled-LL framework becomes inconvenient for describing the physics [32]. We will show, in particular, that at small \hbar_d the interwire hopping amplitudes $g_{\text{sp}}^{(\mathbf{m})}$ obtain a larger scaling dimension than the CDW couplings, which implies that the dimensional crossover potentially preempts the CDW instabilities. However, if the

bare $g_{\text{sp}}^{(\mathbf{m})} \ll g_{\text{CDW}}^{(\mathbf{m})}$, the dimensional crossover scale is pushed below the CDW gap.

The correlation functions of the interwire hopping operators decay as

$$\langle O_{\text{sp}}^{(\mathbf{m})}(0, \Delta x, \mathbf{n}) O_{\text{sp}}^{(\mathbf{m})}(0, 0, \mathbf{n}) \rangle \sim \rho_0 |\lambda_0 \Delta x|^{-2\eta_{\text{sp}}(\hbar_d, \mathbf{m})}, \quad (33)$$

where $\eta_{\text{sp}}(\hbar_d, \mathbf{m}) = (\eta_{\text{CDW}}(\hbar_d, \mathbf{m}) + \eta_{\text{SC}}(\hbar_d, \mathbf{m}))/4$. By defining $\eta_{\text{CDW}}(\hbar_d, \mathbf{m}) = 2 - 4\epsilon_1(\hbar_d, \mathbf{m})$ and $\eta_{\text{SC}}(\hbar_d, \mathbf{m}) = 2 + 4\epsilon_2(\hbar_d, \mathbf{m})$, we express $\eta_{\text{sp}}(\hbar_d, \mathbf{m}) = 1 - \epsilon_1(\hbar_d, \mathbf{m}) + \epsilon_2(\hbar_d, \mathbf{m})$, where $\epsilon_i(\hbar_d, \mathbf{m}) \geq 0$. Since $\lim_{\hbar_d \rightarrow 0} \epsilon_i(\hbar_d, \mathbf{m}) = 0^+$, weak Coulomb repulsion is not sufficient for overcoming the large bare scaling dimension of $g_{\text{sp}}^{(\mathbf{m})}$. Consequently, the system may undergo a dimensional crossover below a momentum scale

$$\lambda_{\text{cross}}^{(\mathbf{m})}(\Upsilon, \mathbf{h}_d) = \lambda_0 (\tilde{g}_{\text{sp}}^{(\mathbf{m})}(\Upsilon, \mathbf{h}_d; \lambda_0))^{-\frac{1}{1+\epsilon_1(\hbar_d, \mathbf{m})-\epsilon_2(\hbar_d, \mathbf{m})}}. \quad (34)$$

Such a crossover from a lower to a higher-dimensional metallic state may be modified by quantum fluctuations that were not considered in this work. Moreover, the higher-dimensional metal itself may become unstable to the formation of a density wave or superconducting state [33], in which case the crossover would get masked by the symmetry broken state.

Within the purview of the present analysis, the CDW transitions discussed in Sec. IV A are present if $\lambda_{\text{cross}}^{(\mathbf{m})} < \lambda_{\text{CDW}}^{(\mathbf{m})}$. In contrast, for $\lambda_{\text{cross}}^{(\mathbf{m})} > \lambda_{\text{CDW}}^{(\mathbf{m})}$, the system crosses over from the CSLL to a d -dimensional Fermi liquid metal which preempts the CDW instabilities. At fixed \mathbf{m} , in the small \hbar_d regime, the former limit is satisfied for $\tilde{g}_{\text{sp}}^{(\mathbf{m})} < \tilde{g}_{\text{CDW}}^{(\mathbf{m})} < 1$. We note that such a limit is physical, since the interwire hopping and backscattering originate from distinct processes, viz. interwire single-particle tunneling and Coulomb interaction, respectively, as a result of which they are independently tunable. In particular, the rate of exponential decay of the interwire hopping amplitude with increasing interwire distance is controlled by a short-distance scale on the order of interwire lattice spacing, while the decay rate of the strength of backscatterings is controlled by the average interparticle distance on each wire as shown in Eq. (29). Since in a dilute system the latter is much smaller than the former, interwire backscatterings dominate over interwire hoppings.

V. CONCLUSION

In this paper, we analyzed the effect of backscatterings at the CSLL fixed point obtained by coupling an infinite number of quantum wires in d dimensions with Coulomb interaction in the forward scattering channel. We showed that in the absence of a dimensional crossover, an infinitesimal Coulomb interaction destabilizes the CSLL towards CDW ordering. This implies that the metallic state discussed in reference [17] is likely unstable. Several CDW states, including Wigner crystals, are shown to compete at linear order in the backscattering couplings. While CDW states with larger wave vectors are favored at large values of the effective fine structure constant \hbar_d and low density, CDW states with smaller wave vectors become dominant in the opposite limit. These properties are

demonstrated by constructing both zero and finite temperature phase diagrams.

Upon the inclusion of quantum fluctuations in terms of interwire backscatterings the relevant (in RG sense) backscattering couplings at quadratic order are expected to renormalize the transition scales [34]. Furthermore, it is in principle possible to obtain critical fixed points where a subset of the $g_{\text{CDW}}^{(\mathbf{m})} \propto \hbar_d$. Both outcomes will modify the phase diagrams obtained here. Since there is a large number of relevant couplings with finely spaced scaling dimensions at small \hbar_d , a general analysis of the higher order effects is complicated. However, by focusing on specific regions of the phase diagram, e.g., fixing \hbar_d , the physics may become more amenable to loop-wise renormalization group analyses. We leave such considerations to future work.

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APPENDIX A: DERIVATION OF ANOMALOUS DIMENSION OF HYDRODYNAMIC FIELDS

In this appendix, we outline the derivation of the scaling dimensions of the hydrodynamic fields $\vartheta_{\mathbf{n}}$ and $\varphi_{\mathbf{n}}$. We consider two different geometries of the first BZ, viz. square and triangular lattices.

1. Square lattice

For a $(d-1)$ -dimensional square lattice, the BZ is a square of sides $\frac{2\pi}{a}$. The correlation functions in Eqs. (12) and (13) are given by

$$\langle e^{i\varphi_{\mathbf{n}}(\tau, \mathbf{x})} e^{-i\varphi_{\mathbf{n}}(\tau, \mathbf{x} + \Delta \mathbf{x})} \rangle = \exp \left\{ -a^{d-1} \int_{\text{BZ}} \frac{d^{d-1}K}{(2\pi)^{d-1}} \int_{-\infty}^{\infty} \frac{dk_0 dk_x}{(2\pi)^2} [1 - \cos(xk_x)] \frac{\pi v_F \Xi_{\lambda_0}(k_x)}{k_0^2 + v_F V_{\varphi}(\mathbf{K}) k_x^2} \right\} = |\lambda_0 x|^{-2\eta_{\varphi}}, \quad (\text{A1})$$

$$\langle e^{i\vartheta_{\mathbf{n}}(\tau, \mathbf{x})} e^{-i\vartheta_{\mathbf{n}}(\tau, \mathbf{x} + \Delta \mathbf{x})} \rangle = \exp \left\{ -a^{d-1} \int_{\text{BZ}} \frac{d^{d-1}K}{(2\pi)^{d-1}} \int_{-\infty}^{\infty} \frac{dk_0 dk_x}{(2\pi)^2} [1 - \cos(xk_x)] \frac{\pi V_{\varphi}(\mathbf{K}) \Xi_{\lambda_0}(k_x)}{k_0^2 + v_F V_{\varphi}(\mathbf{K}) k_x^2} \right\} = |\lambda_0 x|^{-2\eta_{\vartheta}}, \quad (\text{A2})$$

where $\Xi_{\lambda_0}(k_x)$ is a UV regulator for k_x . It is convenient to choose a soft cutoff, e.g., $\Xi_{\lambda_0}(k_x) = \exp(-|k_x|/\lambda_0)$, since a hard cutoff, $\Xi_{\lambda_0}(k_x) = \Theta(\lambda_0 - |k_x|)$ with $\Theta(x)$ being the Heaviside theta function, leads to unphysical oscillations. We note that our choice of cutoff breaks the $1+1$ -dimensional Lorentz invariance of the Gaussian fixed point. Therefore technically this choice is not appropriate, since we do not expect the quantum fluctuations to break the Lorentz invariance. However, in the results presented here the absence of Lorentz invariance does not affect the scaling exponents; it only modifies the prefactors of the scaling terms.

The scaling exponents above are given by

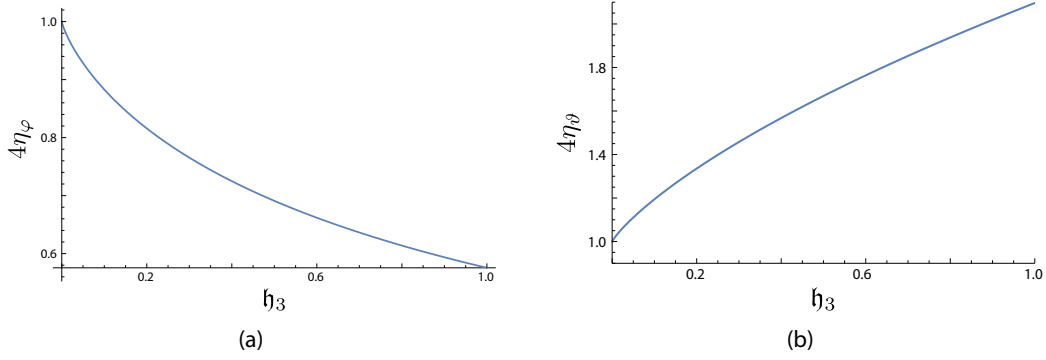
$$\eta_{\varphi} = \frac{a^{d-1}}{4} \int_{-\pi/a}^{\pi/a} \frac{d^{d-1}K}{(2\pi)^{d-1}} \left[1 + \frac{\pi^{d-1} \hbar_d}{a^{d-1} |\mathbf{K}|^{d-1}} \right]^{-\frac{1}{2}}, \quad (\text{A3})$$

$$\eta_{\vartheta} = \frac{a^{d-1}}{4} \int_{-\pi/a}^{\pi/a} \frac{d^{d-1}K}{(2\pi)^{d-1}} \left[1 + \frac{\pi^{d-1} \hbar_d}{a^{d-1} |\mathbf{K}|^{d-1}} \right]^{\frac{1}{2}}. \quad (\text{A4})$$

For $\hbar_d > 0$, $\eta_{\varphi} < 0$, and $\eta_{\vartheta} > 0$, which implies that the phase (ϑ) correlation weakens at large separation, while the density (φ) correlation is enhanced.

2. Triangular lattice

We repeat the same calculation as for the square lattice for a triangular lattice in $d = 3$. For the lattice in coordinate space, we choose the primitive vectors as, $\mathbf{e}_1 = a\hat{y}$ and $\mathbf{e}_2 = a(\hat{y}/2 + \sqrt{3}\hat{z}/2)$, where a is the lattice spacing. The reciprocal vectors, $(\mathbf{e}_1^*, \mathbf{e}_2^*)$, are deduced from the condition $\{\mathbf{e}_1^* \cdot \mathbf{e}_2 = 0, \mathbf{e}_1^* \cdot \mathbf{e}_1 = 2\pi\}$ and $\{\mathbf{e}_2^* \cdot \mathbf{e}_2 = 2\pi, \mathbf{e}_2^* \cdot \mathbf{e}_1 = 0\}$. This leads to $\mathbf{e}_1^* = \frac{2\pi}{a}(\hat{y} - \frac{1}{\sqrt{3}}\hat{z})$ and $\mathbf{e}_2^* = \frac{2\pi}{a}\frac{2}{\sqrt{3}}\hat{z}$. The reciprocal vectors define a hexagonal BZ, where the area enclosed by BZ is $\frac{2\pi}{\sqrt{3}}(\frac{2\pi}{a})^2$. In order for this area to equal $4\pi^2/l_{\text{BZ}}^2$, we need to choose $l_{\text{BZ}} = \sqrt{\frac{\sqrt{3}}{2}}a$. The sides of the hexagonal BZ is $s = \frac{4\pi}{3a}$, and integration of a function, $f(k_y, k_z)$,

FIG. 5. Due to long-range Coulomb interaction the scaling dimension of $e^{i\varphi}$ ($e^{i\vartheta}$) decreases (increases).

over BZ is given by

$$\mathcal{A}[f] = \int_{\text{BZ}}^2 \left[\int_{-s}^{-\frac{s}{2}} \frac{dk_y}{2\pi} \int_{-\sqrt{3}(k_y+s)}^{\sqrt{3}(k_y+s)} \frac{dk_z}{2\pi} + \int_{-\frac{s}{2}}^{\frac{s}{2}} \frac{dk_y}{2\pi} \int_{-\frac{\sqrt{3}}{2}s}^{\frac{\sqrt{3}}{2}s} \frac{dk_z}{2\pi} + \int_{\frac{s}{2}}^s \frac{dk_y}{2\pi} \int_{-\sqrt{3}(s-k_y)}^{\sqrt{3}(s-k_y)} \frac{dk_z}{2\pi} \right] f(k_y, k_z). \quad (\text{A5})$$

By setting $f(k_y, k_z) = 1$, it is easily checked that $\mathcal{A}[1] = 1$, which verifies the required normalization for the integration over BZ. The equal-time correlations between the simplest vertex operators are given by

$$\langle e^{i\vartheta_{\mathbf{n}}(\tau, x)} e^{-i\vartheta_{\mathbf{n}}(\tau, x + \Delta x)} \rangle = e^{-\mathcal{A} \left[\int \frac{dk_0 dk_x}{(2\pi)^2} [1 - \cos(xk_x)] \frac{\pi V_{\varphi}(\mathbf{K}) \Xi_{\lambda_0}(k_x)}{k_0^2 + v_F V_{\varphi}(\mathbf{K}) k_x^2} \right]} = |\lambda_0 x|^{-2\eta_{\vartheta}}, \quad (\text{A6})$$

$$\langle e^{i\varphi_{\mathbf{n}}(\tau, x)} e^{-i\varphi_{\mathbf{n}}(\tau, x + \Delta x)} \rangle = e^{-\mathcal{A} \left[\int \frac{dk_0 dk_x}{(2\pi)^2} [1 - \cos(xk_x)] \frac{\pi v_F \Xi_{\lambda_0}(k_x)}{k_0^2 + v_F V_{\varphi}(\mathbf{K}) k_x^2} \right]} = |\lambda_0 x|^{-2\eta_{\varphi}}. \quad (\text{A7})$$

The exponents are

$$\eta_{\vartheta} = \frac{1}{4} \mathcal{A} \left[\left(1 + \frac{\pi^2 h_3}{a^2 (k_y^2 + k_z^2)} \right)^{-1/2} \right], \quad (\text{A8})$$

$$\eta_{\varphi} = \frac{1}{4} \mathcal{A} \left[\left(1 + \frac{\pi^2 h_3}{a^2 (k_y^2 + k_z^2)} \right)^{1/2} \right]. \quad (\text{A9})$$

Unlike the square BZ, it is hard to find analytical expressions for these exponents. We compute them numerically and plot the results in Fig. 5. We check that in the noninteracting limit (i.e., $h_3 \rightarrow 0$), $\eta_{\vartheta} = \eta_{\varphi} = 1/4$. This reproduces the correct scaling for the left and right moving fermions.

APPENDIX B: SINE-GORDON TERMS

In this Appendix, we deduce the scaling dimension of various “sine-Gordon” terms defined in Eqs. (22), (23), and (24). We use the methods in Appendix A to find

$$\langle O_{\text{CDW}}^{(\mathbf{m})}(\tau, x, \mathbf{n}) O_{\text{CDW}}^{(\mathbf{m})}(\tau, x + \Delta x, \mathbf{n}) \rangle = \frac{2}{|\lambda_0 \Delta x|^{2\eta_{\text{CDW}}(h, \mathbf{m})}}, \quad (\text{B1})$$

$$\langle O_{\text{SC}}^{(\mathbf{m})}(\tau, x, \mathbf{n}) O_{\text{SC}}^{(\mathbf{m})}(\tau, x + \Delta x, \mathbf{n}) \rangle = \frac{2}{|\lambda_0 \Delta x|^{2\eta_{\text{SC}}(h, \mathbf{m})}}. \quad (\text{B2})$$

The exponents are functions of h_d , \mathbf{m} and the geometry of the underlying lattice. For the two dimensional square lattice and the triangular lattice they are, respectively,

$$[\text{square}] \quad \eta_{\text{CDW}}(h_d, \mathbf{m}) = \frac{1}{2} \int_{-1}^1 dw_1 dw_2 [1 - \cos(\pi \mathbf{m} \cdot \mathbf{w})] \left(1 + \frac{h_d}{w_1^2 + w_2^2} \right)^{-\frac{1}{2}}, \quad (\text{B3})$$

$$\eta_{\text{SC}}(h_d, \mathbf{m}) = \frac{1}{2} \int_{-1}^1 dw_1 dw_2 [1 - \cos(\pi \mathbf{m} \cdot \mathbf{w})] \left(1 + \frac{h_d}{w_1^2 + w_2^2} \right)^{\frac{1}{2}} \quad (\text{B4})$$

$$[\text{triangle}] \quad \eta_{\text{CDW}}(\mathbf{h}_d, \mathbf{m}) = 2\mathcal{A} \left\{ \left[1 - \cos \left(\mathbf{a} \left(m_1 + \frac{m_2}{2} \right) k_y + \mathbf{a} \frac{\sqrt{3}}{2} m_2 k_z \right) \right] \left(1 + \frac{\pi^2 \mathbf{h}_3}{\mathbf{a}^2 (k_y^2 + k_z^2)} \right)^{-1/2} \right\}, \quad (\text{B5})$$

$$\eta_{\text{SC}}(\mathbf{h}_d, \mathbf{m}) = 2\mathcal{A} \left\{ \left[1 - \cos \left(\mathbf{a} \left(m_1 + \frac{m_2}{2} \right) k_y + \mathbf{a} \frac{\sqrt{3}}{2} m_2 k_z \right) \right] \left(1 + \frac{\pi^2 \mathbf{h}_3}{\mathbf{a}^2 (k_y^2 + k_z^2)} \right)^{1/2} \right\}. \quad (\text{B6})$$

Here, $\mathbf{m} = \sum_{i=1}^{d-1} m_i \mathbf{e}_i$, where $\{\mathbf{e}_i\}$ are the direct lattice primitive vectors. Since the interwire interaction is rotationally symmetric, in $d = 3$, the behavior of the scaling dimensions, η_{CDW} and η_{SC} , measured along the line $m_2 = 0$ is identical to that along any *equivalent* line obtained by rotating \mathbf{e}_1 by $2\pi/n$ for a lattice with C_n symmetry. New classes of equivalent lines (directions) are constructed by similar transformations of lines along $(1,1), (1,2), (1,3), \dots$ directions.

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- [1] S. Tomonaga, *Prog. Theor. Phys.* **5**, 544 (1950).
 - [2] J. M. Luttinger, *J. Math. Phys.* **4**, 1154 (1963).
 - [3] D. C. Mattis and E. H. Lieb, *J. Math. Phys.* **6**, 304 (1965).
 - [4] F. D. M. Haldane, *Phys. Rev. Lett.* **47**, 1840 (1981); *J. Phys. C* **14**, 2585 (1981).
 - [5] M. Fabrizio and A. Parola, *Phys. Rev. Lett.* **70**, 226 (1993).
 - [6] M. Fabrizio, *Phys. Rev. B* **48**, 15838 (1993).
 - [7] A. M. Finkelstein and A. I. Larkin, *Phys. Rev. B* **47**, 10461 (1993).
 - [8] S. P. Strong, D. G. Clarke, and P. W. Anderson, *Phys. Rev. Lett.* **73**, 1007 (1994).
 - [9] V. J. Emery, E. Fradkin, S. A. Kivelson, and T. C. Lubensky, *Phys. Rev. Lett.* **85**, 2160 (2000).
 - [10] A. Vishwanath and D. Carpentier, *Phys. Rev. Lett.* **86**, 676 (2001).
 - [11] R. Mukhopadhyay, C. L. Kane, and T. C. Lubensky, *Phys. Rev. B* **64**, 045120 (2001).
 - [12] S. L. Sondhi and K. Yang, *Phys. Rev. B* **63**, 054430 (2001).
 - [13] S. A. Kivelson, V. J. Emery, and E. Fradkin, *Nature (London)* **393**, 550 (1998).
 - [14] E. Fradkin and S. A. Kivelson, *Phys. Rev. B* **59**, 8065 (1999).
 - [15] Unless the order parameter is a conserved quantity; see, e.g., K. Yang, *Phys. Rev. Lett.* **93**, 066401 (2004).
 - [16] D. Bulmash, X.-L. Qi, and C.-M. Jian, *Phys. Rev. B* **96**, 045134 (2017).
 - [17] X.-X. Zhang and N. Nagaosa, *Phys. Rev. B* **95**, 205143 (2017).
 - [18] V. Celli and N. D. Mermin, *Phys. Rev.* **140**, A839 (1965).
 - [19] H. Fukuyama, *Solid State Commun.* **26**, 783 (1978).
 - [20] V. A. Miransky and I. A. Shovkovy, *Phys. Rep.* **576**, 1 (2015), and references therein.
 - [21] An effective description in terms of spinless fermions is appropriate for electrons in the presence of an external magnetic field which leads to spin polarization.
 - [22] P. Kopietz, V. Meden, and K. Schönhammer, *Phys. Rev. Lett.* **74**, 2997 (1995); *Phys. Rev. B* **56**, 7232 (1997).
 - [23] H. J. Schulz, *Phys. Rev. Lett.* **71**, 1864 (1993).
 - [24] H. J. Schulz, *J. Phys. C* **16**, 6769 (1983).
 - [25] S. Barišić, *J. Phys.* **44**, 185 (1983).
 - [26] S. Botrić and S. Barišić, *J. Phys.* **45**, 185 (1984).
 - [27] A. H. MacDonald and M. P. A. Fisher, *Phys. Rev. B* **61**, 5724 (2000).
 - [28] C. Castellani, C. D. Castro, and W. Metzner, *Phys. Rev. Lett.* **69**, 1703 (1992).
 - [29] S. Brazovskii and V. Yakovenko, *Zh. Eksp. Teor. Fiz.* **89**, 2318 (1985).
 - [30] C. Bourbonnais and L. G. Caron, *Int. J. Mod. Phys.* **05**, 1033 (1991).
 - [31] D. Boies, C. Bourbonnais, and A.-M. S. Tremblay, *Phys. Rev. Lett.* **74**, 968 (1995).
 - [32] X.-G. Wen, *Phys. Rev. B* **42**, 6623 (1990).
 - [33] J. C. Nickel, R. Duprat, C. Bourbonnais, and N. Dupuis, *Phys. Rev. Lett.* **95**, 247001 (2005); *Phys. Rev. B* **73**, 165126 (2006).
 - [34] G. Sierra and E. H. Kim, *J. Phys. A* **36**, L37 (2003).