

Nonperturbative behavior of the quantum phase transition to a nematic Fermi fluid

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We discuss shape (Pomeranchuk) instabilities of the Fermi surface of a two-dimensional Fermi system using bosonization. We consider in detail the quantum critical behavior of the transition of a two-dimensional Fermi fluid to a nematic state which breaks spontaneously the rotational invariance of the Fermi liquid. We show that higher dimensional bosonization reproduces the quantum critical behavior expected from the Hertz-Millis analysis, and verify that this theory has dynamic critical exponent $z=3$. Going beyond this framework, we study the behavior of the fermion degrees of freedom directly, and show that at quantum criticality as well as in the quantum nematic phase (except along a set of measure zero of symmetry-dictated directions) the quasiparticles of the normal Fermi liquid are generally wiped out. Instead, they exhibit short-ranged spatial correlations that decay faster than any power law, with the law $|x|^{-1}\exp(-\text{const } |x|^{1/3})$ and we verify explicitly the vanishing of the fermion residue utilizing this expression. In contrast, the fermion autocorrelation function has the behavior $|t|^{-1}\exp(-\text{const } |t|^{-2/3})$. In this regime we also find that, at low frequency, the single-particle fermion density of states behaves as $N^*(\omega)=N^*(0)+B\omega^{2/3}\ln\omega+\dots$, where $N^*(0)$ is larger than the free Fermi value, $N(0)$, and B is a constant. These results confirm the non-Fermi liquid nature of both the quantum critical theory and of the nematic phase.

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I. INTRODUCTION

The behavior of interacting Fermi systems near continuous quantum phase transitions is a central problem in the physics of strongly correlated systems. Although much work has been done on this subject, there are still many open and as yet unresolved questions. At present the standard theory of quantum phase transitions¹⁻³ relies primarily on an analysis on the effects of fluctuations perturbatively about the results of Hartree-Fock theory. This analysis proceeds in almost complete analogy with the theory of classical critical phenomena about its upper critical dimension, and its straightforward extension to quantum phase transitions. In practice it consists of an effective theory for a suitable order parameter field while other degrees of freedom, including fermions, are often integrated out at the outset.

In many cases of interest the systems are metallic and have gapless fermionic excitations. In the standard approach, their net effect is to introduce damping in the collective modes associated with the order parameter field. In practice this results in the introduction of dissipative terms in the effective action. While much of this is certainly correct, this approach implicitly assumes that the fermions are largely unaffected by quantum criticality. Why this should be the case is far from obvious.

The assumptions of the Landau theory of the Fermi liquid⁴⁻⁶ are self-consistent and well justified within the Landau phase which has a sizable basin of stability, except in one dimensional systems.⁷⁻¹⁶ However, there is no reason for these assumptions to hold outside the Landau phase. However, there is also mounting evidence that these assumptions may also not hold in a number of phases [and not just at

quantum critical points (QCPs)], including ferromagnetic metals³³ and nematic phases of Fermi fluids.^{19,20} The possibility that quantum criticality may lead to non-Fermi liquid behavior has been a focus of research in recent years, primarily (but not only) in connection with the physics of the “normal phase” of high temperature superconductors,^{3,21-23} and with heavy-fermion systems.²⁴

The simplest example where the Landau assumptions on the behavior of the quasiparticles are violated is the quantum phase transition from a normal (Landau) Fermi liquid phase to a nematic Fermi fluid.¹⁹ A nematic Fermi fluid is a uniform phase of a system of interacting fermions in which the shape of the Fermi surface is distorted *spontaneously*, thus breaking rotational invariance.²⁵ This state is an example of the fate of a Fermi liquid beyond a Pomeranchuk instability.²⁶ In this case, the Landau assumptions appear to be violated throughout this phase, and not just at the quantum critical point.^{19,20}

The clearest experimental evidence to date of a nematic Fermi fluid phase has been found in very clean two-dimensional (2D) electron gases in magnetic fields in ultra-clean samples.^{27,28} The striking resistivity anisotropies that are observed in these experiments can be explained by the onset of nematic order at low temperatures.²⁹ It has also been proposed that phases of this type may play a central role on the physics of high temperature superconductors.^{30,31} This charge-ordered state of a strongly correlated system of fermions is the simplest example of an electronic liquid crystal phase.³⁰

The problem of the fate of the fermions at quantum criticality, and in the “non-Fermi liquid” phases mentioned above, so far has only been considered within perturbative

corrections to Hartree-Fock random-phase approximation (RPA) theory. Oganessian *et al.*¹⁹ found that the quasiparticles are wiped out as well-defined quantum states. This is due to the large fluctuations of (overdamped) quadrupolar collective modes. These authors found, within a Hartree-Fock and RPA theory, an overdamped collective mode with $z=3$ at the critical point. They also found that the fermion self-energy acquires, at the quantum critical point, an imaginary part with a frequency dependence following the law $\omega^{2/3}$. A similar behavior has been found in the case of the Stoner transition and in the antiferromagnetic phase.^{17,33} Oganessian and co-workers also found that this behavior holds inside the nematic phase, except along a set of measure zero of directions determined by the symmetry breaking.³² However, it seems quite likely that such leading order behavior^{19,20} may actually signal the complete failure of the Landau theory of the Fermi liquid. It is clear that to better understand this problem a nonperturbative analysis of the behavior of the fermions at the quantum phase transitions (and beyond) is needed. Chubukov³³ has given arguments which, in the context of the ferromagnetic metallic transition, suggest that this behavior may persist beyond the lowest order in perturbation theory.

In this paper we will consider the nematic quantum phase transition in Fermi fluids using the nonperturbative approach of higher dimensional bosonization.^{10,12,15} We will not discuss the (important) lattice effects here. Bosonization is a powerful tool to study the nonperturbative behavior of one-dimensional gapless Fermi systems, the best understood fermionic quantum critical systems.³⁴ As is well known, the kinematics of one-dimensional systems is so constrained that the bosonic collective modes completely exhaust the spectrum of these fermionic systems, allowing even for a full reconstruction of the fermionic operators entirely in terms of bosons. A striking result in the one-dimensional system is that the electron acquires a nontrivial anomalous dimension and it is no longer the quasiparticle of these systems, even for arbitrarily weak interactions. For these reasons one-dimensional gapless Fermi systems have been termed “Luttinger liquids.” The actual quasiparticles are nontrivial solitons which are orthogonal to a bare electron.³⁵

In dimensions higher than one the physics (and the kinematics) is quite different than in one dimension. Nevertheless bosonization methods still yield the physics of the Landau theory of the Fermi liquid correctly.^{10,12,15} Superficially this may seem surprising since in dimensions higher than one there are no longer strong kinematic constraints, and consequently the bosonic collective modes cannot exhaust the spectrum of an interacting Fermi system. Instead, except for narrow regimes in which the collective modes are stable quantum states, they exhibit Landau damping, reflecting their decay into particle-hole pairs. It is a key check of the validity of higher dimensional bosonization that it gets the physics of Landau damping.¹³

One appealing feature of higher dimensional bosonization is that it is actually a theory of the quantum fluctuations of the shape of the Fermi surface. It is thus a natural approach to study quantum phase transitions associated with Pomeranchuk instabilities, and in particular the nematic state.³⁶ More specifically, we focus on the nematic case for spinless fermi-

ons and compare with the work of Oganessian and co-workers¹⁹ based on RPA and Hartree-Fock. We find that the physics of the bosonic collective modes is the same in bosonization and in RPA, and thus our results agree with those of Ref. 19 in the Landau phase, in the nematic phase, and at the quantum critical point. Perhaps this is not so surprising since at long wavelengths RPA becomes asymptotically exact and this is the regime in which bosonization is correct (for a more thorough discussion, see Ref. 37). In particular we derive the effective action near the quantum critical point and find that it does have a Hertz-Millis form with dynamic critical exponent $z=3$, consistent with the findings Oganessian and co-workers,¹⁹ and by Nilsson and Castro Neto,³⁸ but in disagreement with the results of K. Yang.³⁹

We further use bosonization methods to obtain the fermion propagator. This result is well beyond the Hartree-Fock/RPA theory and thus it allows us to study the fate of the fermions nonperturbatively. We find striking violations of the Landau assumptions for Fermi liquids. Thus, the equal-time behavior of the fermion propagator at the quantum critical point (at zero temperature) is found to fall off faster than any power, decaying instead with a law $1/|x| \exp(-\text{const } |x|^{1/3})$ as a function of distance. The same behavior is found in the nematic phase except along symmetry-determined directions. We also verify explicitly from this expression the vanishing of the fermion residue as expected from this kind of behavior. In contrast to the equal-time behavior, at quantum criticality the fermion autocorrelation function behaves as $1/|t| \exp(-\text{const } |t|^{-2/3})$, with a similar albeit anisotropic law in the nematic phase as well. We also find that the low energy behavior of the one-particle density of states $N^*(\omega)$ exhibits an enhancement to a zero frequency value $N^*(0)$ which we find to be larger than $N(0)$, its noninteracting value. At finite but low frequency we further find that this one-particle density of states behaves as $N^*(\omega) = N^*(0) + B\omega^{2/3} \ln \omega$ (B is a constant), i.e., a cusp at $\omega=0$.

Thus, our bosonization results confirm that the nematic phase of a Fermi fluid is a *non-Fermi liquid*. However, its behavior is more complex than the predictions of the Hartree-Fock/RPA theory. Recently Chubukov³³ has examined the behavior of the fermion self-energy in perturbation theory at the ferromagnetic quantum critical point and found that the frequency dependence is not changed by higher order corrections. Our results for the autocorrelation function are consistent with his results, as well as with Refs. 19, 20, and 40. However, we also find that the equal-time propagator (the “one-particle density matrix”) has a very different behavior than what is predicted from these diagrammatic methods.

The paper is organized as follows: In Sec. II we derive a theory of the nematic QCP via higher dimensional bosonization. Here we present a theory of the quantum phase transition to the nematic Fermi fluid, Sec. II A, and show that it reproduces the analog of Hertz-Millis theory for this problem. In particular we give a detailed analysis of the spectral functions of the collective modes, Sec. II B, and derive the effective action valid in the vicinity of the quantum phase transition, Sec. II C. In Sec. III we use bosonization methods to calculate the fermion propagator. Here we extract the full

diagrammatic perturbation theory of the fermion Green function from bosonization, and use it to calculate the fermion self-energy. Here we check that the bosonization formulas reproduce correctly the non-Fermi liquid behavior found within the Hartree-Fock/RPA theory.¹⁹ We then use the full bosonized expression for the fermion propagator. Here we find large violations to Fermi liquid theory both at quantum criticality and in the nematic Fermi fluid phase. As an application we give a calculation of the fermion one-particle density of states. Finally, in Sec. IV we draw our conclusions. To help keep this paper self-contained, in Appendix A we give a short review the extension of bosonization to D -dimensional Fermi systems. (For a more in-depth review, see Ref. 16). In Appendix B we summarize details of the effective quadrupolar interactions, including fermion screening and Landau damping effects. In Appendix C we discuss the effects of the (uncondensed) s -wave channel on the effective theory for the nematic. The details of the calculation of the boson propagators are given in Appendix D.

II. THE NEMATIC QUANTUM PHASE TRANSITION AND THE ORDER PARAMETER

In this section, we consider the boson theory, obtained via an extension of bosonization to greater than one dimension, near a nematic (Pomeranchuk) instability of a translationally invariant fermion system. In the notation of Appendix A, we take the following action for the bosons:

$$S_0 = \frac{N(0)}{2} \sum_S \int d^2x dt [-\partial_t \varphi_S \mathbf{v}_S \cdot \nabla \varphi_S - (\mathbf{v}_S \cdot \nabla \varphi_S)^2] \quad (2.1)$$

and forward scattering interactions

$$S_{\text{int}} = \frac{N(0)}{2} \sum_{S,T} \int d^2x d^2x' dt \times F_{S-T}(\mathbf{x} - \mathbf{x}') \mathbf{v}_S \cdot \nabla \varphi_S(\mathbf{x}) \mathbf{v}_T \cdot \nabla \varphi_T(\mathbf{x}'). \quad (2.2)$$

Here, S labels the patch defined by coarse graining the Fermi surface, and the density of quasiparticles in a patch may be obtained from the boson field $\varphi_S(\mathbf{x}, t)$ via the relation $\delta n_S(\mathbf{x}, t) = N(0) \mathbf{v}_S \cdot \nabla \varphi_S(\mathbf{x}, t)$. $F_{S-T}(\mathbf{x} - \mathbf{x}')$ is therefore the interaction between particle-hole pairs in patches S and T .

We begin by analyzing our bosonized theory for a constant field configuration and reproduce Pomeranchuk's result. Consider configurations such that δn_S is constant in space and time over some particular range of time T . The resulting action is

$$S_{\text{shape}} = -\frac{VT}{N(0)} (1 + F_0) (m_0^+)^2 - \frac{VT}{2N(0)} \sum_{\ell > 0}^{N/2} (1 + F_\ell) ((m_\ell^+)^2 + (m_\ell^-)^2), \quad (2.3)$$

where we have expanded δn_S as

$$\delta n_S = \sqrt{\frac{2}{N}} \sum_{\ell=0}^{N/2} [m_\ell^+ \cos(\ell \theta_S) + m_\ell^- \sin(\ell \theta_S)] \quad (2.4)$$

and introduced the Fermi liquid parameters via

$$F_{S-T} = \frac{1}{N} F_0 + \frac{2}{N} \sum_{\ell > 0} F_\ell \cos \ell (\theta_S - \theta_T). \quad (2.5)$$

Hence, for arbitrary m_ℓ^\pm , we find that any $F_\ell < -1$ will destabilize the Fermi liquid. The point $F_\ell = -1$ we shall call the Pomeranchuk (nematic for $\ell=2$) quantum critical point (QCP). Though Fermi liquid theory breaks down at this point, Luttinger's theorem is still obeyed.

It should also be noted that in the above analysis we could have included interactions involving large angle scattering (which may lead to charge density/spin density wave instabilities), corrections to the linearized dispersion, three- or four-body interactions, and BCS processes in the bosonized theory. However, except for BCS processes, these effects are irrelevant in the Fermi liquid phase though some become important near the nematic critical point to be discussed below.

For simplicity, in the rest of this section, we shall specialize to the $\ell=2$ instability in two-spatial dimensions, the 2D quantum nematic liquid crystal, though the results generalize easily.

A. Saddle point expansion near the nematic QCP

To study the nematic QCP, originally considered by Oganesyan *et al.*,¹⁹ we set all F_ℓ to zero for $\ell \neq 2$. This is reasonable since these other modes are not critical and their effect is only to introduce finite renormalizations of the parameters in the effective theory of the critical (quadrupolar) modes (see below).

On the broken symmetry side (i.e., $F_2 < -1$), the quadratic action is no longer stable. So, in order to make the theory consistent, we need (at least) a quartic term in the bosonized action. Here, as a specific example, we consider the quartic interaction that arises from corrections to the linearized dispersion in the bosonized form given in Ref. 36,

$$S_4 = \frac{\gamma N(0)}{4!} \sum_S \int d^2r dt (\mathbf{v}_S \cdot \nabla \varphi_S)^4. \quad (2.6)$$

This term can be found by a direct extension of our ϵ_n expansion of (A12) to third order, noticing that it is an exponential series in $i\delta n_S \epsilon_n / [N(0) \hbar v_F]$. An estimate of γ may therefore naturally be obtained from the equations of motion of either the boson or fermion pictures. However, this is not the only quartic contribution to the action since an eight-fermion interaction term (that is quartic in densities) also contributes at this level with a more general form similar to the Fermi liquid interactions. Nevertheless, this will give us a flavor of the broken symmetry phase.

Since this effective theory is no longer quadratic in the bosonic fields, we will examine its behavior within a semiclassical approximation, which means that we will first find the extremal configuration and then expand our action about it. Thus, we write

$$\delta n_S = N(0) \mathbf{v}_S \cdot \nabla \varphi = \delta n_S^{cl} + \xi_S, \quad (2.7)$$

where δn_S^{cl} is the solution of the classical equations of motion at a uniform, mean field level

$$\delta n_S^{cl} + \frac{2}{N} F_2(0) \sum_T \cos[2(\theta_S - \theta_T)] \delta n_T^{cl} + \frac{\gamma}{3! N(0)^2} (\delta n_S^{cl})^3 = 0, \quad (2.8)$$

where as stated earlier, we set $F_\ell(0)=0$ for $\ell \neq 2$. Because the nonlinear term is cubic, we seek solutions of the type

$$\delta n_S^{cl} = \frac{1}{\sqrt{N}} \sum_{\ell=-N/2}^{+N/2} m_\ell e^{i\ell \theta_S}, \quad (2.9)$$

where $m_\ell \neq 0$ for $\ell = \{\pm 2, \pm 6, \pm 10, \dots\}$. In terms of the m_ℓ 's, the equation of motion becomes

$$m_\ell + F_2(m_2 \delta_{\ell,2} + m_{-2} \delta_{\ell,-2}) + \frac{\gamma}{3! N(0)^2 N} \sum_{\ell_1, \ell_2} m_{\ell_1} m_{\ell_2} m_{\ell - \ell_1 - \ell_2} = 0. \quad (2.10)$$

We first observe that for $F_2 \geq -1$, the only possible solution is $m_\ell = 0$ for all ℓ , as expected for the isotropic case. If $F_2 < -1$, there exists a whole set of nontrivial solutions involving, in general, all harmonics $\ell = \{\pm 2, \pm 6, \pm 10, \dots\}$ obeying particle hole symmetry. Nevertheless, when we are near the phase transition, i.e., $F_2 \lesssim -1$, one can find the set of solutions analytically with

$$m_2 m_{-2} = \frac{1}{2} ((m_2^+)^2 + (m_2^-)^2) = \frac{2}{\gamma} N(0)^2 N |1 + F_2(0)| \quad (2.11)$$

using the notation of (4) and for the higher harmonics

$$|m_{4k+2}|^2 \propto N(0)^2 N \left(\frac{2}{\gamma} |1 + F_2(0)| \right)^{2k+1}, \quad (2.12)$$

so that in the limit $F_2 \rightarrow -1$ we can neglect the higher harmonics.

From this calculation, we conclude that near the $F_2 = -1$ nematic QCP, the Fermi surface takes the simple shape

$$\delta n_S^{cl} = N(0) \sqrt{\frac{2|1 + F_2(0)|}{\gamma}} \cos(2\theta_S - \alpha), \quad (2.13)$$

where α picks out the major axis of the ellipse that is spontaneously chosen. Without loss of generality, from now on we will set $\alpha = 0$. We also conclude that further away from the critical point, the anharmonic quartic terms generically introduce higher harmonics to this shape. From this point of view, for example, the Fermi surface in the nematic phase may become increasingly flatter away from the critical point leading to an additional instability towards a smectic phase, breaking translational order in one direction. However, this phase transition cannot be seen within our forward scattering only model and all that happens deeper into the phase here is that the shape becomes more anharmonic. It should also be noted that the other harmonics, $\ell = \{0, \pm 4, \pm 8, \dots\}$, appear when particle-hole symmetry is broken, e.g., when a cubic

term in the action is introduced (S_3), corresponding to the addition of quadratic terms to the fermion dispersion.

B. Theory of the quadrupole moment density

The next step beyond mean field theory is to formulate an order parameter theory. To that end, in this section we are interested in focusing on describing the behavior of the quadrupole moment density, $m_2^\pm(\mathbf{q}, \omega)$. Again, for our present purposes we shall keep only $F_2(q)$. However, in Appendix C we show that by keeping also $F_0(q)$, the additional effects of the noncritical modes do not change our results in any essential way.

Now, our goal here is to obtain an action entirely in terms of m_2^\pm . This can be easily accomplished by following Ref. 16 and using a Hubbard-Stratonovich transformation to aid the diagonalization. In the end, however, both the auxiliary fields and m_ℓ^\pm for $\ell \neq 2$ shall be integrated out.

In momentum space, our free action can be written as

$$S_0 = \frac{1}{2} \sum_S \int \frac{d^2 q d\omega}{(2\pi)^3} (\chi_S^0)^{-1}(q, \omega) \delta n_S(\mathbf{q}, \omega) \delta n_S(-\mathbf{q}, -\omega), \quad (2.14)$$

where

$$\chi_S^0(\mathbf{q}, \omega) = N(0) \frac{\mathbf{v}_S \cdot \mathbf{q}}{\omega - \mathbf{v}_S \cdot \mathbf{q}} = N(0) \frac{\cos \theta_S}{s - \cos \theta_S} \quad (2.15)$$

is the density-density response function in the ‘‘small q ’’ limit of patch S with $s = \omega / q v_F$. (Note: when needed, we regularize the denominator by letting $s \rightarrow s + i\epsilon \text{sign}(s)$ according to the usual time-ordering prescription.)

The interactions, described by the quadratic action S_{int} , cf. Eq. (2.2), become *diagonal* in the angular momentum basis, i.e., in terms of the multipole densities $m_\ell^\pm(q, \omega)$. In particular, the contribution to S_{int} from the $\ell = 2$ (quadrupolar) densities, is

$$S_{\text{int}} = \frac{1}{2} \int \frac{d^2 q d\omega}{(2\pi)^3} f_2(q) [|m_2^+(q, \omega)|^2 + |m_2^-(q, \omega)|^2], \quad (2.16)$$

with $f_2(q)$ defined as usual in Fermi liquid theory through $F_2(q) = N(0) f_2(q)$. The contributions from the other angular momentum channels have a similar form (in terms of the respective Landau parameters).

To aid the diagonalization, we split up the free part of the action, S_0 , using the Hubbard-Stratonovich transformation,

$$S_0 = - \frac{1}{2} \sum_S \int \frac{d^2 q d\omega}{(2\pi)^3} [\chi_S^0(q, \omega) \sigma_S(q, \omega) \sigma_S(-q, -\omega) + (\sigma_S(q, \omega) \delta n_S(-q, -\omega) \sigma_S(-q, -\omega) \delta n_S(q, \omega))] \quad (2.17)$$

and then switch over to the angular momentum basis,

$$S_0 = -\frac{1}{2} \sum_{\eta=\pm} \sum_{\ell, \ell'} \int \frac{d^2 q d\omega}{(2\pi)^3} [(\chi_{\ell-\ell'}^0 + \eta \chi_{\ell+\ell'}^0) \sigma_\ell^\eta \sigma_{\ell'}^\eta + 2(\delta_{\ell-\ell',0} + \eta \delta_{\ell+\ell',0}) (\sigma_\ell^\eta m_\ell^\eta)] \quad (2.18)$$

In the large- N limit, we have

$$\chi_\ell^0 = N(0) \int_0^{2\pi} \frac{d\theta}{2\pi} \frac{\cos \theta}{s - \cos \theta} e^{i\ell\theta}. \quad (2.19)$$

Here we have Fourier transformed with respect to $\hat{\mathbf{q}} = (\cos \phi, \sin \phi)$, that is, $\theta = \theta_s - \phi$.

Now, we will integrate out all the m_ℓ^\pm densities, except for $\ell=2$. This can be easily done since S_0 is a linear function of these fields, while they are absent in S_{int} . This is so for this model with only a quadrupolar interaction, i.e., we have set their corresponding Fermi liquid parameters to zero. (In the vicinity of the nematic transition it is straightforward to include the effects of the $\ell \neq 0$ channels. Their net effect is to give rise to simple renormalizations of the effective theory we are about to derive. A detailed analysis is given Appendix C.) The result is a delta function for the σ_ℓ fields, allowing us to also integrate them out with the net result to simply set $\sigma_\ell^\pm = 0$ for all ℓ except $\ell=2$. This gives us the following simple expression for the effective free action:

$$S_0 = -\frac{1}{2} \sum_{\eta=\pm} \int \frac{d^2 q d\omega}{(2\pi)^3} [(\chi_0^0 + \eta \chi_4^0) |\sigma_2^\eta|^2 + 2(\sigma_2^\eta m_2^\eta)] \quad (2.20)$$

and we now have an action entirely in terms of the quadrupole moment density.

The final step is to integrate out the σ_2^\pm fields. This is easily accomplished and we obtain the Gaussian level of the order parameter theory, including the effects of the interactions S_{int} . The action of the effective theory is

$$S_2[m_2^+, m_2^-] = \frac{1}{2} \int \frac{d^2 q d\omega}{(2\pi)^3} \left[\frac{1}{\chi_2^+} |m_2^+|^2 + \frac{1}{\chi_2^-} |m_2^-|^2 \right], \quad (2.21)$$

where $\chi_2^\pm(s, q)$ is the dynamical correlation function (susceptibility) of the quadrupolar densities

$$\chi_2^\pm(s, q) = \frac{\chi_0^0(s) \pm \chi_4^0(s)}{1 - f_2(q)(\chi_0^0(s) \pm \chi_4^0(s))} \quad (2.22)$$

and

$$\chi_{2\ell}^0 = N(0) \left[-\delta_{\ell,0} + K_0(s) \left(\frac{1 - K_0(s)}{1 + K_0(s)} \right)^\ell \right], \quad (2.23)$$

with

$$K_0(s) = \frac{s}{\sqrt{s-1}\sqrt{s+1}}. \quad (2.24)$$

Naturally, this is just RPA quadrupolar susceptibility of Ref. 19.

We should stress that the effective action S_2 of Eq. (2.21) does not include the effects of the nonlinear interactions rep-

resented in S_4 , cf. Eq. (2.6), which are crucial to stabilize the nematic state past the nematic QCP [for $F_2(0) < -1$]. As we discussed above, these nonlinear terms do mix the different angular momentum channels. However, provided there is no condensation for $\ell \neq 2$ it is still possible to integrate out these degrees of freedom, at least perturbatively. Thus, sufficiently close to the nematic QCP, the effects of the higher angular momentum channels will remain perturbatively small.

Furthermore, should we have decided to integrate out the density fields, m_2^\pm in favor of the auxiliary fields σ_2^\pm , we would have found the propagators of the σ_ℓ^\pm fields to be the RPA effective interaction

$$V_2^\pm(s, q) = \frac{f_2(q)}{1 - f_2(q)[\chi_0^0(s) \pm \chi_4^0(s)]}. \quad (2.25)$$

The action of the σ_ℓ^\pm fields is precisely that found by Oganesyan and co-workers¹⁹ also obtained through a Hubbard-Stratonovich transformation but performed directly in the fermion theory. We shall find that understanding this interaction is the key to understanding the physics of the nematic QCP.

C. Order parameter theory of the nematic QCP

The theory with the action given by Eqs. (2.1) and (2.2) seemingly describes the quantum mechanics of a fluctuating surface. This suggests that the effective degrees of freedom ought to be long-lived bosonic modes of the fluctuations of the shape of the Fermi surface. In other terms, this bosonized theory would seem to be entirely described by stable collective modes. However, these bosonic excitations are not generally stable due to *Landau damping* effects, represented by the branch cut singularities in (2.23).^{19,41}

We will examine this problem more closely. It is useful to introduce the following *quadrupole density spectral functions*:

$$S_2^\pm(q, s) \equiv -2 \text{sign}(s) \text{Im} \chi_2^\pm(q, s) \\ = -\frac{2 \text{sign}(s) \text{Im} V_2^\pm(q, s)}{f_2(q)^2} \equiv \frac{B_2^\pm(q, s)}{f_2(q)^2}, \quad (2.26)$$

where $B_2^\pm(q, s)$ denotes the spectral function of $V_2^\pm(q, s)$, the correlation function of the σ_2^\pm fields.

The analysis is greatly simplified upon recognizing that $V_2^\pm(q, s)$ is a polynomial function of $K_0(s)$. For $\ell=2$ the denominator of $V_2^\pm(q, s)$ is a cubic function of K_0 , whereas for $V_2^-(q, s)$ it is quadratic in K_0 . In Appendix B we show that $V_2^\pm(q, s)$ have the partial fraction expansions

$$V_2^\pm(q, s) = \frac{1}{N(0)} \sum_{\beta} \frac{\mathcal{Z}_\beta^\pm(q)}{\delta_\beta^\pm(q) - K_0(s)}, \quad (2.27)$$

where $\beta=a, b, c$ for V_2^+ , and $\beta=a, b$ for V_2^- . We can therefore view $V_2^\pm(q, s)$ as a sum of terms each of the form of the RPA s -wave channel effective interaction, renormalized by a residue \mathcal{Z}_β^\pm and with an effective interaction $f_\beta^\pm(q) = [1 - \delta_\beta^\pm(q)]^{-1}$. Details of these expansions are given in Appendix B.

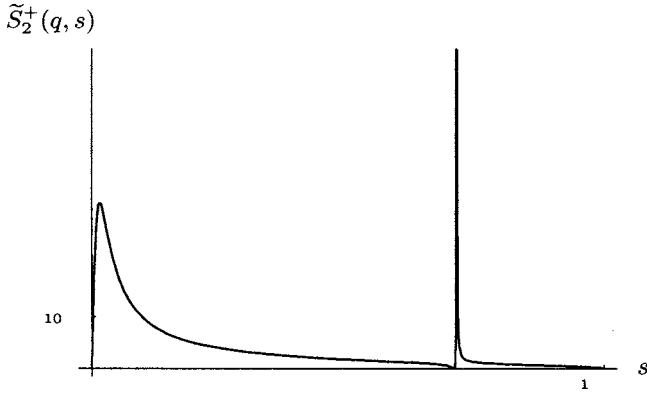


FIG. 1. The (normalized) quadrupole density spectral function $\tilde{S}_2^+(q, s) = S_2^+(q, s)/N(0)$, Eqs. (2.26) and (2.29) for $q \neq 0$, very close to the nematic quantum phase transition from the Fermi liquid phase, $F_2(0) \rightarrow -1^+$. As the QCP is approached, there is a large increase of spectral weight in the $z=3$ overdamped quadrupolar mode at very low frequency. Notice the delta function contribution of the propagating mode with $z=1$ discussed in the text. The propagating mode with $z=2$ gives a similar delta function contribution (with smaller spectral weight) to $S_2^-(q, s)$.

Near the nematic QCP, where $|F_2(0) + 1| \ll 1$, upon defining the quantity $\delta_2(q)$,

$$\delta_2(q) = \frac{1 + F_2(q)}{F_2(q)} = 1 + \frac{1}{F_2(0)} - \kappa q^2, \quad (2.28)$$

the effective interactions $V_2^\pm(q, s)$ become simple and we obtain

$$V_2^+(q, s) \approx \frac{1}{2N(0)} \left[\frac{1}{\frac{\delta_2(q)}{2} - K_0(s)} - \frac{1/4}{s^2 - 1/2} \right], \quad (2.29)$$

$$V_2^-(q, s) \approx \frac{1}{4N(0)} \frac{1}{s^2 + \frac{\delta_2(q)}{4}}. \quad (2.30)$$

Thus, near the critical point we find a propagating mode with dispersion $\omega_q = (1/\sqrt{2})qv_F$ (a $z=1$ mode), represented by the pole in the second term for V_2^+ [see Eq. (2.29)], and another one with dispersion $\omega_q = (\sqrt{\delta_2(q)}/2)qv_F$ (a $z=2$ mode), given by the pole in V_2^- [see Eq. (2.30)]. Furthermore, we also find an overdamped mode, given by the pole in first term of V_2^+ [see Eq. (2.29)], with a dispersion relation of $s = -i\delta_2(q)/2$, and dynamic critical exponent $z=3$. For q sufficiently small, and asymptotically close to the nematic quantum critical point, we find that this overdamped mode dominates the spectral function over the other two (propagating) modes.

In Fig. 1 we present a plot of the spectral function $S_2^+(q, s) \approx B_2^+(q, s) = -2 \text{sign}(s) \text{Im} V_2^+(q, s)$. It shows that, as the QCP is approached from the Fermi liquid side, there is a large transfer of spectral weight in the quadrupolar spectral function to the low frequency end of the spectrum, associated with the emergence of the overdamped $z=3$ mode. This mode thus controls the quantum critical behavior.

Conversely, normal Fermi-liquid behavior is obtained if $\delta_2(0)$ is finite so that $z=1$. On the other hand, if the overdamped mode were to be absent, the dynamic critical behavior would be controlled by propagating mode with $z=2$ discussed above. The trend is thus opposite to what one might naively expect: the higher the z , the higher the effective dimension $D+z$, the stronger the divergence of the overdamped mode. Recently, Yang³⁹ proposed that the transition to the quantum nematic state should have dynamic critical exponent $z=2$. The analysis we just presented shows that this is not the case.

As a result of the above analysis, we obtain the following effective action for the quadrupole density near the nematic QCP:

$$S_{\text{QCP}} = \frac{1}{2N(0)} \int \frac{d^2q d\omega}{(2\pi)^3} [[2i|s| - 1 - F_2(q)] |m_2^+|^2 + [4s^2 - 1 - F_2(q)] |m_2^-|^2] - \frac{\gamma}{8N(0)^3 N} \int d^2x dt (m_2^{+2} + m_2^{-2}). \quad (2.31)$$

The order parameter field is

$$m_2^+(\mathbf{q}, \omega) = \sqrt{\frac{2}{N}} \sum_S \delta n_S(\mathbf{q}, \omega) \cos 2(\theta_S - \phi),$$

$$m_2^-(\mathbf{q}, \omega) = \sqrt{\frac{2}{N}} \sum_S \delta n_S(\mathbf{q}, \omega) \sin 2(\theta_S - \phi), \quad (2.32)$$

where, again, ϕ is the direction of $\mathbf{q} = q(\cos \phi, \sin \phi)$. This action is therefore very similar to Hertz's action for the ferromagnetic quantum phase transition in itinerant fermionic systems,¹ but here within the context of the formation of nematic order (and similar actions may also be obtained for higher ℓ). Recently, Nilsson and Castro Neto³⁸ derived this action using Fermi liquid theory methods.

Now, let us look on the broken symmetry side, in the nematic phase. As in any theory with an $O(2)$ symmetry, here we will find that the order parameter will spontaneously pick a direction. As a result, it is no longer useful to Fourier transform with respect to ϕ in Eq. (2.32). Rotating back and after the saddle point expansion about the classical configuration (2.13), we find the quadratic action on the nematic side

$$S_{\text{QCP}}^{\text{nematic}} = \frac{1}{2N(0)} \int \frac{d^2q d\omega}{(2\pi)^3} \mathbf{m}_2(\mathbf{q}, \omega) \cdot \chi_2^{-1} \cdot \mathbf{m}_2(-\mathbf{q}, -\omega) \quad (2.33)$$

where $\mathbf{m}_2(\mathbf{q}, \omega) = [m_2^+(\mathbf{q}, \omega), m_2^-(\mathbf{q}, \omega)]$ and

$$\chi_2^{-1} = [i|s| - \kappa q^2 - |1 + F_2(0)|] \begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix} + i|s| \times \begin{pmatrix} \cos 4\phi & \sin 4\phi \\ \sin 4\phi & -\cos 4\phi \end{pmatrix} - |1 + F_2(0)| \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix} \quad (2.34)$$

in the reference frame in which the nematic order parameter

is diagonal, i.e., its principal axes, whose orientation is determined spontaneously. The difference between this action and the previous one discussed above for the symmetric phase is the emergence of the last term which originates from the nonlinear (quartic) term in the effective action. This term ruins our ability to rotate ϕ out of the action.

Noting that the off-diagonal terms are higher order in s , we may write this in the simplified form

$$S_{\text{QCP}}^{\text{nematic}} = \frac{1}{2N(0)} \int \frac{d^2q}{(2\pi)^2} \int \frac{d\omega}{2\pi} [[2i|s|\cos^2(2\phi) - 2|1 + F_2(0)| - \kappa q^2]|m_2^+|^2 + [2i|s|\sin^2(2\phi) - \kappa q^2]|m_2^-|^2]. \quad (2.35)$$

We see that now m_2^+ is the amplitude mode, which has $z=1$, while m_2^- is the nematic Goldstone mode *which continues to have dynamic critical exponent $z=3$ even in the nematic phase.*¹⁹

The last point is to discuss how this affects the original boson theory. If we bring back all the integrated out angular momentum channels, we find that the free action on the broken symmetry side has

$$\tilde{\chi}_S^0(\mathbf{q}, \omega) = N(0) \frac{\mathbf{v}_S \cdot \mathbf{q}}{\omega - \mathbf{v}_S \cdot \mathbf{q}}, \quad (2.36)$$

with a weakly renormalized Fermi velocity

$$\tilde{\mathbf{v}}_S = [1 + 4|1 + F_2(0)|\cos^2(2\theta_S)]\mathbf{v}_S. \quad (2.37)$$

Hence, spontaneous symmetry breaking essentially produces a Goldstone mode that continues the critical, $z=3$ behavior into the nematic phase while leaving the rest of the theory virtually untouched until deep into the broken symmetry phase.

III. FERMIONS IN THE CRITICAL REGIME

In the past sections we discussed the behavior of the collective modes near the nematic quantum phase transition. Much of what we discussed in the previous section on the behavior of the collective modes is indeed in complete agreement with the RPA treatment of this theory.¹⁹ This should not be a surprise since RPA is asymptotically exact at low energies and at low frequencies. This is also the reason while bosonization works in the same regime.

We will now turn our attention to the behavior of the fermionic degrees of freedom near the nematic QCP and in the nematic phase. This is very different problem. In Ref. 19 the behavior of the fermion Green function was studied perturbatively and a startling non-Fermi liquid behavior was found already at the lowest (“Fock”) order. However, this very finding raises questions on the applicability of perturbation theory for the fermion propagator. In this section we will use bosonization methods to address this problem.

Here we will use bosonization to compute the fermion propagator. Within this approach one has a theory for the bosonized degrees of freedom and a set of operator identities relating observables of the fermionic theory to those of the

bosonic theory. For a summary see Appendix A. There are two important issues to keep in mind. One is that the bosonized theory is exact for a fermionic theory with a linearized dispersion and forward scattering interactions (i.e., those described by Landau parameters). The other is that one has, within this theory, an operator to represent the fermion. Corrections to the linear dispersion as well as other (nonforward scattering) interactions are represented by nonlinear terms in the action of the bosonized theory. The expressions that we will derive below apply strictly speaking to the fixed point theory, in which these perturbations are not included. It turns out that corrections due to the nonlinearities of the fermion dispersion and other such terms are irrelevant (both in the Landau phase and at the quantum critical point). As such they will affect the results at high energies and at momenta but their effects become negligible in the low energy limit. Please note that one such operator, discussed in the previous section, stabilizes the nematic phase, i.e., it is a prototypical dangerous irrelevant operator.

The boson Green function may be found from our action in a similar way as we found the above density-density correlation functions, that is, using a Hubbard-Stratonovich approach. The result is

$$G_{B(S,T)}(\mathbf{x}, t) = G_{B(S,T)}^0(\mathbf{x}, t) + i \int \frac{d^2k d\varepsilon}{(2\pi)^3} G_{F(S)}^0(k, \varepsilon) V_{S,T}(k, \varepsilon) G_{F(T)}^0(k, \varepsilon) \times (e^{i(\mathbf{k}\cdot\mathbf{x} - \varepsilon t)} - 1), \quad (3.1)$$

where $G_{B(S,T)}^0$ is infrared divergent unless $\mathbf{x} \parallel \mathbf{v}_S$ and $S=T$. On the same patch, $G_{B(S,S)}^0(\mathbf{x}, t)$ is given by the standard expression

$$G_{B(S,S)}^0(\mathbf{x}, t) = \langle \varphi_S(\mathbf{x}, t) \varphi_S(\mathbf{0}, 0) \rangle - \langle \varphi_S(\mathbf{0}, 0)^2 \rangle = -\ln \left(\frac{\hat{n}_S \cdot \mathbf{x} + iv_F t + ia \text{sign } t}{ia} \right), \quad (3.2)$$

where a is a short-distance cutoff. In Eq. (3.1) we have denoted by $V_{S,T}(q, \omega)$ the effective interaction

$$V_{S,T}(q, \omega) = \frac{1}{N(0)} \sum_{\ell, \ell'} e^{i[\ell(\theta_S - \phi) + \ell'(\theta_T - \phi)]} \langle \sigma_\ell(q, \omega) \sigma_{\ell'}(-q, -\omega) \rangle, \quad (3.3)$$

which for the quadrupolar case, becomes simply

$$V_{S,T}(q, \omega) = V_2^+(q, \omega) \cos 2(\theta_S - \phi) \cos 2(\theta_T - \phi) + V_2^-(q, \omega) \sin 2(\theta_S - \phi) \sin 2(\theta_T - \phi), \quad (3.4)$$

where the relative angle, $\theta_S - \phi$ appears here [see Eq. (2.18)]. We also have used in (3.1) the free fermion Green function

$$G_{F(S)}^0(q, \omega) = \frac{1}{\omega - \mathbf{v}_S \cdot \mathbf{q} + i\varepsilon \text{sign } (\omega)}. \quad (3.5)$$

We will now use the bosonized form of the fermion operator [see Eq. (A20) of Appendix A] to give us the fermion Green function of the form

$$G_F(\mathbf{x}, t) = \sum_{S, T} \left(\frac{-i}{N} \right) \langle T \psi_S(\mathbf{x}, t) \psi_T^\dagger(\mathbf{0}, 0) \rangle e^{i\mathbf{k}_S \cdot \mathbf{x}} \quad (3.6)$$

$$\equiv \frac{1}{N} \sum_S G_{F(S)}(\mathbf{x}, t) e^{i\mathbf{k}_S \cdot \mathbf{x}}, \quad (3.7)$$

for which we find the explicit expression

$$G_{F(S)}(\mathbf{x}, t) = G_{F(S)}^0(\mathbf{x}, t) \exp \left[i \int \frac{d^2 k d\varepsilon}{(2\pi)^3} G_{F(S)}^0(k, \varepsilon) V_{S,S}(k, \varepsilon) \times G_{F(S)}^0(k, \varepsilon) (e^{i(\mathbf{k} \cdot \mathbf{x} - \varepsilon t)} - 1) \right]. \quad (3.8)$$

This expression has many similarities with the bosonization formulas usually obtained in one dimension. In particular, the free-fermion prefactor also arises there. However, the exponential factor, which in one dimension yields an anomalous dimension for the fermion operator, plays a very different role in dimensions higher than one.

A. Diagrammatic expansion for the bosonized theory

We will first show that the bosonized formula of Eq. (3.8) is consistent with the perturbative results of Ref. 19. To do that we will expand the exponential and Fourier transform to momentum space to find its diagrammatic expansion. The result will be a series of convolutions since in real space they are products. The first order term is

$$\delta G_F^{(1)}(\mathbf{q}, \omega) = i \int \frac{d^2 k d\varepsilon}{(2\pi)^3} (G_{F(S)}^0(\mathbf{q} - \mathbf{k}, \omega - \varepsilon) - G_{F(S)}^0(\mathbf{q}, \omega)) \times [G_{F(S)}^0(\mathbf{k}, \varepsilon)]^2 V_{S,S}(k, \varepsilon), \quad (3.9)$$

which does not look like it obeys the Feynman rules for the perturbation theory of nonrelativistic fermions. However, we may utilize the following identity:

$$[G_{F(S)}^0(\mathbf{q} - \mathbf{k}, \omega - \varepsilon) - G_{F(S)}^0(\mathbf{q}, \omega)] G_{F(S)}^0(\mathbf{k}, \varepsilon) = G_{F(S)}^0(\mathbf{q} - \mathbf{k}, \omega - \varepsilon) G_{F(S)}^0(\mathbf{q}, \omega) \quad (3.10)$$

and use it again through

$$G_{F(S)}^0(\mathbf{q} - \mathbf{k}, \omega - \varepsilon) = G_{F(S)}^0(\mathbf{q} - \mathbf{k}, \omega - \varepsilon) - G_{F(S)}^0(\mathbf{q}, \omega) + G_{F(S)}^0(\mathbf{q}, \omega) \quad (3.11)$$

to obtain

$$\delta G_F^{(1)}(\mathbf{q}, \omega) = i [G_{F(S)}^0(q, \omega)]^2 \int \frac{d^2 k d\varepsilon}{(2\pi)^3} (G_{F(S)}^0(\mathbf{q} - \mathbf{k}, \omega - \varepsilon) + G_{F(S)}^0(\mathbf{k}, \varepsilon)) V_{S,S}(k, \varepsilon). \quad (3.12)$$

The second term is actually the shift in the chemical potential, $\Sigma(k_F, 0)$, and is zero by the effective particle-hole symmetry of this theory (with a linearized fermion dispersion). We therefore obtain

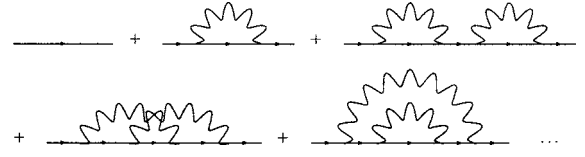


FIG. 2. Bosonization's Feynman diagram series.

$$\delta G_F^{(1)}(\mathbf{q}, \omega) = i [G_{F(S)}^0(q, \omega)]^2 \int \frac{d^2 k d\varepsilon}{(2\pi)^3} \times G_{F(S)}^0(\mathbf{q} - \mathbf{k}, \omega - \varepsilon) V_{S,S}(k, \varepsilon), \quad (3.13)$$

which is the correct result since, as usual, the Hartree term vanishes.

Using the same tricks and assumptions, the second-order contribution can also be worked out (though it is much more work). The series to second order is shown in Fig. 2; please keep in mind that the interaction, $V_{S,S}$ is the full bubble summed RPA interaction. Higher dimensional bosonization therefore keeps all diagrams in perturbation theory that contain up to simple bubbles while neglecting more complicated bubbles as is usual in the RPA (in the Landau theory of the Fermi liquid, these irrelevant operators contribute subdominant potentially nonanalytic temperature and frequency dependent terms to physical quantities¹⁸).

Some time ago, Kopietz and Castilla^{42,43} used a somewhat different (and in principle equivalent) form of bosonization and discussed the effects of a quadratic term in the fermion energy dispersion. However, instead of using an operator identity (and thus not exploiting the nonperturbative character of bosonization) they chose to make contact with perturbation theory, and proceeded to propose a modified form for the fermion propagator directly. It gave an exponential factor similar to the one appearing here except the fermion Green functions that appear in the exponential include the quadratic terms in their energy dispersion. However, these Green functions do not satisfy the criterion $\Sigma(k_F, 0) = 0$ so that in addition to the exponential, they include a preexponential factor that is necessary, for example, to cancel the second term of Eq. (3.12). As a result, order-by-order in $V_{S,S}$, one needs to keep precisely the right preexponential factor to cancel the additional terms. A calculation with their method can therefore only be carried out to a small finite order, and one can no longer think of the exponential factor as separate from the preexponential factor. In contrast, we have seen here that the bosonized expressions, when treated consistently, yield exact results which agree with those of perturbation theory order by order, albeit only in the low energy limit, including the singular behavior.

B. Perturbative results

Before computing the full nonperturbative form of the fermion Green function, let us verify that our bosonized theory, reproduces the perturbative results of Oganessian *et al.*¹⁹ near the nematic QCP. We shall be interested, therefore, in the integral

$$\Sigma_S^{(1)}(\mathbf{q}, \omega) = i \int_{\mathcal{P}_S} \frac{d^2 k d\varepsilon}{(2\pi)^3} G_{F(S)}^0(\mathbf{q} - \mathbf{k}, \omega - \varepsilon) V_{S,S}(k, \varepsilon). \quad (3.14)$$

Here $V_{S,S}(k, \varepsilon)$ is the effective interaction mediated by the collective modes, cf. Eqs. (2.29), (2.30), and (3.4). The main contribution to the self-energy is due to the overdamped Goldstone modes, as noted in Ref. 19. Thus we take a generic interaction of the form

$$V_{S,S}(k, \varepsilon) = \frac{1/N(0)}{\delta(k) - K_0(s)}, \quad (3.15)$$

noting that $\cos^2[2(\theta_S - \phi)]$, which should appear as a coefficient in the nematic case, only introduces irrelevant contributions to the integral and is therefore left out of this analysis. Here, at the QCP, $\delta(k) = -\kappa k^2/2$ but as a check we may characterize the Fermi liquid phase by letting $\delta(k) = \delta$, a constant.

Before computing $\Sigma_S^{(1)}(\mathbf{q}, \omega)$, we should note that this expression gives the clearest definition of the patch. Given the ultraviolet cutoff $k_F - \lambda/2 < |\mathbf{k}| < k_F + \lambda/2$, the patch is defined so that in comparison to the original theory

$$\Sigma^{(1)}(\mathbf{k} = \mathbf{k}_S + \mathbf{q}, \omega) \approx \Sigma_S^{(1)}(\mathbf{q}, \omega), \quad \mathbf{q} \in \mathcal{P}_S, \quad (3.16)$$

and therefore the patch width is of order

$$\Lambda \sim \sqrt{k_F \lambda}, \quad (3.17)$$

dictated by the curvature of the Fermi surface alluded to earlier.

Now, to simplify our calculation, let us focus on the fermion lifetime near the Fermi surface

$$2\Gamma_S(\mathbf{q}, \omega) = -2 \text{sign}(\omega) \text{Im} \Sigma_S(\mathbf{q}, \omega). \quad (3.18)$$

This can most easily be expressed in terms of the spectral function, $B_2(\mathbf{q}, s)$, derived from the imaginary part of $V_{S,S}(q, s)$ as in Eq. (2.26) and Eq. (3.4) (after setting the angular factors to 1),

$$2\Gamma_S(\mathbf{q}, \omega) = \int_{-\Lambda/2}^{\Lambda/2} \frac{dk_t}{2\pi} \int_{q_n - |\omega|/v_F}^{q_n} \frac{dk_n}{2\pi} B_2\left(k, \frac{\omega - v_F(q_n - k_n)}{kv_F}\right), \quad (3.19)$$

where $q_n = \mathbf{v}_S \cdot \mathbf{q} / v_F$. Here we note that this integral is dominated by the contribution of the overdamped mode, which enters in $B_2^+(q, s)$. (The other contributions, associated with the propagating collective modes, only yield regular dependences in the fermion frequency.) Setting $q_n \rightarrow 0$ and looking at the limit $\omega \rightarrow 0$ we obtain

$$\Gamma_S(\omega) = \frac{1}{2\pi^2} \int_0^{\omega/v_F} dk_n \int_0^\Lambda dk_t B_2^+\left(\sqrt{k_n^2 + k_t^2}, \frac{k_n}{k_t}\right), \quad (3.20)$$

using normal and tangential coordinates. From here, we first integrate over the tangential momenta, look at small q_n and obtain the long wavelength behavior. We find that in the Fermi Liquid phase [$\delta(k) = \text{const}$]

$$\Gamma_S(\omega) \sim \frac{1}{N(0)} \left(\frac{\omega}{v_F}\right)^2 |\ln \omega / \Lambda v_F|, \quad (3.21)$$

while at the nematic QCP, using $\delta = -\kappa k^2/2$,

$$\Gamma_S(\omega) \sim \frac{1}{N(0)} \left(\frac{\omega}{\kappa v_F}\right)^{2/3}. \quad (3.22)$$

In the Fermi liquid, the energy of the quasiparticle is proportional to ω and we classify them as long lived. However, at the nematic QCP, one can show via Kramers-Kronig that the real part of the self-energy also goes like $|\omega|^{2/3}$. Hence at the nematic QCP, the lifetime of a quasiparticle is not well defined and to try to understand it as a perturbation about a free theory of long-lived quasiparticles is meaningless.

Thus, the bosonized theory reproduces the results found earlier on perturbatively by Oganessian and co-workers¹⁹ (see also Refs. 20 and 40). It should be noted that the $\omega^{2/3}$ law was also found to appear in the perturbative calculation of the fermion self-energy in a model of holes interacting via a forward scattering $U(1)$ gauge interaction with a similar form to (3.15) in the context of high temperature superconductors,^{44,45} and in the perturbative treatment of the quantum critical point in a ferromagnetic metal.³³

C. Nonperturbative results

We now return to the nonperturbative bosonized expression for the fermion propagator of Eq. (3.8), which we will write as

$$G_{F(S)}(\mathbf{x}, t) = Z_S(\mathbf{x}, t) G_{F(S)}^0(\mathbf{x}, t). \quad (3.23)$$

In the *Fermi liquid phase* and at long distances and low frequencies, the factor Z_S approaches a constant value, $Z_S = Z_F < 1$, i.e., the quasiparticle residue of the Fermi liquid state. Our goal here is to investigate the behavior of $Z_S(\mathbf{x}, t)$ near the nematic QCP and in the nematic phase. However, given the complexity of the full analytic expression, in this paper we will consider only on the equal-time, $t=0$, behavior (sometimes called the ‘‘one-particle density matrix’’) and the equal-position, $\mathbf{x}=0$, dynamical correlation function, and only at zero temperature. We will discuss its full behavior elsewhere.

1. The equal-time fermion propagator

As in the perturbative calculation, it is convenient to express $\ln Z$ as an integral over the spectral functions $B_2^\pm(q, s)$. Once again, at the *nematic QCP*, the important contribution is due to the overdamped collective mode in $B_2^+(q, s)$, and we will neglect all other contributions. [This is an approximation which gives the longtime behavior accurately. An expression valid for all times also includes the contribution of $B_2^-(q, s)$.] Since we have set $t=0$ here, the result is quite simple,

$$\ln Z(x_n, 0) = \int_0^{\Lambda/2} \frac{dk_n}{2\pi} I_2(k_n) [\cos(k_n x_n) - 1], \quad (3.24)$$

where

$$I_2(k_n) = \int_{-\Lambda/2}^{\Lambda/2} \frac{dk_t}{2\pi} \int_0^\infty \frac{d\omega}{2\pi} \frac{B_2^+(k, \omega/kv_F)}{(\omega + v_F k_n)^2}. \quad (3.25)$$

We find, both from a numerical computation and from an analytic estimate, that

$$I_2(k_n) = \frac{1}{N(0)v_F} \frac{f_2(k_n\sqrt{\kappa})}{(k_n\sqrt{\kappa})^{4/3}}, \quad (3.26)$$

where $f_2(k_n\sqrt{\kappa})$ is a regular function of its argument. From this analysis, we conclude after performing the final Fourier transformation,

$$Z_S(x_n, 0) = C_P \exp \left\{ - \frac{b}{N(0)v_F\sqrt{\kappa}} \left| \frac{x_n}{\sqrt{\kappa}} \right|^{1/3} \right\}, \quad (3.27)$$

valid for $|x_n| \gg \sqrt{\kappa}$. Here $b=0.0658$ and C_P is a constant factor resulting from subdominant terms in $I_2(k_n)$. This sharp decay of $Z_S(x_n, 0)$, faster than any power law, introduces a scale (similar to a correlation length arising from a gap in the spectrum) and dominates over the Fermi liquid behavior at low energies. From the above expression, the correlation length is of order $\xi \sim \sqrt{\kappa}(k_F\sqrt{\kappa})^3$ which is much longer than length of the interactions, $\sqrt{\kappa}$.

Let us also compute the behavior of Z_S on the *nematic phase* by focusing on the effect of the Goldstone modes. Recalling our discussion of the order parameter theory, we replace (3.15) with

$$V_{s,s}(k, \varepsilon) = \frac{\sin^2(2\theta_S)}{N(0)} \frac{1}{-\kappa k^2/2 + i|s|\sin^2(2\phi)}, \quad (3.28)$$

which is the contribution from the Goldstone mode ($K_0(s) \approx -i|s|$). In the limit, $k_n \rightarrow 0$, that is on the Fermi surface, $\phi \rightarrow \theta_S + \pi/2$ and therefore, we simply have

$$V_{s,s}(k, \varepsilon) = \frac{1}{N(0)} \left(- \frac{\kappa k^2}{\sin^2(2\theta_S)} - K_0(s) \right)^{-1}, \quad (3.29)$$

so that the difference between this and the symmetric side is simply that $\kappa \rightarrow \kappa/\sin^2(2\theta_S)$. Hence, we may directly write down Z_S

$$Z_S(x_n, 0) = C_{P(S)} \exp \left\{ - \frac{b|\sin(2\theta_S)|^{4/3}}{N(0)v_F\sqrt{\kappa}} \left| \frac{x_n}{\sqrt{\kappa}} \right|^{1/3} \right\}, \quad (3.30)$$

where b is the same constant of Eq. (3.27). In Eq. (3.30) we have not included the subdominant contributions which become the leading terms along the symmetry-dictated directions, the ‘‘nematic axes,’’ along which the angular factor vanishes. Along the nematic axes the behavior of the equal-time correlation function has a more Fermi liquid-like long distance behavior as shown by Oganessian *et al.*, but due to the introduction of patches, we cannot accurately capture this behavior here.

Thus, we see that on the broken symmetry side, we have special points at $\theta_S = n\pi/2$ where the Goldstone mode weakens and subdominant behavior takes over. At these points, $Z_S = C_{P(S)} = Z_F < 1$ and the quasiparticles become long lived. It is interesting to note that the angular dependence

shown here with a power of 4/3 is similar to that of the perturbative calculation using the same transformation of $\kappa \rightarrow \kappa/\sin^2(2\theta_S)$ on Eq. (3.22) and it agrees with the results of Ref. 19.

2. The fermion residue

One simple calculation we can do with the above result for $Z_S(x_n, 0)$ is the fermion residue following Migdal:⁴⁶

$$Z_q = n(\mathbf{k}_F - \mathbf{q}) - n(\mathbf{k}_F + \mathbf{q}) \quad (3.31)$$

$$= \int \frac{d\omega}{2\pi} (G_F(\mathbf{k}_F - \mathbf{q}, \omega) - G_F(\mathbf{k}_F + \mathbf{q}, \omega)) e^{i\varepsilon\omega}, \quad (3.32)$$

where the exponential factor tells us to close the contour in the upper half plane. This expression may be written in terms of the real-space, real-time fermion Green function, at time $t = -\varepsilon$. Inserting our expression for G_F within Bosonization and neglecting interpatch scattering (which should produce analytic in q/Λ contributions here), we obtain

$$Z_{q_n} = \frac{2}{\pi} \int_0^\infty dx_n \frac{\sin q_n x_n}{x_n} Z_S(x_n, 0). \quad (3.33)$$

In the *Fermi liquid phase*, we find $Z_S(x_n, 0) = Z < 1$ in the long distance limit and this leads directly to $Z_{q_n} = Z$ when $q_n \rightarrow 0$ as expected.

At the *nematic QCP* and into the nematic phase away from the nodal points discussed earlier, $Z_S(x_n, 0)$ is short ranged and we may expand to leading order in q_n and obtain

$$Z_{q_n} \approx \frac{12}{\pi} q_n \xi \quad (3.34)$$

with the correlation length $\xi = \sqrt{\kappa}(\sqrt{\kappa}k_F/2\pi b)^3$. Thus the fermion residue vanishes linearly similar to how it would in a Fermi liquid at finite temperature where temperature makes the correlations short ranged. It should be noted, however, that because $Z_S(x_n, 0)$ decays slower than $e^{-\alpha x_n}$ in the long distance limit, this series expansion that we have used is poorly defined at higher order with the coefficient of q_n^{2j+1} growing so rapidly that the Taylor expansion has zero radius of convergence in the complex- q_n plane. Thus Z_{q_n} is not analytic in q_n and the Fermi surface may still be defined as a singular point in $n(\mathbf{k})$. The same behavior occurs in the *nematic phase* for generic momenta, except along the directions of the nematic principal axes where a finite residue is obtained.

3. The fermion autocorrelation function

This case turns out to be more complicated than the equal-time expression and we present a full analysis in Appendix D. By a simple integration by parts, we found that the double pole of Eq. (3.25) may be reduced to a single pole and that the integrals involved were less singular by a full power [diverging like $1/k_n^{1/3}$, unlike Eq. (3.26)] but with logarithmic corrections. From that analysis, we found the general form of Z_S to be

$$Z_S(0,t) = C_P \exp \left\{ -A(1-i\sqrt{3}) \frac{\ln(v_F t / \sqrt{\kappa})}{(v_F t / \sqrt{\kappa})^{2/3}} \right\}. \quad (3.35)$$

In contrast with our result for the equal-time correlation function, Eq. (3.35) approaches a constant at long times. However, it decays to that constant much more slowly than in a Fermi liquid where we would expect the exponent 2/3 to become 2. As a result, the nonperturbative effects are here less important and, consequently, this time dependence appears to exhibit the same power-law behavior as the lifetime calculated perturbatively in Sec. III B.

The coefficient of the exponential was found to be $A \propto 1/(k_F \lambda \kappa)$ and deserves some attention. In a Fermi liquid, we would find $A \propto \lambda/k_F \ll 1$ and the time dependence would present a small correction from the Fermi liquid behavior. However, A need not be small if $1/\kappa \gg k_F \lambda \sim \Lambda^2$, where Λ is the patch width [cf. Eq. (3.17)]. Due to the emergence of Λ here, this limit occurs precisely where the Fermi-surface curvature begins to matter, and where the interaction length scale is still quite small ($k_F^{-1} \ll \sqrt{\kappa} \ll \Lambda^{-1}$).

In the nematic phase, again letting $\kappa \rightarrow \kappa/\sin^2 2\theta_S$, we find

$$Z_S(0,t) = C_{P(S)} \exp \left\{ -A(1-i\sqrt{3}) \frac{\ln \left(\frac{v_F t}{\sqrt{\kappa}} |\sin 2\theta_S| \right)}{(v_F t / \sqrt{\kappa})^{2/3}} \right\}, \quad (3.36)$$

so that the angular dependence is similar to that of the equal-time behavior.

4. The one-particle density of states

We close this section with an application of these results to the calculation of the one-particle density of states (DOS) in both the Fermi liquid and the nematic phases, and at the nematic (Pomeranchuk) quantum critical point.

The one-particle DOS is defined by the standard expression

$$N^*(\omega) = -\text{sign}(\omega) \frac{1}{\pi} \text{Im} G_F(x,x;\omega). \quad (3.37)$$

In *Fermi liquid phase*, using $Z_S(0,t) = C_F$, a constant dependent upon the Fermi liquid parameters that goes to 1 for the noninteracting case, as expected we find

$$\begin{aligned} G_F(x,x,\omega) &= \sum_S \frac{1}{N} C_F \int dt G_{F(S)}^0(0,t) e^{i\omega t} \\ &= C_F \int dt \frac{N(0)v_F}{-v_F t + ia \text{sign}(t)} e^{i\omega t} \\ &= -i\pi \text{sign}(\omega) C_F N(0), \end{aligned} \quad (3.38)$$

so that here, $N^*(\omega) = C_F N(0)$ as expected. Note: C_F is not the fermion residue and may be greater than 1.

At the *nematic QCP*, we have instead

$$\begin{aligned} G_F(x,x,\omega) &= \sum_S \frac{1}{N} \int_{-\infty}^{\infty} dt C_P e^{i\omega t} \\ &\times \exp \left\{ -A(1-i\sqrt{3}) \frac{\ln(v_F t / \sqrt{\kappa})}{(v_F t / \sqrt{\kappa})^{2/3}} \right\} G_{F(S)}^0(0,t), \end{aligned} \quad (3.39)$$

so that $N^*(\omega) = C_P N(0) I(\bar{\omega})$ with

$$I(\omega) = \frac{2}{\pi} \mathcal{R}e \int_0^{\infty} \frac{du}{u} \sin(\bar{\omega}u) \exp \left\{ -A(1-i\sqrt{3}) \frac{\ln u}{u^{2/3}} \right\}, \quad (3.40)$$

with $\bar{\omega} = \sqrt{\kappa}\omega/v_F$, valid for $\omega \ll v_F \min\{\kappa^{-1/2}, \lambda\}$. Notice that we have used throughout these expressions only the long time limit of the exponential factor. At shorter times the behavior of the exponential should be dominated by high energy effects which are insensitive to whether the system is in a Fermi liquid, a quantum critical point, or in a nematic phase. Thus, the time integrals have an implicit short distance cutoff (which we have denoted by "0"). In any event, we are only interested in the low frequency behavior which is dominated by the long time part of the integration range.

By inspection we see that the exponential factor approaches unity (quite rapidly) for large $u \gg 1$. Thus, the main effect of this factor is a correction to $I(\omega)$ away from its value at zero frequency i.e., $I(0) = 1$. The leading finite frequency behavior, as $\omega \rightarrow 0$, is obtained by expanding the exponential factor in Eq. (3.40)

$$\begin{aligned} I(\omega) &= 1 - \frac{2A}{\pi} \int_0^{\infty} \frac{du}{u} \frac{\sin \bar{\omega}u}{u^{2/3}} \ln u + \dots \\ &= 1 + A \frac{3\sqrt{3}}{2\pi} \Gamma(1/3) \left(\frac{\sqrt{\kappa}\omega}{v_F} \right)^{2/3} \ln \left(\frac{\sqrt{\kappa}\omega}{v_F} \right) + \dots, \end{aligned} \quad (3.41)$$

where the last line is accurate for $\sqrt{\kappa}\omega/v_F \ll 0.1$. The ellipsis in Eq. (3.41) represents subdominant contributions at low frequencies, which vanish faster than $\omega^{2/3} \ln \omega$ as $\omega \rightarrow 0$. As a result, the $\omega^{2/3}$ behavior of the inverse lifetime $\Gamma(\omega)$ calculated perturbatively, appears here as a cusp in the DOS (with a logarithmic correction). Unlike the lifetime, though, here A depends on the product of the patch width cutoff $\sqrt{k_F \lambda} \sim \Lambda$ and $\sqrt{\kappa}$ while $\Gamma(\omega)$ depends only on the product of k_F and $\sqrt{\kappa}$.

Thus we find that the low frequency oneparticle density of states has the form

$$N^*(\omega) = N^*(0) + B\omega^{2/3} \ln \omega + \dots, \quad (3.42)$$

where $N^*(0) = C_P N(0) > N(0)$, since we found the constant $C_P > 1$ [see Eq. (D17)], and $B = A C_P N(0) (3\sqrt{3}/2\pi) \Gamma(1/3)$ [see Eq. (D18)]. Hence, the zero frequency value of the density of states is *larger* than the Fermi liquid value. As the frequency increases, the one-particle density of states decreases from its zero frequency value according to the $\omega^{2/3} \ln \omega$ correction term. This is a cusp singularity at $\omega = 0$. It is important to stress that we obtained these expressions

upon expanding the exponential factor in the autocorrelation function. This is consistent since this factor asymptotically (and rather rapidly) approaches 1 at long times. This also implies that, in this regime, our results should be consistent with the behavior of the fermion Green function found in perturbation theory around Hartree-Fock/RPA.^{19,20,33,40}

In the nematic phase, an angular average of Eq. (3.36) enters our expression for $N^*(\omega)$. In the long time limit, when the argument of the exponential is much less than one, we expect that for low frequencies

$$A \rightarrow \langle |\sin 2\theta_s|^{4/3} \rangle A \approx 0.58A. \quad (3.43)$$

Hence, in the nematic phase, the Goldstone modes continue the critical behavior of this function in only a mildly weaker form.

In summary, in this section we used bosonization to compute the nonperturbative behavior of the fermion propagator using the bosonized form of the fermion operator. We first checked that the non-Fermi liquid behavior of the nematic QCP and in the nematic phase, which were obtained earlier using conventional diagrammatic (perturbative) methods,^{19,20} is recovered here upon expanding the bosonized expression to leading order in V_{SS} , i.e., a single boson exchange. However, upon a closer examination of the full bosonized result we found that the equal-time fermion correlation function has a much more singular behavior that could have been predicted in perturbation theory. In contrast, the fermion autocorrelation function (and hence the oneparticle density of states) is seemingly consistent with the perturbative analysis of the quantum critical behavior.

We note here that Chubukov³³ has analyzed the quantum critical behavior of ferromagnetic Fermi liquids and claims that the $\omega^{2/3}$ behavior, found at lowest order, persists to all orders in perturbation theory. The results of this section for the fermion autocorrelation function appear to agree with those of Chubukov and co-workers. However our results for the equal-time fermion correlator apparently disagree with these results. Clearly, a more detailed analysis of the bosonized expression for this propagator is warranted. We will discuss this problem in a separate publication.

IV. CONCLUSION

In this paper, we have utilized the method of high dimensional bosonization to study nonperturbatively the quantum phase transition from a Landau Fermi liquid state to a nematic phase, a nematic (Pomeranchuk) instability. For this purpose, we have constructed an order parameter theory from the boson theory by integrating out noncritical modes and verified that this boson theory is equivalent to RPA. We then turned to studying the bosonization form of the fermion propagator and found its diagrammatic expansion, proving the correctness and clarifying the arguments leading up to that expression. This diagrammatic expansion keeps all diagrams up to the simple bubble in the spirit of RPA as applied to the density-density propagator and shows, in specific, that bosonization goes beyond the self-consistent Born approximation to include vertex corrections. We then found explicitly that bosonization reproduces the results of Hartree-Fock

with an RPA interaction by showing that the lifetime computed in this limit also has an $|\omega|^{2/3}$ dependence as originally found for the case of the nematic QCP by Oganesyan and co-workers.¹⁹

Lastly, we calculated the fermion propagator nonperturbatively and found the dramatic effect of the overdamped critical mode that induces short-ranged spatial correlations that decay nearly exponentially $(1/|x|)e^{-\text{const}|x|^{1/3}}$, while the autocorrelation function exhibits a milder non-Fermi liquid behavior of the form $(1/|t|)\exp(-\text{const}|t|^{-2/3})$. From this short-ranged behavior of the equal time Green function, we verify that the fermion residue vanishes at the critical point and into the nematic phase except at four special points. We also calculated the one-particle (fermion) density of states $N^*(\omega)$. We found the low frequency behavior $N^*(\omega) = N^*(0) + B\omega^{2/3}\ln\omega$ [with $N^*(0) > N(0)$] both at the quantum critical point and into the nematic phase. Thus, the fermion propagator exhibits unexpected behaviors which could not have been anticipated by the existing perturbative results.^{17,19,20,33,38} In a separate publication we will present a more detailed analysis of the fermion spectral function in both phases and at finite temperature.

Two recent papers, one by Yang³⁹ and another by Nilsson and Castro Neto,³⁸ also study Pomeranchuk nematic instabilities in two-dimensional Fermi systems. Yang also derives an order parameter theory within high dimensional bosonization. However, contrary to our results, concludes that the critical mode has $z=2$ and is undamped. While we agree that a $z=2$ propagating mode does exist at the critical point, we find that the spectral function is completely dominated by the overdamped $z=3$ mode. This effect is due to Landau damping, a consequence of the curvature of the Fermi surface, and dominates the low energy behavior of the theory at the critical point. The reason for this disagreement is that in Ref. 39 the effects of Landau damping are ignored. We find that these effects are crucial.

On the other hand, Nilsson and Castro Neto³⁸ approach the nematic quantum phase transition within the more traditional methods found in the Fermi liquid theory literature. They first find an order parameter theory by constructing a path integral, whose classical equations of motion give the collisionless Boltzmann equation as in high dimensional bosonization, and integrating out the noncritical modes. Their results, however, agree with ours in all essential details, including the existence of a $z=3$ mode. They then calculate the fermion lifetime using the Bethe-Salpeter equations and Fermi's Golden rule, finding that $\tau^{-1} \propto \epsilon^{2/3}$, as in (3.22), and agree with us, in the perturbative regime, in concluding that this represents both a breakdown of Fermi liquid theory and perturbation theory.

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APPENDIX A: SUMMARY OF BOSONIZATION IN D-DIMENSIONAL FERMI SYSTEMS

Consider the Fermi liquid theory of spinless fermions interacting via a short but perhaps finite range forward scattering interaction living in a translationally invariant D -dimensional world. This is a low energy theory and as such, following Landau, we shall linearize the energy dispersion near the Fermi surface, though corrections to this may be considered when necessary. To this end, let us build a construction in which we linearize within N equally sized patches approximating the Fermi surface. For this construction to be reasonable, our end result should be relatively insensitive to the details belonging to this partitioning. Keep in mind, however, that a remarkable property of Fermi liquids is that they only require a few of the lowest angular momentum Fermi liquid parameters to understand a wide range of phenomena.

The number of patches, N , approximating the Fermi surface will naturally be inversely proportional to its curvature. As such, if the density, $n \rightarrow \infty$ then $N \rightarrow \infty$ and the construction becomes exact. An exact solution to leading order in N of this linearized theory is therefore equivalent to the asymptotic low energy limit as dictated by the renormalization group. More specifically, the number of patches N and the patch width $\Lambda \sim \sqrt{k_F \lambda}$, where λ is the energy cutoff, must be related by the condition $2\pi k_F = N\Lambda$, required for the Fermi system to have a finite density of states and curvature (see below). Consequently, the number of patches N must scale as $N \sim 2\pi\sqrt{k_F/\lambda}$. Clearly $N \rightarrow \infty$ in the infrared limit $\lambda \rightarrow 0$. Many of the results of this paper can be understood in light of this basic reasoning.

Under our construction, the fermion annihilation operator becomes

$$\hat{c}(\mathbf{x}) = \sum_{\mathbf{k}} \hat{c}_{\mathbf{k}} \frac{e^{i\mathbf{k}\cdot\mathbf{x}}}{\sqrt{L^D}} = \sum_S \sum_{\mathbf{q} \in \mathcal{P}_S} \hat{c}_{\mathbf{k}_S + \mathbf{q}} \frac{e^{i(\mathbf{k}_S + \mathbf{q})\cdot\mathbf{x}}}{\sqrt{L^D}} = \sum_S \hat{c}_S(\mathbf{x}) \frac{e^{i\mathbf{k}_S\cdot\mathbf{x}}}{\sqrt{N}}, \quad (\text{A1})$$

where S labels the patch, \mathcal{P}_S is the volume in \mathbf{k} space around the point \mathbf{k}_S and the new fermion operators, $\psi_S(\mathbf{x})$, obey the canonical commutation relations

$$\{c_S(\mathbf{x}), c_T^\dagger(\mathbf{x}')\} = \delta_{S,T} \delta^D(\mathbf{x} - \mathbf{x}'). \quad (\text{A2})$$

It is also important to note the Fourier transform normalizations within the construction

$$\begin{aligned} \delta^D(\mathbf{x} - \mathbf{x}') &= \delta_t^{D-1}(\mathbf{x}_t - \mathbf{x}'_t) \delta(x_n - x'_n) \\ &= \left(N \sum_{|\mathbf{q}_t| < \frac{\Lambda}{2}} \frac{e^{i\mathbf{q}_t \cdot (\mathbf{x}_t - \mathbf{x}'_t)}}{L^{(D-1)}} \right) \left(\sum_{|\mathbf{q}_n| < \frac{\Lambda}{2}} \frac{e^{iq_n(x_n - x'_n)}}{L} \right), \end{aligned} \quad (\text{A3})$$

where we have introduced Λ to characterize the tangential width of the patch ($N\Lambda$ is the area of the Fermi surface) and λ as an ultraviolet cutoff about the Fermi surface.

Now, in terms of our fermion operators, the linearized Hamiltonian is

$$:\hat{H}:= = \sum_S \int d^x (\mathcal{H}_S^0(\mathbf{x}) + \mathcal{H}_S^{\text{Int}}(\mathbf{x})), \quad (\text{A4})$$

where the free Hamiltonian density is

$$\mathcal{H}_0(\mathbf{x}) = \frac{\hbar}{2i} \mathbf{v}_S \cdot (:\nabla \hat{c}_S^\dagger(\mathbf{x})]: \hat{c}_S(\mathbf{x}) - \hat{c}_S^\dagger(\mathbf{x}) \nabla \hat{c}_S(\mathbf{x})): \quad (\text{A5})$$

and the forward scattering interactions are described by

$$\mathcal{H}_S^{\text{Int}}(\mathbf{x}) = \sum_T \int d\mathbf{x}' F_{S-T}(\mathbf{x} - \mathbf{x}') \delta \hat{n}_S(\mathbf{x}) \delta \hat{n}_T(\mathbf{x}'). \quad (\text{A6})$$

Here, the density fluctuations are defined by

$$\delta \hat{n}_S(\mathbf{x}) \equiv : \hat{n}_S(\mathbf{x}) : = : \hat{c}_S^\dagger(\mathbf{x}) \hat{c}_S(\mathbf{x}) :, \quad (\text{A7})$$

and throughout this description we have been using the usual normal ordering procedure for any operator \mathcal{O}

$$:\mathcal{O}:|G\rangle = 0 \rightarrow \mathcal{O} = : \mathcal{O} : - \langle G | \mathcal{O} | G \rangle, \quad (\text{A8})$$

where we take the filled Fermi sea as our ground state

$$|G\rangle = \prod_S \prod_{\{\mathbf{q} \in \mathcal{P}_S | \mathbf{v}_S \cdot \mathbf{q} < 0\}} \hat{c}_{\mathbf{q},S}^\dagger |0\rangle. \quad (\text{A9})$$

It was shown by Haldane,^{10,11} Castro Neto and Fradkin,¹²⁻¹⁴ and Houghton and Marston^{15,16} that this Hamiltonian can be entirely described in terms of the electron density operators, $\delta \hat{n}_S(\mathbf{x})$ in the high density limit and it is quadratic in these operators. A fermion operator may then be constructed following well-known one-dimensional (1D) bosonization techniques and so the theory can be solved exactly. (This represented a major step forward since the introduction of RPA by Bohm and Pines.⁴) Here we shall outline the proof of this solution, but with the traditional approach of point-splitting regularization, commonly used in the one-dimensional case (see, for example Ref. 47).

The expectation value of the density operator $\hat{n}_S(\mathbf{x})$ in the ground state of the Fermi sea is clearly divergent if we send the density of fermions to infinity. As a result, in the high density limit we are interested in, it is poorly defined. To control this divergence, we introduce the point-split operator

$$\hat{n}_S^\epsilon(\mathbf{x}) \equiv \hat{c}_S^\dagger(\mathbf{x} + \epsilon/2) \hat{c}_S(\mathbf{x} - \epsilon/2) = : \hat{n}_S^\epsilon(\mathbf{x}) : - \langle G | \hat{n}_S^\epsilon(\mathbf{x}) | G \rangle. \quad (\text{A10})$$

Here, by short distance it is meant a length scale short compared with the separation of all operators of interest but long compared with physical short length scales, i.e., It should be

noted that physically point-splitting can be thought of as a means of $|\mathbf{x}| \gg |\epsilon| \gg \lambda^{-1}$.

The divergent part may be computed explicitly

$$\begin{aligned} \langle G | \hat{n}_S^\epsilon(\mathbf{x}) | G \rangle &= \frac{N}{L^D} \sum_{\mathbf{q}_t} \sum_{q_n < 0} e^{-i\mathbf{q}_t \cdot \epsilon} \\ &= \delta_t^{D-1}(\epsilon_t) \int_{-\infty}^0 \frac{dq_n}{2\pi} e^{-iq_n \epsilon_n + q_n/\lambda} \\ &= \frac{i \delta_t^{D-1}(\epsilon_t)}{2\pi(\epsilon_n + i\lambda^{-1})}, \end{aligned} \quad (\text{A11})$$

where we have implemented the ultraviolet cutoff, λ , as a soft cutoff $e^{-|q_n|/\lambda}$. This is a highly anisotropic expression in ϵ . It vanishes if we first send $\epsilon_n \rightarrow 0$, looking along a tangential direction ($|\epsilon_t| \gg \Lambda^{-1} \rightarrow 0$) but diverges if we first send $\epsilon_t \rightarrow 0$ looking along $\epsilon_n \gg \lambda^{-1} \rightarrow 0$. Therefore, to capture the basic physics of the density operator we must choose the latter limit as the definition of the point-split operator. Keeping in mind that $\delta_t(0) = A_F / (2\pi)^{D-1} = 2\pi \hbar v_F N(0)$, where A_F is the $(D-1)$ -dimensional Fermi surface area and $N(0)$ is the density of states at the Fermi surface, we obtain

$$\hat{n}_S^\epsilon(\mathbf{x}) = -\frac{iN(0)\hbar v_F}{\epsilon_n} + \delta \hat{n}_S(\mathbf{x}) + \frac{i\epsilon_n}{\hbar v_F} \mathcal{H}_S^0(\mathbf{x}) + \dots \quad (\text{A12})$$

to leading order in $\epsilon_n \gg \lambda^{-1}$ and where we kept the expansion of $:\hat{n}_S^\epsilon(\mathbf{x}):$ to first order noticing the useful emergence of the free Hamiltonian density operator.

Now that we have a controlled definition of the density operator, we proceed with computing its commutator,

$$\begin{aligned} &[\hat{n}_S^{\epsilon_n}(\mathbf{x}), \hat{n}_T^{\epsilon'_n}(\mathbf{x}')] \\ &= \frac{iN(0)\hbar v_F}{\epsilon_n + \epsilon'_n} \delta_{S,T} \delta_t^{D-1}(\mathbf{x}_t - \mathbf{x}'_t) (\delta[x_n - x'_n - (\epsilon_n + \epsilon'_n)/2] \\ &\quad - \delta[x_n - x'_n + (\epsilon_n + \epsilon'_n)/2]). \end{aligned} \quad (\text{A13})$$

Expanding both sides of this equation in powers of ϵ_n and equating like powers gives us the following result:

$$[\delta \hat{n}_S(\mathbf{x}), \delta \hat{n}_T(\mathbf{x}')] = -i \hbar N(0) \delta_{S,T} \mathbf{v}_S \cdot \nabla \delta^D(\mathbf{x} - \mathbf{x}'), \quad (\text{A14})$$

$$[\delta \hat{n}_S(\mathbf{x}), \mathcal{H}_T^0(\mathbf{x}')] = -i \hbar \mathbf{v}_S \cdot \nabla \delta \hat{n}_S(\mathbf{x}) \delta^D(\mathbf{x} - \mathbf{x}'). \quad (\text{A15})$$

Using these commutators, we may compute the equation of motion for the Heisenberg operator $\delta \hat{n}_S(\mathbf{x}, t)$,

$$\begin{aligned} &\partial_t \delta \hat{n}_S(\mathbf{x}, t) + \mathbf{v}_S \cdot \nabla \delta \hat{n}_S(\mathbf{x}, t) + \mathbf{v}_S \\ &\cdot \nabla \sum_T \int d\mathbf{x}' F_{S-T}(\mathbf{x} - \mathbf{x}') \delta \hat{n}_T(\mathbf{x}', t) = 0, \end{aligned} \quad (\text{A16})$$

where $F_{S-T} = N(0) f_{S-T}$. This is the linearized collisionless Boltzmann equation in operator form found by Castro Neto and Fradkin in the context of a coherent state formalism.¹² We also notice through this derivation that we may let

$$\mathcal{H}_S^0(\mathbf{x}) = \frac{1}{2N(0)} \delta \hat{n}_S^2(\mathbf{x}) \quad (\text{A17})$$

and obtain exactly the same answer. Hence, the Hamiltonian may be expressed entirely in terms of the density operator $\delta \hat{n}_S(\mathbf{x})$.

A natural consequence of (A14) is that the density operator may be expressed in terms of a chiral boson field,

$$\delta \hat{n}_S(\mathbf{x}) = N(0) \mathbf{v}_S \cdot \nabla \hat{\phi}_S(\mathbf{x}), \quad (\text{A18})$$

with a canonically conjugate momentum

$$\hat{\pi}_S(\mathbf{x}) = -N(0) \mathbf{v}_S \cdot \nabla \hat{\phi}_S(\mathbf{x}). \quad (\text{A19})$$

These chiral bosons are a direct extension of the right/left chiral bosons in the context of 1D bosonization. Following this extension then, we may express the fermion operator as a vertex operator

$$\hat{\psi}_S(\mathbf{x}) = \eta_S(\mathbf{x}_t) \sqrt{N(0)v_F\lambda} e^{-i\hat{\phi}_S(\mathbf{x})/\hbar}, \quad (\text{A20})$$

where $\eta_S(\mathbf{x}_t)$ is a set of Klein factors responsible for ensuring that $\hat{\psi}_S(\mathbf{x})$ obey the proper anticommutation relations within the patch and on different patches. The relation between this fermion operator and the original $\hat{c}_S(x)$ operators will be made precise in Sec. III A via direct comparison of the perturbation series in $F_{S-T}(\mathbf{x} - \mathbf{x}')$ obtained in the bosonized theory. $\hat{\psi}_S(\mathbf{x})$, therefore, is equivalent to $\hat{c}_S(\mathbf{x})$ in the free case and in the interacting case, it is this operator projected onto the high-density subspace. Hence, this is an effective low energy theory of a dense Fermi system and $\hat{\psi}_S^\dagger(\mathbf{x})$ can be viewed as actually creating Landau quasiparticles in the Fermi liquid phase.

The canonical structure of the bosonized theory also allows us to directly write down a path integral formulation of the problem, including interactions described by a set of Landau parameters (for a derivation using coherent states, see Ref. 12). The action for the bosonized theory has the general form

$$\begin{aligned} S &= \frac{N(0)}{2} \sum_S \int d^2x dt [-\partial_t \varphi_S \mathbf{v}_S \cdot \nabla \varphi_S - (\mathbf{v}_S \cdot \nabla \varphi_S)^2] \\ &\quad + \frac{N(0)}{2} \sum_{S,T} \int d^2x d^2x' dt \\ &\quad \times F_{S-T}(\mathbf{x} - \mathbf{x}') \mathbf{v}_S \cdot \nabla \varphi_S(\mathbf{x}) \mathbf{v}_T \cdot \nabla \varphi_T(\mathbf{x}'). \end{aligned} \quad (\text{A21})$$

This action is a quadratic form in the Bose fields. Here we have not included higher order terms (such as those discussed in the body of the paper, in the context of the nematic instability). Such terms are generally present due to nonlinearities in the fermion dispersion relation, as well as many-body effective interactions.³⁶ We have also not included ‘‘vertex operators’’ such as those associated with pairing (BCS) interactions.^{12,16} We shall find it more convenient to work within this path-integral formulation when constructing a theory of the nematic quantum critical point.

Before leaving our discussion of high dimensional bosonization, we should make a final comment on its validity. As discussed in the main body of the paper, the boson theory here completely recovers the random phase approximation (RPA) in the long wavelength limit. This should not be surprising since, in the asymptotic low energy limit, both RPA and bosonization saturate the f -sum rule and are (formally) exact. This is a well-known established property of bosonization, extensively discussed in the literature since the 1970's.

APPENDIX B: Analysis of V_2^+ and V_2^-

Here we present the partial fraction expansion of the effective interactions V_2^+ and V_2^- and their behavior near quantum criticality.

1. Partial fraction expansion of $V_2^+(q, s)$

In terms of the function $K_0(s)$ we may write V_2^+ as

$$V_2^+(q, s) = \frac{1}{N(0)} \left[\delta_2(q) - K_0(s) \left(1 + \left(\frac{1-K_0(s)}{1+K_0(s)} \right)^2 \right) \right]^{-1} \\ = \frac{1}{2N(0)} \left[\frac{(1+x)^2}{x^3 - \frac{\delta_2}{2}x^2 + [1 - \delta_2(q)]x - \frac{\delta_2(q)}{2}} \right], \quad (\text{B1})$$

where $x=K_0(s)$. The denominator is thus a cubic polynomial in $K_0(s)$ and solving for the poles, $x=\delta_\beta^+(q)$, we find

$$\delta_\beta^+(q) = \left\{ \frac{1}{6} \left[\delta_2(q) + \frac{\delta_2(q)^2 + 12\delta_2(q) - 12}{e^{in\pi/3}h(\delta_2(q))} \right. \right. \\ \left. \left. + e^{in\pi/3}h(\delta_2(q)) \right] \right\} \Big| n \in \{-1, 1, 3\}, \quad (\text{B2})$$

where

$$h(x) = [12\sqrt{12 - 36x + 42x^2 + 3x^3} - x(x^2 + 18x + 36)]^{1/3}, \quad (\text{B3})$$

and we assign $\beta=\{a, b, c\}$ to each of these three poles. The partial fraction expansion is

$$V_2^+(q, s) = \frac{1}{N(0)} \sum_{\beta=a, b, c} \frac{\mathcal{Z}_\beta^+(q)}{\delta_\beta^+(q) - K_0(s)}, \quad (\text{B4})$$

with residues

$$\mathcal{Z}_a^+(q) = \frac{1}{2} \frac{(1 + \delta_a^+)^2}{(\delta_a^+ - \delta_b^+)(\delta_a^+ - \delta_c^+)}, \\ \mathcal{Z}_b^+(q) = \frac{1}{2} \frac{(1 + \delta_b^+)^2}{(\delta_b^+ - \delta_a^+)(\delta_b^+ - \delta_c^+)}, \\ \mathcal{Z}_c^+(q) = \frac{1}{2} \frac{(1 + \delta_c^+)^2}{(\delta_c^+ - \delta_a^+)(\delta_c^+ - \delta_b^+)}. \quad (\text{B5})$$

For $|\delta_2(q)| \ll 1$ these formulas simplify to

$$\delta_a^+ \approx i, \quad \delta_b^+ \approx -i, \quad \delta_c^+ \approx \delta_2/2, \quad (\text{B6})$$

$$\mathcal{Z}_a^+ \approx i/2, \quad \mathcal{Z}_b^+ \approx -i/2, \quad \mathcal{Z}_c^+ \approx 1/2, \quad (\text{B7})$$

as a result we may write near the nematic QCP

$$V_2^+(q, s) \approx \frac{1}{2N(0)} \left[\frac{1}{\frac{\delta_2(q)}{2} - K_0(s)} - \frac{1/4}{s^2 - 1/2} \right]. \quad (\text{B8})$$

2. Partial fraction expansion of $V_2^-(q, s)$

We may write V_2^- as

$$V_2^-(q, s) = \frac{1}{N(0)} \left[\delta_2(q) - K_0(s) \left(1 - \left(\frac{1-K_0(s)}{1+K_0(s)} \right)^2 \right) \right]^{-1} \\ = \frac{1}{[4 - \delta_2(q)]N(0)} \left[\frac{1}{(1+x)^2} x^2 - 2 \frac{\delta_2(q)}{4 - \delta_2} x - \frac{\delta_2}{4 - \delta_2} \right]. \quad (\text{B9})$$

The denominator is simply quadratic and we find poles at $x=\delta_\beta^-$ with

$$\delta_a^-(q) = -\frac{\sqrt{\delta_2(q)}}{\sqrt{\delta_2(q)} + 2}, \quad \delta_b^-(q) = -\frac{\sqrt{\delta_2(q)}}{\sqrt{\delta_2(q)} - 2}. \quad (\text{B10})$$

Hence, we can now expand in partial fractions to obtain

$$V_2^-(q, s) = \frac{1}{N(0)} \left(\frac{\mathcal{Z}_a^-(q)}{\delta_a^-(q) - K_0(s)} + \frac{\mathcal{Z}_b^-(q)}{\delta_b^-(q) - K_0(s)} \right), \quad (\text{B11})$$

with residues

$$\mathcal{Z}_a^-(q) = \frac{1}{4 - \delta_2} \left[\frac{(1 + \delta_a^-)^2}{\delta_a^- - \delta_b^-} \right], \quad \mathcal{Z}_b^-(q) = \frac{1}{4 - \delta_2} \left[\frac{(1 + \delta_b^-)^2}{\delta_b^- - \delta_a^-} \right]. \quad (\text{B12})$$

Again, these formulas simplify for $|\delta_2| \ll 1$,

$$\delta_a^- \approx -\frac{\sqrt{\delta_2}}{2}, \quad \delta_b^- \approx \frac{\sqrt{\delta_2}}{2}, \quad (\text{B13})$$

$$\mathcal{Z}_a^- \approx -\frac{1}{4\sqrt{\delta_2}}, \quad \mathcal{Z}_b^- \approx \frac{1}{4\sqrt{\delta_2}}. \quad (\text{B14})$$

Thus, near the nematic QCP we can write

$$V_2^-(q, s) \approx \frac{1}{4N(0)} \left(\frac{1}{s^2 + \frac{\delta_2(q)}{4}} \right). \quad (\text{B15})$$

APPENDIX C: INCLUSION OF $F_0(q)$

Here we calculate $V_{S,T}$ for the case when both F_0 and F_2 are present and we let F_2 approach the nematic QCP

($F_2 \rightarrow -1$). We shall find, however, that the critical behavior is utterly independent of F_0 ! Now, let

$$F_{S-T}(q) = \frac{2}{N}(F_0 + F_2(q)\cos[2(\theta_S - \theta_T)]). \quad (C1)$$

Returning to the σ_ℓ^α theory, whose correlators are the RPA interaction $V_{S,T}$, we find in this case

$$S_\sigma = -\frac{1}{2} \int \frac{d^2 q d\omega}{(2\pi)^2} [\sigma^+(\mathbf{q}, \omega) \cdot (\mathbf{V}^+)^{-1} \cdot \sigma^+(-\mathbf{q}, -\omega) + V_2^- |\sigma_2^-|^2], \quad (C2)$$

where $\sigma^+ = (\sigma_0^+, \sigma_2^+)$ and

$$(\mathbf{V}^+)^{-1} = \begin{pmatrix} 2\left(\chi_0^0 - \frac{1}{F_0}\right) & 2\chi_2^0 \\ 2\chi_2^0 & \chi_0^0 + \chi_4^0 - \frac{1}{F_2} \end{pmatrix}. \quad (C3)$$

As a result, V_2^- is completely unaffected by the presence of F_0 due to the fact that there is no σ_0^- field and opposite “signs” decouple.

Again, we may write this expression as a function of $x=K_0(s)$ and taking the inverse, we find

$$\mathbf{V}^+ = \frac{1}{x^3 + ax^2 + bx + c} \mathbf{W}(x), \quad (C4)$$

where

$$\mathbf{W}(x) = \frac{1}{2\delta_0 + \delta_2 - 4} \times \begin{pmatrix} -x^3 + (\delta_2/2)x^2 - (1 - \delta_2)x + \delta_2/2 & x(1 - x^2) \\ x(1 - x^2) & \delta_0(1 + x^2) \end{pmatrix} \quad (C5)$$

and

$$a = \frac{(2 - \delta_0)\delta_2}{2\delta_0 + \delta_2 - 4}, \quad b = \frac{2\delta_0 + \delta_2 - 2\delta_0\delta_2}{2\delta_0 + \delta_2 - 4}, \quad c = \frac{-\delta_0\delta_2}{2\delta_0 + \delta_2 - 4}, \quad (C6)$$

with $\delta_0 = 1 + 1/F_0$ and $\delta_2 = 1 + 1/F_2$. Again we find that the polynomial is cubic. Hence, the addition of F_0 only complicates the algebra, but not the general structure of the solution. We therefore continue as in Appendix B.

Solving the cubic equation in (C4) leads to an algebraically complicated result. However, it simplifies near the nematic QCP and we find to lowest order in δ_2 ,

$$\delta_a^+ = i\sqrt{\frac{\delta_0}{\delta_0 - 2}}, \quad \delta_b^+ = -i\sqrt{\frac{\delta_0}{\delta_0 - 2}}, \quad \delta_c^+ = \delta_2/2, \quad (C7)$$

which return to our previous result if we let $F_0 \rightarrow 0$ or $\delta_0 \rightarrow \infty$. Notice that, to lowest order δ_c^+ is independent of F_0 , as it should since the s -wave mode is noncritical and this “pole” precisely represents the critical behavior. Expanding in partial fractions leads to the expression

$$\mathbf{V} = \frac{1}{N(0)} \sum_{\beta=a,b,c} \frac{\mathbf{Z}_\beta^+}{\delta_\beta - K_0(s)}, \quad (C8)$$

with residue matrices

$$\mathbf{Z}_a^+ = \frac{i}{2} \sqrt{\frac{\delta_0}{\delta_0 - 2}} \begin{pmatrix} \frac{1}{\delta_0(\delta_0 - 2)} & \frac{\delta_0 - 1}{\delta_0(\delta_0 - 2)} \\ \frac{\delta_0 - 1}{\delta_0(\delta_0 - 2)} & 1 + \frac{i}{\sqrt{\delta_0(\delta_0 - 2)}} \end{pmatrix},$$

$$\mathbf{Z}_b^+ = \frac{-i}{2} \sqrt{\frac{\delta_0}{\delta_0 - 2}} \begin{pmatrix} \frac{1}{\delta_0(\delta_0 - 2)} & \frac{\delta_0 - 1}{\delta_0(\delta_0 - 2)} \\ \frac{\delta_0 - 1}{\delta_0(\delta_0 - 2)} & 1 - \frac{i}{\sqrt{\delta_0(\delta_0 - 2)}} \end{pmatrix},$$

$$\mathbf{Z}_c^+ = \frac{1}{2} \begin{pmatrix} 0 & 0 \\ 0 & 1 \end{pmatrix}, \quad (C9)$$

which again returns to the previous result if we let $F_0 \rightarrow 0$. Notice that \mathbf{Z}_c^+ is independent of F_0 to this order in δ_2 . Consequently, the critical behavior of the fermions, (3.27), is unaffected by the inclusion of a noncritical mode such as F_0 .

APPENDIX D: Equal-Position Boson Propagator

Here we are interested in the quantity

$$G_{B(S,S)}(\mathbf{0}, t) = G_{B(S,S)}^0(\mathbf{0}, t) + i \int \frac{d^2 k d\varepsilon}{(2\pi)^3} G_{F(S)}^0(k, \varepsilon) V_{S,S}(k, \varepsilon) \times G_{F(S)}^0(k, \varepsilon) (e^{-i\varepsilon t} - 1), \quad (D1)$$

the equal position part of Eq. (3.1). In particular, we focus on the second term, labeling it $\ln Z_S(\mathbf{0}, t)$. We begin by writing the interaction in terms of its spectral function:

$$V_{S,S}(k, \varepsilon) = \int \frac{d\varepsilon'}{2\pi} \frac{B(k, \varepsilon')}{\varepsilon - \varepsilon' + i\varepsilon' \text{sign } \varepsilon'}, \quad (D2)$$

where

$$B(\mathbf{k}, \omega) = B_2^+(k, \omega) \cos^2[2(\theta_S - \phi)] + B_2^-(k, \omega) \sin^2[2(\theta_S - \phi)] \quad (D3)$$

and ϕ the direction of \mathbf{k} . Note: in the appropriate scaling within a patch, $\phi \rightarrow \theta_S + \pi/2$ so that the cosine factor scales to 1 while the sine factor scales to 0 driving the $B_2^-(k, \omega)$ contribution irrelevant (this has been checked explicitly).

This allows us to do the ε integration immediately and rewrite our expression in a more physical form. The result is (after letting $\varepsilon' \rightarrow 0$),

$$\begin{aligned} \ln Z_S(\mathbf{0}, |t|) = & \int_{-\Lambda/2}^{\Lambda/2} \frac{dk_t}{2\pi} \int_{-\Lambda/2}^{\Lambda/2} \frac{dk_n}{2\pi} \int_0^\infty \frac{d\varepsilon}{2\pi} B(k, \varepsilon) R_1(k_n, \varepsilon, |t|) \\ & + \int_{-\Lambda/2}^{\Lambda/2} \frac{dk_t}{2\pi} \int_0^\infty \frac{d\varepsilon}{2\pi} B(k, \varepsilon) R_2(k_n, \varepsilon, |t|), \end{aligned} \quad (\text{D4})$$

which we have written directly in terms of the patch coordinates. The residue of the single pole is

$$R_1(k_n, \varepsilon, |t|) = \frac{e^{-i\varepsilon|t|} - 1}{(\varepsilon - v_F k_n + i\varepsilon \text{sign } \varepsilon)^2} \quad (\text{D5})$$

$$= \frac{1}{v_F} \frac{d}{dk_n} \left[\frac{e^{-i\varepsilon|t|} - 1}{\varepsilon - v_F k_n + i\varepsilon \text{sign } \varepsilon} \right], \quad (\text{D6})$$

and the residue of the double pole is

$$R_2(k_n, \varepsilon, |t|) = -\frac{1}{v_F} \frac{d}{dk_n} \left[\frac{e^{-ik_n|t|} - 1}{\varepsilon - v_F k_n + i\varepsilon \text{sign } \varepsilon} \right], \quad (\text{D7})$$

so that it is clear that an integration by parts removes the double pole altogether, leaving us with an integral only over a single pole. One can view this as a cancellation of the double pole's contribution to the integral and expect the result to be less singular than the equal-time case.

Using the symmetries of the spectral function, we may rewrite Eq. (D4) as

$$\begin{aligned} \ln Z_S(\mathbf{0}, |t|) = & \int_{-\infty}^{\infty} \frac{dk_t}{2\pi} \int_0^\infty \frac{dk_n}{2\pi} \int_0^\infty \frac{d\varepsilon}{2\pi} B(k, \varepsilon) [R_1(k_n, \varepsilon, |t|) \\ & + R_1(-k_n, \varepsilon, |t|) + R_2(k_n, \varepsilon, |t|) - R_2(k_n, -\varepsilon, |t|)]. \end{aligned} \quad (\text{D8})$$

Performing the integration by parts, and a quick change of variables allows us to separate the time dependence in terms of the following integrals:

$$\begin{aligned} \ln Z_S(\mathbf{0}, |t|) = & \frac{1}{N(0)v_F\sqrt{\kappa}} \left[-2 \int_0^\infty d\nu \left(I'_\nu(-\nu) - \frac{1}{\lambda\sqrt{\kappa}} [I_\nu(-\nu) + I'_\nu^{(1)}(-\nu)] \right) \right. \\ & + \int_0^\infty du_n \left(I'_u(-u_n) + I'_u(u_n) - \frac{1}{\lambda\sqrt{\kappa}} [I_u(-u_n) + I_u(u_n) + u_n I'_u(-u_n) + u_n I'_u(u_n)] \right) e^{-iu_n\bar{t}} \\ & \left. + \int_0^\infty d\nu \left(I'_\nu(-\nu) - I'_\nu(\nu) - \frac{1}{\lambda\sqrt{\kappa}} [I_\nu(-\nu) - I_\nu(\nu) + I'_\nu^{(1)}(-\nu) - I'_\nu^{(1)}(\nu)] \right) e^{-i\nu\bar{t}} \right], \end{aligned} \quad (\text{D9})$$

where we have sent the cutoffs to infinity, keeping the lowest order correction in λ (having checked that higher orders are non-singular), and where the various integrals are

$$I_\nu(\nu) = \frac{1}{4\pi^3} \mathcal{P} \int_0^\infty du_t \int_0^\infty du_n \frac{B(u, \nu)}{\nu - u_n}, \quad (\text{D10})$$

$$I'_\nu(\nu) = \frac{1}{4\pi^3} \mathcal{P} \int_0^\infty du_t \int_0^\infty du_n \frac{\frac{d}{du_n} B(u, \nu)}{\nu - u_n}, \quad (\text{D11})$$

$$I'_\nu^{(1)}(\nu) = \frac{1}{4\pi^3} \mathcal{P} \int_0^\infty du_t \int_0^\infty du_n \frac{u_n \frac{d}{du_n} B(u, \nu)}{\nu - u_n}, \quad (\text{D12})$$

$$I_u(u_n) = \frac{1}{4\pi^3} \mathcal{P} \int_0^\infty du_t \int_0^\infty d\nu \frac{B(u, \nu)}{\nu - u_n}, \quad (\text{D13})$$

$$I'_u(u_n) = \frac{1}{4\pi^3} \mathcal{P} \int_0^\infty du_t \int_0^\infty d\nu \frac{\frac{d}{d\nu} B(u, \nu)}{\nu - u_n}. \quad (\text{D14})$$

In this latest form of $\ln Z_S$, we have defined $\mathbf{u} = \mathbf{k}\sqrt{\kappa}$, $\nu = \varepsilon\sqrt{\kappa}/v_F$, $\bar{t} = v_F t/\sqrt{\kappa}$ due to the form of the spectral function at a nematic instability

$$B(k, \varepsilon) = \frac{\varepsilon k v_F}{\varepsilon^2 + \kappa^2 v_F^2 k^6/4} = \frac{\nu u}{\nu^2 + u^6/4}, \quad (\text{D15})$$

where we have only included the effects of the overdamped mode since it is responsible for the leading singular behavior. Here we have replaced the angular factor $\cos^2(2\theta_S)$ by a constant of order unity. (This is consistent since we will be interested in the angular-averaged fermion Green function and the angular factor in the exponent vanishes only on a set of measure zero.)

Table I shows the low frequency-long wavelength limit of these integrals computed numerically. Two basic forms emerge: a $\ln^2(\nu)$ divergence and a $\nu^{-1/3}$ power-law divergence (with logarithmic corrections), each form occurring at

TABLE I. Low frequency/long wavelength limit of the required integrals.

	$\lim_{\{\nu, u_n\} \rightarrow 0^-}$	$\lim_{\{\nu, u_n\} \rightarrow 0^+}$
$I_\nu(\nu)$	$-0.0053 \ln(\nu)/\nu^{1/3}$	$0.0053 \ln(\nu)/\nu^{1/3}$
$I'_\nu(\nu)$	$-0.0018 \ln^2(\nu)$	$-0.0018 \ln^2(\nu)$
$I''_\nu(\nu)$	$0.0077/\nu^{1/3}$	$-0.0077/\nu^{1/3}$
$I_u(u_n)$	$0.0280/u_n^{1/3}$	$0.0140/u_n^{1/3}$
$I'_u(u_n)$	$0.0052 \ln^2(u_n)$	$-0.0052 \ln^2(u_n)$

different scales. However, the $\ln^2(\nu)$ contributions all vanish. With these limits in mind, we perform the final integral and obtain

$$Z_S(\mathbf{0}, |t|) = C_P \exp \left\{ -A(1 + \sqrt{3}i) \frac{\ln(b|\bar{t}|)}{|\bar{t}|^{2/3}} \right\}, \quad (\text{D16})$$

valid for $|\bar{t}| \gg 1$. In this expression,

$$C_P = \exp \left\{ \frac{1}{N(0)v_F\sqrt{\kappa}} \left(0.0476 + \frac{0.0391}{\lambda\sqrt{\kappa}} + \dots \right) \right\} > 1, \quad (\text{D17})$$

$$A = \frac{0.00724}{N(0)v_F\kappa\lambda} \propto \frac{1}{k_F\lambda\kappa}, \quad (\text{D18})$$

$$b = -4.95 + 1.51i, \quad (\text{D19})$$

where the numbers calculated for C_P were also computed numerically.

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