

Two-point function of a $d = 2$ quantum critical metal in the limit $k_F \rightarrow \infty$, $N_f \rightarrow 0$ with $N_f k_F$ fixedPetter Sätterskog,^{*} Balazs Meszена,[†] and Koenraad Schalm[‡]*Institute Lorentz Δ ITP, Leiden University, P.O. Box 9506, Leiden 2300 RA, The Netherlands*

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We show that the fermionic and bosonic spectrum of $d = 2$ fermions at finite density coupled to a critical boson can be determined nonperturbatively in the combined limit $k_F \rightarrow \infty$, $N_f \rightarrow 0$ with $N_f k_F$ fixed. In this double scaling limit, the boson two-point function is corrected but only at one loop. This double scaling limit therefore incorporates the leading effect of Landau damping. The fermion two-point function is determined analytically in real space and numerically in (Euclidean) momentum space. The resulting spectrum is discontinuously connected to the quenched $N_f \rightarrow 0$ result. For $\omega \rightarrow 0$ with k fixed the spectrum exhibits the distinct non-Fermi-liquid behavior previously surmised from the RPA approximation. However, the exact answer obtained here shows that the RPA result does not fully capture the IR of the theory.

DOI: [10.1103/PhysRevB.96.155125](https://doi.org/10.1103/PhysRevB.96.155125)**I. INTRODUCTION**

The robustness of Landau's Fermi liquid theory relies on the protected gapless nature of quasiparticle excitations around the Fermi surface. Wilsonian effective field theory then guarantees that these protected excitations determine the macroscopic features of the theory in generic circumstances [1,2]. Aside from ordering instabilities, there is a poignant exception to this general rule. These are special situations where the quasiparticle excitations interact with other gapless states. This is notably so near a symmetry breaking quantum critical point. The associated massless modes should also contribute to the macroscopic physics. In $d \geq 3$ dimensions this interaction between Fermi surface excitations and gapless bosons is marginal/irrelevant and these so-called quantum critical metals can be addressed in perturbation theory as first discussed by Hertz and Millis [3–6]. In $2 + 1$ dimensions, however, the interaction is relevant and the theory is presumed to flow to a new interacting fixed point [6–9]. This unknown fixed point has been offered as a putative explanation of exotic physics in layered electronic materials near a quantum critical point such as the Ising-nematic transition. As a consequence, the deciphering of this fixed point theory is one of the major open problems in theoretical condensed matter physics. There have been numerous earlier studies of Fermi surfaces coupled to gapless bosons, but to be able to capture their physics one has almost always been required to study certain simplifying limits [10–20].

In this paper we show that the fermionic and bosonic spectrum of the most elementary $d = 2$ quantum critical metal can be computed nonperturbatively in the double limit where the Fermi momentum k_F is taken large, $k_F \rightarrow \infty$, while the number of fermion species N_f is taken to vanish, $N_f \rightarrow 0$, with the combination $N_f k_F$ held constant. This is an extension of previous work [21] where we studied the purely quenched limit $N_f \rightarrow 0$ followed by the limit $k_F \rightarrow \infty$. In this pure quenched $N_f \rightarrow 0$ limit the boson two-point function does not receive any corrections and the fermion two point function

can be found exactly. However, it is well known that for finite N_f and k_F the boson receives so-called Landau damping contributions that dominate the IR of the theory.

These Landau damping corrections are always proportional to N_f , and a subset of these are also proportional to k_F . These terms in particular influence the IR as the large scale, low energy behavior should emerge when k_F is large. Studying the double scaling limit where the combination $N_f k_F$ is held fixed gives a more complete understanding of the small N_f and/or large k_F limit and their interplay. In particular, this new double scaling limit makes precise previous results in the literature on the RPA approximation together with the $N_f \rightarrow 0$ limit and the strong forward scattering approximation [22–25]. Importantly, we shall show that the RPA results qualitatively capture the low energy at fixed momentum regime but not the full IR of the theory in the double scaling limit. The idea of this limit is similar to the limit taken in Ref. [26] where they study a similar model but in a matrix large N limit. In this limit they keep the quantity k_F^{d-1}/N fixed while taking both N and k_F large.

All the results here refer to the most elementary quantum critical metal. This is a set of N_f free spinless fermions at finite density interacting with a free massless scalar through a simple Yukawa coupling. Its action reads (in Euclidean time)

$$S = \int dx dy d\tau \left[\psi_j^\dagger \left(-\partial_\tau + \frac{\nabla^2}{2m} + \mu \right) \psi^j + \frac{1}{2} (\partial_\tau \phi)^2 + \frac{1}{2} (\nabla \phi)^2 + \lambda \phi \psi_j^\dagger \psi^j \right], \quad (1)$$

where $j = 1 \dots N_f$ sums over the N_f flavors of fermions and $\mu = \frac{k_F^2}{2m}$. We will assume a spherical Fermi surface, meaning k_F both sets the size of the Fermi surface, $2\pi k_F$, and the Fermi surface curvature, $1/k_F$. We will study the fermion and boson two-point functions of this theory in the double scaling limit $N_f \rightarrow 0$, $k_F \rightarrow \infty$. By this we mean that we take $k_F \rightarrow \infty$ while keeping the external momenta (measured from Fermi surface), energies, the coupling scale λ^2 , and the Fermi velocity $v = k_F/m$ fixed. We shall not encounter any UV divergences, but to address any ambiguities that may arise the usual assumption is made that the above theory is an effective theory below an energy and momentum scale Λ_0, Λ_k ,

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each of which is already much smaller than k_F ($\Lambda_0, \Lambda_k \ll k_F$). We do not address fermion pairing instabilities in this paper. They have been studied and found for similar models in other limits, outside of the particular double scaling limit studied here [26–30].

II. REVIEW OF THE QUENCHED APPROXIMATION

($N_f \rightarrow 0$ first, $k_F \rightarrow \infty$ SUBSEQUENTLY)

Let us briefly review the earlier results of Ref. [21] as they are a direct inspiration for the double scaling limit. Consider the fermion two-point function for the action above, Eq. (1). Coupling the fermions to external sources and integrating them out, and taking two derivatives w.r.t. the source, the formal expression for this two point function is

$$\begin{aligned} G_{\text{full}}(\omega, k) &= \langle \psi^\dagger(-\omega, -k) \psi(\omega, k) \rangle \\ &= \int \mathcal{D}\phi \det^{N_f} (G^{-1}[\phi]) G(\omega, k)[\phi] e^{-\int \frac{1}{2}(\partial_\tau \phi)^2 + \frac{1}{2}(\nabla \phi)^2}, \end{aligned} \quad (2)$$

where $G(\omega, k)[\phi]$ is the fermion two-point function in the presence of a background field ϕ , defined by

$$\left(-\partial_\tau + \frac{\nabla^2}{2m} + \mu + \lambda \phi \right) G(t, x)[\phi] = \delta(t - t') \delta^2(x - x'). \quad (3)$$

In the limit $k_F \rightarrow \infty$, for external momentum k close to the Fermi surface, we may approximate the derivative part with $-\partial_\tau + i v \partial_x$. The defining equation for the Green's function can then be solved in terms of a free fermion Green's function dressed with the exponential of a linear functional of ϕ . In the quenched $N_f \rightarrow 0$ limit this single exponentially dressed Green's function can be averaged over the background scalar with the Gaussian kinetic term. The result in real space is again an exponentially dressed free Green's function

$$G_{R, N_f \rightarrow 0}(r, t) = G_{R, \text{free}}(r, t) e^{I(t, r)} \quad (4)$$

with the exponent $I(r, t)$ given by

$$I(\tau, r) = \lambda^2 \int \frac{d\omega dk_x dk_y}{(2\pi)^3} \frac{\cos(\tau\omega - r k_x) - 1}{(i\omega - k_x v)^2} G_B(\omega, k), \quad (5)$$

and r conjugate to momentum measured from the Fermi surface (k_x), not the origin. Here $G_B(\omega, k)$ is the free boson Green's function determined by the explicit form of the boson kinetic term in the action Eq. (1). This is of course a known function and due to this simple dressed expression the retarded Green's function and therefore the fermionic spectrum of this model can be determined exactly in the limit $N_f \rightarrow 0$. The retarded Green's function in momentum space reads [21] (here ω is Lorentzian)

$$G_{R, N_f \rightarrow 0}(\omega, k_x) = \frac{1}{\omega - k_x v + \frac{\lambda^2}{4\pi\sqrt{1-v^2}} \sigma(\omega, k_x)}, \quad (6)$$

where σ is the solution of the equation

$$\begin{aligned} \frac{\lambda^2}{4\pi\sqrt{1-v^2}} (\sinh(\sigma) - \sigma \cosh(\sigma)) + v\omega - k_x \\ - \cosh(\sigma)(\omega - k_x v + i\epsilon) = 0, \end{aligned} \quad (7)$$

with k_x the distance from the Fermi surface, $v = k_F/m$ is the Fermi velocity, and $\epsilon \rightarrow 0^+$ is an $i\epsilon$ prescription that selects the correct root.

This nonperturbative result already describes interesting singular fixed point behavior: the spectrum exhibits non-Fermi liquid scaling behavior with multiple Fermi surfaces [21]. Nevertheless, it misses the true IR of the theory as the quenched limit inherently misses the physics of Landau damping. This arises from fermion loop corrections to the boson propagator that are absent for $N_f \rightarrow 0$. Below the Landau damping scale $\omega < \sqrt{\lambda^2 N_f k_F}$ the physics is expected to differ from the quenched approximation.

III. LOOP CANCELLATIONS AND BOSON TWO-POINT FUNCTION

It is clear from the review of the quenched derivation that finite N_f , i.e., fermion loop corrections, that only change the boson two-point function, can readily be corrected for by replacing the free boson two-point function $G_B(\omega, k)$ by the (fermion-loop) corrected boson two-point function in Eq. (5) (valid at large k_F). This is the essence of many RPA-like approximations previously studied. A weakness is that finite N_f corrections will also generate higher-order boson interactions and these can invalidate the simple dressed expression obtained here.

At the same time, it has been known for some time that finite density fermion-boson models with simple Yukawa scalar-fermion-density interactions as in Eq. (1) have considerable cancellations in fermion loop diagrams for low energies and momenta after symmetrization [24,31,32]. These cancellations make loops with more than three interaction vertices $V \geq 3$ finite as the external momenta and energies are scaled uniformly to zero. We will now argue that this result also means that in the $N_f \rightarrow 0, k_F \rightarrow \infty$ limit with $N_f k_F$ fixed, these $V \geq 3$ loops vanish. In this limit *only* the boson two-point function is therefore corrected and only at one loop and we can directly deduce that in this double limit the exact fermion correlation function is given by the analog of the dressed Green's function in Eq. (4). We comment on the limitations of considering this limit for subdiagrams in perturbation theory later in this section.

Consider the quantum critical metal before any approximations; i.e., we have a fully rotationally invariant Fermi surface with a finite k_F . The Yukawa coupling shows that the boson couples to the density operator $\psi^\dagger(x)\psi(x)$. All corrections to the boson therefore come from fermionic loops with fermion density vertices. These loops always show up symmetrized in the density vertices. Consider a fermion loop with a fixed number V of such density vertices, dropping the overall coupling constant dependence, and arbitrary incoming energies and momenta. Such a loop (ignoring the overall momentum conserving δ function) has energy dimension $3 - V$. These fermionic loops are all UV finite so they are independent of the scale of the UV cutoffs. There are only two important scales, the external bosonic energies and momenta ω_i, k_i , the fermi momentum k_F . A symmetrized V -point loop can by dimensional analysis be written as

$$I(\{\omega_i\}, \{k_i\}) = k_F^{3-V} f(\{\omega_i/k_F\}, \{k_i/k_F\}). \quad (8)$$

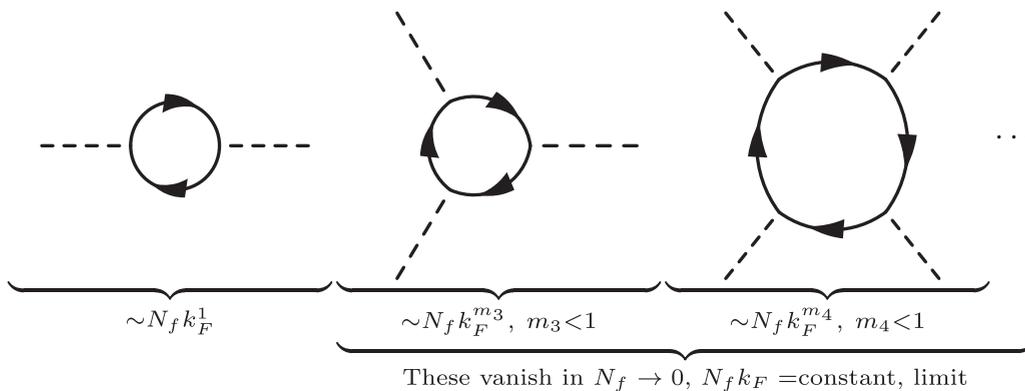


FIG. 1. Here we show the dominant scaling of fermion loops with different numbers of vertices in the limit of $N_f \rightarrow 0$ with $N_f k_F$ constant. This is the scaling after symmetrizing the external momenta. The two-vertex loop on the left does not get symmetrized and is the only loop that does not vanish in this limit.

Since fermion loops of our theory with $V \geq 3$ vertices have been shown to be finite as external energies and momenta are uniformly scaled to 0 [31], we thus have that f is finite as k_F is taken to infinity. This in turn means that $I(\{\omega_i\}, \{k_i\})$ scales as k_F^n with $n \leq 0$ for large k_F when $V \geq 3$.¹ Note that the use of the small external energies and momenta limit from Ref. [31] was merely a way of deducing the large k_F limit. We do not rely on the physical IR scaling to be the same as in Ref. [31], indeed we will find it not to be the same. All single fermionic loops additionally contain a sum over fermionic flavors so are therefore proportional to N_f . Combining this we see that a fermionic loop with $V \geq 3$ density vertices comes with a factor of $N_f k_F^{m_V}$ where $m_V \leq 0$ after symmetrizing the vertices. By now considering the combined limit of $N_f \rightarrow 0$ and $k_F \rightarrow \infty$ with $N_f k_F$ constant we see that these $V \geq 3$ loops all vanish. See Fig. 1.

We have now concluded that for a fixed set of external momenta, all symmetrized fermion loops vanish in our combined limit, except the $V = 2$ loop. There is still a possibility that diagrams containing $V > 2$ loops are important when taking the combined limit *after* performing all bosonic momentum integrals and summing up the infinite series of diagrams. In essence, the bosonic integrals and the infinite sum of perturbation theory need not commute with the combined $N_f \rightarrow 0, k_F \rightarrow \infty$ limit. What the IR of the full theory (finite N_f) looks like is not known so taking the $N_f \rightarrow 0$ limit last is currently out of reach. In Ref. [33] the authors show that divergence of fermionic loops does not cancel under a nonuniform low-energy scaling of energies and momenta where the momenta are additionally taken to be increasingly collinear. The scaling they use is motivated by the perturbative treatment in Ref. [10], and if this is the true IR scaling and it persists at small N_f , then there will be effects unaccounted for in the above.

¹Naively corrections to the boson would be expected to scale as k_F since it receives corrections from a Fermi surface of size $2\pi k_F$. However k_F also sets the curvature of the Fermi surface and for a large k_F we approach a flat Fermi surface for which $V \geq 3$ loops completely cancel. This is shown in more detail in Appendix A.

Regardless of the above mentioned caveat, in keeping the $V = 2$ fermion loops we take the combined limit *after* performing the fermionic loop integrals and thus move closer than in our previous work [21] to the goal of understanding the IR of quantum critical metals. To summarize: the ordered limit we consider is:

- (1) We first perform rotationally invariant finite k_F fermionic loop integrals.
- (2) Then we take the limit $N_f \rightarrow 0, k_F \rightarrow \infty$ with $N_f k_F$ fixed; this only keeps $V = 2$ loops.
- (3) Next we perform bosonic loop integrals.
- (4) Finally we sum all contributions at all orders of the coupling constant.

The result above means the fully quantum corrected boson remains Gaussian in this ordered limit and only receives corrections from the $V = 2$ loops.

A. Boson two-point function

We now compute the one-loop correction to the boson two-point function; in our ordered double scaling limit this is all we need. We then substitute the Dyson summed one-loop corrected boson two-point function into the dressed fermion Green's function to obtain the exact fermionic spectrum.

The one-loop correction—the boson polarization—in the double scaling limit is given by the large k_F limit of the two-vertex fermion loop. This can be calculated using a linearized fermion dispersion:

$$\Pi_1(Q) = \lambda^2 N_f k_F \int \frac{dk_0 dk d\theta}{(2\pi)^3} \times \frac{1}{(ik_0 - vk)(i(k_0 + q_0) - v(k + |\vec{q}| \cos \theta))}. \quad (9)$$

Note from the $\cos \theta$ dependence in the numerator that we are not making a “patch” approximation. In the low energy limit this angular dependence is the important contribution of the rotationally invariant fermi surface, whereas the subleading terms of the dispersion can be safely ignored. As stated earlier, the result of these integrals is finite. However, it does depend

on the order of integration. The difference is a constant C

$$\begin{aligned} \Pi_1(Q) &= \frac{\lambda^2 N_f k_F}{2\pi v} \left(\frac{|q_0|}{\sqrt{q_0^2 + v^2 \vec{q}^2}} + C \right) \\ &\equiv M_D^2 \left(\frac{|q_0|}{\sqrt{q_0^2 + v^2 \vec{q}^2}} + C \right). \end{aligned} \quad (10)$$

As pointed out in for instance Refs. [19,20], the way to think about this ordering ambiguity is that one should strictly speaking first regularize the theory and introduce a one-loop counterterm. This counterterm has a finite ambiguity that needs to be fixed by a renormalization condition. Even though the loop momentum integral happens to be finite in this case, the finite counterterm ambiguity remains. The correct renormalization condition is the choice $C = 0$. This choice corresponds to the case when the boson is tuned to criticality since a nonzero C would mean the presence of an effective mass generated by quantum effects.

A more physical way to think of the ordering ambiguity is as the relation between the frequency (Λ_0) and momentum (Λ_k) cutoff. We will assume that $\Lambda_k \gg \Lambda_0$, which means that we evaluate the k integral first and then the frequency k_0 integral. In this case $C = 0$ directly follows.

IV. FERMION TWO-POINT FUNCTION

With the single surviving one-loop correction to the boson two-point function in hand, we can immediately write down the expression for the full fermion two-point function. This is the same dressed expression Eq. (5) as in Ref. [21] but with a modified boson propagator $G_B = 1/(G_{B,0}^{-1} + \Pi)$, with Π the one-loop polarization of Eq. (10). Substituting this in we thus need to calculate the integral

$$\begin{aligned} I(\tau, r) &= \lambda^2 \int \frac{d\omega dk_x dk_y}{(2\pi)^3} \\ &\times \frac{\cos(\tau\omega - rk_x) - 1}{(i\omega - k_x v)^2 (\omega^2 + k_x^2 + k_y^2 + M_D^2 \frac{|\omega|}{\sqrt{v^2(k_x^2 + k_y^2) + \omega^2}})}. \end{aligned} \quad (11)$$

At this moment, we can explain clearly how our result connects to previous approaches. A similarly dressed propagator can be proposed based on extrapolation from 1d results [22,24]. An often used approximation in the literature is to now study this below the scale M_D , see e.g., Refs. [6,23]. This is the physically most interesting limit since in the systems of interest N_f is order one and we are considering large k_F . In this limit the polarization term will dominate over the kinetic terms, but since the rest of the integrand in (11) has no k_y dependence, it is necessary to keep the k_y term in the boson propagator. The ω and k_x momenta will suppress the integrand when they are of order λ^2 whereas the k_y term will do so once it is of order $\lambda^{2/3} M_D^{2/3}$. This means that for $M_D \gg \lambda^2$, the relevant k_y will be much larger than the relevant ω and k_x .

This argues that we can truncate to the large M_D propagator

$$G_{B, M_D \rightarrow \infty}(\omega, k_x, k_y) = \frac{1}{k_y^2 + M_D^2 \frac{|\omega|}{v|k_y|}}. \quad (12)$$

This Landau-damped propagator has been used extensively, for instance, Refs. [6,23]. In Ref. [23] this propagator was used for the type of nonperturbative calculation we are proposing here. We discuss this here, as we will now show that using this simplified propagator has a problematic feature. This propagator only captures the leading large M_D contribution, but the nonperturbative exponential form of the exact Green's function sums up powers of the propagator which then are subleading in M_D .

Using the large M_D truncated boson Green's function the integral $I(\tau, r)$ to be evaluated simplifies to

$$I_{M_D \rightarrow \infty} = \lambda^2 \int \frac{d\omega dk_x dk_y}{(2\pi)^3} \frac{\cos(\tau\omega - xk_x) - 1}{(i\omega - k_x v)^2 (k_y^2 + M_D^2 \frac{|\omega|}{v|k_y|})}. \quad (13)$$

Writing the cosine in terms of exponentials we can perform the k_x integrals using residues by closing the contours in opposite half planes. Summing up the residues gives

$$I_{M_D \rightarrow \infty} = -\lambda^2 \int d\omega dk_y \frac{k_y^2 |r| e^{-|\omega|(\frac{|r|}{v} + i \operatorname{sgn}(r)\tau)}}{8\pi^2 M_D^2 |\omega k_y| + k_y^4 v}. \quad (14)$$

The k_y integral can be performed next to yield

$$I_{M_D \rightarrow \infty} = -\lambda^2 \int d\omega \frac{|r| e^{-|\omega|(\frac{|r|}{v} + i \operatorname{sgn}(r)\tau)}}{12\sqrt{3}\pi (M_D^2 v^5 |\omega|)^{1/3}}. \quad (15)$$

The primitive function to this ω integral is the upper incomplete gamma function, with argument $2/3$. Evaluating this incomplete gamma function in the appropriate limits and substituting the final expression for $I_{M_D \rightarrow \infty}(\tau, r)$ into the expression for the fermion two-point function gives us:

$$\begin{aligned} G_f(\tau, r)_{M_D \rightarrow \infty} &= \frac{1}{2\pi(i\tau - v\tau)} \\ &\times \exp\left(-\frac{|r|}{l_0^{1/3} (|r| + i v \operatorname{sgn}(r)\tau)^{2/3}}\right), \end{aligned} \quad (16)$$

where the length scale l_0 is given by

$$l_0^{1/3} = \frac{6\sqrt{3}\pi v M_D^{2/3}}{\Gamma(\frac{2}{3})\lambda^2}. \quad (17)$$

This result has been found earlier in Ref. [22] (see also Ref. [24]). However, this real space expression hides the inconsistency of the approach. This becomes apparent in its momentum space representation. The Fourier transform of the real space Green's function

$$\begin{aligned} G_f(\omega, k)_{M_D \rightarrow \infty} &= \int d\tau dr \frac{e^{i(\omega\tau - kr)}}{2\pi(i\tau - v\tau)} \\ &\times \exp\left(-\frac{|r|}{l_0^{1/3} (|r| + i v \operatorname{sgn}(r)\tau)^{2/3}}\right) \end{aligned} \quad (18)$$

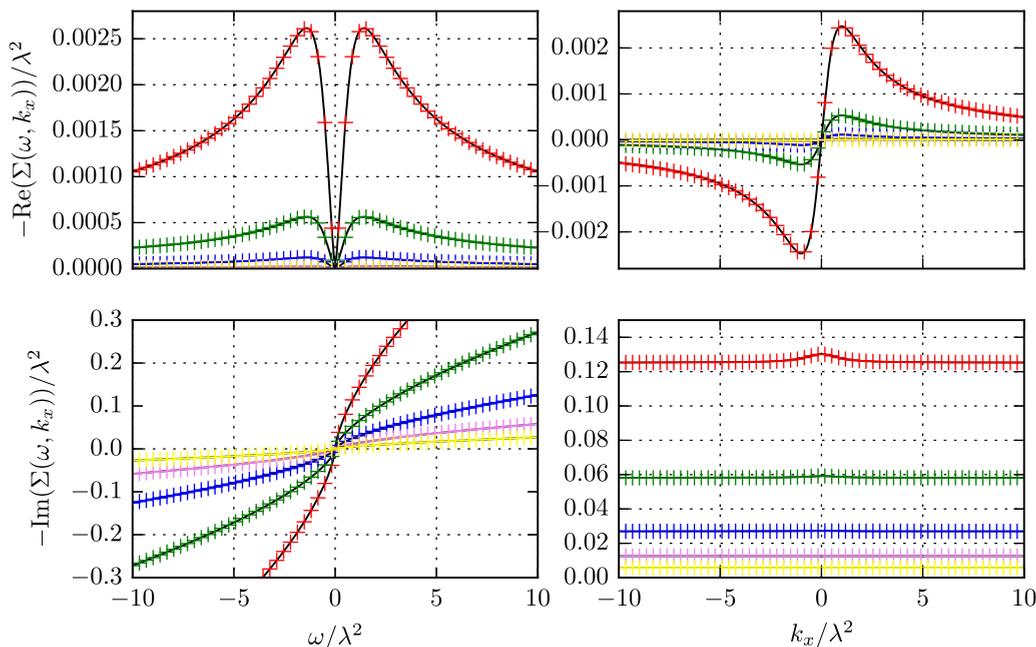


FIG. 2. Real and imaginary parts of the self energy obtained using the large- k_F Landau-damped propagator. This plot shows the agreement between the numerics and the analytical solution, verifying that both solutions are correct. Notice the difference in magnitude between the real and imaginary part. The agreement of the real parts shows that the numerical procedure has a very small relative error. All plots are for the $k_x, \omega = \lambda^2$ slice with $v = 1$.

is tricky, but remarkably can be done exactly. We do so in Appendix B. The result is

$$\begin{aligned}
 G_{f, M_D \rightarrow \infty}(\omega, k_x) &= \frac{1}{i\omega - k_x v} \cos\left(\frac{\omega}{v l_0^{1/2}(\omega/v + i k_x)^{3/2}}\right) \\
 &+ \frac{6\sqrt{3}i\Gamma(\frac{1}{3})\omega^{2/3}}{8\pi l_0^{1/3}v^{5/3}(\omega/v + i k_x)^2} \\
 &\times {}_1F_2\left(1; \frac{5}{6}, \frac{4}{3}; -\frac{\omega^2}{4l_0v^2(\omega/v + i k_x)^3}\right) \\
 &+ \frac{3\sqrt{3}i\Gamma(-\frac{1}{3})\omega^{4/3}}{8\pi l_0^{2/3}v^{7/3}(\omega/v + i k_x)^3} \\
 &\times {}_1F_2\left(1; \frac{7}{6}, \frac{5}{3}; -\frac{\omega^2}{4l_0v^2(\omega/v + i k_x)^3}\right). \quad (19)
 \end{aligned}$$

This expression has been compared with numerics to verify its correctness; see Fig. 2.

We can now show the problematic feature. Recall that Eq. (19) is the Green's function in Euclidean signature. Continuing to the imaginary line, $\omega = -i\omega_R$, this becomes the proper retarded Greens function, $G_R(\omega_R, k_x)$, and from this we can obtain the spectral function $A(\omega_R, k_x) = -2\text{Im} G_R(\omega_R, k_x)$. As it encodes the excitation spectrum, the spectral function ought to be a *positive* function that moreover equals 2π when integrated over all energies ω_R , for any momentum k . This large M_D spectral function contains an oscillating singularity at $\omega_R = vk_x$. We are free to move the contour into complex ω_R plane by deforming $\omega_R \rightarrow \omega_R + i\Omega$ where Ω is positive but otherwise arbitrary. Upon doing this it is easy to numerically verify that indeed the integral over ω_R gives 2π . However, if we look at the behavior close to the essential singularity, the

function oscillates rapidly and does not stay positive as one approaches the singularity; see Fig. 3. This reflects that the large M_D approximation done in this way is not consistent.

Even though the approximation for the exponent $I(\tau, r) \equiv \frac{\tilde{I}(\tau, r)}{(M_D)^{2/3}}$ is valid to leading order in $1/M_D$, this is not systematic after exponentiation to obtain the fermion two-point function

$$G_{f, M_D \rightarrow \infty}(\tau, r) = \frac{1}{2\pi(ir - v\tau)} \exp\left(\frac{\tilde{I}(\tau, r)}{M_D^{2/3}} + \mathcal{O}\left(\frac{1}{(M_D)^{4/3}}\right)\right). \quad (20)$$

Reexpanding the exponent one immediately sees that keeping only the leading term in $I(\tau, r)$ mixes at higher order with the

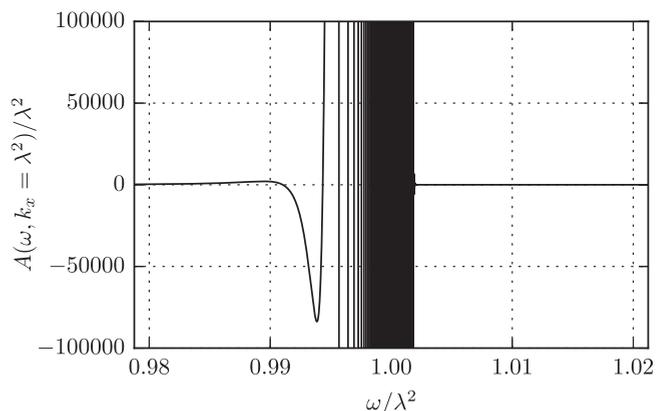


FIG. 3. Exact fermion spectral function based on the large- M_D approximation for the exact boson propagator. Notice that the function is not positive everywhere. Here $k = \lambda^2$, $v = 1$ and $M_D = 2\pi\lambda^2$.

subleading terms at lower order in $1/M_D$

$$G_f(\tau, x) \underset{M_D \rightarrow \infty}{=} \frac{1}{2\pi(ir - v\tau)} \left(1 + \frac{\tilde{I}(\tau, r)}{M_D^{2/3}} + \mathcal{O}(M_D^{-4/3}) \right. \\ \left. + \frac{1}{2} \left(\frac{\tilde{I}(\tau, r)}{M_D^{2/3}} + \dots \right)^2 \right). \quad (21)$$

Despite this problematic feature, we will show from the exact result that in the IR G_f (with a small modification) does happen to capture the correct physics.

A. Exact fermion two-point function for large k_F with M_D fixed; $v = 1$

We therefore make no further assumption regarding the value of M_D and we return to the full integral Eq. (11) to determine the real space fermion two-point function. Solving this in general is difficult, and to simplify mildly we consider the special case $v = 1$. In our previous studies of the quenched $M_D = 0$ limit we saw that this choice for value of v is actually not very special, even though it appears that there is an enhanced symmetry. In fact, nothing abruptly happens as $v \rightarrow 1$, except that the quenched $M_D = 0$ solution can be written in closed form for this value of $v = 1$. Nor for the case of large M_D is the choice $v = 1$ in any way special. As can be seen above in Eq. (16) for large M_D all v are equivalent up to a rescaling of τ versus r and a rescaling of the single length scale l_0 . We may therefore expect that for a finite M_D , the physics of $0 < v < 1$ is qualitatively the same as the (not-so-) special case $v = 1$.

After setting $v = 1$ and changing to spherical coordinates we have

$$I = \lambda^2 \int d\tilde{r} d\phi d\theta e^{2i\phi} \frac{\cos(\tilde{r} \sin(\theta)(\tau \sin(\phi) - r \cos(\phi))) - 1}{8\pi^3 \sin(\theta)^2 (M_D^2 |\sin(\phi)| + \tilde{r}^2 / \sin(\theta))}. \quad (22)$$

Performing the \tilde{r} integral gives us

$$I = \pi^4 \lambda^2 \int d\phi d\theta e^{2i\phi} \frac{e^{-M_D |\tau \sin(\phi) - r \cos(\phi)| \sqrt{|\sin(\phi)| \sin(\theta)^3}} - 1}{16M_D \sqrt{|\sin(\phi)| \sin(\theta)^3}}. \quad (23)$$

Note that if the signs of both τ and r are flipped, then this is invariant. Changing the sign of only τ , and simultaneously

making the change of variable $\phi \rightarrow -\phi$, then the (real) fraction is invariant but the exponent in the prefactor changes sign. Thus, I goes to I^* as the sign of either τ or r is changed. Without loss of generality, we can assume that both of them are positive from now on. We further see that the integrand is invariant under $\phi \rightarrow \phi + \pi$, so we may limit the range of ϕ to $(0, \pi)$ by doubling the value of integrand. Similarly we limit θ to $(0, \pi/2)$ and multiply by another factor of 2. We then make the changes of variables:

$$\phi = \tan^{-1}(s) + \pi/2 \quad \theta = \sin^{-1}(u^{2/3}) \quad (24)$$

with $s \in \mathbb{R}$ and u is integrated over the range $(0, 1)$. For convenience we introduce the function

$$z(s) = M_D |sr + \tau| (1 + s^2)^{-3/4}. \quad (25)$$

Now the two remaining integrals can be written as

$$I = \frac{\lambda^2}{M_D} \int ds du \frac{(e^{-uz(s)} - 1)(s - i)}{6\pi^2 (s + i)(1 + s^2)^{3/4} u^{4/3} \sqrt{(1 - u^{4/3})}}. \quad (26)$$

After expanding the exponential we can perform the u integral term by term. We are left with

$$I = \frac{\lambda^2}{M_D} \int ds \sum_{n=1}^{\infty} \frac{(1 + s^2)^{1/4} (-z(s))^n \Gamma(\frac{3n-1}{4})}{8\pi^{3/2} n! (i + s)^2 \Gamma(\frac{3n+1}{4})}. \quad (27)$$

This can be resummed into a sum of generalized hypergeometric functions, but this is not useful at this stage. Instead we once again integrate term by term. Collecting the prefactors and introducing the constant $a = \tau/r$, the n -th term can be written as

$$I = \sum_{n=1}^{\infty} c_n \int ds (s - i)^2 |s + a|^n (1 + s^2)^{-(7+3n)/4}. \quad (28)$$

This can be written as

$$I = \sum_{n=1}^{\infty} c_n \int ds dw (s - i)^2 |s + a|^n \frac{e^{-w(s^2+1)} w^{3(1+n)/4}}{(3(1+n)/4)!}, \quad (29)$$

where w is integrated on $(0, \infty)$. After splitting the integral at $s = -a$ to get rid of the absolute value we can calculate the s integrals in terms of confluent hypergeometric functions ${}_1F_1(a, b; z)$. Adding the two halves $s < -a$ and $s > -a$ of the integral we have

$$I = \sum_{n=1}^{\infty} c_n \int dw \frac{\Gamma(\frac{1+n}{2}) e^{-(1+a^2)w}}{2(3(1+n)/4)!} \left(2w(i+a)^2 {}_1F_1\left(\frac{2+n}{2}, \frac{1}{2}; a^2 w\right) \right. \\ \left. + (2+n) {}_1F_1\left(\frac{4+n}{2}, \frac{1}{2}; a^2 w\right) - 4aw(2+n)(i+a) {}_1F_1\left(\frac{4+n}{2}, \frac{3}{2}; a^2 w\right) \right). \quad (30)$$

It may look like we have just exchanged the s integral for the w integral, but by writing the hypergeometric functions in series form,

$$I = \sum_{n=1}^{\infty} c_n \int dw \sum_{m=0}^{\infty} \frac{2^{2m-1} a^{2m} e^{-(1+a^2)w} w^{\frac{n-3}{4}+m} \Gamma(\frac{1+n}{2} + m)}{\Gamma(\frac{7+3n}{4}) \Gamma(2+2m)} (n(1+2m-4a(i+a)w) + (1+2m)(1+2m-2(1+a^2)w)), \quad (31)$$

the w integral can now be performed. The result is

$$I = \sum_{n=1, m=0}^{\infty} c_n \frac{(a+i)4^{m-1}a^{2m}(a^2+1)^{-\frac{1}{4}(4m+n+5)}\Gamma(m+\frac{n+1}{4})\Gamma(m+\frac{n+1}{2})}{\Gamma(2m+2)\Gamma(\frac{7+3n}{4})} \times (a(2m-6mn-2n^2-n+1)-i(2m+1)(n+1)). \quad (32)$$

The sum over m can be expressed in terms of the ordinary hypergeometric function, ${}_2F_1(a_1, a_2; b; z)$:

$$I = \sum_{n=1}^{\infty} c_n \frac{(n+1)(a^2+1)^{-\frac{n}{4}-\frac{1}{4}}\Gamma(\frac{n+1}{4})\Gamma(\frac{n+1}{2})}{24(a-i)^2(a+i)\Gamma(\frac{3n}{4}+\frac{7}{4})} \quad (33)$$

$$\times \left(a^2(n+1)(-3an+a-i(n+1)) {}_2F_1\left(\frac{n+3}{2}, \frac{n+5}{4}; \frac{5}{2}; \frac{a^2}{a^2+1}\right) \right) \quad (34)$$

$$-6(a^2+1)(a(2n-1)+i) {}_2F_1\left(\frac{n+1}{4}, \frac{n+1}{2}; \frac{3}{2}; \frac{a^2}{a^2+1}\right). \quad (35)$$

The space-time dependence in this expression is in $a = \tau/r$ and with additional r dependence in the coefficients c_n . The result above is the value for both τ and r positive. Using the known symmetries presented above, the solution can be extended to all values of τ and r by appropriate absolute value signs. Then changing variables to

$$\begin{aligned} \tau &= R \cos(\Phi) \\ r &= R \sin(\Phi) \end{aligned} \quad (36)$$

we have

$$I = \frac{\lambda^2 f(RM_D, \Phi)}{M_D} \quad (37)$$

with the function $f(\tilde{R}, \Phi)$ given by

$$\begin{aligned} f(\tilde{R}, \Phi) &= \sum_{n=1}^{\infty} f_n \tilde{R}^n \\ f_n &= \frac{e^{i\Phi} 2^{-1-n} (-1)^n \Gamma(\frac{n+1}{4}) |\sin(\Phi)|^{\frac{1+3n}{2}}}{9\pi(3n-1)\Gamma(\frac{n}{2}+1)\Gamma(\frac{1+3n}{4})} \cdot \left({}_2F_1\left(\frac{n+3}{2}, \frac{n+5}{4}; \frac{5}{2}; \cos^2(\Phi)\right) (n+1) \cdot \cos^2(\Phi) \right. \\ &\quad \left. \times ((1-3n)\cos(\Phi) - i(n+1)\sin(\Phi)) + {}_2F_1\left(\frac{n+1}{4}, \frac{n+1}{2}; \frac{3}{2}; \cos^2(\Phi)\right) 6((1-2n)\cos(\Phi) - i\sin(\Phi)) \right). \end{aligned} \quad (38)$$

This exact infinite series expression for the exponent $I(\tilde{R}, \Phi)$ gives us the exact fermion two-point function in real (Euclidean) space (time). We have not been able to find a closed form expression for this final series. Note that $f_n \sim 1/n!$ for large n , and the series therefore converges rapidly. Moreover, numerically the hypergeometric functions are readily evaluated to arbitrary precision (e.g., with Mathematica), and therefore the value of $f(\tilde{R}, \Phi)$ can be robustly evaluated to any required precision.

As a check on this result, we can compare it to the exact result in the quenched $M_D = 0$ limit in Ref. [21], where the exact answer was found in a different way. In the limit where $M_D \rightarrow 0$ we see that only the first term of this series gives a contribution and the expression for the exponent collapses to

$$\lim_{M_D \rightarrow 0} I(R, \Phi) = \lambda^2 f_1 R = \lambda^2 \frac{e^{2i\Phi}}{12\pi} R. \quad (39)$$

In Cartesian coordinates this equals

$$\lim_{M_D \rightarrow 0} I(\tau, r) = \lambda^2 \frac{(\tau + ir)^2}{12\pi \sqrt{\tau^2 + r^2}}. \quad (40)$$

This is the exact same expression as found in Ref. [21] for $v = 1$.

There is one value of the argument for which $f(\tilde{R}, \Phi)$ drastically simplifies. For $r = 0$ ($\Phi = 0, \pi$) we have

$$f_n(\Phi = 0) = -\frac{(-1)^n}{6\pi\Gamma(n+2)} \quad (41)$$

and thus

$$f(\tilde{R}, \Phi = 0) = \frac{1}{6\pi} + \frac{e^{-\tilde{R}} - 1}{6\pi\tilde{R}}. \quad (42)$$

Further numerical analysis shows that the real part of $f(\tau, r)$ is maximal for $r = 0$.

B. The IR limit of the exact fermion two-point function compared to the large- M_D expansion

With this exact real space answer, we can now reconsider why the large M_D (large $N_f k_F$) limit fails and which expression does reliably capture the strongly coupled IR

physics of interest. The expression obtained above, Eq. (38), is not very useful for extracting the IR Green's function or the Green's function at a large M_D as the expression is organized in an expansion around $RM_D = 0$. To study the limit where $RM_D \gg 1$ we can go back to Eq. (26). With this expression we see that the exponential in the integrand, $e^{-uz(s)}$ with $z \sim M_D|sr + \tau| \sim \tilde{r}$, is generically suppressed for large $\tilde{R} = RM_D$. The exceptions are when either $sr + \tau$ is small, s is large, or u is small. The first two cases are

also unimportant in the $\tilde{R} \gg 1$ limit. In the first case we restrict the s integral to a small range of order $1/\tilde{R}$ around $-\tau/r$; this contribution therefore becomes more and more negligible in the limit $\tilde{R} \gg 1$. In the second case we will have a remaining large denominator in s outside the exponent that also suppresses the overall integral. Thus for large \tilde{R} , the only appreciable contribution of the exponential term to the integral in $I(\tau, r)$ arises when u is small. To use this, we first write the integral as

$$\begin{aligned}
 I_{\text{IR}} &= I_{\text{IR,exp}} + I_{\text{IR,-1}}, \\
 I_{\text{IR,exp}}(\tau, r) &= \lambda^2 \int_{-\infty}^{\infty} ds \frac{s-i}{6\pi^2(s+i)(1+s^2)^{3/4}M_D} \left(\int_0^1 du \frac{e^{-4uz(s)} - 1}{u^{4/3}\sqrt{1-u^{4/3}}} - \int_1^{\infty} du \frac{1}{u^{4/3}} \right) \\
 &\simeq \lambda^2 \int_{-\infty}^{\infty} ds \frac{s-i}{6\pi^2(s+i)(1+s^2)^{3/4}M_D} \left(\int_0^1 du \frac{e^{-4uz(s)} - 1}{u^{4/3}} - \int_1^{\infty} du \frac{1}{u^{4/3}} \right), \\
 I_{\text{IR,-1}}(\tau, r) &= \lambda^2 \int_{-\infty}^{\infty} ds \frac{s-i}{6\pi^2(s+i)(1+s^2)^{3/4}M_D} \left(\int_1^{\infty} du \frac{1}{u^{4/3}} + \int_0^1 du \left(\frac{-u^{-4/3}}{\sqrt{1-u^{4/3}}} + u^{-4/3} \right) \right). \quad (43)
 \end{aligned}$$

We have added and subtracted an extra term to each to ensure convergence of each of the separate terms. Since the important contribution to $I_{\text{IR,exp}}$ is from the small u region we can extend its range from $(0,1)$ to $(0,\infty)$. This way, the integrals can then be done

$$I_{\text{IR,exp}} = \int_{-\infty}^{\infty} ds \frac{-\lambda^2|sr + \tau|^{1/3}}{3^{3/2}\pi M_D^{2/3}(s+i)^2\Gamma(\frac{4}{3})} = -\frac{\Gamma(\frac{2}{3})\lambda^2|r|^{1/3}}{3^{3/2}\pi M_D^{2/3}(1+\frac{i\tau}{r})^{2/3}}, \quad (44)$$

$$I_{\text{IR,-1}} = \int_{-\infty}^{\infty} ds \frac{\lambda^2(s-i)}{6\pi^2(s+i)(1+s^2)^{3/4}M_D} \left(\int_0^{\infty} du \left(\frac{1}{u^{4/3}} - \frac{\theta(1-u)}{u^{4/3}\sqrt{1-u^{4/3}}} \right) \right) = \frac{\lambda^2}{6\pi M_D}. \quad (45)$$

In total we have for large \tilde{R} :

$$I = -\frac{\Gamma(\frac{2}{3})\lambda^2|r|}{3^{3/2}\pi M_D^{2/3}(|r|+i\text{sgn}(r)\tau)^{2/3}} + \frac{\lambda^2}{6\pi M_D} + \mathcal{O}(\lambda^2 M_D^{-4/3} R^{-1/3}). \quad (46)$$

We see that the leading order term in R is the same as was obtained from the large M_D approximation of the exponent. The first subleading term is just a constant. This is good news because we already have the Fourier transform of this expression. This result is valid for length scales larger than $1/M_D$ with a bounded error of the order $R^{-1/3}$. Defining this approximation as G_{IR} , i.e.,

$$G_{\text{IR}} = G_0 \exp \left(-\frac{\Gamma(\frac{2}{3})\lambda^2|r|}{3^{3/2}\pi M_D^{2/3}(|r|+i\text{sgn}(r)\tau)^{2/3}} + \frac{\lambda}{6\pi M_D} \right), \quad (47)$$

the error of this approximation follows from:

$$\Delta G_{\text{IR}} = G - G_{\text{IR}} = G_{\text{IR}}(\exp(\mathcal{O}(\tilde{R}^{-1/3})) - 1). \quad (48)$$

Since the exponential in G_{IR} is bounded we have that $\Delta G_{\text{IR}} = \mathcal{O}(r^{-4/3})$. After Fourier transforming this translates to an error of order $\mathcal{O}(k^{-2/3})$.

C. The exact fermion two-point function in momentum space: Numerical method

Having understood the shortcomings of the naive large M_D answer, the way to derive the *exact* answer in real space, and the

correct IR approximation, we can now analyze the behavior of the quantum critical metal at low energies. For this we need to transform to frequency-momentum space. As our exact answer is in the form of an infinite sum, this is not feasible analytically. We therefore resort to a straightforward numerical Fourier analysis.

To do so we first numerically determine the real space value of the exact Green's functions. To do so accurately, several observations are relevant:

(1) The coefficients f_n in the infinite sum for $I(\tau, r)$ decay factorially in n so once n is of order \tilde{R} , convergence is very rapid.

(2) The hypergeometric functions for each n are costly to compute with high precision, but with the above choice of polar coordinates the arguments of the hypergeometric functions are independent of R and M_D . We therefore numerically evaluate the series over a grid in \tilde{R} and Φ . We can then reuse the hypergeometric function evaluations many times and greatly decrease computing time.

(3) The real space polar grid will be limited to a finite size. The IR expansion from Eq. (47) can be used instead of the exact series for large enough \tilde{R} . To do so, we have to ensure an overlapping regime of validity. It turns out that a rather large value of \tilde{R} is necessary to obtain numerical

agreement between these two expansions, i.e., one needs to evaluate a comparably large number of terms in the expansion. For the results presented in this paper it has been necessary to compute coefficients up to order 16 000 in \tilde{R} , for many different angles Φ . The function is bounded for large τ and \tilde{R} but each term grows quickly. This means that there are large cancellations between the terms that in the end give us a small value. We therefore need to calculate these coefficients to very high precision in evaluating the polynomial. For these high precision calculations, we have used the Gnu Multiprecision Library [34].

(4) On this polar grid we computed the exact answer for $\tilde{R} < \tilde{R}_0 \approx 1000$ and used cubic interpolation for intermediate values. For larger \tilde{R} we use the asymptotic expansion in Eq. (47).

We then use a standard discrete numerical Fourier transform (DFT) to obtain the momentum space two-point function from this numeric prescription for $G(R, \Phi)$. Sampling $G(R, \Phi)$ at a finite number of discrete points, the size of the sampling grid will introduce an IR cutoff at the largest scales we sample and a UV cutoff set by the smallest spacing between points. These errors in the final result can be minimized by using the known asymptotic values analytically. Rather than Fourier transforming $G(\tau, r)$ as a whole, we Fourier transform $G_{\text{diff}}(\tau, r) = G(\tau, r) - G_{\text{IR}}(\tau, r)$ instead. Since both these functions approach the free propagator in the UV, the Fourier transform of its difference will decay faster for large

ω and k . This greatly reduces the UV artifacts inherent in a discrete Fourier transform. These two functions *also* approach each other for large τ and x . In fact, with the numerical method we use to approximate $f(\tilde{R}, \Phi)$ described above, they will be identical for $\tilde{R}_0 < M_D \sqrt{\tau^2 + r^2}$. This means that we only need to sample the DFT within that area. With a DFT we will always get some of the UV tails of the function that gives rise to folding aliasing artifacts. Now our function decays rapidly so one could do a DFT to very high frequencies and discard the high frequency part. This unfortunately takes up a lot of memory so we have gone with a more CPU intensive but memory friendly approach. To address this we perform a convolution with a Gaussian kernel, perform the DFT, keep the lowest 1/3 of the frequencies and then divide by the Fourier transform of the kernel used. This gives us a good numeric value for $G_{\text{diff}}(\omega, k)$. To this we add our analytic expression for $G_{\text{IR}}(\omega, k)$.

V. THE PHYSICS OF 2+1 QUANTUM CRITICAL METALS IN DOUBLE SCALING LIMIT

With the exact real space expression and the numerical momentum space solution for the full nonperturbative fermion Green's function, we can now discuss the physics of the elementary quantum critical metal in the double scaling limit. Let us emphasize right away that all our results are in Euclidean space. Although a Euclidean momentum space Green's

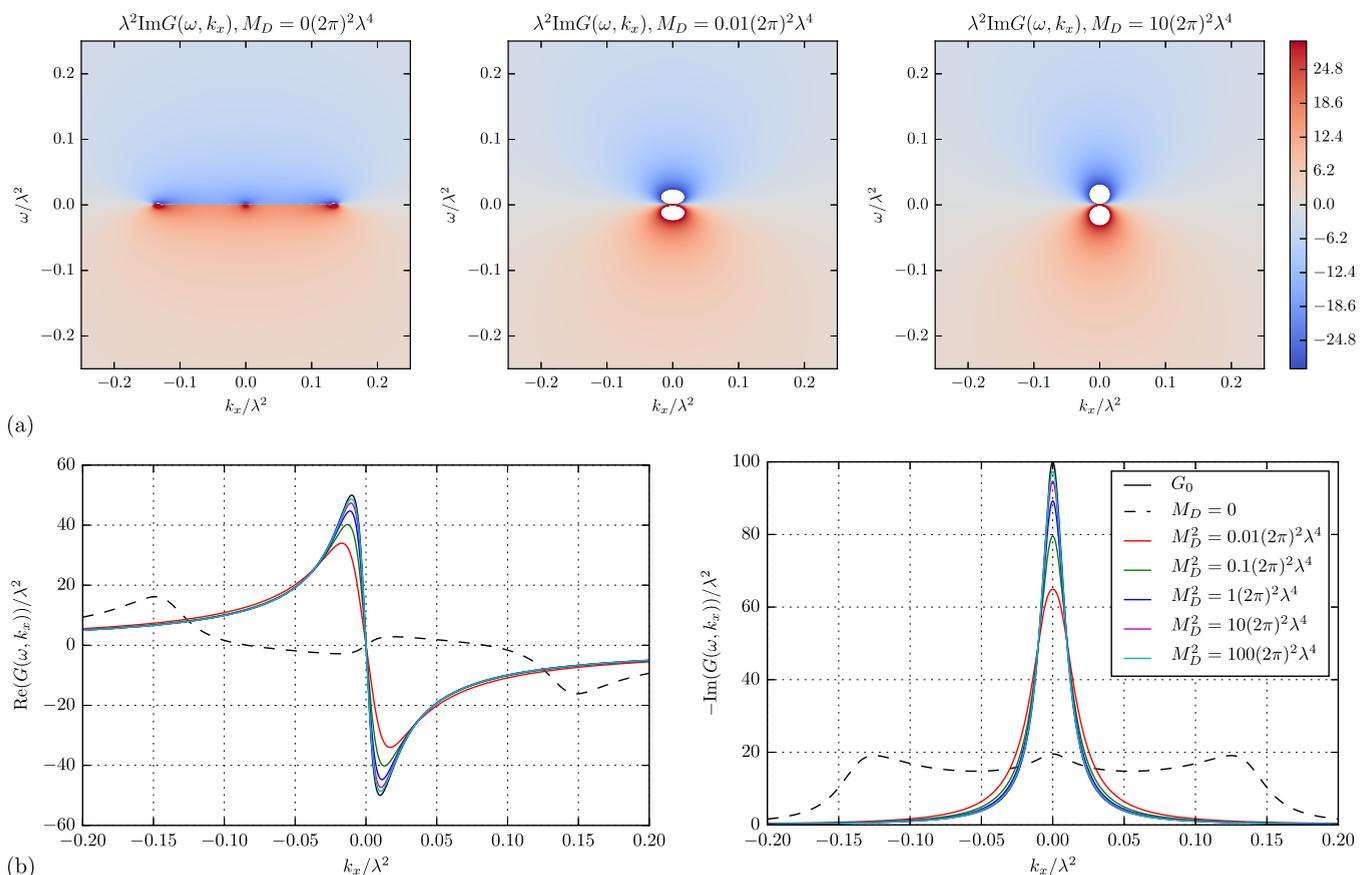


FIG. 4. (a) Density plots of the imaginary part of the exact (Euclidean) fermion Green's function $G(\omega, k_x)$ for various values of M_D . In the quenched limit $M_D = 0$ the three Fermi surface singularities are visible. For any appreciable finite M_D the Euclidean Green's function behaves as a single Fermi surface non-Fermi liquid. (b) Real and imaginary parts of $G(\omega, k_x)$ for very small $\omega = 0.01\lambda^2$.

function can be used to find a good Lorentzian continuation with a well-defined and consistent spectral function, this function is not easily obtainable from our numerical Euclidean result. We leave this for future work. The Euclidean signature Green's function does not visually encode the spectrum directly, but for very low energies/frequencies the Euclidean and the Lorentzian expressions are nearly identical, and we can extract much of the IR physics already from the Euclidean correlation function.

In Fig. 4 we show density plots of the imaginary part of $G(\omega, k_x)$ for different values of M_D as well as cross sections at fixed low ω . For the formal limit $M_D = 0$ we detect three singularities near $\omega = 0$ corresponding with the three Fermi surfaces found in Lorentzian signature in our earlier work [21]. However, for any appreciable value of the dimensionless ratio M_D/λ^2 one only sees a single singularity. As the plots for $G(\omega, k_x)$ at low frequency show, its shape approaches that of the strongly Landau-damped $M_D \rightarrow \infty$ result, Eq. (19), as one increases M_D/λ^2 , though for low M_D it is still distinguishably different. Recall that our results are derived in the limit of large k_F and therefore a realistic ($N_f \sim 1$) value for M_D is $M_D \sim \lambda\sqrt{N_f k_F} \gg \lambda^2$.

This result is in contradistinction to what happens to the bosons. When the bosons are not affected in the IR, i.e., the quenched limit, the fermions are greatly affected by the

boson: There is a topological Fermi surface transition and the low-energy spectrum behaves as critical excitations [21]. However, once we increase M_D to realistic values, the bosonic excitations are rapidly dominated in the IR by Landau damping but we now see that this *reduces* the corrections to the fermions. As M_D is increased for fixed ω, k_x , the deep IR fermion two-point function approaches more and more that of the simple RPA result with self-energy $\Sigma \sim i\omega^{2/3}$.

A more careful analysis of the IR reveals that there are several distinct $\omega \rightarrow 0$ limits of the two-point function:

$$\lim_{\substack{\omega \rightarrow 0, \\ k_x \text{ fixed}}} G_{\text{IR}}(\omega, k_x) \approx e^{\frac{\lambda^2}{6\pi M_D}} \frac{1}{i\omega - k_x - \Sigma_{\text{RPA}}}$$

$$\lim_{\substack{\omega \rightarrow 0, \\ k_x/\omega \text{ fixed}}} G_{\text{IR}}(\omega, k_x) \approx e^{\frac{\lambda^2}{6\pi M_D}} \frac{1}{i\omega - k_x - \frac{4\pi}{3\sqrt{3}} \Sigma_{\text{RPA}}}.$$

However in the case of $\omega^2 \sim l_0 k_x^3$, the full expression for G_{IR} is necessary to describe the low energy limit,

$$\begin{aligned} \lim_{\substack{\omega \rightarrow 0, \\ l_0 k_x^3 \omega^{-2} \text{ fixed}}} G_{\text{IR}}(\omega, k_x) \\ = e^{\frac{\lambda^2}{6\pi M_D}} \left[\frac{1}{i\omega - k_x} \cos\left(\frac{\omega}{l_0^{1/2}(\omega + ik_x)^{3/2}}\right) \right] \end{aligned}$$

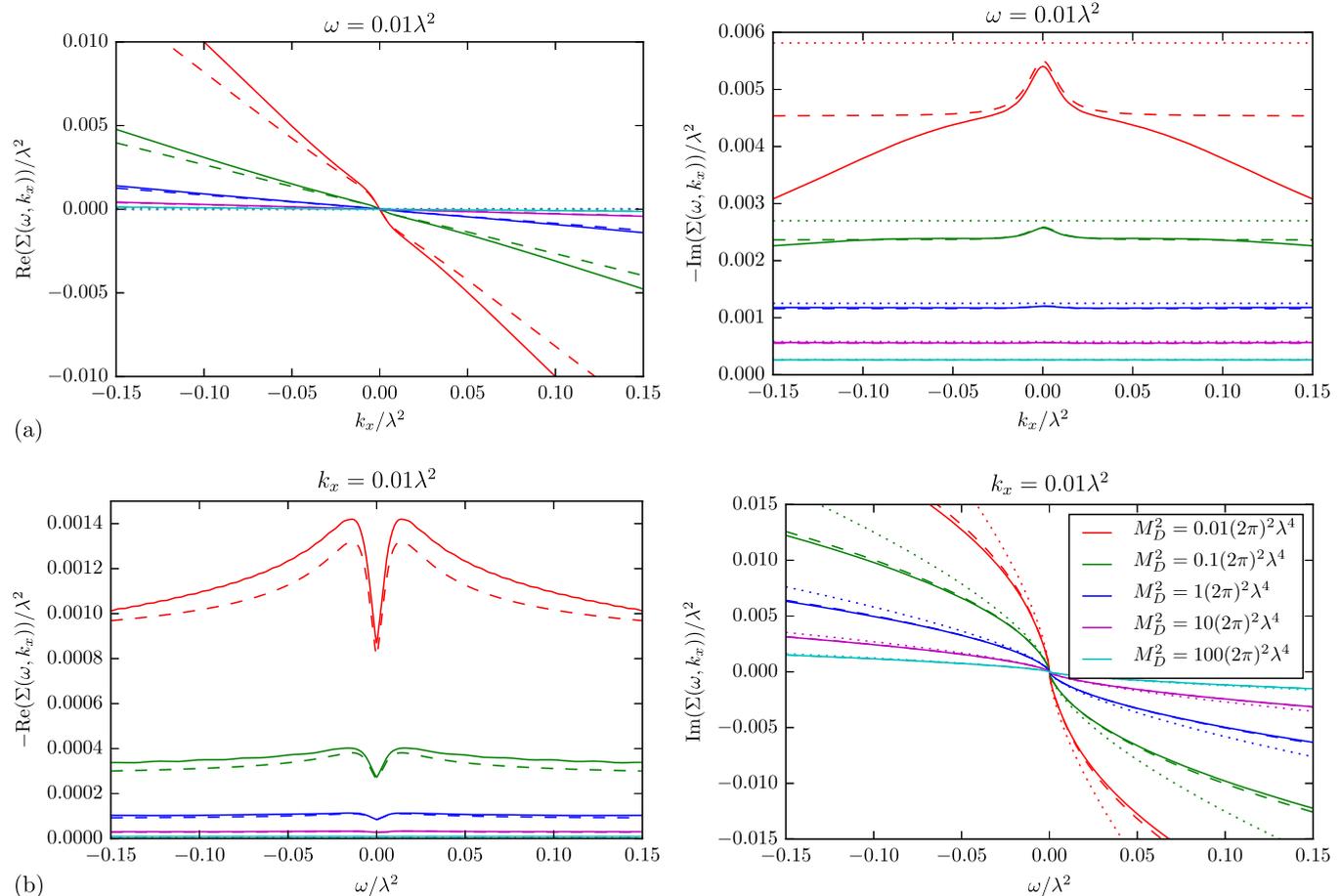


FIG. 5. Real and imaginary parts of the fermion self-energy for (a) $\omega = 0.01\lambda^2$ and for (b) $k_x = 0.01\lambda^2$. Dashed lines show the G_{IR} approximation; dotted lines show the RPA result.

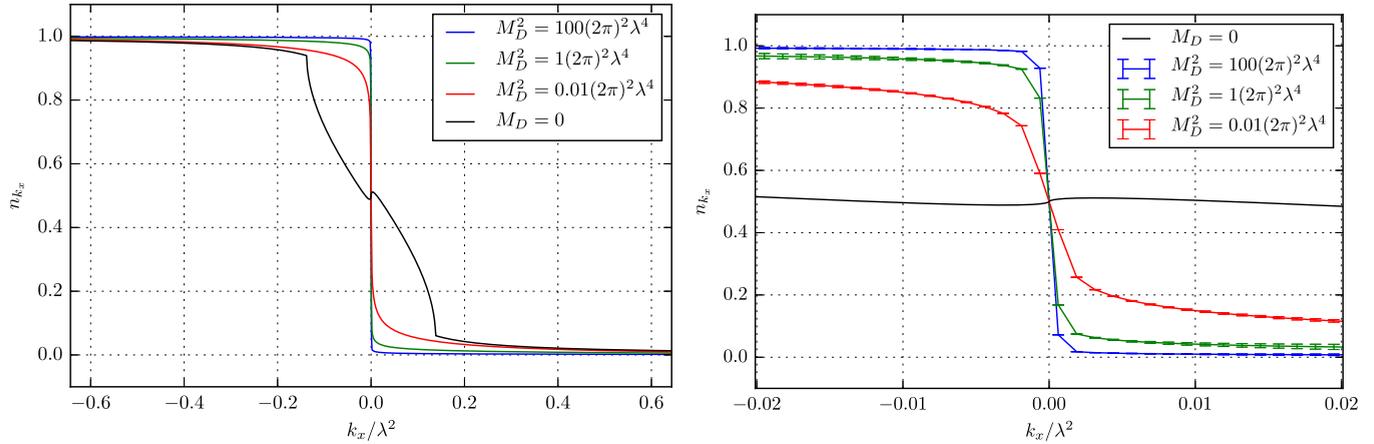


FIG. 6. Momentum distribution function. The two plots show the same function for different ranges. The error bars of the lower figure show an estimate of the error due to using the free Green's function close to the UV. The error bars are exaggerated by a factor 100.

$$\begin{aligned}
 & + \frac{6\sqrt{3}i\Gamma(\frac{1}{3})\omega^{2/3}}{8\pi l_0^{1/3}(\omega + ik_x)^2} {}_1F_2\left(1; \frac{5}{6}, \frac{4}{3}; -\frac{\omega^2}{4l_0(\omega + ik_x)^3}\right) \\
 & + \frac{3\sqrt{3}i\Gamma(-\frac{1}{3})\omega^{4/3}}{8\pi l_0^{2/3}(\omega + ik_x)^3} {}_1F_2\left(1; \frac{7}{6}, \frac{5}{3}; -\frac{\omega^2}{4l_0(\omega + ik_x)^3}\right) \Big].
 \end{aligned}
 \tag{49}$$

This existence of multiple limits shows in fact that the RPA result is never a good low energy (less than M_D) approximation for any value of M_D/λ^2 . We can illustrate this more clearly by studying the self-energy of the fermion $\Sigma(\omega, k_x) = G(\omega, k_x)^{-1} - G_0(\omega, k_x)^{-1}$. It is shown in Fig. 5 that the naive large M_D RPA result (dotted lines) does agree for large M_D at $\omega = 0, k_x = 0$ and the leading ω dependence of the imaginary part is captured. The leading k_x dependence is not captured by RPA. On the other hand, our improved approximation for the low energy regime G_{IR} (dashed lines) captures these higher order terms in the low energy expansion of $G(\omega, k_x)$ very well and also works for finite values of M_D/λ^2 .

In all cases it is clear that the Fermion excitation, though qualitatively sharp, is not a Fermi liquid quasiparticle. We can calculate the occupation number and check whether it is consistent with the non-Fermi-liquid nature of the Green's function. With a Fermi liquid by definition is meant a spectrum with a discontinuity in the zero-temperature momentum distribution function $n_k = \int_{-\infty}^0 d\omega_R A(\omega_R, k_x)/2\pi$ with $A(\omega_R, k_x)$ the spectral function. As the spectral function is the imaginary part of the retarded Green's functions and the latter is analytic in the upper half plane of ω_R we can move this contour to Euclidean ω and use the fact that $G(\omega, k_x)$ approaches $G_0(\omega, k_x)$ in the UV to calculate the momentum distribution function from our Euclidean results. In detail

$$\begin{aligned}
 n_{k_x} & = - \int_{-\infty}^0 d\omega_R \text{Im} \frac{G_R(\omega_R, k_x)}{\pi} \\
 & = \text{Im} \left[\int_0^\Lambda d\omega i \frac{G(\omega, k_x)}{\pi} + \int_C dz i \frac{G(z, k_x)}{\pi} \right]; \tag{50}
 \end{aligned}$$

the first integral can be done with the numerics developed in the preceding section. The contour C goes from Λ to $-\infty$ and for

large enough Λ this is in the UV and can well be approximated by the free propagator. The resulting momentum distribution $n(k)$ is shown in Fig. 6. Within our numerical resolution, these curves are continuous as opposed to a Fermi liquid. This is of course expected; the continuity reflects the absence of a clear pole in the IR expansions in the preceding subsection. Note also that as M_D is lowered, the finite M_D curves approach the quenched result for $|k_x| > k_x^*$ where k_x^* is the point of the discontinuity of the derivative of the quenched occupation number. At k^* the (derivative) of the quenched momentum distribution number does have a discontinuity (reflecting the branch cut found in Ref. [21]).

VI. CONCLUSION

We have presented a nonperturbative answer for the (Euclidean) fermion and boson two-point functions of the elementary quantum critical metal in the double limit of small N_f and large k_F with $N_f k_F$ fixed. This limit was taken order by order in perturbation theory and additionally before performing bosonic momentum integrals.

Our exact results clarify how approximations that have been made in the past hang together. The $N_f \rightarrow 0, k_F \rightarrow \infty$ theory is characterized by two energy scales λ^2 and M_D . For very large ω/λ^2 and ω/M_D^2 one can use perturbation theory in λ to understand the theory. This is the perturbation around the UV-fixed point of a free fermion plus a free boson. In the deep IR region, $\omega \ll M_D^2$, we have an analytic expression for the two-point function, G_{IR} . This shows a $\omega^2 \sim l_0 k_x^3$ scaling. In the intermediate regime, the full numerical results that we have presented are necessary.

The quenched result we obtained earlier [21] does not seem to have a useful regime of validity. As the momentum occupation number $n(k)$ indicates, its precise regime of applicability depends discontinuously on the momentum k/λ^2 . The discontinuity is surprising, but it can be explained analytically as an order of limits ambiguity. Although it is hard to capture the deep IR region for very small N_f in the full numerics we find from G_{IR} that in the ω - k plane there is a region where the limit $N_f \rightarrow 0$ and $\omega, k_x \rightarrow 0$ do not commute. We show an indication of this in Appendix C. Thus for scales below

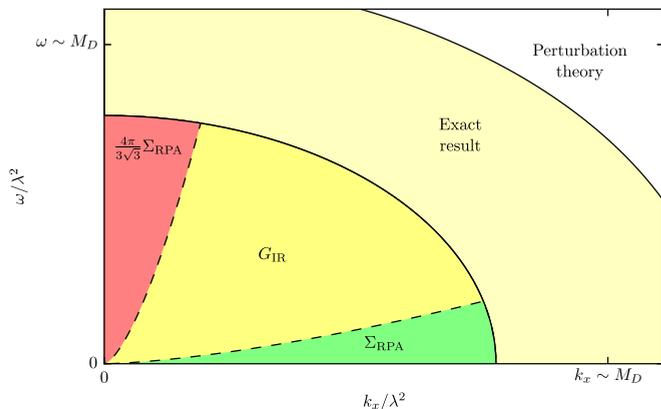


FIG. 7. A sketch of the regimes of applicability of various approximations to the exact fermion Green's function of the elementary quantum critical metal in the double scaling limit. The G_{IR} approximation is applicable both in the deep yellow region and the red and green regions.

λ^2 , the quenched result is not useful. For scales above λ^2 with M_D small, we can use perturbation theory so also in this regime the quenched result is not useful. For small ω and k_x , below M_D , the IR approximation we presented above gives a good description of the physics. We have presented a pictorial overview of how the various approximations are related in Fig. 7.

As argued before the spectrum of the elementary quantum critical metal in the ordered double scaling limit obtained here is not the definitive one. Taking this limit after the bosonic integrals and the perturbative sum may give a different result. Another estimate for the true IR spectrum of the elementary quantum critical metal has been postulated before based on large N_f approximations [25,35,36], but our result at small N_f gives a different IR. Our nonperturbative answer follows from the insight that in the limit of large k_F all fermion loops with more than two external boson legs have cancellations upon symmetrization such that their scaling in k_F is reduced. This higher loop cancellation has been previously put forward as a result of the strong forward scattering approximation. This approximation is engineered such that the Schwinger-Dyson equations combined with the fermion number Ward identity collapse to a closed set of equations. The solution to this closed set is the same dressed fermion correlator that we have presented. However the exact connection between the strong forward scattering approximation and the double scaling limit is not clear. In the double scaling limit here, it is manifest that only the boson propagator is corrected and this in turn implies that the exact fermion Green's function in real space is an exponentially dressed version of the free Green's function.

At the physics level, an obvious next step is therefore to explore the spectrum of the elementary quantum critical metal including corrections in N_f that are not proportional to k_f . Since this necessarily involves higher-point boson correlations, the role of self-interactions of the boson needs to be considered. These are also relevant in the IR and may therefore give rise to qualitatively very different physics than found here. We leave this for future work.

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APPENDIX A: MULTILoop CANCELLATION

The robustness of the one-loop result in case of linear fermionic dispersion was recognized before under the name of multiloop cancellation [24]. The technical result is that for a theory with a simple Yukawa coupling and linear dispersion around a Fermi surface, a symmetrized fermion loop with more than two fermion lines vanishes. In our context the linear dispersion is a consequence of the large k_F limit. In other words all higher loop contributions to the polarization Π should be subleading in $1/k_F$. This was explicitly demonstrated at two loops in Ref. [37].

We will give here a short derivation of this multiloop cancellation in the limit $k_F \rightarrow \infty$. We may assume in this limit that the momentum transfer at any fermion interaction is always much smaller than the size of the initial (\vec{k}) and final momenta (\vec{k}') which are of the order of the Fermi momentum, i.e., $|\vec{k}' - \vec{k}| \ll k_F$ with $|\vec{k}|, |\vec{k}'| \sim k_F$. The free fermion Green's function then reflects a linear dispersion

$$G_0(\omega, k) = \frac{1}{i\omega - vk}. \quad (\text{A1})$$

We now Fourier transform back to real space, as multiloop cancellation is most easily shown in this basis. The real space transform of the “linear” free fermion propagator above is

$$G_0(\tau, r) = -\frac{i}{2\pi} \frac{\text{sgn}(v)}{r + i\tau}, \quad (\text{A2})$$

where as before r is the conjugate variable to $k = |\vec{k}| - k_F$. The essential step in the proof is that real space Green's function manifestly obeys the identity [21]

$$G_0(z_1)G_0(z_2) = G_0(z_1 + z_2)(G_0(z_1) + G_0(z_2)) \quad (\text{A3})$$

with $z \equiv r + i\tau$. Consider then (the subpart of any correlation function/Feynman diagram containing) a fermion loop with $n \geq 2$ vertices along the loop connected to indistinguishable scalars (i.e., no derivative interactions and all interactions are symmetrized). The corresponding algebraic expression in a real space basis will then contain the expression

$$F(z_1, \dots, z_n) = \sum_{(i_1, \dots, i_n) \in S_n} G_0(z_{i_1} - z_{i_2})G_0(z_{i_2} - z_{i_3}) \dots \times G_0(z_{i_{n-1}} - z_{i_n})G_0(z_{i_n} - z_{i_1}), \quad (\text{A4})$$

where S_n is the set of permutations of the numbers 1 through n . Using the “linear dispersion” identity Eq. (A3) and the

shorthand notation $G_0(z_{i_1} - z_{i_2}) = G_{12}$ we obtain

$$F(z_1, \dots, z_n) = \sum_{(i_1, \dots, i_n) \in S_n} G_{12} G_{23} \dots G_{n-1,1} (G_{n-1,n} + G_{n,1}). \quad (\text{A5})$$

Next we cyclically permute the indices from 1 to $n - 1$: $1 \rightarrow 2$, $2 \rightarrow 3, \dots, n - 1 \rightarrow 1$ in the sum

$$\begin{aligned} & \sum_{(i_1, \dots, i_n) \in S_n} G_{12} G_{23} \dots G_{n-1,1} G_{n-1,n} \\ &= \sum_{(i_1, \dots, i_n) \in S_n} G_{23} G_{34} \dots G_{12} G_{1,n}. \end{aligned} \quad (\text{A6})$$

This gives us

$$F(z_1, \dots, z_n) = \sum_{(i_1, \dots, i_n) \in S_n} G_{12} G_{23} \dots G_{n-1,1} (G_{1,n} + G_{n,1}). \quad (\text{A7})$$

Then since $G_{i,j}$ corresponds to a (spinless) fermionic Green's function, it is antisymmetric $G_{i,j} = -G_{j,i}$, and we can conclude that F vanishes for $n \geq 3$. For $n = 2$ it is not possible to use the identity Eq. (A3) since we would need to evaluate $G(0)$ which is infinite.

APPENDIX B: THE FOURIER TRANSFORM OF THE FERMION GREEN'S FUNCTION IN THE LARGE M_D APPROXIMATION

We will now show how to perform this Fourier transform. We need to calculate the following integral:

$$\begin{aligned} G_f(\omega, k) &= \int d\tau dx \frac{e^{i(\omega\tau - kx)}}{2\pi(ix - v\tau)} \\ &\times \exp\left(-\frac{|x|}{l_0^{1/3}(|x| + iv \operatorname{sgn}(x)\tau)^{2/3}}\right). \end{aligned} \quad (\text{B1})$$

First we note that the integrand is τ analytic in the region $\{\tau \in \mathbb{C} : \min(0, vx) < \operatorname{Im}(\tau) < \max(0, vx)\}$. Since the integrand necessarily goes to zero at ∞ we can thus shift the τ contour, $\tau \rightarrow \tau + ix/v$. We now have

$$\begin{aligned} G_f(\omega, k) &= - \int \frac{d\tau dx}{2\pi v\tau} \exp \\ &\times \left(i\omega\tau - x \left(ik + \omega/v + \frac{\operatorname{sgn}(x)}{l_0^{1/3}(iv \operatorname{sgn}(x)\tau)^{2/3}} \right) \right). \end{aligned} \quad (\text{B2})$$

We have allowed ourselves to choose the order of integration, shift the contour, and then change the order. We see that the x integral now is trivial but only converges for

$$-\frac{v^{1/3}}{2l_0^{1/3}|\tau|^{2/3}} < \operatorname{Re}(\omega) < \frac{v^{1/3}}{2l_0^{1/3}|\tau|^{2/3}}. \quad (\text{B3})$$

This is fine since we know that the final result is ω analytic in both the right and left open half planes. As long as we can

obtain an answer valid within open subsets of both of these sets we can analytically continue the found solution to the whole half planes. We thus proceed assuming $\operatorname{Re}(\omega)$ is in this range. One can further use symmetries of our expression to relate the left and right ω half planes, so to simplify matters, from now on we additionally assume ω to be positive. Let us now consider a negative x ; we then see that the τ integral can be closed in the upper half plane and since it is holomorphic there the result will be 0. We can thus limit the x integrals to \mathbb{R}^+ . We then have

$$G_f(\omega, k) = \int d\tau \frac{e^{i\omega\tau}}{2\pi v\tau} \frac{-1}{a + \frac{1}{l_0^{1/3}(iv\tau)^{2/3}}}, \quad (\text{B4})$$

where $a = ik + \omega/v$. The integrand has a pole at $(i\tau)^{2/3} = -1/(al_0^{1/3}v^{2/3})$. Now break the integral in positive and negative τ and write it as

$$G_f(\omega, k) = \frac{h((-i)^{2/3}u_0^*)^* - h((-i)^{2/3}u_0)}{2\pi va}, \quad (\text{B5})$$

where

$$\begin{aligned} h(u) &= \int_0^\infty d\tau \frac{e^{i\tau}}{\tau + u\tau^{1/3}}, \\ u_0 &= \frac{\omega^{2/3}}{al_0^{1/3}v^{2/3}}. \end{aligned} \quad (\text{B6})$$

Using Morera's theorem we can prove that h is holomorphic on \mathbb{C}/\mathbb{R}^- . Consider any closed curve C in \mathbb{C}/\mathbb{R}^- . We need to show that

$$\int_C du h(u) = 0. \quad (\text{B7})$$

We do this by rotating the contour slightly counterclockwise. For any curve C there is clearly a small $\epsilon > 0$ such that we will still not hit the pole $\tau + u\tau^{1/3} = 0$.

$$\int_C du \int_0^{(1+i\epsilon)\infty} d\tau \frac{e^{i\tau}}{\tau + u\tau^{1/3}}, \quad (\text{B8})$$

The piece at ∞ converges without the denominator and thus goes to 0. Since the integral now converges absolutely we can use Fubini's theorem to change the orders of integration

$$\int_0^{(1+i\epsilon)\infty} d\tau \int_C du \frac{e^{i\tau}}{\tau + u\tau^{1/3}} = 0. \quad (\text{B9})$$

Since also the integrand is holomorphic on a connected open set containing C the proof is finished.

The function h can for $0 < \arg(u) < 2\pi/3$ be expressed as a Meijer G function:

$$h(u) = \frac{3}{8\pi^{5/2}} G_{3,5}^{5,3} \left(0, \frac{1}{3}, \frac{2}{3} \middle| -\frac{u^3}{4} \right). \quad (\text{B10})$$

The G function has a branch cut at \mathbb{R}^- and because of that we can not easily write $h(u)$ in terms of it since the argument u appears cubed in the argument of the G function and we know that h is holomorphic on all of \mathbb{C}/\mathbb{R}^- . We can however express h as an analytic continuation of the G function past this branch cut, onto the further sheets of its Riemann surface (Fig. 8).

We will write the G function as a function with two real arguments, first the absolute value and second the phase of its otherwise complex argument. We will allow the phase to be any real number and when outside the range $[-\pi, \pi]$, we let the function be defined by its analytical continuation to the corresponding sheet. We hereafter omit the constant parameters of the G function. We then have:

$$\begin{aligned} G_f(\omega, k) &= 3 \frac{G_{3,5}^{5,3}(\frac{|u_0|^3}{4}, -2\pi - 3 \arg(u_0))^* - G_{3,5}^{5,3}(\frac{|u_0|^3}{4}, -2\pi + 3 \arg(u_0))}{16\pi^{7/2}va} \\ &= 3 \frac{G_{3,5}^{5,3}(\frac{|u_0|^3}{4}, 3 \arg(u_0) + 2\pi) - G_{3,5}^{5,3}(\frac{|u_0|^3}{4}, 3 \arg(u_0) - 2\pi)}{16\pi^{7/2}va}. \end{aligned} \quad (\text{B11})$$

In the last step we used the fact that the G function commutes with complex conjugation. We see that the Green's function is given by a certain monodromy of the G function. It is given by the difference in its value starting at a point $u_0^3/4$ on the sheet above the standard one and then analytically continuing clockwise around the origin twice to the sheet below the standard one and there return to $u_0^3/4$. Since we only need this difference, we might expect this to be a, in some sense, simpler function as would happen for, e.g., monodromies of the logarithm. To see how to simplify this we look at the definition of the G function. It is defined as an integral along L :

$$G_{p,q}^{m,n} \left(\begin{matrix} a_1, \dots, a_p \\ b_1, \dots, b_q \end{matrix} \middle| z \right) = \frac{1}{2\pi i} \int_L \frac{\prod_{j=1}^m \Gamma(b_j - s) \prod_{j=1}^n \Gamma(1 - a_j + s)}{\prod_{j=m+1}^q \Gamma(1 - b_j + s) \prod_{j=n+1}^p \Gamma(a_j - s)} z^s ds. \quad (\text{B12})$$

There are a few different options for L and which one to use depends on the arguments. In our case L starts and ends at $+\infty$ and circles all the poles of $\Gamma(b_j - s)$ in the negative direction. Using the residue theorem we can recast the integral to a series. We have double poles at all negative integers and some simple poles in between. The calculation to figure out the residues of all these single and double poles is a bit too technical to present here but in the end the series can be written as

$$G_{3,5}^{5,3}(z) = \sum_{n=0}^{\infty} (a_n z^n + b_n z^n \log(z) + c_n z^{n+1/3} + d_n z^{n+1/2} + e_n z^{n+2/3}). \quad (\text{B13})$$

Now we perform the monodromy term by term and a lot of these terms cancel out.

$$G_{3,5}^{5,3}(|z|, \arg(z) + 2\pi) - G_{3,5}^{5,3}(|z|, \arg(z) - 2\pi) = \sum_{n=0}^{\infty} (4\pi i b_n z^n + i\sqrt{3}c_n z^{n+1/3} - i\sqrt{3}e_n z^{n+2/3}) \quad (\text{B14})$$

The coefficients a_i contain both the harmonic numbers and the polygamma function whereas the other coefficients are just simple products of gamma functions. This simplification now lets us sum this series to a couple of generalized hypergeometric functions. Inserting the expressions for a and u_0 we have

$$\begin{aligned} G_f(\omega, k) &= \frac{1}{i\omega - kv} \cos\left(\frac{\omega}{v l_0^{1/2}(\omega/v + ik)^{3/2}}\right) + \frac{6\sqrt{3}i\Gamma(\frac{1}{3})\omega^{2/3}}{8\pi l_0^{1/3}v^{5/3}(\omega/v + ik)^2} {}_1F_2\left(1; \frac{5}{6}, \frac{4}{3}; -\frac{\omega^2}{4l_0v^2(\omega/v + ik)^3}\right) \\ &+ \frac{3\sqrt{3}i\Gamma(-\frac{1}{3})\omega^{4/3}}{8\pi l_0^{2/3}v^{7/3}(\omega/v + ik)^3} {}_1F_2\left(1; \frac{7}{6}, \frac{5}{3}; -\frac{\omega^2}{4l_0v^2(\omega/v + ik)^3}\right). \end{aligned} \quad (\text{B15})$$

Note that this expression is ω holomorphic for ω in the right half plane so our previous assumptions on ω can be relaxed as long as ω is in the right half plane. We note from expression (B1) that if we change sign on both ω and k and do the changes of variables $\tau \rightarrow -\tau$ and $x \rightarrow -x$ we end up with the same integral up to an overall minus sign. We can thus get the left half plane result using the relation

$$G_f(-\omega, k) = - G_f(\omega, -k). \quad (\text{B16})$$

As mentioned in the main text, this expression has been compared with numerics to verify that we have not made any mistakes. See Fig. 2. We have also done the two integrals for the Fourier transform in the opposite order, first obtaining a different Meijer G function then using the G-function convolution theorem to do the second integral. In the end one obtains the same monodromy of the G function as above. This

expression can also be found in Appendix A.2 of Ref. [25], however it was not entirely clear from the phrasing of the last paragraph that this function is actually the exact Fourier transform, Eq. (B1).

APPENDIX C: THE DISCONTINUOUS TRANSITION FROM THE QUENCHED TO THE LANDAU-DAMPED REGIME

We show here why including Landau damping (finite M_D) physics starting from the quenched $N_f \rightarrow 0$ result is discontinuous in the IR. To do so we calculate the Green's function by imposing the $N_f \rightarrow 0$ limit from the beginning. We need to evaluate the Fourier transform integral:

$$G_L(\omega, k) = \int_{-\infty}^{\infty} dx \int_{-L}^L d\tau G_0 \exp(I_0 + i\omega \cdot \tau - ik \cdot x), \quad (\text{C1})$$

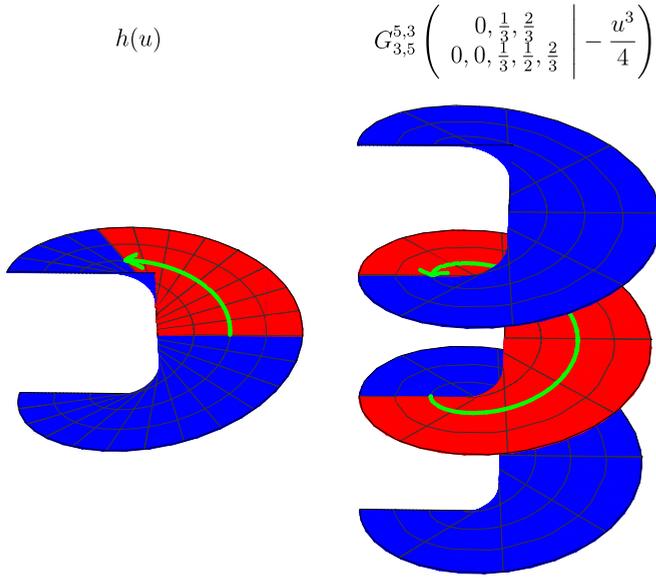


FIG. 8. The left image shows part of the Riemann surface of the function h and the right image shows part of the Riemann surface of the G function. The part in red of the left image can be expressed as the G function evaluated on its first sheet, the red part of the right image. The green arrow shows how points are mapped from one Riemann surface to the other by the mapping $u \mapsto -u^3/4$.

where as before the free propagator is

$$G_0(\tau, x) = -\frac{i}{2\pi} \frac{1}{x + i \cdot \tau} \quad (C2)$$

and the $N_f \rightarrow 0$ limit of the exponent of the real space Green's function:

$$I_0 = \frac{(\tau + i \cdot x)^2}{12\pi \sqrt{\tau^2 + x^2}}. \quad (C3)$$

Note however that the τ integral is divergent in this limit. Therefore in (C1) we have introduced a cutoff L in this direction. By looking at the full expansion of I we can see that the natural value of L is of the order $1/M_D$. For larger values of τ the asymptotic expansion describes I better. We expect that for large enough momenta and frequencies the asymptotic region does not contribute to the Fourier transform and therefore the cutoff can be removed. This naive expectation however is only partly true. We will shortly see that the region in the ω - k plane where the cutoff can be removed is more complicated and asymmetric in terms of the momenta and frequency.

Let us turn now to the evaluation of (C1). After making the coordinate change $\tau \rightarrow \tau, x \rightarrow u \cdot \tau$ one of the integrals (τ) can be evaluated analytically:

$$G_L(\omega, k) = \int_{-\infty}^{\infty} du (I_1(u) + I_2(u) + I_3(u)), \quad (C4)$$

where

$$I_1(u) = \frac{6 \exp\left(ikLu\sqrt{u^2 + 1} - iL \cdot \omega \cdot \sqrt{u^2 + 1} - \frac{L(u-i)^2}{12\pi}\right)}{12\pi(u+i)(ku-\omega) + i\sqrt{u^2 + 1}u + \sqrt{u^2 + 1}}, \quad (C5)$$

$$I_2(u) = \frac{6i \exp\left(-ikLu\sqrt{u^2 + 1} + iL \cdot \omega \cdot \sqrt{u^2 + 1} - \frac{L(u-i)^2}{12\pi}\right)}{12i\pi k(u+i)u + \sqrt{u^2 + 1}u - i\sqrt{u^2 + 1} + 12\pi(1-iu)\omega}, \quad (C6)$$

$$I_3(u) = -\frac{144\pi^2 k^2 (u+i) + u - 3i}{u(-3 + u(144\pi^2 k^2 (u+i) + u - 3i)) - 288\pi^2 ku(u+i)\omega + 144\pi^2 (u+i)\omega^2 + i}. \quad (C7)$$

By numerically performing the single integral u we can obtain $G_{M_D \rightarrow 0}$. Since it is easier than evaluating the Fourier transform

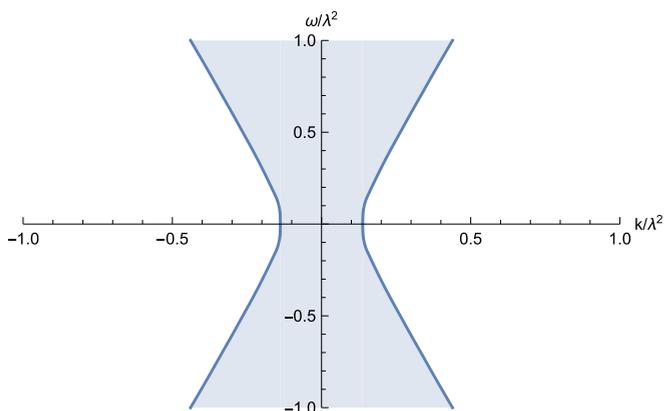


FIG. 9. The region of convergence. In the shaded area the $L \rightarrow \infty$ ($N_f \rightarrow 0$) limit is not convergent while outside this area $\lim_{L \rightarrow \infty} G_L = G_{\text{quenched}}$.

of the true real-space version of the Green's function it is worth understanding how the $L \rightarrow \infty$ (which is equivalent to $M_D \rightarrow 0$) works. The result is depicted on Fig. 9. In the shaded region (which corresponds to small k) the limit is not well defined while outside of this region the limit is equal to the quenched result. The numerics shows that for zero frequency, the edge of this region is at k^* , where the Green's function is singular.

We can qualitatively determine the line separating the convergent and divergent region. For this we assume that when L is large one can expand the exponent in u

$$I_1(u) \sim \exp\left[u^2\left(-\frac{1}{12\pi} - \frac{1}{2}i\omega\right)L + \frac{i u(6\pi k + 1)}{6\pi}L - iL\omega + \frac{L}{12\pi} + \mathcal{O}(Lu^3)\right]. \quad (C8)$$

We see that because of the term $-L \cdot u^2/(12\pi)$, the integrand is nonzero only in a narrow region around $u = 0$. For the same

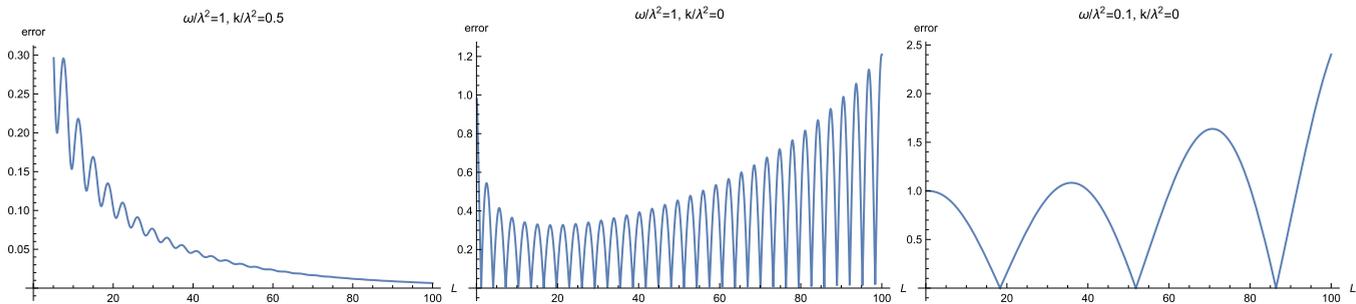


FIG. 10. The relative difference between G_L and the quenched result as a function of L for different frequencies and momenta. Left: for a point in the (ω, k) plane which is outside of the shaded region of Fig. 9 the “error” goes to zero for large L . Middle: inside the shaded region the error is oscillating with diverging amplitude as L goes to larger values. However, for larger values of ω there is an intermediate range of L , where the relative error is smaller than 0.3. Therefore the quenched approximation is qualitatively correct. Right: for small ω the amplitude of the oscillation is large.

reason we approximate the denominator of (C5) by replacing u by zero there. With these simplifications we arrive to a gaussian integral which can be evaluated analytically:

$$\int_{-\infty}^{\infty} I_1(u) du \approx \frac{12i\sqrt{3}\pi \exp\left(\frac{iL(6\pi k^2 + 2k + \omega(-12\pi\omega + i))}{12\pi\omega - 2i}\right)}{(12\pi\omega + i)\sqrt{L(1 + 6i\pi\omega)}}. \quad (\text{C9})$$

The real part of (C9) is

$$L \frac{3(\omega^2 - k^2)\pi - k}{36\pi^2\omega^2 + 1}. \quad (\text{C10})$$

It is clear that if this value is positive (i.e., $3(\omega^2 - k^2)\pi - k > 0$) than the $L \rightarrow \infty$ limit is divergent. Looking at the numerical result in Fig. 9 we indeed see that the boundary of the shaded region is indeed a hyperbola. Note, however, that the exact location of this hyperbola obtained from expanding the exponent is slightly off.

It is interesting to note that if we are in the divergent region G_L is not convergent for large L but for some intermediate values it can still get close to the quenched result G_{quenched} . To quantify this let us introduce a relative “error” function

$$\text{error}(L) = \left| \frac{G_L - G_{\text{quenched}}}{G_{\text{quenched}}} \right|. \quad (\text{C11})$$

In Fig. 10 we show the behavior of this function for various points in the ω - k plane. For a point which is outside the shaded region the error approaches zero when L is large. For $\omega = 1, k = 0$ which is inside the divergent region the error is oscillating but there is an interval of L where the amplitude of this oscillation has a minimum. If the frequency is small ($\omega = 0.1$), the amplitude is larger.

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