Anomalous quasiparticle lifetime in geometric quantum critical metals

arXiv: 2310.07539

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Collaborators



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Reference:

Song, Ma, Kallin, Lee, "Anomalous quasiparticle lifetime in geometric quantum critical metals", arXiv: 2310.07539.

Outline

□ Motivation and our main results

Derivations and details of our results

Gamma Summary and outlook

Transition of fermi surface **topology**



Transition of fermi surface geometry



• cubic inflection point $E_{\mathbf{K}+\mathbf{k}} = v_F k_n + k_t^3$ generic point on Fermi Surface $E_{\mathbf{K}+\mathbf{k}} = v_F k_n + a k_t^2 + o(|k_n| + k_t^2)$

Questions

- Are there any *qualitatively* different low energy physics caused by the appearance of Fermi Surface *inflection* points?
- If yes, can we make any *simple* and *universal* predictions about it?



 $E_{\mathbf{K}+\delta\mathbf{k}} = v_F \delta k_n + \left(\delta k_t\right)^m$

Our main results

Qualitatively different low energy physics due to the Fermi Surface inflection points?

Our results (based on random phase approximation & patch approximation): • Near **K** with $E_{\mathbf{K}+\delta\mathbf{k}} = v_F \delta k_n + (\delta k_t)^m$, the quasiparticle decay rate \propto



Our main results

Qualitatively different low energy physics due to the Fermi Surface inflection points?

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$$\left|\Im\Sigma_{\mathbf{k}}\left(E_{\mathbf{k}}\right)\right| \stackrel{\mathbf{k} \xrightarrow{m} \mathbf{K}}{\propto} \begin{cases} E_{\mathbf{k}}^{2} \ln E_{\mathbf{k}}, & m = 2, \\ E_{\mathbf{k}}^{\frac{m}{m-1}}, & m > 2. \end{cases}$$

$$\frac{d \ge 3}{\Im \Sigma_{\mathbf{k}} (E_{\mathbf{k}})} \stackrel{E_{\mathbf{k}} \to 0}{\sim} E_{\mathbf{k}}^{2} \qquad E_{\mathbf{k}}^{2} \ln E_{\mathbf{k}} \qquad E_{\mathbf{k}}^{1-\alpha} d$$

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Outline of our calculation

Bare Hamiltonian

$$H = \int \frac{d^2 \mathbf{k}}{(2\pi)^2} \xi_{\mathbf{k}} \psi_{\mathbf{k}}^{\dagger} \psi_{\mathbf{k}} + \int \frac{d^2 \mathbf{q}}{(2\pi)^2} \int \frac{d^2 \mathbf{k}_1}{(2\pi)^2} \int \frac{d^2 \mathbf{k}_2}{(2\pi)^2} \frac{V_{\mathbf{q}}}{2} \psi_{\mathbf{k}_1 + \mathbf{q}}^{\dagger} \psi_{\mathbf{k}_2 - \mathbf{q}}^{\dagger} \psi_{\mathbf{k}_2} \psi_{\mathbf{k}_1}, \ V_{\mathbf{q}} = \frac{e^2}{|\mathbf{q}|}.$$

Self-consistent equations

$$\begin{split} G\left(\mathbf{p},\varepsilon\right) &= \frac{1}{-\varepsilon\left(1+i0^{+}\right)+\xi_{\mathbf{p}}+\Sigma\left(\mathbf{p},\varepsilon\right)},\\ \Sigma\left(\mathbf{p},\varepsilon\right) &= -i\int_{\mathbf{q},\omega}\mathcal{V}\left(\mathbf{q},\omega\right)G\left(\mathbf{p}-\mathbf{q},\varepsilon-\omega\right),\\ \mathcal{V}\left(\mathbf{q},\omega\right) &= \frac{1}{V_{\mathbf{q}}^{-1}-\Pi\left(\mathbf{q},\omega\right)},\\ \Pi\left(\mathbf{q},\omega\right) &= -i\int_{\mathbf{p},\varepsilon}G\left(\mathbf{p},\varepsilon\right)G\left(\mathbf{p}+\mathbf{q},\varepsilon+\omega\right). \end{split}$$

$$E_{\mathbf{p}} \implies \operatorname{Im}\Sigma\left(\mathbf{p}, E_{\mathbf{p}}\right) \sim \frac{1}{\tau_{\mathbf{p}}}$$

Dyson's equation for Green's function

leading order of self-energy

screened interaction

polarization in RPA

p+q

Self-consistency computation

$$G(\mathbf{p},\varepsilon) = \frac{1}{-\varepsilon(1+i0^{+}) + \xi_{\mathbf{p}} + \Sigma(\mathbf{p},\varepsilon)}$$

$$\simeq \frac{Z_{\mathbf{p}}}{-\varepsilon + E_{\mathbf{p}} - i \operatorname{sgn}(\varepsilon) \tau_{\mathbf{p}}^{-1}}$$

$$G(\mathbf{p},\varepsilon) \approx \frac{1}{-\varepsilon(1+i0^{+}) + E_{\mathbf{p}}}$$

$$E_{\mathbf{p}} = \xi_{\mathbf{p}} + \operatorname{Re}\Sigma(\mathbf{p}, E_{\mathbf{p}})$$

$$\tau_{\mathbf{p}}^{-1} \sim \left|\operatorname{Im}\Sigma\left(\mathbf{p}, E_{\mathbf{p}}\right)\right|$$

$$\Sigma(\mathbf{p},\varepsilon) = -i \int_{\mathbf{q},\omega} \mathcal{V}(\mathbf{q},\omega) G(\mathbf{p} - \mathbf{q},\varepsilon - \omega)$$

$$\Pi(\mathbf{q},\omega) = -i \int_{\mathbf{p},\varepsilon} G(\mathbf{p},\varepsilon) G(\mathbf{p} + \mathbf{q},\varepsilon + \omega)$$

$$\mathcal{V}(\mathbf{q},\omega) = \frac{1}{V_{\mathbf{q}}^{-1} - \Pi(\mathbf{q},\omega)}$$

Step 1: compute $\Pi(q, \omega)$

$$\Pi(\mathbf{q},\omega) = -i \int_{\mathbf{p},\varepsilon} G(\mathbf{p},\varepsilon) G(\mathbf{p}+\mathbf{q},\varepsilon+\omega) = \int_{\mathbf{p}} \frac{\Theta(-E_{\mathbf{p}+\mathbf{q}}) - \Theta(-E_{\mathbf{p}})}{-\omega(1+i0^+) + E_{\mathbf{p}+\mathbf{q}} - E_{\mathbf{p}}}$$

For small
$$\omega \to 0$$
, $\Pi(\mathbf{q}, 0) = \int_{\mathbf{p}} \frac{\Theta(-E_{\mathbf{p}+\mathbf{q}}) - \Theta(-E_{\mathbf{p}})}{E_{\mathbf{p}+\mathbf{q}} - E_{\mathbf{p}}} = -\int_{\mathbf{p}} \delta(E_{\mathbf{p}})$

density of quasiparticles at Fermi energy

$$\operatorname{Im}\Pi(\mathbf{q},\omega) = \pi\operatorname{sgn}(\omega)\int_{\mathbf{p}} \left(\Theta(-E_{\mathbf{p}+\mathbf{q}}) - \Theta(-E_{\mathbf{p}})\right)\delta\left(E_{\mathbf{p}+\mathbf{q}} - E_{\mathbf{p}} - \omega\right) \sim \sum_{\mathbf{N}\in\mathcal{N}_{-\mathbf{q}}} \frac{-|\omega|}{\left|\mathbf{v}_{\mathbf{N}+\mathbf{q}}\times\mathbf{v}_{\mathbf{N}}\right|}$$

$$\Pi(\mathbf{q},\omega) \sim -1 - i \sum_{\mathbf{N} \in \mathcal{N}_{-\mathbf{q}}} \frac{|\omega|}{|\mathbf{v}_{\mathbf{N}+\mathbf{q}} \times \mathbf{v}_{\mathbf{N}}|} \text{ for the purpose of determining scaling behavior}$$

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Step 1: compute $\Pi(q, \omega)$



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Step 2: compute $\mathcal{V}(q, \omega)$

For the purpose of determining scaling behavior, $\Pi(\mathbf{q}, \omega) \sim -1 - i \sum_{\mathbf{N} \in \mathscr{N}_{-\mathbf{q}}} \frac{|\omega|}{|\mathbf{v}_{\mathbf{N}+\mathbf{q}} \times \mathbf{v}_{\mathbf{N}}|}$

$$\mathcal{V}(\mathbf{q},\omega) = \frac{1}{V_{\mathbf{q}}^{-1} - \Pi(\mathbf{q},\omega)} \sim \frac{1}{\frac{|\mathbf{q}|}{2\pi e^2} + 1 - i \operatorname{Im}\Pi(\mathbf{q},\omega)} \sim \frac{1}{1 - i \operatorname{Im}\Pi(\mathbf{q},\omega)}$$

Coulomb potential is "screened".

Step 3: compute imaginary part of $\Sigma(p, \varepsilon)$

Use the spectral representation $f^{R}(\mathbf{p}, \varepsilon) = \int \frac{d\varepsilon'}{\pi} \frac{\operatorname{Im} f^{R}(\mathbf{p}, \varepsilon')}{-(\varepsilon + i0^{+}) + \varepsilon'}$ for the retarded Green's functions.

$$\Sigma(\mathbf{p},\varepsilon) = -i \int_{\mathbf{q},\omega} \mathcal{V}(\mathbf{q},\omega) G(\mathbf{p}-\mathbf{q},\varepsilon-\omega)$$

$$\Sigma(\mathbf{p},\varepsilon) = \int \frac{d^2q}{(2\pi)^2} \int \frac{d\omega'}{\pi} \int \frac{d\varepsilon'}{\pi} \frac{\operatorname{Im} \mathcal{V}^R(\mathbf{q},\omega') \operatorname{Im} G^R(\mathbf{p}-\mathbf{q},\varepsilon')}{-\varepsilon (1+i0^+)+\varepsilon'+\omega'} \left[\Theta(\varepsilon')-\Theta(-\omega')\right]$$

$$\operatorname{Im} \Sigma(\mathbf{p},\varepsilon) \approx -\operatorname{sgn}(\varepsilon) \int \frac{d^2p'}{(2\pi)^2} \left|\operatorname{Im} \mathcal{V}^R\left(\mathbf{p}-\mathbf{p}',\varepsilon-E_{\mathbf{p}'}\right)\right| \Theta\left(E_{\mathbf{p}'}\left(\varepsilon-E_{\mathbf{p}'}\right)\right)$$

$$\tau_{\mathbf{p}}^{-1} \sim \left| \operatorname{Im} \Sigma \left(\mathbf{p}, E_{\mathbf{p}} \right) \right| \sim \int_{\mathbf{p}'} \left| \operatorname{Im} \mathcal{V} \left(\mathbf{p} - \mathbf{p}', E_{\mathbf{p}} - E_{\mathbf{p}'} \right) \right| \Theta \left(E_{\mathbf{p}'} \left(E_{\mathbf{p}} - E_{\mathbf{p}'} \right) \right)$$

Step 4: compute τ_p

$$\frac{1}{\tau_{\mathbf{p}}} \sim \int_{\mathbf{p}'} \left| \operatorname{Im} \mathcal{V} \left(\mathbf{p} - \mathbf{p}', E_{\mathbf{p}} - E_{\mathbf{p}'} \right) \right| \Theta \left(E_{\mathbf{p}'} \left(E_{\mathbf{p}} - E_{\mathbf{p}'} \right) \right) \sim \int_{\mathbf{p}'} \left| \operatorname{Im} \frac{\Theta \left(E_{\mathbf{p}'} \left(E_{\mathbf{p}} - E_{\mathbf{p}'} \right) \right)}{1 + i \frac{|\omega|}{|\mathbf{v}_{\mathbf{p}} \times \mathbf{v}_{\mathbf{p}'}|}} \right|$$

The decay rate due to scattering between unparalleled patches: $\tau_{\mathbf{p}}^{\#} \sim E_{\mathbf{p}}^{2}$ **p**' Focusing on one patch $\frac{1}{\tau_{\mathbf{p}}^{\parallel}} \sim \begin{cases} \frac{1}{a_{2}} E_{\mathbf{p}}^{2} \ln \left| \frac{1}{E_{\mathbf{p}}} \right|, & n = 2\\ a_{n}^{-\frac{1}{n-1}} E_{\mathbf{p}}^{\frac{n}{n-1}}, & n > 2 \end{cases}$ Focusing on one patch *n*: order of inflection point on a parallel patch

Step 4: compute τ_p

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An <u>exact</u> computation is possible for scattering within a single patch:

$$\frac{1}{\tau_{\mathbf{p}}^{\parallel}} \sim \int_{E_{\mathbf{p}'},k_{\parallel}'} \left| \operatorname{Im} \frac{\Theta\left(E_{\mathbf{p}'}\left(E_{\mathbf{p}}-E_{\mathbf{p}'}\right)\right)}{1-i\frac{1}{a_{n}}\frac{E_{\mathbf{p}}-E_{\mathbf{p}'}}{\left|k_{\parallel}^{\prime n-1}\right|}} \right|$$
$$\sim \begin{cases} \frac{1}{a_{2}}E_{\mathbf{p}}^{2}\ln\frac{1}{E_{\mathbf{p}}}, & n=2, \\ \frac{1}{a_{n}^{1/(n-1)}}E_{\mathbf{p}}^{\frac{n}{n-1}}, & n>2, \end{cases}$$
for $\mathbf{p} = \mathbf{K} + (k_{\perp}, 0).$

$$\mathbf{p}' = \mathbf{K} + \mathbf{k}'$$

$$\mathbf{K} \quad \mathbf{k}' \quad \mathbf{p} = \mathbf{K} + \mathbf{k}, E_{\mathbf{p}}$$

$$E_{\mathbf{K}+\mathbf{k}} = v_F k_{\perp} + a_n k_{\parallel}^n$$

Determine the crossover

$$\frac{1}{\tau_{\mathbf{p}}} \sim \int_{\mathbf{p}'} \left| \operatorname{Im} \mathcal{V} \left(\mathbf{p} - \mathbf{p}', E_{\mathbf{p}} - E_{\mathbf{p}'} \right) \right| \Theta \left(E_{\mathbf{p}'} \left(E_{\mathbf{p}} - E_{\mathbf{p}'} \right) \right) \sim \int_{\mathbf{p}'} \left| \operatorname{Im} \frac{\Theta \left(E_{\mathbf{p}'} \left(E_{\mathbf{p}} - E_{\mathbf{p}'} \right) \right)}{1 + i \frac{|\omega|}{|\mathbf{v}_{\mathbf{p}} \times \mathbf{v}_{\mathbf{p}'}|}} \right|$$

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$$\sim \begin{cases} \frac{1}{a_{2}}E_{\mathbf{p}}^{2}\ln\frac{1}{E_{\mathbf{p}}}, & n=2, \\ \frac{1}{a_{n}^{1/(n-1)}}E_{\mathbf{p}}^{\frac{n}{n-1}}, & n>2, \end{cases}$$



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• Near **K** with $E_{\mathbf{K}+\delta\mathbf{k}} = v_F \delta k_n + (\delta k_t)^m$, the quasiparticle decay rate \propto

$$\frac{1}{\tau_{\mathbf{p}}} \sim \begin{cases} E_{\mathbf{p}}^2 \ln \left| \frac{1}{E_{\mathbf{p}}} \right|, & m = 2\\ E_{\mathbf{p}}^{\frac{m}{m-1}}, & m > 2 \end{cases}$$



Outlook

> The leading non-analytic correction to the renormalized dispersion E_p at the inflection points.

How the geometric quantum criticality manifests itself in collective modes that describe fluctuations of Fermi surface shape?

>Try to explicitly test our predictions by photoemission spectroscopy.

Thank you for your attention!