Superconducting critical temperature in the extended diffusive SYK model

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Models for strongly interacting fermions in disordered clusters forming an array, with electron hopping between sites, reproduce the linear dependence on temperature of the resistivity, typical of the strange metal phase of High Temperature Superconducting materials (Extended Sachdev-Ye-Kitaev (SYK) models). We identify the low energy collective excitations as neutral, energy excitations, diffusing in the lattice of the thermalized, non Fermi liquid phase. However, the diffusion is heavily hindered by coupling to the pseudo Goldstone modes of the conformal broken symmetry SYK phase, which are local in space. The imaginary time evolution of the extended model in the strong interaction and $1/N$ expansion limit is presented, in the incoherent non chaotic regime. On the other hand, a Fermi electronic liquid at low energy becomes marginal when perturbed by the SYK dots. A critical temperature for superconductivity is derived, which is not BCS-like, in case the collective excitations are assumed to mediate an attractive Cooper-pairing.

I. INTRODUCTION

Understanding the physics of copper-oxide materials, which undergo the superconducting transition at higher temperature, is still an unsettled topic of Condensed Matter Physics. Recent work suggests the breakdown of the Fermi Liquid (FL) theory at intermediate temperatures in these metals, while FL is the conventional starting point for low critical temperature superconductivity. New approaches to study high-temperature superconductivity are recently investigated, in particular lattice fermionic models with a strong local interaction.

Recently a $(0 + 1)−d$ dimensional model, the Sachdev-Ye-Kitaev (SYK) model, describing random all-to-all $J$-interaction between $N$ Majorana fermions, has been extensively studied. In the infrared (IR) limit, when $N$ is large and the temperature $T$ is low, the model has an emergent approximate conformal symmetry and has become quite popular for its large-$N$ "melons" diagrammatics, which allows for a simple representation of the power-law decay in time of the correlation functions and for the analysis of the thermodynamic and chaotic properties, providing a holographic dual for gravity theories.

Generalized SYK models have been proposed with extension to higher space dimensions also having in mind applications to High Critical Temperature ($HT_c$) superconducting materials. Indeed, there seems to be a widespread consensus that inhomogeneity and strong coupling could be distinguished factors for the cuprates and their $2−d$ CuO planes. Moreover, universal features emerge in the high temperature "strange metal" phase, which is recognized as a Non Fermi Liquid (NFL) phase. The most striking of these is the linear increase with temperature of the electrical conductivity.

The conformal symmetry of the SYK model is spontaneously broken down to the $SL(2, \mathbb{R})$ group symmetry and Goldstone modes arise which are only approximately gapless, when ultraviolet (UV) corrections are taken into account. The nature and the role of these collective excitations has not been satisfactorily investigated, to our knowledge, up to now, in phenomenological approaches for the description of the low temperature metal phases of extended SYK models.

FIG. 1: A cartoon of the SYK model extended to include hopping between $2−d$ square lattice sites. The neutral fermions are depicted as small blue circles, grouped on the lattice sites (their number is $N$ on each site). In the right upper part of the picture, a magnification of four sites with just four Majorana particles at each site, to represent the all-to-all fermion interaction $J_{ijkl}$ [red arrows] and the residual interaction $U_c$ between quasiparticles (complex superposition of intra-cluster pairs of majoranas) [dashed arrow] hopping between sites (with hopping matrix element $t_0$ [full dark green arrows]).

We consider a lattice of $(0 + 1)−d$ SYK clusters (or

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dots), each composed of strongly correlated $N$ neutral fermions, via the SYK interaction. A sketch of the lattice, in two space dimensions, is depicted in Fig.1. The first part of this work discusses the collective bosonic excitations, in two space dimensions, is depicted in Fig.1. The fermions, via the SYK interaction. A sketch of the lattice sites of the lattice as if their capacity were infinite. we ignore the possibility of charging of the clusters at the clusters are non-conserving and neutral. In this respect, energies, while the (“intradot”) excitations in the SYK symmetry for the (“interdot”) dynamics in the lattice. Disorders is a distinct feature of the SYK model. Disorder averages make the SYK and its generalizations solvable. We assume random hopping in the lattice and we assume that self-averaging restores space translational invariance. In our description, the bilocal auxiliary fields $G_x(\tau_1, \tau_2)$ and $\Sigma_x(\tau_1, \tau_2)$, in imaginary time $\tau$, acquire a slowly varying phase $\varphi_p(\tau)$. Here the subscript $x$ denotes the space coordinate and the wavevector $p = k\vec{a}$ ($\vec{a}$ is the lattice parameter) is used as a quantum number in the continuum space limit. $G_x(\tau_1, \tau_2) \equiv G_x(\tau_1, \tau_2, \tau_+)$ acts as an order parameter which characterizes the SYK-phase in the scaling to strong interaction $J \to \infty, N \to \infty$ with finite $\beta J / N$ ratio. Here $\tau_{12}$ is the relative time coordinate, which takes care of the "intradot" dynamics, while $\tau_{+}$ is the center of mass coordinate of the "slow" interdot dynamics. The first task (Section III) is to study the correlations of the $pGm$'s, when minimal coupling to the compact dynamical $U(1)$ gauge boson $\varphi_z(\tau_+)$ is established. Fig.3 displays a "dressed" correlator, $\langle \delta G_z(\tau_{12}, \tau_+) \delta G_z^*(\tau_{34}, 0) \rangle \approx \langle \delta G_z(0^+, \tau_+) \delta G_z^*(0, 0) \rangle$ compared with a zero order, "naked" one, continued to real time and in the limit $k\vec{a} \ll 1$. The naked correlation can be derived with a real time approach in Appendix C. Both correlators decay with real time, but the dressed one decays by far faster. This confirms that the extended SYK model at hand describes incoherent dynamics. However, Fig.3 proves that, as long as the gauge boson lacks its own dynamics, correlations cannot be said to be diffusive over the lattice. Actually, diffusivity on a temperature dependent (and scaling-dependent) space distance $\tilde{a}_d(T)$, much larger than the lattice parameter, is expected. In fact, the presence of impurities, low dimensionality, strong interaction and disorder, usually makes the collective excitations diffusive at low frequencies and small momentum$^{20}$. The $pGm$ fermionic excitations of the SYK dots generate fluctuations of the chemical potential in the lattice $\varphi_z(\tau_+) \varphi_p(0)$, driven by quasiparticle hopping between lattice sites, parametrized by the matrix element $t_0$ and produce the bosonic Q-excitations.

Our aim is twofold. On the one hand we want to caracterize the quantum diffusion of the Q-excitation in the lattice$^{11}$. On the other hand we want to study the response $D^\beta(p, \Omega_n)$, of these modes to interdot tunnelling, a $J_Q - J_Q$ response, where $J_Q$ is an energy flux density which is somehow canonical conjugate to $\partial_\tau \varphi_p(\tau_+)$, the latter takes the role of a space dependent chemical potential across the lattice$^{32}$. The probability of quantum diffusion, involving retarded and advanced Green’s function in real time, $G^R$ and $G^A$ respectively, is written in the form:

$$ P(r, r'; \Omega) \propto \int d\omega \, G^R(r, r'; \omega)G^A(r', r; \omega - \Omega), $$

where overline denotes disorder average. On the other hand the retarded density response function $D^{R}_{Q, Q}$ involves the retarded $G^R$ and the Keldysh $G^K$ Green’s function. Our approach will be in imaginary time, but correct time ordering is crucial to guarantee a correct analytial continuation to real times. A relevant quantity typical of the diffusion processes is its Fourier transform in time, denoted as the heat kernel$^{32}$, which is defined as the probability $z(t)$ to return to the origin, integrated over the point of departure. In Section V we derive a form of it, $z(t) \sim e^{\frac{D_Q}{2} t^2}$, after the Q-excitations have been integrated out (Eq.(36)) and we determine how the diffusion parameter $D_Q$ depends on the scaling to strong interaction.

In dealing with the "bad" metal at finite temperature $T$, we concentrate on two temperature scales in-
olved in the extended SYK model, $T_0$ and $T_{coh}$. At $T \lesssim T_{coh} \sim t_0^2/N^2$, transport in the lattice is assumed to acquire coherence. This crossover is out of reach in the present work. We expect that the Q-excitations merge into the particle hole (p-h) continuum of the low energy MFL. A derivation of the Landau damped acoustic plasmon embedded in the p-h continuum, is reported in the appendix E. At $T \gtrsim T_0$, thermalization in the system is very effective and diffusion is incoherent. pGMs are the intradot excitations which drive the incoherence. This is a feature of the SYK model and is attained in the present extended version of the model. The relaxation time is $\sim h\beta^{34}$. At later times the system evolves toward the scrambled phase and the chaotic dynamics, as the analysis of the out-of-time ordered correlator (OTOC) shows. The fate of the chaotic single dot regime in the extended model deserves specific concern\textsuperscript{21,28,35,36} beyond the present paper.

Usual hydrodynamical approaches to the response function $D^R$ do not involve the role of the pGMs at energies $\sim \kappa g T_0$. This is highly questionable, because the diffusion constant, $D_Q$, is strongly renormalized by the inverse of four point function of the SYK dots, $F^{-1}$. Indeed the first UV correction plays the role of keeping the pGM propagator $F$ finite. A brief presentation of this approximation to $D^R$, which does not go beyond the conformal limit\textsuperscript{13} and uses the real time Keldysh contour, is reported in the appendix C. By contrast, our approach is quite simple and even naive, but it aims to stress the parameter renormalization in the scaling process. In fact, the separation in energy of $T_0$ and $T_{coh}$ allows us to perform a kind of adiabatic factorization, between the "fast" intradot pGMs and the "slow" interdot Q-fluctuations. We discuss the UV local space-time correction and show how they influence the time correlation of the Q-excitations.

Physically, we concentrate in distinguishing the two regimes $T \lesssim T_0$. The $T \gtrsim T_0$ regime, being characterized by strong thermalization, is governed by the order parameter of the SYK model which, in the UV corrected form, is described by a complex field $\phi(x,\tau_+)$ in Section IV. The Q-excitations, arising from the minimal coupling with the gauge mode, are interpreted as energy excitations induced by the fluctuations of the chemical potential. Energy density $\tilde{N}$ and energy flux density $\tilde{N}^\tau / J_Q$ are the physical dynamical variables\textsuperscript{12}. The corresponding parameters which rule the response are thermal capacitance $C_V$ and the thermal conductivity $\kappa$.

The structure of the paper is as follows.

In Section II the extended SYK model is presented. In the conformal symmetry limit of our approach, the SYK clusters acquire an hopping dependent selfenergy of the kind $\sim t_0^2 G_c G_{c\tau}$, where $G_c(\tau_1,\tau_2)$ is the fermionic propagator of the SYK model\textsuperscript{13}. A term of this kind is suggested by a simple derivation of the hopping between two neighbouring SYK sites. The local correlations arising from the kinetic term are obtained by gaussian integration of the $\delta g_{mn}$ fluctuations in presence of a source term, the chemical potential $\partial_\tau \varphi_\pm (\tau_+)$. They are derived in Section III. In Section IV we clarify that the proper dynamics of the chemical potential fluctuations should be added to account for the UV corrections which, by giving mass to the pGMs, make the partition functional convergent. This implies a renormalization of the correlations provided by the pGM propagator $F$, in which the first UV correction is included. To this end we introduce a complex local order parameter $\phi(x,\tau_+)$, which is promoted to a bosonic coherent field in Section IV, by means of a more conventional model for the Q-excitations. The inclusion of the dynamics via the local action $\tilde{S}_2[\partial_\tau \varphi(x,\tau_+)]$ of Eq.(28,29) implies that the short range, exponentially decaying dependence on real time $t_+$, of the correlators turns into a diffusive dynamics for $T \gtrsim T_0$, the energy window in which our approximations are justified (Section V.A). Section V.B discusses qualitatively how the transport parameters evolve with scaling in the incoherent and coherent energy ranges. They can be used to qualify the diffusion parameter $D_Q$ by means of the Einstein relation. In Section VI we show how a coherent low energy FL, when perturbed by a higher energy SYK-type environment, becomes marginal. A conventional Eliashberg\textsuperscript{37,38} approach to the gap equation is presented in Section VI, where the Q-excitations constitute a bosonic virtual pairing mechanism but with diffusive dynamics. The self-consistent equation for the non-BCS critical temperature $T_c$ is derived. Additional remarks and a summary are reported in the Conclusions (Section VII). The Appendices give details of the derivations.

II. THE EXTENDED SYK MODEL

Let the Hamiltonian for the extended model be $H_0 + H_K$. $H_0$ is the sum of the neutral fermion Hamiltonians of uncoupled 0+1-d SYK dots, $H_d$, in a two-dimensional lattice with intradot random interaction, labeled by the lattice site $a$, and $H_K$ adds the kinetic energy of electrons with interdot random hopping between neighbouring dots. $H_K$ (given by Eq.(6)) is derived in this Section. The Hamiltonian $H_0$ for the uncoupled 0+1-d SYK dots is:

$$H_0 = \sum_a H_a = \frac{1}{4!} \sum_{a,klnm} J_{aklnm} \chi_{a,k} \chi_{a,l} \chi_{a,m} \chi_{a,n}.$$  \hspace{1cm} (2)

where $\chi_{a,t}$ are Majorana fermion operators on site $a \ (klmn \in 1, .., N)$.

Electronic quasiparticles hop from site $a$ to a neighbouring site $b$. $c^\dagger_a c_b$ and $c^\dagger_b c_a$ are the complex fermionic spinless operators for the electrons, which can be represented in terms of two flavours of the neutral fermions on the same site:

$$c_b = \frac{1}{\sqrt{2}} (\chi_{b1} + i \chi_{b2}) \ , \ c_b^\dagger = \frac{1}{\sqrt{2}} (\chi_{b1} - i \chi_{b2}).$$  \hspace{1cm} (3)
The kinetic term describing the hopping can be written as \( h_K = t_0 \delta_i^c \delta_i^a + \text{h.c.} \), where \( t_0 \) is a constant hopping energy.

The time dependence of the operator \( \delta_i^c \) in the interaction picture is:

\[
-\frac{\partial}{\partial \tau} \delta_i^c(\tau) = e^{(H_b + H_a)} \left[ \delta_i^c, H_0 \right] e^{-(H_b + H_a)}. \tag{4}
\]

The commutator with the Hamiltonian can be performed by applying the commutation relations for neutral fermions: \( \chi_a,k \chi_{a,l} + \chi_{a,l} \chi_a,k = \delta_{l,k} \) and \( \chi_a,k \chi_{b,l} + \chi_{b,l} \chi_a,k = 0 \) for \( a \neq b \), exploiting the antisymmetry of \( J_{bklnm} \) in the permutation of the \( klmn \) indices. From Eq.(4) we get:

\[
\frac{\partial}{\partial \tau} \delta_i^c(\tau) = i \frac{1}{3!} \sum_{lm} J_{b12lm} \chi_{b,l}(\tau) \chi_{b,m}(\tau) \delta_i^c(\tau). \tag{5}
\]

\( \delta_i^c \) commutes with \( H_b \) so that it can be added afterwards. The hermitian conjugate term \( \chi^c_b \delta_i^c \) gives the same result with \( b \rightarrow a, i \rightarrow -i \).

This allows to identify the hopping Hamiltonian term in the interaction representation, from the evolution operator in a single hopping process, \( \delta U(\tau,0) \), to lowest order:

\[
\mathcal{H}_K(\tau) = i \frac{1}{3!} \sum_{lm} J_{j12lm} \chi_{j,l}(\tau) \chi_{j,m}(\tau) + \text{h.c.}. \tag{6}
\]

Here \( J_{j12lm} \) is random interdot hopping for hopping onto site \( j \). Eq.(6) shows that, starting from the neutral fermions of the SYK model, a symmetric description of conserving and non-conserving charge processes is provided. This feature sets charge (and spin) dynamics free with respect to energy dynamics, which is the premise for NFL behavior.

The disorder average of the standard SYK model includes here the gaussian average of \( J_{j12lm} \). The next step is the integration over the Majorana fields \( \chi_{j,l}(\tau) \), with the help of Hubbard-Stratonovich fields which become complex due to an additional \( U(1) \) minimal coupling. The final result is the action in terms of the complex bilocal auxiliary fields \( G_x(\tau_1, \tau_2) \) and \( \Sigma_x(\tau_1, \tau_2) \), with a phase \( \varphi_x \) introduced in the next Section.

\[
\frac{I_{ex}}{N} = \sum_x \left[ -\ln \text{Det} [\partial_{\tau} - \Sigma_x] + \int \! d\tau_1 d\tau_2 \left\{ -\frac{J^2}{4} \left[ G_x(\tau, \tau') \right]^2 + \Sigma_x(\tau, \tau') G^*_x(\tau, \tau') \right\} \right]. \tag{7}
\]

The last term of the action is the interdot kinetic term. The expansion up to quadratic terms of this action in \( \delta G_x, \delta \tau_\varphi \) is discussed in Section III and in the appendix A. The single dot 0–1-d SYK action can be recovered by dropping \( \delta \Sigma_x \), the last term and the sum over sites. The auxiliary fields are now real and the Det has to be substituted with a Pfaffian. In this case the IR limit corresponds to the dropping of \( \partial_{\tau} \) in the Pfaffian. On the contrary, \( \partial_{\tau} \) plays an important role in the extended model.

### III. KINETIC CORRELATIONS OF THE EXTENDED SYK MODEL

The single particle Green’s function of the SYK model, in the conformal symmetry limit, is local in space (i.e. wavevector independent) and, assuming particle-hole (p-h) symmetry and low temperature, it is given by:

\[
G_x(i\omega_n) = i \frac{\text{sign}(\omega_n)}{\sqrt{\sqrt{\omega_n}}}, \tag{8}
\]

where \( \omega_n \) are fermionic frequencies.

Our aim is to include correlations between sites of the lattice, here denoted by the subscript \( x \). The Green function and the self-energy become complex fields, \( G_x(\vartheta_1, \vartheta_2), \Sigma_x(\vartheta_1, \vartheta_2) \). They include space dependent fluctuations of the modulus and of the phase, close to the saddle point \( G_x(\vartheta_{12}), \Sigma_x(\vartheta_{12}) \):

\[
G_x(\vartheta_1, \vartheta_2) = [G_x(\vartheta_{12}) + \delta G(x, \vartheta_{12}, \vartheta_+) \ e^{i \varphi_x(\vartheta_+)}, \Sigma_x(\vartheta_1, \vartheta_2) = [\Sigma_x(\vartheta_{12}) + \delta \Sigma(x, \vartheta_{12}, \vartheta_+)] \ e^{i \varphi_x(\vartheta_+)}, \tag{9}
\]

where \( \vartheta_{12} = \vartheta_1 - \vartheta_2 \) and \( \vartheta_+ = (\vartheta_1 + \vartheta_2)/2 \). We have moved to the center of mass time coordinate \( \vartheta_+ \) and the relative time coordinate \( \vartheta_{12} \) of the incoming particles and of the outgoing ones. Here \( \vartheta = 2\pi / 3 \) is a dimensionless time and the Green’s functions and selfenergy are also dimensionless, everywhere, except when explicitly stated. Nevertheless we will most of the times denote the dimensionless time as \( \tau \) unless differently specified. To spell out the structure of the kinetic term, we calculate the correlator of the \( \delta G \) fluctuations between neighbouring sites and Fourier transform it with respect to space. Ignoring the relevant role of the \( p \bar{G} \bar{m} \)’s, we neglect, in the IR limit, the local correction \( \delta \bar{G}(x, \tau_1, \tau_2, \tau_+) \ e^{i \varphi_x(\tau_2)} \) appearing in Eq.(9) and we consider just nearest neigh-
bour $x, x'$ terms in a lattice of spacing $\tilde{a}$. We get:

$$
\delta G_{c,x}(\tau_{12}, \tau_+) \delta G_{c,x}^*(\tau_{34}, \tau'_+) = [G_x(\tau_1, \tau_2) G_x^*(\tau_3, \tau_4) - G_c(\tau_1 - \tau_2) G_c(\tau_3 - \tau_4)]
$$

$$
\approx \frac{1}{2} G_c(\tau_{12}) \left(e^{-i\tilde{a} \cdot \nabla_x [\varphi_x(\tau_+) - \varphi_x(\tau'_+) - 1]} - 1\right) G_c(\tau_{34}),
$$

where we have qualified the lowest order, originating from the conformal Green's functions, with the label $c$. Only the quadratic terms of the exponential are included in the expansion, to account for the additional complex conjugate contribution, giving

$$
\approx -\frac{1}{2} G_c(\tau_{12}) \left(\tilde{a} \cdot \nabla_x [\varphi_x(\tau_+) - \varphi_x(\tau'_+)]\right)^2 G_c(\tau_{34}).
$$

We now approximate (A9) $[\varphi_x(\tau_+) - \varphi_x(\tau'_+)] \approx (\tau_+ - \tau'_+) \partial_+ \varphi_x(\tau_+)$ and define

$$
R_c^{-1} \Lambda_c R_c^{-1} = \frac{1}{2}(\tilde{a} \cdot \nabla_x) G_c(\tau_{12})(\tau_+ - \tau'_+) \frac{1}{2} G_c(\tau_{34})(\tilde{a} \cdot \nabla_{\tau'} x').
$$

Owing to the self-averaging established for the SYK model at large $N$, translational invariance allows space Fourier transform:

$$
\frac{1}{N} FT_p \left[\delta G_{c,x}(\tau_{12}, \tau_+) \delta G_{c,x}^*(\tau_{34}, \tau'_+)\right] = \frac{\delta^2}{\delta \partial_+ \varphi_p(\tau_+)} \frac{1}{N} (\partial_+ \varphi_p) \left[ R_c^{-1} \Lambda_c R_c^{-1} | \partial_+ \varphi_p \right] (12)
$$

($FT_p$ denotes Fourier Transform with respect to space coordinate of lattice spacing $\tilde{a}$, with $p = k\tilde{a}$). We now express Eq.(12) in the frequency space. The matrix elements of the kernel are labeled by $m, m', \ell$ indices. $m, m'$ indices refer to the intradot fluctuations $\delta g$ which are fermionic in the origin, while $\ell$ labels bosonic frequencies $\Omega_\ell$, corresponding to the spectrum of the Q-fluctuations. We get

$$
\frac{1}{N} FT \left[\delta G_{c,x}(\tau_{12}, \tau_+ \delta G_{c,x}^*(\tau_{34}, 0)\right] |_{k \neq 0} = \frac{\beta \tilde{a}^2}{N} k^2 \tilde{a}^2 \times \sum_{n} \sum_{m, m'} e^{i\Omega_\ell \tau_+} \left( R_c^{-1} \Lambda_c R_c^{-1} \right)_{m, m'} e^{i\omega_m \tau_{12}} e^{i\omega_m \tau_{34}} (13)
$$

Restricting ourselves to the IR limit, we plot in Fig.2 the time Fourier transform, keeping just the dependence on the relative coordinate $\tau_{12} - \tau_{34}$ mod. $2\pi$ ($\omega_m = -\omega_n$).

$$
\frac{1}{N} \sum_{\ell, m} \left( R_c^{-1} \Lambda_c R_c^{-1} \right)_{m, m'} \frac{1}{\pi(2m + 1)} e^{i\omega_m \tau},
$$

where Eq.(8) has been used. It is denoted as $\langle \delta G_{c,x}(\tau, 0)\delta G_{c,x}(0, 0)\rangle_k$ in Fig.2. This quantity, together with the dressed correlator of Eq.(35) (blue curves), is plotted for $\tau_+ \rightarrow 0^+$. The prefactors $k^2 \tilde{a}^2 \beta \tilde{a}^2/(2\pi \tilde{a})$ have been dropped in the plots. The real part of the continuation of Eq.(14) to real time $t_+$, when $\tau \rightarrow 0^+$, $\Re \langle \delta G_{c,x}(0, t_+) \delta G_{c,x}(0, 0)\rangle_k$ is plotted in Fig.3. Note the difference in the scale of decay between this correlation derived from the naked kinetic term and the one of Eq.(35), including UV corrections, which we are going to discuss in detail in the next Section.

Integrating out the $\delta \Sigma$ fluctuations (see (A10)), the functional integral in terms of the fluctuations $\delta g(\tau_{12}, \tau_+)$ is

$$
Z \left[\partial_+ \varphi_p(\tau_+)\right] = \int \left(\Pi \delta g_{\tau_{12}, \tau_+}^{\ast}(\tau_+) \Pi \delta g_{\tau_{12}, \tau_+}\right) e^{\frac{\beta}{2} \left[\delta g [R_c^{-1} \Lambda_c R_c^{-1} \delta g]\right] - \frac{\beta}{2} \Re \left\{ -i R_c^{-1} \partial_r \varphi_p [\partial_r \varphi_p] \right\}} \times e^{\frac{\beta}{2} \left[ -i R_c^{-1} \partial_r \varphi_p [\partial_r \varphi_p] \right]} (15)
$$

The forks $(\ldots)$ in Eq.(15) include integration over $\tau_{12}$ and $\tau_+$. Here $\langle \delta g(\tau_{12}, \tau_2) \rangle = R_c(\tau_{12}, \tau_2) G_c(\tau_{12}, \tau_2)$ and $R_c(\tau_{12}, \tau_2) = \beta \tilde{a} (q - 1) [G_c(\tau_{12}, \tau_2)]$ (with $q = 4$ in the usual notation).

Integrating out $\delta g_{\tau_{12}, \tau_+}, \delta g_{\tau_{12}, \tau_+}^{\ast}$, the generating func-

$$
Z \left[\partial_+ \varphi_p(\tau_+)\right] = e^{\frac{\beta}{2} \left[ -i \partial_r \varphi_p (\langle \partial_+ \varphi_p \rangle + \frac{\partial^2}{\partial^2_r \varphi_p R_c^{-1} \Lambda_c R_c^{-1}} \right)]} (17)
$$
\[ F(\tau_1, \tau_2, \tau_3, \tau_4) = R_c^{-1} K_c |1 - K_c|^{-1} R_c^{-1} \] (18)

is the four point function of the 0+1−d SYK model. Integration over intermediate times is intended. Here \( F \) is \( O(1) \), with the meaning of \( O \left( \frac{\delta g}{N} \right)^0 \). As \( R_c^{-1} \Lambda_e R_c^{-1} \propto G_c G_c \) is \( O(N/\beta J) \), it appears from Eq.(17) that we can define a physical parameter \( \beta \delta g / N \beta J \) of \( O(1) \), to guarantee that the hopping across the lattice is not irrelevant in the scaling. It turns out however that both \( F \) and \( \beta \delta g / N \beta J \) become of \( O \left( \frac{\delta g}{N} \right)^0 \) when the UV correction is included, which is crucially important to give sense to the functional integration of Eq.(15), as we explain here below.

Actually, the functional integral of Eq.(15) includes a divergent contribution due to the Goldstone modes \( \delta g_c \) corresponding to eigenvalues of \( K_c \rightarrow 1 \), which has to be regularized resorting to the first UV correction \( (\delta g) |K_c^{-1} - 1| \delta g \sim \beta J \). The Faddeev Popov regularization provides an integration performed in the orthogonal space with respect to the \( pGm \), while the smallest eigenvalue of the kernel \( 1 - K_c \) is approximated with its UV correction, given by \( 1 - k_c (h = 2, n) \approx \frac{\delta g}{N} |n| + \ldots (\alpha K \approx 3 \text{ is a constant}) \). It follows that the large, but finite contribution to \( F \) in Eq.(18) with this UV correction (i.e. for \( \tau \sim 1/J \)) is not \( O(1) \) as stated here above, but \( O \left( \frac{\delta g}{N} \right)^0 \) and the same has to occur for \( \frac{\beta g}{N \beta J} \). We will discuss this point in the next Section. The temperature threshold for coherence defined here, \( T_{coh} = \frac{\beta}{2\pi} \), is recurrent in the next.

If we ignore this matter for the time being, the generating functional of Eq.(17) provides the correlator
\[ \frac{1}{2} \int \mathcal{D} \varphi \int \mathcal{D} \varphi (\tau) \int \mathcal{D} \varphi (\tau') \int \mathcal{D} \varphi (\tau'') \int \mathcal{D} \varphi (\tau''') \int \mathcal{D} \varphi (\tau''') (\delta g_c (\tau_1, \tau_2) \delta g_c (\tau_1, \tau_4)) |_{k=\delta g} \] in imaginary time, inclusive of the hopping in the lattice:
\[ \frac{\delta^2}{\delta \partial \varphi (\tau_+)} \frac{\delta^2}{\delta \partial \varphi (\tau_+)} (\frac{1}{N} \sum_{\nu} \ln Z [\partial \varphi (\tau_+)] \times \left( 1 + \frac{t_0^2}{2N} \mu^2 R_c^{-1} \Lambda_e R_c^{-1} F^{-1} \right) F. \] (19)

This result adds to the naked correlator \( F \) the contribution coming from Eq.(13), so that the two dynamics are just added together in this approximation. However, one can envisage the present one as the lowest order of a ladder resummation which will appear more clearly in the next Section. The operator \( G_c G_c F^{-1} \) appearing in the Kernel of Eq.(19) is the inverse matrix of
\[ F \{G_c, G_c\}^{-1} = \frac{6 \alpha_0 \beta J}{\pi^2 a K} \times \sum_{n \geq 2, even} e^{i n (y_1 - y_4)} \frac{\pi^2 (n^2 - 1)}{f_n(t_1) f_n(t_4)}, \] (20)

where \( y_i \equiv (\tau_i + \tau_4)/2 \) in unit of \( h \beta / 2 \pi \). The basis functions \( f_n(t_1) \) are defined in Appendix B, together with the spectral representation of the kernel \( K_c^{-1} [1 - K_c] \), as well as with their Fourier transform.

IV. DRESSED CORRELATOR OF \( pGm \) MODES

The derivation of the previous Section has assumed that \( \varphi (\tau_+) \) is given as an external source. However, continuation to real time requires that \( \varphi (\tau_+) \) acquires a dynamics. Meanwhile, the symmetry breaking induced by the UV perturbation source \( \partial \varphi \) couples to \( G_{IR} \) and \( G_{IR} \) is derived from a time reparametrization under the diffeomorphism \( e^{i \varphi} \rightarrow e^{i \varphi (\varphi)} \) of the conformal Green’s function \( \Delta = 1/4 \):
\[ G_{IR}(\varphi_1, \varphi_2) = G_c (\varphi_1 (\varphi_1), \varphi_2 (\varphi_2)) \varphi_1 (\varphi_1) \varphi_2 (\varphi_2) \Delta, \] (21)

where \( \varphi = 2 \pi \tau / \beta \) and \( \varphi' \equiv \partial \varphi \).

The leading correction to the conformal action arising from this reparametrization (apart for a shift of the ground state energy) is the Schwarzian:
\[ I_{local} = -2 \pi \alpha \varepsilon e^{i \varphi}. \] (22)

Here \( \alpha \varepsilon \) is a constant and \( \varepsilon = 1/\beta J \). Hence, the full action in place of the one appearing in Eq.(17) reads:
Now the field \( \partial_\tau \phi_x \) has its own dynamics and could be integrated out, possibly after adding a source term to get a generating functional of \( \langle \partial_\tau \phi \partial_\tau \phi \rangle \) correlators. However, the action of Eq. (23) is essentially a "phase only" model for hopping of \( \partial_\tau \phi_x \) across the lattice. This is so, because we have neglected \( \delta G(x, \tau_2, \tau_+ \) appearing in Eq. (9). So as it stands, \( I_{\text{local}}[\partial_\tau \phi] \sim O(N/\beta J) \), while the other term is \( O(1) \). Hence, the local action \( I_{\text{local}}[\partial_\tau \phi] \) is irrelevant in the \( \beta J \rightarrow \infty \) limit and the phase lacks its own dynamics. Writing down a partition function for the order parameter given by Eq. (9), with inclusion of its modulus, gives the chance of extracting correlation functions with include the UV correction and can be extended to real time. Let us denote the complex order parameter in two space dimension \( \phi(x, \tau_1, \tau_+) = \sqrt{\rho_0 + i \rho} e^{i \theta} \), for each degree of freedom. The functional integral, with \( \tau \) of the dimension time in the following, reads:

\[
\frac{G_c(\tau_2)}{a} \int D\phi(x, \tau_1, \tau_+) D\phi(x, \tau_1, \tau_+) e^{-\tilde{S}[\phi^*, \phi]}.
\]

The action leading to the one of Eq. (23) can involve also space derivatives:

\[
\tilde{S} = \int \beta \tau_1 \int d^2 x \int d^2 \tau \left[ -\pi^2 \partial_\tau \partial_\tau \phi \partial_\tau \phi + (\phi^* \partial_\tau \phi - \phi \partial_\tau \phi^*) - \frac{1}{2\pi \alpha s \epsilon} |\phi|^4 \right] (25)
\]

which can be identified with \( \frac{G_c}{a^2} \partial_\tau \phi \) of Eq. (22) provided we also introduce space non locality there, by trading \( \frac{G_c(\tau_2)}{a} \partial_\phi \phi \), which appears in Eq. (28), for \( \partial_\tau \phi \). Identification requires that

\[
\rho_0 = |\phi_0|^2 = \frac{\pi \alpha s \epsilon}{a^2} |G_c(\tau_2)|^2, \quad \partial_\tau \phi \equiv \partial_\tau \phi
\]

(an extra factor \( N \) pops up from the number of flavours in the first equality). Introducing \( \partial_\phi = |G_c(\tau_2)| \) and substituting \( -i \partial_\tau \phi \rightarrow e^{-i \theta} \partial_\phi \partial_\tau \phi \), the functional integral becomes

\[
\int D\phi(x, \tau_1, \tau_+) D\phi(x, \tau_1, \tau_+) e^{-\tilde{S}[\phi^*, \phi]} e^{-\frac{1}{2\pi \alpha s \epsilon} \sum \phi^* \tilde{R}_c^{-1} \phi \tilde{R}_c^{-1}} \sum \langle \phi \tilde{R}_c^{-1} \tilde{G}_c \left( F + \frac{i}{2} \pi^2 \partial_\tau \phi \right) \phi \tilde{R}_c^{-1} \rangle \rangle (30)
\]
where $\tilde{S}$ is given by Eq.(25). It is useful to redefine the $\phi = \phi R_c^{-1}$ in the functional integral. In the change of
the integration field, the action $\tilde{S}[\phi^*,\phi]\rightarrow \tilde{S}[\phi^*,\phi]$ acquires a factor $(G_c(\tau_{12}))^2$, except for the $|\phi|^4$ term which acquires the fourth power. Note, however, that in the
UV domain is $\tau_{12}\sim J^{-1}$, so that, with $8b^{-2} = \sqrt{4\pi J}$,
\begin{equation}
(G_c(\tau_{12}))^2 = \frac{b^2}{|\tau_{12}|} \sim \mathcal{O}(1).
\end{equation}
Hence, the last term of the full action from Eq.(30) is
$\mathcal{O}(\beta J)$ in the large $\beta J$ limit, while the first contribution
to the full action, given by $\tilde{S}$, is $\mathcal{O}(1)$. Actually the $|\phi|^4$ term in $\tilde{S}$ is also $\mathcal{O}(\beta J)$, but we stick to zero order in the
anharmonic functional integration. The evolution of Eq.(30) is characterized by an interplay between the dynamics of the intradot fluctuations and the dynamics of the interdot $\partial_r \varphi_p$ fluctuations, which is mostly represented by the action $\tilde{S}$. If $\tilde{S}$ is dropped altogether, because it becomes irrelevant in the scaling, the gaussian
integration of Eq.(30) can be easily performed, giving rise to a density matrix $\rho_{\phi \phi}$ of the intradot fluctuations at each given time $\tau_r$. When Fourier transformed with respect to the intradot times $\tau_{12}, \tau_{34}$, stripping off the unperturbed evolution $\mathcal{F}$, the result of the functional integration of Eq.(30), in the absence of $\tilde{S}$ is (again, $y_{ij}$ are center-of-mass times in unit of $h/\beta/2\pi$ and here $\tau \equiv \tau_r = y_{12} - y_{34}$ in unit of $h\beta/2\pi$):
\begin{equation}
\rho_{\delta \delta}(y_{12} - y_{34}; p)_{m,m'} \sim \frac{1}{N} \left\{ 1 + \frac{t_0^2 p^2}{N} \partial_r \left( [R_c^{-1} \Lambda_c R_c^{-1}] \dot{\hat{f}}^{-1} \right) \partial_r \right\}^{-1}_{\tau_r,m,m'}.
\end{equation}
We define
\begin{equation}
\bar{B}_{m,m'}(\tau_r) \equiv \partial_r \left( [R_c^{-1} \Lambda_c R_c^{-1}] \dot{\hat{f}}^{-1} \right) \partial_r \right\}_{m,m'} \sim \tau_r^{-1}.
\end{equation}
which will be used in the following.

The correlation function $\langle FT[\delta G_c(\tau_{12},\tau_r)\delta G_c^*(\tau_{34},0)] \rangle_{k \neq 0}$, corresponding to
Eq.(13), but including the ladder resummation, is obtained by tracing on the density matrix of Eq.(32), after the $p = 0$ term has been subtracted. To lowest order, we get:
\begin{equation}
\langle FT[\delta G_c(0^+,\tau_r)\delta G_c^*(0,0)] \rangle_{k \neq 0} \approx \sum_{m,m'} G_c(\omega_m) \rho_{\delta \delta}(\tau_r;k)_{m,m'} G_c(\omega_m^*).
\end{equation}
where $G_c(\omega_m)$ is given by Eq.(8). When $k\tilde{a} << 1$, the contribution of the ladder can be dropped and the un-

normalized correlator $\langle \delta G(\tau_{12},\tau_r)\delta G(\tau_{34},0) \rangle_{k \neq 0}$ reads:
\begin{equation}
= \frac{1}{N} \sum_m \frac{t_0^2}{N} k^2 \tilde{a}^2 \left[ \overline{G_c}(\tau_r) \right]_{y_m,m} \mathcal{E}(\omega_m(\tau_{12} - \tau_{34})).
\end{equation}
Only the dependence on the relative coordinate $\tau_{12} - \tau_{34}$ mod. $2\pi$ has been retained.

In Fig.2, the correlator $\langle \delta G(\tau,0^+)\delta G(0,0) \rangle_k$ from
Eq.(35) is plotted and compared with the naked $\langle \delta G_c(\tau_r)\delta G_c^*(\tau_r) \rangle$ correlator given by Eq.(14). The main panel of Fig.2 displays the modulus while the phase appears in the inset of Fig.2. The prefactor $k^2\tilde{a}^2 \beta t_0^2/(2\pi J)$ has been dropped.

The correlator $\langle \delta G(\tau,0^+)\delta G(0,0) \rangle_k$ has been calculated as reported in appendix B, using the Fourier Trans-
form of Eq.20, with the inclusion of $F^{-1}$ in the evolution. We had to truncate the sum over the (even) indices $n$
up to $n = 12$, and consequently the sum over internal (odd) indices just includes up to $m, m' = 5$. Its modulus and phase, compared to those of the corresponding naked $\langle \delta G_c(\tau_r)\delta G_{c'}^* \rangle_k$ correlator, are plotted in Fig.2. The modulus of the naked correlator is exponentially decaying at the intradot time $\tau \sim 0$, mod $[2\pi]$, while the dressed one is powerlaw, highlighting the criticality of the phase, when
the UV correction is included. The Fourier transform of the sawtooth phase oscillations of $\langle \delta G_{c/c} \rangle_k$ (blue curves) appearing in Fig.2 is not simply $\propto 1/\omega_m$, revealing the "fast" intradot time scale induced by the UV correction, with respect to the phase of the naked correlator. They could have acquired further structure, if larger $n, m$ values had been retained.

The real part of the analytic continuation to real times of the center of mass coordinate $t_+$ in
$\langle \delta G(0^+,\tau_r)\delta G(0,0) \rangle_k$ is plotted in the main panel of
Fig.3 and compared to the corresponding naked correla-
tion of Eq.(14). The same correlators, but keeping the dependence on the relative imaginary time coordinate $\tau_{12} - \tau_{34}$, mod $[2\pi]$ as in Eq.(35), are plotted for various values of $\tau = \tau_{12} - \tau_{34}$ in the inset panel. The dependence on $\tau$ is oscillating and we have chosen values for $\tau$
within a single oscillation. The prefactor $k^2\tilde{a}^2 \beta t_0^2/(2\pi J)$ has been dropped again. The $\langle \delta G(0^+,\tau_r)\delta G(0,0) \rangle_k$'s appearing in Fig.3 are scaled by $\times 10$ with respect to the corre-
larators arising from the naked kinetic term of Eq.(14). The $t_+$ dependence in the presence of UV corrections appears very localized and highly variable with the intradot time, as compared with the naked one. The UV corrections squeeze the interdot correlations in time, in-
creasing their "local" nature. This drastic drop in time of the correlations cannot guarantee quantum diffusion on an extended space scale, much larger than the lattice spacing, and we have to resort to a better approximation which retains the dynamics entailed by the action $\bar{S}$, which was lost in this result.

Besides, the strong dependence of the dressed correla-
A. Partition function of the Q-excitations

In Section IV we have shown that, to improve the \( t_i \)-dependence of the correlator \( \Re \langle \delta G(\tau, t_i) \delta G(0, 0) \rangle_k \) of Eq.(34), the UV local time corrections should be included more carefully. In fact, the result of the previous Section is unsatisfactory, because, in the flowing to the fixed point of the partition function of Eq.(30), we had to drop the order parameter dynamics entailed by the action \( S \) of Eq.(25). A semiclassical approach to the diffusion process can be still envisaged, however, in the results of the previous Section. When the trace over the intradot frequencies is performed, the density matrix of Eq.(32), appropriately continued to real time, \( \rho_{\delta \delta} (y_1 - y_2; k \rightarrow \mathcal{P}(r, r', t_i)) \), takes the form of a heat kernel \( z(t) \), typical of a diffusion process\(^{33} \), defined as the probability to return to the origin, integrated over the point of departure. From Eq.(32), in the \( k\tilde{a} << 1 \) limit, we have:

\[
z(t_i) = \frac{1}{N} \sum_{m} e^{-\phi(v, t_i)} d^{2}r
\]

where we have restored the free intradot evolution.

Now that we know what the drawback is, we reconsider the UV correction to the action given by Eq.(22). Its variation with respect to \( \varphi' \) gives a simple equation of motion \( \partial_{\tau}^{2} \varphi' = -\varphi' \). When derived from the action of Eq.(28), this motion equation is rewritten in the form of lattice space oscillations. In the following we quantize these space extended excitations by means of a phenomenological 2-d Lagrangian with canonical conjugate variables, introduced in appendix B:

\[
\dot{\theta} = \left( \frac{\kappa C}{\hbar T} \right)^{1/2} \frac{JQ T_{Q}}{kB}, \quad \nabla \theta = \left( \frac{\kappa}{\hbar C T} \right)^{1/2} \frac{\kappa}{T} \nabla T.
\]

Here \( JQ \) is the thermal energy current density. The corresponding Lagrangian is

\[
\mathcal{L} = \frac{1}{2} \int d^{2}x \left[ \frac{k_{B}}{T} \left( \frac{JQ T_{Q}}{kB} \right)^{2} + \frac{h}{\kappa C} \left( \frac{\kappa}{T} \nabla T \right)^{2} \right]
\]

\[
\equiv \frac{1}{2} \int d^{2}x \left[ \frac{h k_{B}}{\kappa C} \hat{\theta}^{2} + T \left( \nabla \theta \right)^{2} \right].
\]

The terms in the square brackets have dimension \( \mathcal{E}/\ell^{2} \) (\( \mathcal{E} \cong \text{energy, } \ell \cong \text{length} \)).

This Lagrangian is of course conserving, but we have introduced the relaxation time \( T_{Q} \), so that we can reproduce a diffusive motion equation of the form \( \dot{J}_{Q} = -\kappa \nabla T \), if we approximate the time derivative of the energy current fluctuations \( J_{Q} \propto J_{Q}/T_{Q} \). Here \( \kappa = \frac{C_{v} \ell}{\alpha_{T}} \) is the thermal conductivity in \( 2 - d \), where \( \ell \) and \( v \) are typical mean free path and velocity, respectively, while
\[ \tilde{\alpha}_t^2 \sim \tilde{D}_Q T_Q \] is the area over which the thermal capacitance \( C \) is defined and will be introduced here below.

We quantize the corresponding Hamiltonian, in terms of the creation and destruction bosonic operators \( a_k^\dagger, a_k \):

\[
\begin{align*}
\pi_k &= -i T^{1/2} \frac{1}{(2\Omega_k)^{1/2}} |k| \left( a_{-k} - a_k^\dagger \right), \\
\theta_k &= T^{-1/2} \frac{(2\Omega_k)^{1/2}}{|k|} \left( a_k + a_k^\dagger \right), \\
\Omega_k &= \tilde{a} \left[ \frac{\kappa}{2} \frac{C T_p}{\hbar} \right]^{1/2} |k| \equiv v |k|, \\
H_0^D &= \sum_k \Omega_k a_k^\dagger a_k + \text{cnst}. \quad (40)
\end{align*}
\]

\( \Omega_k \) is the linear dispersion law of these modes with velocity \( v \) defined in Eq.(39). From the damped fluctuations of these modes, the response function \( D_\beta(\omega) \) is derived in eq (D12), within this Lagrangian approach. On the contrary, here in the incoherent phase, we aim to derive the quantum diffusion probability, stressing the interplay between intradot \( \delta g_m \) modes and the kinetics of the Q-fluctuations in the lattice.

From Eq.(30), we recognize the coupling Hamiltonian \( \hat{H}_D \), which, in the interaction representation of \( H_0^D \), takes the form:

\[
\hat{H}_D(\tau) = -\frac{\tilde{\alpha}^2}{\pi \alpha_S N \varepsilon} \sum_{p \neq 0} p^2 \beta \tau^2 2N \times \sum_k \hat{B}(\tau) a_{k+p}^\dagger(\tau) a_k(\tau). \quad (41)
\]

\( \hat{H}_D(\tau) \) of Eq.(41) represents an "effective interaction Hamiltonian" for energies in the incoherent phase. We remind that \( \hat{B}(\tau) \), defined in Eq.(33), is \( \sim \mathcal{O} \left( \frac{N}{\tau^2} \right) \) and that the hat denotes the \( m \times m \) matrix structure. As \( \frac{\tilde{\alpha}^2}{2N} \sim \mathcal{O} \left( \frac{\tilde{a}^2}{N} \right) \), the additional factor \( (\pi \alpha_S N \varepsilon)^{-1} \) in Eq.(41) makes39 \( \hat{H}_D \) of \( \left( \frac{\tilde{a}^2}{N} \right) \) and allows us to define a scaled length \( \tilde{\alpha}_t \sim \tilde{a} (\pi \alpha_S N \varepsilon)^{-1/2} \gg \tilde{a} \), which is the length scale for diffusion in the lattice.

The partition function of Eq.(30), represented in the bosonic coherent field \( \phi = \phi R e^{-1} \), can be expressed as

\[
\text{Tr} e^{-\beta H_{\text{SYK}}^D} \left\{ \text{tr}_{\phi} e^{-\beta \tilde{H}_{\text{D}}(\phi^*, \phi, \tau_+)} \right\} \quad (42)
\]

and the full quantum dynamics, is included (we drop the tilde on \( \phi \) henceforth). Here \( \text{tr}_{\phi} \) denotes the trace of a time ordered functional integral (\( \text{Tr} \tau_+ \) is time ordering in \( \tau_+ \)), while we keep the symbol \( \text{Tr} \) for the trace of the \( m \times m \) matrices. \( \hat{H}_D(\phi^*, \phi, \tau_+) \) is the matrix element derived from Eq.(41), in the coherent basis representation. In performing the trace, we assume \( \hat{H}_D(\phi^*_p, \phi_p, \tau) \) to be diagonal in the \( k \) label.

As we are dropping the \( |\phi|^4 \) term appearing in the original action \( \tilde{S} \) of Eq.(30), our toy model involves non-interacting bosonic fields only. The partition function can be written down straightforwardly by slicing the trace \( \text{tr}_{\phi} \) into \( \frac{\beta}{M} \) time slices (\( M \) integer)40:

\[
\mathcal{Z} = \lim_{M \to \infty} \text{Tr} \left\{ e^{-\beta H_{\text{SYK}}^D} \right\} \times \prod_k \left[ 1 - \left\{ e^{\frac{\beta}{M} k^2 \frac{2\varepsilon}{\pi \alpha_S N \varepsilon} \frac{2\varepsilon}{\pi} \theta \left( \frac{2\varepsilon}{\pi} \right)} \right]^M \right]. \quad (43)
\]

In Eq.(43) the dynamics of the intradot fluctuations \( \delta g_m \) and their interdot extension to the lattice are fully tangled. In view of some simplification, we limit ourselves to the regime in which the inverse timescale of the Q-fluctuations in the lattice, \( \tau^{-1} \equiv -i T_Q^{-1} \), is much smaller than the typical frequency scale of the intradot evolution (which includes the dominant term of the UV corrections). In this regime we factorize the \( \left( \frac{2\varepsilon}{\pi} \right) \) slices of the intradot propagator generated by \( H_{\text{SYK}}^D \) which is \( \tilde{F} \), while \( \tilde{B} \) includes the kernel \( G_0 G_1 \tilde{F}^{-1} \) of the Q-fluctuations. The factorization amounts to a kind of "non-interacting blip approximation"41,42 and can be justified as long as the thermalization is very effective.

With this approximation, the functional integration of the partition function of Eq.(43) can be cast in the form:

\[
\lim_{M \to \infty} \prod_k \text{Tr} \left\{ \tilde{F}(\beta) \left[ 1 - \left( 1 - \frac{\beta}{M} k_B T_0 \tilde{f}(k) \right)^M \right]^{-1} \right\}, \quad (44)
\]

where \( \frac{\beta}{M} k_B T_0 \tilde{f}(k) \) is a linearized \( m \times m \) matrix, for a small increment \( \frac{\beta}{M} \) of \( \tau_+ \), arising from the correspondence

\[
-k^2 \tilde{a}_t^2 \frac{\beta t^2}{\pi \alpha_S N \varepsilon} B \left( \frac{\beta}{M} \right) \rightarrow \frac{\beta}{M} k_B T_0 \tilde{f}(k). \quad (45)
\]
The subscript  \( \tau \) is to remind that the factorization of the traces is only justified in a limited temperature range in which the separation of the time scales holds. We have extracted a temperature scale \( T_0 \) from the left hand side, of \( O\left( \frac{\beta T}{N} \right) \), and introduced the function \( f_\tau \) of \( \mathcal{O}(1) \).

As we are on a closed time contour \(^{33}\), the partition function should be unity. The intradot propagation should be periodic in \( \tau_+ \), as well: \( \text{Tr} \left( \hat{f}(\beta) \right) = 1 \). As both \( \hat{F}(\beta) \) and \( \hat{f}_\tau \) in Eq.(44) are \( m \times m \) matrix of rank \( \tilde{r}_m \), the limit of the trace is costly from the numerical point of view. It can be done straightforwardly if we trade \( 1/\tilde{r}_m \) for the stripping of \( \hat{F}(\beta) \) off the trace. Once done this, we have checked what is the minimal \( M \) value, \( \tilde{M} \), which fulfills unitarity, at a given approximation order:

\[
z_{\tilde{M}}(k) = \frac{1}{\tilde{r}_m} \text{Tr} \left\{ \frac{1}{1 - \frac{1}{\tilde{r}_m} \left( e^{-\frac{2\pi}{\tilde{M}} k \cdot \tilde{r}_m} \right)^3} \right\} \approx 1. \quad (46)
\]

In Fig.4, we plot an interpolated smoothed curve of the (approximate) lowest \( \tilde{M} \) value, which satisfies Eq.(46), vs. \( T_0/T \), for \( k^2\tilde{a}_t^2 = 1 \). Precision is up to \( > 10^{-5} \). \( \tilde{M} \) is practically constant, when \( \frac{T_0}{T} \geq 1 \), but it increases strongly when \( T \) takes values \( T > T_0 \). The trend is only meaningful for \( T_0/T \sim 1 \), because \( T \) values larger than \( T_0 \) require \( n > 12 \) in the spectral representation of Eq.(20) and matrices \( m \times m \) of rank \( \tilde{r} > 3 \), i.e. higher than the ones used here. Fig.4 is the numerical proof that \( T_0 \) represents the temperature above which the thermalization is more efficient and our factorization between evolutions breaks down. The threshold temperature scale \( T_0 \) introduced in Eq.(45) and the space scale \( \tilde{a}_t \) defined after Eq.(41) are discussed in Subsection C.

**B. \( (\delta G^{\mu \nu} \delta G^\lambda) \) diffusion probability**

The generating functional to obtain the correlator of the field \( \phi_\rho(\tau) \) at different times \( y' - y'' \) ( \( y'' \equiv 2\pi \tau/M; \ r, s \text{ integers} \) ) can be derived from Eq.(42) by adding a source term. Its \( r, s \) matrix element can be denoted as \( \text{Tr} \left( \left[ \hat{f}_r + \hat{H}_D \right]^{-1} \right) \) :

\[

\lim_{M \to \infty} \text{Tr} \left\{ \left[ \hat{f}_r + \hat{H}_D \right]^{-1} \right\} \approx \lim_{M \to \infty} \text{Tr} \left\{ \left[ \hat{f} \left( \frac{\beta}{M} \right) \right]^{-r-s} \right\} \frac{\left( 1 - \frac{1}{M} \hat{f}_r \right)^{(r-s)}}{1 - \left[ \hat{f} \left( \frac{\beta}{M} \right) \left( 1 - \frac{1}{M} \hat{f}_r \right)^M \right]}
\]

\[

\approx \lim_{M \to \infty} \text{Tr} \left\{ \left[ \hat{f} \left( \frac{\beta}{M} \right) \right]^{-r-s} \right\} \text{Tr} \left\{ \left[ e^{-\frac{2\pi}{\beta M} \hat{\tau}} \right]^{(y''-y')} \right\}, \quad (47)
\]

To sum up, the steps of the logical inference starting from Eq.(45) are

\[

- k^2 \tilde{a}_t^2 \frac{\alpha_s N_c}{\pi} \delta_{\beta N_0} \beta M \left\{ \frac{\beta T_0}{M} \right\} \to \frac{\beta}{M} kBT_0 \hat{f}_r \left( \frac{\beta}{M} \right) \\

\to \ln \left( \text{Tr} \left\{ e^{-\frac{2\pi}{\beta M} \hat{\tau}} \right\}^{\delta_{\beta N_0}^2} \right) \to \frac{T_0}{MT} \hat{f}_r \to \tilde{D}_Q k^2 T_0^2 \quad (49)
\]

As the left hand side is \( \mathcal{O} \left( \frac{\delta \beta}{N} \right) \), the product \( \tilde{D}_Q T_0^2 \) is \( \mathcal{O} \left( \frac{\delta \beta}{N} \right) \), as well. In fact we will put \( \tilde{D}_Q T_0^2 = \tilde{a}_t^2 = \tilde{a}_t^2 \frac{T_0^2}{T} \). The diffusion coefficient \( \tilde{D}_Q \) and the relaxation time scale of the diffusion process \( i T_0^2 \equiv \tau \) are discussed
FIG. 5: Plot of the logarithm of the propagator appearing in Eq. (48) for $y^r - y^s = 1$ (i.e., $M = 1$),\[ \ln \left( \text{Tr} \left[ e^{-T_k y^r} \right] \right) \], vs $(k\tilde{a})^2$, for $T_h = 0.005, 0.01, 0.03$, displaying an approximately linear behavior as a function of $(k\tilde{a})^2$, for $(k\tilde{a})^2 > 0.5$.

in Subsection C.

Our simplified approach to the diffusive constant $D_Q$ provides an analytical approximate expression for $\langle \delta G(0^+, \tau_+) \delta G(0, 0) \rangle_k$. From the left hand side of the second line of Eq. (47) we write:

$$\langle \delta G(0^+, \tau_+) \delta G(0, 0) \rangle_{k\neq 0} \sim \frac{e^{-T_k y^r}}{1 - e^{-2\pi T_k}}, \tag{50}$$

with $T_k = \beta D_Q k^2 / 2\pi$. By Fourier transforming to Matsubara Bose frequencies ($n$ stands here for $\beta \Omega_n / 2\pi$) we obtain:

$$\int_0^{2\pi} dy \frac{e^{-T_k y^r}}{1 - e^{-2\pi T_k}} e^{in y} = \frac{1 - e^{-2\pi T_k} e^{2\pi i n}}{T_k - i n} \frac{1}{1 - e^{-2\pi T_k}},$$

so that:

$$\{ \langle \delta G(0^+, \tau_+) \delta G(0, 0) \rangle \}_{k\neq 0, n} \sim \frac{2\pi}{\beta} \frac{1}{D_Q k^2 - i \Omega_n}. \tag{51}$$

This result highlights the diffusive pole in the Fourier transform of the Q-fluctuation correlator.

Identifications of the threshold temperature, $T_0$, and of the space parameters, $\tilde{a}_k, D_Q T_0$, introduced as scales in the previous derivation, require a modelization of the damped dynamics of the Q-excitation, which we derive in the next Subsection C. While these parameters, in the course of the derivation, have been recognized as marginal in a renormalization group sense, as they are $O\left( \frac{\beta T_0}{\tilde{N}} \right)$ (in the limit $J, N \to \infty, J/N \to \text{const}$), they should rest on phenomenological fundamental quantities, like the thermal capacitance $C$ per unit mass and the thermal conductivity $\kappa$. These quantities will be related to two parameters, i.e. the damping of the neutral Q-excitation, $\Gamma$, and their propagation velocity $\nu$, given by Eq. (39) in our model. The velocity $\nu$, appears in the linear spectrum of the Q-excitation given by Eq. (40), while $\Gamma$ is introduced as a broadening of their spectral peak. Subsection C is devoted to the presentation and discussion of these relations.

C. Thermalized and coherent energy processes

In the previous Sections we have shown how the $pGm$ within each SYK dot, $\delta g_m$, generate energy modes diffusing in the lattice of the extended SYK model. The validity of our approach, involving the partial factorization that we have adopted in our traces, rests on different temporal dependence scales of the center of mass times $y^r - y^s$ on one side and of the intradot fluctuations on the other. The "interdot" dynamical time scale is discussed phenomenologically in this Subsection.

The "interdot" time scale is the thermalization time $T_Q$, introduced in Eq. (49). $T_Q$, the threshold temperature scale for thermalization $T_0$, and the space scale $\tilde{a}_s$, are connected with the velocity $\nu$, given by Eq. (39), and with the phenomenological damping $\Gamma$, which is the inverse lifetime of the Q-excitation. In turn, these quantities depend on the thermal capacitance $C$ per unit mass and the thermal conductivity $\kappa$, which are the phenomenological, experimentally measurable quantities.

When adopting our rough approximations, we cannot ignore that these parameters depend on temperature. In particular, the two sets of scales should be considered, $(T_{coh}, T_Q)$ related to particle transport in the coherent phase and $(T_0, \frac{\tilde{N}}{\tilde{m}})$ for a thermalized system in the incoherent phase, when $T > T_0 > T_{coh}$. The parameters $T_{coh}, T_Q, T_0$ and $\tilde{a}_s^2$ are $O\left( \frac{\beta T_0}{\tilde{N}} \right)$. We will discuss these two regimes in this Subsection. At finite $T$ the small gap of the $Q-$excitations can be disregarded. We assume that both regimes have gapless and chargeless bosonic excitation modes of energy $\Omega_k = \hbar v k$, given by Eq. (39). Indeed, we exclude charging effects in transport. In the incoherent regime the gaussian action of Eq. (28) involves energy density $N \propto \partial \theta \partial \theta$ fluctuations and energy flux density $\tilde{N} \propto J_Q$ fluctuations$^{12}$, the Q-excitation. In the coherent phase, bosonic excitations are particle-hole excitations with fluctuations of the particle number $N_0$ and first sound excitations. In the appendix E we show that the sound mode survives when the interaction with the SYK dots is turned on perturbatively, embedded in the p-h continuum. We attribute an inverse lifetime $\Gamma \propto T$ to these excitations.

We proceed with the incoherent regime first, at $T \approx T_0$. The thermal conductivity, in presence of a damping $\Gamma$, is derived in Eq. (D14) from the $J_Q - J_Q$ response:

$$\kappa = k_B \Gamma \left( \frac{h \Gamma}{C T} \right)^{1/2}. \tag{52}$$
From one of the Einstein relations, the diffusivity $\tilde{D}_Q$ is related to the thermal capacitance $C$ and to the density $\rho_0$, according to (p-h symmetry is assumed)
\[
\tilde{D}_Q = \frac{k}{C\rho_0}. \tag{53}
\]
As the chemical potential $\mu$ is assumed to vanish, the 2-d particle density $\rho_0$ involved in these excitations, given by Eq.(29), is not well phenomenologically defined. We will estimate it as $\rho_0 = \Gamma^2/v^2$, a choice that will turn out to be consistent with our results of this Subsection. We proceed now by deriving an estimate of $T_0$. In this case, energy diffusion is mainly due to heat transport in a highly thermalizable environment and we use the first temperature dependent correction to the energy of the SYK model\(^8\): $\delta E = c/(2\beta^2) = CT$ in Eq.((52), where $c = 4\pi^2\alpha_S N/J$. From Eq.(52), we get:
\[
\kappa = k_B \Gamma^{3/2} (h\beta)^{1/2} \left( \frac{\beta J}{\pi \alpha_S N} \right)^{1/2}, \tag{54}
\]
which, inserted in Eq.(53), with $\rho_0 = \Gamma^2/v^2$ gives:
\[
\tilde{D}_Q = \Gamma \left( \frac{\beta J}{\pi \alpha_S} \right) \tilde{a}^2, \tag{55}
\]
where Eq.(39) has been used. On the other hand, the last inference in Eq.(49), together with Eq.(45), suggests that $\tilde{T} \propto k^2\tilde{a}^2$, with $\tilde{a}^2 \sim \tilde{a}^2 T_0/T$. We conclude from Eq.(55) that $\tilde{D}_Q \propto \Gamma \tilde{a}^2$ and, as $\tilde{D}_Q T_Q = \tilde{a}^2$, the relaxation time $^{34} T_Q \sim \Gamma^{-1} \sim h\beta$. Thermalization is better handled in euclidean time. Putting $T_Q \rightarrow h\beta$ in $\tilde{a}^2 T_0 - \tilde{D}_Q T_Q$ and using Eq.(55), we conclude that
\[
k_B T_0 \sim 2\pi h \Gamma \left( \frac{\beta J}{N\pi \alpha_S} \right). \tag{56}
\]
This equation qualifies $k_B T_0$ as a threshold energy for efficient thermalization and confirms that $T_0$ is $O\left( \frac{\alpha_S N}{N\pi \alpha_S} \right)$, if just the zero order for $\Gamma$ is retained. As we have assumed that $\Gamma \propto T$, both $\tilde{D}_Q$ of Eq.(55) and $T_0$ of Eq.(56) are temperature independent.

Our approximations, which involve some kind of adiabatic factorization, do not allow us to discuss the coherent carrier transport regime, $T \lesssim \Omega_0 \lesssim T_{coh} < T_0$, except for a very qualitative bird’s eye. Indeed, the convergence of the ‘normalization’ of Eq.(46) in Fig.4 is misleading, as one should keep in mind that just the dominant UV contribution of $\mathcal{F}$ has been retained and all the regular contributions (belonging to the fluctuation domain orthogonal to the pGm’s) have been neglected. These include low energy contributions and their evolution cannot be factorized. Anyhow, back to Eq.(53) for this case, an approximated expression for the specific heat arising from the gapless modes of the model given by Eqs.(39,40) is given by Eq (D15):
\[
C_V = k_B \left( \frac{k_B T}{h\nu} \right)^2 6 \zeta[3], \tag{57}
\]
where $\zeta[n]$ is the Riemann function. When the velocity $v$ is inserted in this expression, we get an equation for $1/C_V^2$, which can be related to Eq.(52) to give:
\[
\left( \frac{k}{k_B T} \right)^4 \frac{1}{(h\Gamma)^2} = \frac{1}{C_V^2 T^2} = \frac{2\pi}{k_B} \frac{\hbar}{(k_BT)^4} \frac{\kappa}{6 \zeta[3]} \tag{58}
\]
Inserting this result in Eq.(53), with $v$ given by Eq.(39) and $\rho_0 = \Gamma^2/v^2$ we obtain:
\[
\tilde{D}_Q = \frac{k^2}{\Gamma^2} \frac{T}{\hbar k_B} \tilde{a}^2 = \frac{\hbar^2 T^2}{h \Gamma} \frac{1}{6 \zeta[3]} \tilde{a}^2. \tag{59}
\]
Assuming again $\Gamma \sim T$, Eq.(59) shows that the diffusion constant is in this case $\propto T$ as in the Einstein-Smoluchowski formula. At least formally, it can be put in the form of a bound on the diffusion rate, which has been conjectured for strongly interacting systems at zero chemical potential:\(^{34,44}\):
\[
\tilde{D}_Q \geq \frac{\hbar^2 \tilde{a}^2}{k_BT}. \tag{60}
\]
In this case the velocity which arises here is not $v_F$ but $\tilde{v} \sim \tilde{a} \Gamma$. Eq.(60) is non universal.

Now we proceed just by analogy with the previous case and we assume that, just by replacing $T_0$ with $T_{coh}$, we can put here
\[
\tilde{a}_0^2 \sim \tilde{D}_Q T_Q = \tilde{a}^2 T_{coh}/T. \tag{61}
\]
From Eq.(59) it follows that:
\[
\tilde{a}_0^2 T_{coh}^2 \sim \frac{\hbar^2 T^2}{h} \frac{1}{k_BT} \frac{1}{6 \zeta[3]} \tilde{a}^2 T_{coh}^2 T_Q, \tag{62}
\]
what implies:
\[
T_{Q}^{coh} \sim (6 \zeta[3])^{1/3} \frac{k_B T_{coh}}{\hbar \Gamma^2}. \tag{63}
\]
Given $T_{coh} \propto l_0^2/NJ \sim J/N$, is $T_{Q}^{coh} \propto T^{-2}$ as in the Fermi liquid case.

VI. SUPERCONDUCTIVE COUPLING AT LOW TEMPERATURE

In this Section we present an Eliashberg approach to the superconducting instability of a quantum electron liquid that contains the Q-excitations in its energy spectrum. As explained in the Introduction, we consider a model with two components: a lattice of local $0 + 1 - d$ SYK dots and an underlying FL which interacts with the
The isotropic self-energy arising from the interaction, for a quasiparticle residue $\nu$, and a single particle energy $\epsilon_\nu = \nu_F k$ in the continuum limit, with a renormalized physical velocity $\tilde{v}_F = Z\nu_F$ and a residual local interaction of strength $U_c$, which is dealt with perturbatively. The isotropic self-energy arising from the interaction, for $k$ on the Fermi surface, is:

$$\Sigma(k_F,\omega) = -\omega Z^{-1} - i \alpha n_0 |\omega|^2 \ln \left( \frac{Z\nu_F k_F}{|\omega|} \right) \text{sign}(\omega)$$

For $T > \Omega^* = W^2/U_c$, we should put $2 \ln(\mathcal{J}/W) \to \ln(\mathcal{J}/T)$ in Eq.(67). In Eq.(64), $\theta$ is the angle between $\vec{q}$ and $\vec{k}$ and, for $|q| \ll k_F$, we have approximated $\epsilon_{k_F+\vec{q}} - \epsilon_{k_F} \approx \tilde{v}_F \cos \theta$.

$$\Pi(q,i\omega) = \sum_p \sum_{\omega_n} G(p,i\omega_n) G(p+q,i\omega_n+i\omega_m).$$

In the range of frequencies $\Omega < \Omega^* = W^2/U_c$, where $W$ is the bandwidth, there are two contributions to the polarization, one (labeled by $i = 1$) coming from the residual FL interaction and a second one ($i = 2$) coming from hybridization with the incoherent disordered SYK clusters of $0+1-d$ neutral fermions, interacting at energy $\mathcal{J}$, one at each lattice site (see Fig.1). While $\Pi^1(q,i\Omega)$ uses the Green’s function which appears in Eq.(64) with a simple pole, $\Pi^2(i\Omega)$ is evaluated from the single particle Green’s of the SYK model, in the conformally symmetric limit, which is local in space (i.e. $q-$independent) and reported in Eq.(8). Approximately, is:

$$\Pi^1(q,i\Omega) \approx Z\nu_0 \left[ 1 - \frac{\Omega \text{sign}(\Omega)}{\sqrt{\Omega^2 + (Z\nu_F k)^2}} \right],$$

$$\Pi^2(i\Omega) \approx -\frac{4}{\mathcal{J}} \ln \left( \frac{\mathcal{J}}{\max[\Omega,\Omega^*]} \right) \to -\frac{8}{\mathcal{J}} \ln \left( \frac{\mathcal{J}}{W} \right).$$

Here $\nu_0 = k_F/(2\pi\hbar\tilde{v}_F)$ is the density of states at the Fermi surface and $Z\nu_0 \sim U_c^{-1}$. In performing the integral over momenta $p$, we have assumed that, at low temperatures $T << \Omega^*$, the difference in occupation numbers $n_F \left( Z\epsilon_{p+q} - n_F \left( Z\epsilon_p \right) \right) \approx -\delta(\epsilon_q) \tilde{v}_F q \cos \theta$.

Moving to real frequencies we get:

$$\Sigma(k_F,\omega) = -\omega Z^{-1} - i \alpha n_0 |\omega|^2 \ln \left( \frac{Z\nu_F k_F}{|\omega|} \right) \text{sign}(\omega)$$

(67)

$$= \left[ |\omega| \frac{\epsilon_F}{\Omega^*} \ln \left( \frac{\mathcal{J}}{W} \right) + \frac{1}{\tau} \right] - Z \Im \Sigma(k_F,\omega)$$

(68)

($\alpha$ is a parameter of order one), which shows that to the lowest approximation, the perturbed FL is a Marginal Fermi Liquid. The interaction of the electronic quantum liquid (qL) delocalized over the $2-d$ lattice with the SYK clusters, makes the quasiparticles not well defined, but still with a well defined Fermi surface. In the appendix E, we derive the lowest lying collective excitations in the present perturbative frame. The hydrodynamic collective excitation, the would-be acoustic plasmon, is also rather well defined. At strong coupling, in the limit $U_c \to \mathcal{J}$, its dispersion tends to the boundary of the p-h continuum and the imaginary part, which blurs the mode, vanishes. The acoustic plasmon is on the verge to emerge as a bound state at low energies, split off the p-h continuum.

**B. Superconductive critical Temperature**

We outline here the derivation of the superconducting critical temperature $T_{c}$, of a $2-d$ qL in interaction with the SYK lattice, using the Eliashberg approach. Although we are unable to discuss the nature of the microscopic low temperature electron-electron interactions driven by the virtual Q-fluctuations, we assume that Cooper pairing is induced in a qL of bandwidth $W$, by virtual coupling with the diffusive energy Q-modes in the lattice, which, in turn, are generated by the pGn of the SYK clusters, as discussed in the previous Sections. Three energy scales come into play in this context. The
energy scale \( t_0^2/\mathcal{J} \sim W^2/\mathcal{J} \), associated with the temperature threshold \( T_{coh} \), below which coupling between the SYK clusters and the qL is perturbative. Two more energies associated to the coupling between the qL and the Q-excitation, the coupling strength \( g \) and the energy cutoff of the interaction \( U > U_c \), which also appears in the typical frequency for the attractive interaction \( \Omega' = t_0^2/U \sim W^2/U \) (in this Subsection is \( h = 1 \)). This assumptions immediately implies an electronic energy scale, as the reference scale for the superconducting transition. Our standard approach to the superconducting transition within the Eliashberg theory\(^3\) gives rise to a non BCS-like phase transition. The non BCS critical temperature is a direct consequence of the quantum liquid to be marginal and of the excitation modes to be diffusive.

In a mean field superconducting Hamiltonian, in the Nambu representation, the one electron Green’s function and the electronic self-energy \( \Sigma(p, i\omega_n) \) are \( 2 \times 2 \) matrices defined by the Dyson equation

\[
[G(p, i\omega_n)]^{-1} = [G_0(p, i\omega_n)]^{-1} - \Sigma(p, i\omega_n),
\]

where \( G_0(p, i\omega_n) \) is the one-electron Green’s function for the non interacting system \( [G_0(p, i\omega_n)]^{-1} = i\omega_n - \xi_p \sigma_3 \) and the approximation used for the self-energy is (see appendix F):

\[
\Sigma(p, i\omega_n) = -\frac{1}{\beta N_q} \sum_q \sum_{p'\nu'} \sigma_3 G(p', i\omega_n) \sigma_3 \, d\Omega_q \left( g_{p'p}(\Omega_q) \right)^2 B(\Omega_q, \Omega) \left\{ \frac{1}{i\omega_n - i\omega_n - \Omega} - \frac{1}{i\omega_n - i\omega_n + \Omega} \right\}.
\]

(70)

Here \( \vec{q} = \vec{p} - \vec{p}' \) is the transferred momentum and \( \Omega_q = \bar{D}_Q q^2 \) is the energy of the collective excitations. An isotropic coupling density \( g(p, p') \) is assumed and, in place of the sum over \( p' \) vectors, we integrate over \( \Omega_q \), with the energy density of the \( q \) momenta \( \frac{1}{q^2} \). The imaginary part of the retarded energy flux density response function is

\[
B(\Omega_q, \Omega) = -\frac{1}{\pi} \Im \left\{ D^R(\Omega_q, \Omega) \right\} = -\frac{1}{\pi} \left[ \frac{\Omega_q \Omega}{\Omega^2 + \Omega_q^2} \right] T_Q.
\]

(71)

Note the difference, due to diffusivity, with respect to the usual Eliashberg approach, in which \( B(q, \Omega) \sim \frac{\Omega_q}{\Omega_q + \Omega} \).

We take \( \bar{D}_Q T_{coh} = \bar{a}_q^2 = \bar{a}_q^2 T_{coh}/T \), as in Eq.(61) (we drop the label \( coh \) from \( T_{coh} \) in the following).

Using Eq.(67), Eq.(69), we write \([G(p, i\omega_n)]^{-1} = i Z^{-1} \omega \left( \xi_p - i Z^{-1} \left[ \frac{|\omega| \varepsilon_F}{\Omega_q} \ln \left( \frac{\mathcal{J}}{W} \right) + \frac{\alpha}{Z} v_0 |\omega|^2 \ln \left( \frac{Z \tilde{v}_F k_F}{|\omega|} \right) \right] \right) \sigma_3 - \Xi \sigma_1 \), in which the mean pairing field \( \Xi \) has to be self-consistently determined:

\[
[G(p, \omega)]^{-1} = Z^{-1} \omega \mathbf{1} - \left\{ \xi_p - i Z^{-1} \left[ \frac{|\omega| \varepsilon_F}{\Omega_q} \ln \left( \frac{\mathcal{J}}{W} \right) + \frac{\alpha}{Z} v_0 |\omega|^2 \ln \left( \frac{Z \tilde{v}_F k_F}{|\omega|} \right) \right] \right\} \sigma_3 - \Xi(\omega) \sigma_1.
\]

(72)

The final result for Eq.(70) is:

\[
\Sigma(k_F, \omega) = v_0 \int_{-\infty}^{\infty} d\omega' \Re \left\{ \frac{Z^{-1} \omega' \mathbf{1} - \Xi(\omega') \sigma_1}{|\mathcal{P}(\omega')|^{1/2}} \right\} \int_{0}^{\infty} d\Omega_q \left\{ \sigma_3 \right\} - \Xi(\omega) \sigma_1, \quad \Xi(\omega) \sigma_1.
\]

(73)
where $N(\Omega)$ and $f(\omega)$ are the Bose and Fermi occupation probabilities. The term in curly brackets arises from $\Im \{\nu_0 \int_{-\infty}^{+\infty} d\xi \sigma_3 G(p', \omega) \sigma_3\}$ which turns into a real part by working out the inverse of Eq.(72). A limited region contributes to the integral $\int d\xi'$, but we can extend the integration limits to infinity with no big error.

Following McMillan, we want to find an approximate solution to the gap equation ($\Delta(\omega) = Z(\omega)\Xi(\omega)$). At the critical temperature, $\Delta \sim 0$ and can be dropped in the denominator, but the gap equation has to be satisfied. From Eq.s(72), the term multiplied by $\sigma_1$ gives

$$\Delta(\omega) = Z\nu_0 \int_0^{\infty} \frac{d\omega'}{\omega'^2 + \left(\frac{1}{\tau(\omega')}\right)^2} \frac{\text{Re} \{\Delta(\omega')\}}{\omega'} \int_0^\infty d\Omega \frac{\tilde{\alpha}_c^2}{DQ} \int_0^\infty d\Omega' |g_{kF,\omega'}(\Omega')|^2 \mathcal{B}(\Omega, \Omega)$$

$$\times \left\{ [N(\Omega) + f(-\omega')] \left[ \frac{1}{\omega + \omega' + \Omega} + \frac{1}{-\omega + \omega' + \Omega} \right] - [N(\Omega) + f(\omega')] \left[ \frac{1}{\omega - \omega' + \Omega} + \frac{1}{-\omega - \omega' + \Omega} \right] \right\}. \quad (74)$$

$\tau(\omega')$ is the lifetime of the quasiparticles from Eq.(68).

In the rest of the calculation we neglect the thermal excitations and drop $N(\Omega)$. Two energy ranges contribute to $\Delta(\omega)$:

$$\Delta(\omega) = \begin{cases} \Delta_0 & 0 < \omega < \Omega^* \\ \Delta_\infty & \Omega^* < \omega \end{cases}. \quad (75)$$

The first, $\Delta_a(\omega)$, arises from integration over $0 < \omega' < \Omega^*$ and the second, $\Delta_b(\omega)$, from the integration over $\Omega^* < \omega' < \omega < U$ ($U$ is the cutoff energy). Hence $\Delta(\omega) = \Delta_a + \Delta_b$.

While $\Delta_0$, can be assumed as the usual order parameter in the lattice, it is unclear what $\Delta_\infty$ is, when $\omega > \Omega^*$ and incoherence is established at these energies. In the mean field approach, $\Delta_\infty$ can be thought of some kind of intradot field induced by the ordering of the low energy system. Of course we concentrate on the ordering transition for $\omega < \Omega^*$, but both $\Delta$’s should be non vanishing.

Observing that the integration variable $\Omega_q$ has the meaning of the diffusive energy (see Eq.(71)), it is clear that it cannot be integrated at energies above $\Omega$. We also use the parameter equality $Z\nu_0 = U_c^{-1}$ and we take $|g_{kF,\omega'}(\Omega_q)|^2 = g^2$ constant ($[g]^{-1} \sim \text{time} (h = 1$ here):
two contributions together \( \Delta_0 = \Delta_a(0) + \Delta_b(0) \) we have:

\[
\Delta_0 = \frac{8\pi}{|g|^2 \ln 2} \frac{\tilde{D}_Q}{a_T^2} \frac{1}{T_Q} \frac{W}{U_c \ln \beta \Omega^*} \ln \beta \Omega^* \\
= \frac{\Delta \alpha}{\beta} \left( \frac{U^2}{W^2} \right)^2 \frac{1}{\ln \beta \Omega^*}.
\]

(78)

Using the definition of \( T_Q \equiv T_Q^{\text{coh}} \) given by Eq.(63), the pairing parameter takes the form:

\[
\frac{\tilde{D}_Q}{|g|^2 a_T^2} \frac{1}{T_Q} \sim \frac{1}{|g|^2 T_Q^2} = \left[ \frac{\Gamma}{|g|} \frac{h \Gamma}{k_B T_Q^{\text{coh}}} \frac{1}{(6 \zeta[3])^{1/3}} \right]^2.
\]

(79)

As \( T_Q, T_Q^{\text{coh}} \sim O \left( \frac{k_B T_Q^{\text{coh}}}{N \pi a_S} \right) \), it follows that \( |g| \sim O \left( \frac{N \pi a_S}{k_B T_Q^{\text{coh}}} \right) \), so that \( |g| T_Q^{\text{coh}} \sim O(1) \). Assuming both \( \Delta_0 \) and \( \Delta \alpha \) to be non zero, Eq.(78) gives:

\[
\frac{k_B T_c}{\Omega^*} = \left( \frac{J}{W} \right) \frac{1}{\ln \beta \Omega^*} \frac{\Delta \alpha}{\beta} \left( \frac{U^2}{W^2} \right)^2 \frac{1}{\ln \beta \Omega^*} \frac{1}{(6 \zeta[3])^{1/3}}.
\]

(80)

Eq.(80) provides the value of \( T_c \) on a scale of \( \Omega^* \), which is a power of \( J/W \), which is difficult to determine, because it requires the full quantitative characterization of the model. However, qualitatively, the non-BCS behavior is fully apparent. Indeed, \( T_Q \) itself is a function of the temperature, because the energy width of the mode relaxation, \( \Gamma \), appearing in Eq.(79), is expected to be \( \sim T \). In this case, Eq.(80) defines \( T_c \) only implicitly. Dropping the first negative exponent and writing the second exponent as \( \omega^4 / \lambda \) where \( \omega = \frac{k_B T_c}{\hbar} \), the zeros of the function \( F[\omega] = u - \Theta(u^4 / \lambda) \) give the \( T_c \) value. In the prefactor \( \Theta \sim W / J \), all the unknown features of the pairing interaction is lumped. \( \Theta \) strongly depends on the cutoff energy \( U_c / W \) and on \( J/W \), as well as on the lifetime of the quasiparticles in \( 2 - d \) at higher energy \( \sim W \) (see Eq.(72)). \( F[\omega] \) is plotted in Fig.6 vs \( u \), at \( \Theta = 0.1 \), for \( \lambda = 5, 0.5, 0.2 \). Increasing the pairing strength \( \propto |g|^2 \), \( \lambda \) increases, and so does \( T_c \).

VII. CONCLUSIONS

Hopefully, the intriguing high temperature "strange metal" phase of materials undergoing a HTc superconducting transition is at a turning point, since attention was drawn to the "universal" linear dependence on \( T \) of the resistivity and to features like the possible violation of the Wiedemann and Franz law. The Wiedemann and Franz universal ratio unambiguously rests on the coexistence of heat and charge transport typical of weakly interacting electronic Fermi Liquids. It is accepted that interactions in these systems are strong and play a nonperturbative role. This gives credit to a Non Fermi Liquid (NFL) perspective for the high temperature normal "bad metal" phase. Consensus in the physics community increases on the use of the doped Mott insulator paradigm as an interpretation ground for the copper oxide HTc materials. On the other hand, strong crystal anisotropy and doping tend to privilege the role of copper-oxide planes and the role of collective fluctuations. Even when clean single crystals are available, the doping and the chemistry of the charge reservoir layers separating the CuO2 plane(s) from one another could induce inhomogeneities.

It was really a twist the discovery that the Su-Ye-Kitaev model, in the limit of strong interactions and strong disorder, can be solved exactly in \( 0 + 1 \) dimension displaying a NFL incoherent toy system, with non trivial properties as a zero temperature finite entropy and a chaotic behavior at long times. Moreover, hydrodynamical extensions to higher space dimensions provides the linear dependence of the resistivity which has attracted a flurry of interest from the condensed matter community. By contrast, a conventional Eliashberg approach, typical of FL systems has not been seriously questioned.

On one side we enquire on the influence of a hopping term between neighbouring SYK clusters organized in a \( 2 - d \) lattice. Hopping is assumed to be marginal in a \( 1/N \) expansion and strong coupling \( J \) limit, with \( \beta J / N \) kept finite. On the other side we study the perturbative effect that the SYK lattice exerts on a FL with delocalized electrons (in the continuum hydrodynamical limit), displaying a well defined Fermi surface and a large Fermi
energy. The aim is to characterize the collective excitations of the SYK system in view of identifying the latter as responsible for the superconducting instability via an unidentified mechanism.

There are various ways to extend the SYK model to a lattice and we use one of them\textsuperscript{11,13,18}. All of them rest on a disorder average and we assume that self-averaging allows for a translationally invariant approach with wavevector $k \bar{a} \ll 1$, where $\bar{a}$ is the lattice spacing. We focus on the role of the collective fermionic excitations of a SYK dot $\delta g_m$ of Matsubara frequency $\omega_m$. Among these, there are also incipient Goldstone Modes, which originate from the spontaneous breaking of the conformal symmetry. However, they acquire mass when the first UV correction $\sim O \left( \frac{N}{\pi} \right)$ is included. They are denoted as pseudo Goldstone modes (pGM’s) in the text\textsuperscript{6,8}. The UV correction forces locality in space and time. The real fermionic propagator $\delta g_m$ of the IR limit acquires a complex local phase in the extended SYK action, due to minimal compact $U(1)$ coupling. The energy fluctuations driven by these dressed excitations across the lattice can be monitored by investigating the correlations of a local space-time UV "order parameter" of the incoherent phase, an extension of the bilocal two-point propagator $G_c$ of the conformal symmetric limit. In a NFL system the energy can be interpreted as energy density excitations, better than chemical potential fluctuations. In this work our focus was on the nature of these dressed bosonic fluctuations which we nickname as "Q-excitation" and on the response $D^3$ of the lattice system to perturbations which excite them. In the recent past, the scaling of $U(1)$ RVB models with a gap to both charge and spin excitations has been studied\textsuperscript{51}.

We take advantage of the fact that two time (or temperature) scales come into play, the "fast" intradot fermionic $\delta g_m$ modes and the "slow" interdot bosonic energy density fluctuations originating from the Q-excitation. This allows us to characterize the dynamics in a range of energies $\sim T_0$, where $T_0$ is a threshold temperature for efficient thermalization. In this temperature range the Q-excitation are proved to be diffusive when the dynamics induced by the UV correction is appropriately accounted for. Diffusivity arises from the combination of disorder in the SYK dots and hopping in the superlattice. We find the mode-mode correlations in imaginary time $\propto G_cG_cF^{-1}$ where $F$ is the bilocal 4-point propagator, which diverges in the conformal limit, but is made finite when the UV dominant correction is included. The presence of $F^{-1}$ in the diffusion parameter is the signature of the presence of the pGM and is the main result of this work. The corresponding retarded response function in real time can be derived from the correlations, by analytic continuation to real frequencies $i \omega_m \rightarrow \omega + i 0^+$. A similar result was derived directly in real time\textsuperscript{12}, but without including the role of the pGM and is reproduced in the appendix C. In the real time approach, the factor $F^{-1}$ does not appear as part of the diffusive pole. The scaling rinormalizes the thermalization temperature $T_0$ and the diffusion constant $D_Q$, by introducing a lattice length $\tilde{a}_t \sim \tilde{a} T_0/T$ and a diffusion time $T_Q$, such that $D_Q T_Q \sim \tilde{a}_t$. A simple quantum approach to the dynamics of the energy fluctuations in presence of damping $\Gamma$ allows for their explicit determination, $\Gamma \sim T$ is the energy broadening of the Q-fluctuation excitation due to relaxation in the lattice. If we resort to the Einstein relations which connects the diffusion coefficient $D_Q$ to the transport coefficients\textsuperscript{34}, we derive the temperature dependence of these quantities and obtain Eq.(60) which refers to $\tilde{v} \sim \tilde{a} \Gamma$ as a physical (non universal) diffusion velocity. Eq.(60) has to be contrasted with a bound for incoherent systems that has been conjectured\textsuperscript{44}.

In the study of the correlations, it emerges clearly (see Fig.4) that our approach to the partition function and to the generating functional is only valid for $T \sim T_0$, an energy range which we conclude to be well separated with respect to the one $\sim T_{coh}$, the temperature which marks the prevail of the low energy Fermi Liquid. For $T < T_0$, entanglement of the dynamics of the pGM's in the SYK dots with the dynamics of the energy fluctuations across the lattice require more sophisticated methods than the factorization used here, in the calculation of the thermodynamic functionals. Still, some qualitative hint is presented in Subsection V.B.

In Section VI.B we assume that the Q-excitation has a role in the superconducting instability at low temperature. A dispersive self-energy for an electronic quantum liquid perturbed by a SYK dot with a local interaction $\mathcal{J}$ turns the FL into a marginal FL\textsuperscript{52}, with inverse lifetime of the quasiparticles close to the Fermi surface $\propto \omega$. The quasiparticle lifetime influences the mean field superconductive order parameter $\Delta_{\omega}$ at energies above $\Omega^* \sim W^2/U_c$, where $U_c$ is cutoff energy for the pairing interaction. The topic, whether the Q-excitation could really play the role of virtual excitations inducing pairing, provided an appropriate attractive coupling is active\textsuperscript{53,54}, is beyond the present state of the art. It is an old idea that an incipient Goldstone mode of an ordered phase can accomplish this task. This possibility was examined in the past and it was concluded that the fluctuations involved would lead to a depression of $T_c$\textsuperscript{55}. We think that this pattern may not work here for various reasons. Here, indeed, the vertex corrections vanish to lowest approximation order. However, the fluctuations driving the transition do not arise from an incipient order but are non local in time, in a fully disordered system. What we call "order parameter" here is energy relaxation modes which are effectively non local in space and non number conserving in nature, as phonons would be. We have also omitted the influence of long-range Coulomb interactions, which certainly modifies the spectrum of boson density fluctuations\textsuperscript{49}.

Of course, if the Q-modes play a role, the temperature scale of the superconducting $T_c$ is of electronic origin, $\sim \Omega_{\epsilon}^*$, defined in Eq.(80). $T_c$, as derived using the Eliashberg\textsuperscript{45} and McMillan\textsuperscript{46} approaches, is not BCS-like
and appears as the zeros of a function $F \left[ \frac{k_B T_c}{\Omega \nu} \right]$, which is plotted in Fig.6. It also depends on the "low" energy scale $\sim T_Q^{-1}$, on the lifetime at higher energies of the Cooper-pairing electron charges and on the diffusion length of the Q-excitations. Indeed, the correlation length of the pairs $\xi$ depends on the effective mean square length $\tilde{a}_\ell^2 \sim D_Q T_Q$ which identifies the $2 - d$ range of the pairing attractive potential. In our model, its temperature dependence is $\tilde{a}_\ell^2 \sim T_{coh}/T$. This suggests a possible experimental check for the surmise that the superconducting instability is driven by the Q-modes in the CuO$_2$ planes. Two possibilities arise: a) multiple order parameters could provide different intervortex interactions for different magnetic field strengths in lowering the temperature. However this possibility requires a two-component Ginzburg-Landau formalism, even when only one divergent length scale is associated with the transition at $T_c$.$^{56,57}$ b) a second superconducting phase transition to Type I superconductivity takes place, a rather unlikely possibility. Discussion related to case a) arose in connection with superconductivity in the two band $MgB_2$. $^{58-60}$

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Appendices

Appendix A: Expansion of the action up to second order

We expand Eq.(7) of the Main Text (MT),

$$I_a = \sum_x \left[ -\ln \text{Det}(G_o^{-1} - \Sigma_x) + \frac{1}{2} \int d\theta_1 d\theta_2 \left( \Sigma_x(\theta_1, \theta_2) \Sigma_x^*(\theta_1, \theta_2) - \frac{(\beta \mathcal{J})^2}{4} (G_x(\theta_1, \theta_2))^4 \right) \right]$$

to second order in $\delta \Sigma_x$, $\delta G_x$, $\partial_\tau \varphi_x$:

$$G_x(\theta_1, \theta_2) = (G_c(\theta_1 - \theta_2) + \delta G(x, \theta_1 - \theta_2, \theta_+)) e^{i \varphi_x(\theta_+)}$$

$$\Sigma_x(\theta_1, \theta_2) = (\Sigma_c(\theta_1 - \theta_2) + \delta \Sigma(x, \theta_1 - \theta_2, \theta_+)) e^{i \varphi_x(\theta_+)}$$

where $\theta_+ = (\theta_1 + \theta_2)/2$. Gauge invariance is exploited, transforming $\Sigma_x(\tau_1, \tau_2)$ in such a way that the time derivative $\partial_\tau \varphi_x(\tau)$ appears in the Det, so that the variation of the Det term reads:

$$\frac{1}{2} \ln \text{Det} (G_o^{-1} - \Sigma_c) = \frac{1}{4} \text{Tr} \left( [G_o^{-1} - \Sigma_c]^{-1} [i \partial_\tau \varphi_x + \delta \Sigma] \right)^2 \rightarrow \frac{1}{2} \ln \det (G_o^{-1} - \Sigma_c) = \frac{1}{4} \text{Tr} \left[ 2 \Re \{ G_c i \partial_\tau \varphi_x G_c \delta \Sigma \} - (G_c \partial_\tau \varphi_x)^2 + (G_c \delta \Sigma)^2 \right]$$

Integrals in time are intended in the second term. Close to the conformal symmetry point (using the saddle point equality $\Sigma_c = \mathcal{J}^2 G_c^2$) the second term in the action of Eq.(A1) gives:

$$(\Sigma_c + \delta \Sigma)(G_c + \delta G) - \frac{\mathcal{J}^2}{4} (G_c + \delta G)^4 \sim \frac{3 \mathcal{J}^2}{4} G_c^4 + \delta \Sigma \delta G - \frac{3 \mathcal{J}^2}{2} G_c^2 \delta G^2.$$ (A3)

Only second order terms will be retained.

We introduce a dimensionless approach: $\theta = 2\pi \tau / \beta$ with the substitutions proposed by Kitaev\(^6\) and define $K_c(\theta_1, \theta_2, \theta_3, \theta_4)$:

$$g(\tau_1, \tau_2) = R_c(\tau_1, \tau_2) G(\tau_1, \tau_2),$$

$$f(\tau_1, \tau_2) = R_c^{-1}(\tau_1, \tau_2) \Sigma(\tau_1, \tau_2)$$

$$R_c(\theta_1, \theta_2) = \beta J \sqrt{(q - 1)} \left| \tilde{G}_c(\theta_1, \theta_2) \right|^{\frac{q-2}{2}},$$

$$K_c = R_c(\theta_1, \theta_2) \tilde{G}_c(\theta_1, \theta_3) \tilde{G}_c(\theta_4, \theta_2) R_c(\theta_3, \theta_4)$$ (A4)

For $q = 4$, the saddle point bilocal field is:

$$G_c(\tau) = b \left[ \frac{\pi}{\beta \sin \frac{\pi \tau}{\beta}} \right]^{1/2} \text{sign}(\tau), \quad b^2 = \frac{1}{2 J \pi^{1/2}}.$$ (A5)

Excluding for the time being the hopping term, the action $I_a$ of Eq.(A1), expanded up to second order, reads:

$$\frac{\tilde{I}_a}{N} = -\frac{1}{2} (\delta f | \delta g) - \frac{1}{4} \left[ (\delta f | K_c | \delta f) + 2 \Re \{ (\delta f | K_c | i R^{-1}_c \partial_\tau \varphi_x) \} + (R^{-1}_c \partial_\tau \varphi_x | K_c | R^{-1}_c \partial_\tau \varphi_x) \right] - \frac{1}{4} (\delta g | \delta g)$$ (A6)
We integrate out $\delta f' = \delta f + i \, R_c^{-1} \partial_{\tau} \varphi_x$:

\[
\sim e^{-\frac{N}{2} \sum_x \left[ (\delta f' [K_c] \delta f') + 2 (\delta f' [\delta g]) - 2 (i \, R_c^{-1} \partial_{\tau} \varphi_x [\delta g]) \right]} e^{\frac{N}{2} \partial_{\tau}^2} [\delta g [K_c^{-1} - 1] [\delta g]].
\]  
(A7)

The second variation of the hopping term (third term in Eq.(A1) ), to lowest order can be expanded as follows:

\[
\delta G_{c,x}(\tau_{12}, \tau_{+}) \delta G^*_{c,x}(\tau_{34}, \tau_{+}'') = [G_x(\tau_{12}, \tau_{+}) G_x^*(\tau_{34}, \tau_{+}'') - G_c(\tau_{12} - \tau_{+}) G_c(\tau_{34} - \tau_{+}'')] 
\approx \frac{1}{2} G_c(\tau_{12}) \left( e^{-i \alpha \tau_{x} \left[ \varphi_x(\tau_{+}) - \varphi_x(\tau_{+}'') \right]} - 1 \right) G_c(\tau_{34}) + c.c.
\]

\[
\rightarrow - \frac{1}{2} G_c(\tau_{12}) (\bar{\alpha} \cdot \nabla_{\tau} \left[ \varphi_x(\tau_{+}) - \varphi_x(\tau_{+}'') \right])^2 G_c(\tau_{34}),
\]  
(A8)

where only quadratic terms of the exponential have been included, to account for the additional complex conjugate contribution. We approximate $[\varphi_x(\tau_{+}) - \varphi_x(\tau_{+}'')] = (\tau_{+} - \tau_{+}'') \partial_{\tau} \varphi_x(\tau_{+})$ and define $R_c^{-1} \Lambda_c R_c^{-1} = \frac{1}{2} (\bar{\alpha} \cdot \nabla_{\tau}) G_c(\tau_{12})(\tau_{+} - \tau_{+}'')^2 G_c(\tau_{34})(\bar{\alpha} \cdot \nabla_{\tau'})$, so that, with relative time integrals traced out,

\[
\int d\tau \int d\tau' \partial_{\tau} \varphi_x(\tau) R_c^{-1} \Lambda_c(\tau - \tau') R_c^{-1} \partial_{\tau} \varphi_x(\tau').
\]  
(A9)

Fourier transform of the time dependences gives $\Lambda_c(\omega_{m}, \omega_{m'}; \Omega_{t})$. Fourier transforming the hopping term w.r.t $x$ we obtain the kinetic term added to the action of Eq.(A6) so that the full functional integral is:

\[
\sim \prod_p \left( \Pi_m \delta g_m^* \Pi_m \delta g_m \right) \right) e^{\frac{N}{2} \left( i \, R_c^{-1} \partial_{\tau} \varphi_x [\delta g] \right)} e^{\frac{N}{2} \left[ [\delta g [K_c^{-1} - 1] [\delta g]] \right]} e^{-\frac{\alpha^2}{2} \tau^2 \partial_{\tau}^2 \left[ \left( -i \, R_c^{-1} \partial_{\tau} \varphi_x [\Lambda_c] - i \, R_c^{-1} \partial_{\tau} \varphi_{x'} \right) \right]}
\]  
(A10)

where the forks (...) denote integration over $(\tau(12)_+ - \tau(34)_+)$.

In the conformal limit $i \, \omega_{m}$, can be dropped in $G_c^{-1} \ln$ Eq.(A1) and the functional integral of Eq.(A10) is just a gaussian form, so that, integrating out the $\delta g$’s we have:

\[
\sim \prod_p e^{-\frac{\alpha^2}{2} \left( -i \, R_c^{-1} \partial_{\tau} \varphi_{x}[K_c^{-1} - 1]^{-1} - i \, R_c^{-1} \partial_{\tau} \varphi_{x'} \right) e^{\frac{\alpha^2}{2} \frac{\alpha^2}{2} \tau^2 \left[ \left( -i \, R_c^{-1} \partial_{\tau} \varphi_x [\Lambda_c] - i \, R_c^{-1} \partial_{\tau} \varphi_{x'} \right) \right]}
\]

\[
\sim \prod_p e^{-\frac{\alpha^2}{2} \left( -i \, \partial_{\varphi_x} \left( 1 - \frac{\alpha^2}{2} \frac{\alpha^2}{2} \tau^2 \partial_{\tau}^2 R_c^{-1} \Lambda_c R_c^{-1} \varphi_x \right) \right) - \partial_{\varphi_{x'}} \left( 1 - \frac{\alpha^2}{2} \frac{\alpha^2}{2} \tau^2 \partial_{\tau}^2 R_c^{-1} \Lambda_c R_c^{-1} \varphi_{x'} \right) \right).}
\]  
(A11)

Here we have defined $F = R_c^{-1} K_c [K_c - 1]^{-1} R_c^{-1}$, and in the Kernel of Eq.(A11) we write:

\[
R_c^{-1} K_c [1 - K_c]^{-1} R_c^{-1} - \frac{t_2^0}{N} R_c^{-1} \Lambda_c [1 - K_c] K_c^{-1} K_c [1 - K_c]^{-1} R_c^{-1} = \left( 1 - \frac{t_2^0}{N} \frac{1}{2} R_c^{-1} \Lambda_c R_c^{-1} \right) F.
\]  
(A12)

The numerical evaluation of this kernel is discussed in B.

Appendix B: Numerical evaluation of the Kernel of Eq.(A12)

The correlator $F = R_c^{-1} K_c [K_c - 1]^{-1} R_c^{-1}$ has been defined after Eq.(A11). It diverges in the conformal limit because $[K_c - 1]$ has zero eigenvalues due to the spontaneous symmetry breaking. We will keep just the series of these eigenvalues (usually denoted by $h = 2$), which make the largest contribution to the correlator functions. The spectral representation in imaginary time of the regularized form of the kernel, obtained by shifting the zero eigenvalues,
\( k(h = 2, n) \approx 1 - \frac{\alpha_K}{\beta^2} |n| + \ldots (\alpha_K \approx 3) \), thus including the UV correction at \( O(N/\beta J) \) is:

\[
K_c[1 - K_c]^{-1} = \frac{1}{2} R_c(\tau_{12}) F(\tau_{12}, \tau_{34}) R_c(\tau_{34})
\]

\[
= \sum_{h, n} \Psi_{h, n}(\tau_{12}) \frac{k(h, n)}{1 - k(h, n)} \Psi^*_{h, n}(\tau_{34}).
\]

The basis functions for \( h = 2 \) are

\[
\Psi_{2, n}(x, y) = \gamma_n e^{-in\pi y/2} f_n(x),
\]

\[
f_n(x) = \sin \frac{nx}{\tan \frac{x}{2}} - n \cos \frac{nx}{2}, \quad \gamma_n = \frac{3}{\pi^2 |n(n^2 - 1)|},
\]

with \( x_{12} = \tau_1 - \tau_2 \), \( y_{12} = \frac{\tau_1 + \tau_2}{2} \). We Fourier transform the variables \( \tau_{12} \) and \( \tau_{34} \). In full generality:

\[
O_{m, m'} = \sum_{n \geq 2} e^{in(y_{12} - y_{34})} \Phi_n(\omega_m) (2, n | O | 2, n) \Phi^*_n(\omega_{m'}),
\]

where \( \omega_m, \omega_{m'} \) are the fermionic Matsubara frequencies. We redefine variables in such a way that \( \omega_m \to m \) with \( m \) integer. The basis functions are:

\[
\Phi_n(m) = \int_0^{2\pi} d\tau \frac{1}{2\pi} f_n(\tau) e^{im\tau}.
\]

It can be shown that the Fourier transforms \( \Phi_n(m) \) have a factor \( -\frac{1}{2} (1 + e^{2im\pi}) \) or \( -\frac{1}{2} (1 - e^{2im\pi}) \), depending on \( n \) being even or odd, respectively. It follows that odd \( m \) imply even \( n \) as expected because the \( \tau_n \) time dependence has to be with \( n \) even, i.e. bosonic-like. \( \Phi_n(m) \) have a maximum at increasing values of \( m \) when \( n \) increases and eventually go to zero.

The largest contribution \( O(\beta J/N) \) of \( [1 - K_c]^{-1} \) appearing in Eq.(A11), gives \( R_c^{-1} K_c [1 - K_c]^{-1} R_c^{-1} \sim O(1) \) and, as \( G_c G_c \sim 1/\beta J \), to have the same order of the kinetic term \( t_0 \sim O(\beta J N)^{1/2} \).

In the definition of the matrix function \( R_c^{-1} \Lambda_c R_c^{-1} \) from Eq.(11), the matrix \( G_c G_c \hat{F}^{-1} \) appears, which is the inverse of \( \frac{F}{G_c G_c} \).

The dominant expression for \( \frac{F}{G_c G_c} (\tau_1 \ldots \tau_4) \) in imaginary times, on the subspace orthogonal to the \( pGm \) fluctuations is:

\[
\frac{F}{G_c G_c} (\tau_1 \ldots \tau_4) = \frac{6\alpha_0}{\pi^2 \alpha K} \beta J \sum_{n \geq 2} e^{in(y_{12} - y_{34})} f_n(\tau_{12}) f_n(\tau_{34}).
\]

Its Fourier transform requires the transformed basis functions:

\[
\phi_n(m) = \int_0^{2\pi} d\tau \frac{1}{2\pi} f_n(\tau) e^{im\tau}.
\]

All of them have a factor sin \( m\pi \) which vanishes for \( m \) odd integer. However, this zero can be compensated by a zero in the denominator. Consider the case \( n = 2 \) for example:

\[
\phi(n = 2, m) = \frac{\sin m\pi}{\pi m(m - 1)(m + 1)} e^{im\pi},
\]

We give a finite expression to this vector element using the limit:

\[
\lim_{x \to \pm 1} \frac{\sin x\pi}{(x - 1)(x + 1)} = \frac{1}{2} (\mp \pi),
\]

However, all \( m > 1 \) give zero for \( n = 2 \), because there is no other factor of the kind \( (m - 3), (m - 5), \ldots \) (only odd \( m \))
are considered) in the denominator. We get \( \phi_n(m) = \text{sign}(m) \times \) :

\[
\begin{align*}
\text{for } n = 2 & \Rightarrow \\
& \begin{cases} m = 1 & \frac{1}{2m^2} e^{im\pi}, \\
m = 3, 5, 7, ... & 0 \\
\end{cases} \\
\text{for } n = 4 & \Rightarrow \\
& \begin{cases} m = 1 & -2 \frac{(5m^2 - 2)}{(m - 2)(m + 2)} \frac{1}{2m^2} e^{im\pi}, \\
m = 3, 5, 7, ... & 0 \\
\end{cases} \\
\text{for } n = 6 & \Rightarrow \\
& \begin{cases} m = 1 & -\frac{(36 - 119m^2 + 35m^4)}{(m - 3)(m - 2)(m + 2)(m + 3)} \frac{1}{2m^2} e^{im\pi}, \\
m = 3 & -\frac{1}{(m - 2)(m + 1)(m + 2)} \frac{1}{2m^2} e^{im\pi}, \\
m = 5, 7, 9, ... & 0 \\
\end{cases} \\
\text{for } n = 8 & \Rightarrow \\
& \begin{cases} m = 1 & -2 \frac{(2 - 288 + 1068m^2 - 462m^4 + 42m^6)}{(m - 4)(m - 3)(m - 2)(m + 2)(m + 3)(m + 4)} \frac{1}{2m^2} e^{im\pi}, \\
m = 3 & -\frac{2(2 - 288 + 1068m^2 - 462m^4 + 42m^6)}{(m - 4)(m - 2)(m + 1)(m + 2)(m + 3)(m + 4)} \frac{1}{2m^2} e^{im\pi}, \\
m = 5, 7, 9, ... & 0 \\
\end{cases} \\
\text{for } n = 10 & \Rightarrow \\
& \begin{cases} m = 1 & -3 \frac{(3 - 4800 + 18964m^2 - 9735m^4 + 1386m^6 + 55m^8)}{(m - 5)(m - 4)(m - 3)(m - 2)(m + 2)(m + 3)(m + 4)(m + 5)} \frac{1}{2m^2} e^{im\pi}, \\
m = 3 & -3 \frac{(3 - 4800 + 18964m^2 - 9735m^4 + 1386m^6 + 55m^8)}{(m - 4)(m - 3)(m - 2)(m + 1)(m + 2)(m + 3)(m + 4)(m + 5)} \frac{1}{2m^2} e^{im\pi}, \\
m = 5 & -3 \frac{(3 - 4800 + 18964m^2 - 9735m^4 + 1386m^6 + 55m^8)}{(m - 4)(m - 3)(m - 2)(m - 1)(m + 1)(m + 2)(m + 3)(m + 4)(m + 5)} \frac{1}{2m^2} e^{im\pi}, \\
m = 7, 9, 11, ... & 0 \\
\end{cases} \\
\text{for } n = 12 & \Rightarrow \\
& \begin{cases} m = 1 & -2 \frac{(2(-259200 + 1066164m^2 - 603746m^4 + 1059636m^6 - 6864m^8 + 143m^10)}{(m - 5)(m - 4)(m - 3)(m - 2)(m + 2)(m + 1)(m + 4)(m + 5)} \frac{1}{2m^2} e^{im\pi}, \\
m = 3 & -2 \frac{(2(-259200 + 1066164m^2 - 603746m^4 + 1059636m^6 - 6864m^8 + 143m^10)}{(m - 4)(m - 3)(m - 2)(m - 1)(m + 1)(m + 2)(m + 3)(m + 4)(m + 5)} \frac{1}{2m^2} e^{im\pi}, \\
m = 5 & -2 \frac{(2(-259200 + 1066164m^2 - 603746m^4 + 1059636m^6 - 6864m^8 + 143m^10)}{(m - 4)(m - 3)(m - 2)(m - 1)(m + 1)(m + 2)(m + 3)(m + 4)} \frac{1}{2m^2} e^{im\pi}, \\
m = 7, 9, 11, ... & 0 \\
\end{cases}
\end{align*}
\]

(B6)

Also the polynomials in the numerator could have been factorized but the roots are non-integer. We normalize each \( n = 10 \) \( \Rightarrow \) row, \( \text{for } n = 12 \) \( \Rightarrow \) column \( \times \) row each \( n \)-vector. To the elements with \( m, m' = 1 \), all vectors \( n = 2, 4, 6, 8, 10, 12 \) contribute. To the elements with \( m, m' = 3 \), vectors with \( n = 4, 6, 8, 10, 12 \) contribute. To the elements with \( m, m' = 5 \) only vector with \( n = 10, 12 \) contributes. The final result for \( n \leq 12 \) is a \( 3 \times 3 \) matrix. Each \( 3 \times 3 \) matrix \( W^n \) has eigenvalues \( 1, 0, 0 \). Computations leading to Fig.2,3,4 of the MT have been performed with \( 3 \times 3 \) matrices up to \( n = 12 \). The starting point is Eq.(20) of the text:

\[
\frac{F}{G_c} G_c (y_{12} - y_{34}) |_{m,m'} = \frac{4\alpha \omega}{\pi^2 \alpha K} \beta J \sum_{n \geq 2, even}^{10} \cos n(y_{12} - y_{34}) \frac{\cos n}{n^2(n^2 - 1)} W^n_{m,m'}
\]

where \( y \) are center-of-mass times: \( y = y_{12} - y_{34} \).

In Fig.4 of the MT we plot an interpolated smoothed curve of the (approximate) lowest \( \tilde{M} \) value which fulfills unitarity of the partition function of Eq.(46) of MT vs. \( T_0/T \) for \( \tilde{a} = 1 \). Precision is up to \( > 10^{-5} \). The trend is only meaningful for \( T_0/T \sim 1 \), because larger values require \( n > 12 \) in the spectral representation of Eq.(B3) and matrices \( m \times m \) of rank \( r > 3 \), i.e. higher than the ones used here.

Appendix C: Real time approach to the correlation functions

The response function in real time is derived from extension of the imaginary time along the Keldysh contour. The present approach\(^{13}\) is only semiclassical. In setting up the functional we disregard the \( q - q \) term and neglect variation with respect to the saddle point solutions for the \( G \) and \( \Sigma \) of the SYK dots. This means that, going back to the action in Eq.(A6), terms with \( \delta q \) and \( \delta f \) are disregarded, so that only an extra term is added here, taken from the imaginary time action, that is \(-N \left( R_c^{-1} \partial_{\tau} \varphi \right) K_c R_c^{-1} \partial_{\tau} \varphi \) = \(-N \left( \partial_{\tau} \varphi \right) G_c G_c \left( \partial_{\tau} \varphi \right)\), which we denote as the kinetic energy term.

As for the non local phase fluctuations, Fourier transforming with respect to time the non local fluctuation part of
the action (i.e. the "hopping" term of Eq.(A9)) is of a similar form:

\[
\int dp \bar{D} \frac{p^2}{\omega^2} \sum_{s,s'} \int d\omega \varphi_s(\omega) \overline{\varphi_{s'}} G_{s,s'}(\omega) \varphi_{s'}(-\omega).
\]  

(C1)

With respect to the Fourier transform of the kinetic energy term, the present one has a factor \(\omega^2\) lacking, so that we merge the two together, by defining a function

\[
\left[ \frac{\bar{D} p^2}{\omega^2} h(\omega) - 1 \right] G_c G_c \equiv \zeta G_c G_c.
\]  

(C2)

which defines the function \(\zeta(\omega)\). \(h(\omega)\) excludes the \(\omega = 0\) term. Its retarded form is defined as

\[
\bar{D} h_R(\omega) G^R K_{-\omega} = \bar{D} \int_{-\infty}^{\infty} dt G^R(t) G^K(-t) \left[ e^{i \omega t} - 1 \right] 
\]

\[
\sim i \bar{D} \omega \int_{-\infty}^{\infty} dt G^R(t) G^K(-t) \approx i \bar{D} \omega G^R_0 G^K_0.
\]

When writing \(h(\omega)\), we will not specify the label \(R/A\) for the retarded or advanced form, in the following, as long as no ambiguity arises. Transforming from the branches \(s, s' = +, -\) to the combined \(\alpha, \beta \equiv \epsilon l, q\)\(^{43,61}\), we get:

\[
\varphi^{\epsilon l/q}(t_j) = \frac{1}{\sqrt{2}} (\varphi_+(t_j) \pm \varphi_-(t_j)),
\]

\[
\frac{1}{2} \begin{pmatrix}
  1 & 1 \\
  1 & -1
\end{pmatrix} \begin{pmatrix}
  G_{++} & G_{+-} \\
  G_{+} & G_{--}
\end{pmatrix} \begin{pmatrix}
  1 & 1 \\
  1 & -1
\end{pmatrix} = \begin{pmatrix}
  0 & G^A_{-} \\
  G^A_{+} & G^A_{++}
\end{pmatrix}.
\]  

(C3)

The \(\epsilon l - \epsilon l\) component is zero in the matrices on the right. It reflects the fact that for a pure classical field configuration \((\varphi^q = 0)\), the action is zero. Indeed, in this case \(\varphi_+ = \varphi_-\) and the action on the forward part of the contour is canceled by that on the backward part (safe for the boundary terms, that may be omitted in the continuum limit), because the circuit is closed\(^{43}\).

The integrand of Eq.(C1) becomes\(^{13}\):

\[
(\varphi^\epsilon_\omega \ \varphi^q_\omega) \begin{pmatrix}
  G^A G^R_{\omega} \\
  G^A G^R_{\omega} + G^K_{-\omega} G^K_{-\omega}
\end{pmatrix} \begin{pmatrix}
  \varphi^{\epsilon \omega} \\
  \varphi^{q \omega}
\end{pmatrix}
\]

The resulting matrix can be rewritten as matrix of the self-energies \(\Sigma_D\) due to the \(\bar{D}\) coupling, which shows the same causality structure:

\[
\begin{pmatrix}
  G^A G^R_{\omega} \\
  G^A G^R_{\omega} + G^K_{-\omega} G^K_{-\omega}
\end{pmatrix} = \begin{pmatrix}
  G^A G^R_{\omega} \\
  G^K_{-\omega} G^K_{-\omega}
\end{pmatrix} \begin{pmatrix}
  \Sigma^A_D \ \\
  \Sigma^K_D
\end{pmatrix}.
\]

We neglect the \(qq\) term and write the functional:

\[
\int D\varphi^\epsilon_\omega D\varphi^q_\omega e^{-\frac{i}{\hbar} \int d\omega (\omega \varphi^\epsilon_\omega \ \omega \varphi^q_\omega) \begin{pmatrix}
  G^A G^R_{\omega} \\
  G^K_{-\omega} G^K_{-\omega}
\end{pmatrix} \begin{pmatrix}
  \varphi^{\epsilon \omega} \\
  \varphi^{q \omega}
\end{pmatrix}},
\]

(C4)

where we have also added the kinetic energy term of the semiclassical approach.

The field \(\hat{N}(x,t)\) is the source of \(\partial_t \varphi(x,t)\). We want the response written along the Keldysh contour:

\[
D_{KN}(x,t) = i \theta(t) \{ \hat{N}(x,t), \hat{N}(0,0) \}
\]

\[
= i \{ \hat{N}(x,t) \hat{N}(0,0) + \hat{N}(0,0) \hat{N}(x,t) \}.
\]  

(C5)

To get the generating functional of the \(\varphi - \varphi\) fluctuations we invert the kernel of Eq.(C4), obtaining the matrix:

\[
\begin{pmatrix}
  0 & (\zeta^R G^K_{-\omega} G^R_{\omega})^{-1} \\
  (\zeta^A G^A_{\omega} G^K_{-\omega})^{-1} & (G^R_{\omega} G^K_{-\omega})^{-1}
\end{pmatrix}
\]
The $[G^{-1}]^K$ component for the free field is only a regularization factor, originating from the (time) boundary terms. It is, in general, non-local in $x$ and $x'$, however, being a pure boundary term, it is frequently omitted\textsuperscript{13}. In our case this should apply because $[G^{-1}]^K(t,t') = [G^{-1}]^K \circ F - F \circ [G^{-1}]^A = [G^R]^{-1} \circ F - F \circ [G^A]^{-1} = [G^K]^{-1}$. Integrating out the $\varphi$ fields and ignoring again the $q-q$ term, we get:

$$\propto \exp \left\{ -\frac{1}{2} \int d\omega \left( \hat{N}^\text{c}(\omega) \: \hat{N}^\text{q}(\omega) \right) \left( \begin{array}{cc} 0 & (\zeta^R G^K_{\omega} G^R_{\omega})^{-1} \\ (\zeta^A G^K_{\omega} G^A_{\omega})^{-1} & 0 \end{array} \right) \left( \begin{array}{c} \hat{N}^\text{c}(\omega) \\ \hat{N}^\text{q}(\omega) \end{array} \right) \right\}. \quad (C6)$$

Functional derivation with respect to the sources provides the cross contributions (we keep just the lowest order in $\omega$). Using the definition of Eq.\textsuperscript{(C2)}:

$$N \omega^2 \langle \varphi^c(\omega) \: \varphi^q(-\omega) \rangle \approx (G^K_0 G^R_0)^{-1} \frac{\omega^2}{i D p^2 \omega - \omega^2}$$

$$N \omega^2 \langle \varphi^q(\omega) \: \varphi^c(-\omega) \rangle \approx (G^K_0 G^A_0)^{-1} \frac{\omega^2}{-i D p^2 \omega - \omega^2}.$$ 

Now, the retarded energy flux density response of Eq.\textsuperscript{(C5)} can be estimated, considering that $\delta \varphi^c \text{ and } \hat{N}^c/q$ are conjugate variables $\hat{N}^c/q(t) = \frac{\delta \varphi^c}{\delta q}$, so that, keeping just the $\omega^2$-term in Eq.\textsuperscript{(C4)},

$$\hat{N}^c(\omega) = -N G^K_0 G^R_0 \omega \varphi_q(-\omega),$$

$$\hat{N}^q(\omega) = -N G^K_0 G^A_0 \omega \varphi_c(-\omega)$$

we get

$$\langle \hat{N}^c(\omega) \: \hat{N}^c(-\omega) \rangle = -N^2 G^K_0 G^R_0 G^K_0 G^R_0 \omega^2 \times \langle \varphi^c(-\omega) \: \varphi^c(\omega) \rangle$$

$$= -N^2 \frac{\omega^2}{i D p^2 \omega - \omega^2} G^K_0 G^R_0,$$

$$\langle \hat{N}^q(\omega) \: \hat{N}^c(-\omega) \rangle = -N^2 G^K_0 G^R_0 G^K_0 G^A_0 \omega^2 \times \langle \varphi^c(-\omega) \: \varphi^q(\omega) \rangle G^K_0 G^A_0$$

$$= -N^2 \frac{\omega^2}{-i D p^2 \omega - \omega^2} G^K_0 G^R_0.$$ 

The symmetrized correlation is:

$$\frac{1}{2} \left( \langle \hat{N}^c(\omega) \: \hat{N}^c(-\omega) \rangle + \langle \hat{N}^q(\omega) \: \hat{N}^c(-\omega) \rangle \right) = N^2 \text{Re} \left\{ \frac{\omega^2}{i D p^2 \omega + \omega^2} G^K_0 G^R_0 \right\}$$

$$\approx N^2 \frac{\left( D p^2 \right)^2}{(D p)^2 + \omega^2} \Im m\left[G^K_0 G^R_0\right]. \quad (C7)$$

This result can be rewritten as

$$\text{sign}(\omega) \Im m \left\{ \frac{\omega}{i D p^2} - \omega \right\} \Im m\left[G^K_0 G^R_0\right].$$

Subtracting the $p = 0$ term, we recognize $\Im m\left(D^R(p,\omega)\right)$, the imaginary part of the density response function\textsuperscript{62}

$$\Im m\left\{ \frac{D^R_{NN}(p,\omega)}{D^R_{NN}(p,\omega)} \right\} = -\text{sign}(\omega) \Im m \left\{ \frac{D p^2}{i \omega - D p^2} \right\} \Im m\left[G^K_0 G^R_0\right]. \quad (C8)$$

This result should be compared with Eq.\textsuperscript{(51)} of the MT. Apart for the matrix structure of the function in Eq.\textsuperscript{(47)} of the MT, the important point is that $F^{-1}$ is absent here in the definition of the diffusion parameter.

We add here the important consequence on the electrical conductivity. In the conformal limit, the electrical
conductivity is\textsuperscript{13}:

\begin{equation}
\Re \{ \sigma \} = -\lim_{\omega \to 0} \frac{1}{\omega} \Im \{ D^{R3} (\omega) \} \propto -\frac{\beta}{2\pi^2} \frac{t_0^2}{N} \int d\omega \sech^2 \left( \frac{\beta \omega}{2} \right) \left( \Im \{ G^{R3} (\omega) \} \right)^2
\end{equation}

Resistivity is \( \propto T \) in this approach.

Appendix D: Quantization of gapless diffusive excitation mode

\( J_Q = -\kappa \nabla T \) is a classical diffusion equation of a non conserving system. We now construct a Hamiltonian of the excitation modes which is conserving but we ask that, introducing a relaxation time \( \tau_0 = \hbar/\Gamma \) for these modes, the equation of motion reproduces \( J_Q = -\kappa \nabla T \). We will quantize this Hamiltonian and derive the response function from the fluctuations of these modes. The canonical conjugate variables and the corresponding Lagrangian (in 2-d) are:

\begin{equation}
\dot{\theta} = \left( \frac{\kappa C}{\hbar T} \right)^{1/2} J_Q \tau_0, \quad \nabla \theta = \left( \frac{\hbar}{\kappa C T} \right)^{1/2} \kappa \nabla T
\end{equation}

Here \( C \) is the thermal capacitance, \( \kappa \) is the thermal conductance and \( J_Q \) is the thermal energy current. The corresponding Lagrangian is

\begin{equation}
\mathcal{L} = \frac{1}{2} \int d^2 x \left[ \frac{k_B}{\kappa C} \left( \frac{J_Q \tau_0}{k_B} \right)^2 + \frac{\hbar}{\kappa C} \left( \frac{\kappa}{T} \nabla T \right)^2 \right] \equiv \frac{1}{2} \int d^2 x \left[ A \dot{\theta}^2 + B (\nabla \theta)^2 \right]
\end{equation}

with \( A = \frac{\hbar k_B}{\kappa C} \) and \( B = T \). These choices provide terms in the square brackets which have dimension \( \mathcal{E}/\ell^2 \) (\( \mathcal{E} \equiv \text{energy} \)).

With the approximation \( \tau_0 \dot{J}_Q \approx J_Q \), the equation of motion,

\begin{equation}
\frac{d}{dt} \left( \frac{\partial L}{\partial \dot{\theta}} \right) - \frac{\partial L}{\partial \theta} = A \ddot{\theta} + B \nabla \theta = 0,
\end{equation}

boils down to the diffusion equation:

\begin{equation}
\frac{\hbar}{\kappa C} J_Q = -\frac{\hbar}{\kappa C} \kappa \nabla T.
\end{equation}

Although \( \hbar \) is already in the Lagrangian, we proceed with quantization of the theory\textsuperscript{62}. Fourier transforming, the canonical momentum for \( \theta_k \) is

\begin{equation}
\pi_k = \frac{\partial \mathcal{L}}{\partial \dot{\theta}_k} = A \dot{\theta}_{-k}, \quad [\theta_k, \pi_{k'}] = i \delta_{kk'}.
\end{equation}

The Hamiltonian

\begin{equation}
\mathcal{H} = \frac{1}{2} \sum_k \left[ \frac{1}{A} \pi_{-k} \pi_k + B k^2 \theta_{-k} \theta_k \right].
\end{equation}
is second quantized according to \( \hat{a}_k = u_k \theta_k + i v_k \pi_{-k} \), with

\[
u_k = \frac{1}{\sqrt{2}} (AB)^{1/4}, \quad v_k = \frac{1}{\sqrt{2}} (AB)^{-1/4},
\]

\[
\epsilon_k = \sqrt{\frac{Bk^2}{A}} = \left[ \frac{\kappa}{k_B} \right]^{1/2} \left( k \right) \| v \| v, \quad |k| \equiv v |k|,
\]

\[
\pi_k = -i T^{1/2} (2\epsilon_k)^{1/2} \left| k \right| \left( \hat{a}_k - \hat{a}^\dagger_{-k} \right),
\]

\[
\theta_k = T^{-1/2} (2\epsilon_k)^{1/2} \left| k \right| \left( \hat{a}_k + \hat{a}^\dagger_{-k} \right),
\]

\[
\mathcal{H} = \sum_k \epsilon_k \hat{a}_k^\dagger \hat{a}_k + \text{cnst}
\]

In Eq.(D5) we have defined the velocity \( v \) of these modes. The approach is similar to the one for phonons. \( \pi(x) \) plays the role of the space displacement \( d(x,t) \), while \( \nabla \theta \) plays the role of the phonon impulse \( \Pi(x,t) \). The thermal conductance used in the text is given by

\[
\Re \{ \kappa(\omega) \} = -\frac{1}{\omega} \Im \left\{ \frac{1}{2} \left\{ \left\{ J_{-k}^Q(-\omega), J_k^Q(\omega) \right\} \right\} \right\}_\omega \right\},
\]

where

\[
\frac{1}{2} \left\{ \left\{ J_{-k}^Q, J_k^Q \right\} \right\}_\omega = \frac{1}{\tau_0} \kappa C T \left\{ \left\{ \pi_{-k}, \pi_k \right\} \right\}_\omega.
\]

The symmetrized correlation on the right hand side, \( D^3(\omega) = \left\{ \left\{ \pi_{-k}, \pi_k \right\} \right\}_\omega \), apart of the prefactor \( T \), can be evaluated at zero temperature in a standard way. Eq.(D8) gives:

\[
\Re \{ \kappa(\omega) \} = \pi^2 k_B^2 \omega T_0 \frac{\omega}{v^2 T_0}.
\]

If \( \Re \{ \kappa(\omega) \} \sim \kappa \tau_0 \) we get, from Eq.(D5),

\[
v^2 \tau_0^2 = \frac{k \tau_0}{k_B} \frac{CT}{\hbar} \Rightarrow \kappa = \frac{\pi k_B}{\tau_0} \left( \frac{\hbar}{CT} \right)^{1/2}.
\]

However, introducing the damping of the mode in \( D^3 \), by adding an energy broadening \( \Gamma \), we get

\[
D^3(\omega) = \pi T \int_{-\infty}^{\infty} \frac{2 e^{-\epsilon^0} e}{\epsilon^2 - (\omega + i \Gamma)^2} \epsilon^2 d\epsilon
\]

and

\[
\Im \left\{ D^3(\omega) \right\} = -\frac{2\pi T \omega^2}{v^2 \Gamma} \left\{ \left( 1 - \frac{\Gamma^2}{\omega^2} \right) \text{arctan} \left( \frac{\Gamma}{\omega} \right) + \frac{\Gamma}{\omega} \ln \left( 1 + \frac{\Gamma^2}{\omega^2} \right) \right\},
\]

which, in the limit \( \Gamma/\omega \ll 1 \) gives:

\[
\Im \left\{ D^3(\omega) \right\} = -\frac{2\pi T \omega}{v^2}.
\]

Posing again \( \Re \{ \kappa(\omega) \} \sim \kappa \tau_0 \), in place of Eq.(D11) we have:

\[
v^2 \tau_0^2 = \frac{k \tau_0}{k_B} \frac{CT}{\hbar} \Rightarrow \kappa = k_B \frac{\Gamma}{CT} \left( \frac{\hbar}{CT} \right)^{1/2}.
\]

Using the \( \kappa = C v_F \ell = C v_F^2 \tau_0 \), the gapless bosonic excitations of energy \( \hbar \nu k \) generate a specific heat at fixed \( 2 - d  \)
volume:

\[ \frac{C_V}{a^2} = \frac{d}{dT} \frac{1}{2\pi} \left( \frac{k_B T}{h \nu} \right)^3 \hbar \int_0^\infty \frac{z^2 \, dz}{e^z - 1} = k_B \frac{1}{2\pi} \left( \frac{k_B T}{h \nu} \right)^2 \frac{6 \zeta[3]}{\pi} \]

and the thermal conductivity

\[ \kappa = k_B \frac{1}{\pi} \left( \frac{k_B T}{h \nu} \right)^2 3 \zeta[3] v \ell. \]

Here \( \zeta[n] \) is the Riemann function\(^6^3\). This is the Stefan-Boltzmann relation in two dimensions\(^6^4\).

In the case of the SYK model, based on the saddle point contribution to energy\(^8\):

\[ \ln Z = -\beta E_0 + S_0 + \frac{c}{2\beta} + ..., \quad \text{with} \quad \frac{c}{2} = \frac{2\pi^2 \alpha S N}{\mathcal{J}}, \]

the first energy correction in temperature is \( E = c/(2\beta^2) \) \( (c \approx 0.396 N/\mathcal{J}) \), so that, by taking \( CT \sim c/(2\beta^2) \) in Eq.(D14), we get:

\[ \kappa = k_B \Gamma \left( \frac{h \Gamma}{e^T} \right)^{1/2} \to k_B \Gamma^{3/2} (h \beta)^{1/2} \left( \frac{\beta \mathcal{J}}{\pi \alpha S N} \right)^{1/2} \]

\[ = k_B \Gamma^{3/2} (h \beta)^{1/2} \left( \frac{\beta \mathcal{J} N}{\pi \alpha S} \right)^{1/2} \]

\[ \quad \text{(dimensions are } [C/t] \text{ as always).} \]

The thermal conduction response in the conformal limit\(^13\) requires the energy current response function \( G^{R\beta}_{cQ}(\omega) \)

\[ \frac{\Re \{ \kappa \} }{NT} = - \lim_{\omega \to 0} \frac{1}{\omega NT} \Im \{ D^{R\beta}_{cQ}(\omega) \} \]

\[ = \frac{1}{NT \pi^2} \int d\omega \tanh \frac{\beta \omega}{2} \frac{\partial}{\partial \omega} \left( \omega \Im \left\{ G^{R\beta}_{cQ}(\omega) \right\} \right)^2 \propto \left( \frac{\beta \mathcal{J} N}{\pi \alpha S} \right). \]

In the Fermi liquid case, \( \tau_0 \sim T^{-2} \) and \( C \sim T \), so that \( \kappa \sim T^{-1} \).

**Appendix E: The acoustic plasma mode in the marginal Fermi Liquid**

To characterize the MFL phase, it is important to check the nature of the collective excitations, in particular the particle-hole continuum, under the action of the increasing coupling to the high energy localized modes. We will show that, within our approximations, the real part of the \( \omega(q) \) dispersion of the density excitations is linear, but with a small reduction of the physical velocity \( d\omega/dq \) at small \( q \), and, most of all, a peculiar imaginary part. We also find that, at large couplings, the interaction pulls a linearly dispersed, well defined acoustic plasmon mode out of the particle-hole continuum.

When the residual interaction is turned on, the vertex function \( \Gamma(p, p - q; q, i \Omega) \) satisfies the Bethe-Salpeter equation\(^6^5\),

\[ \Gamma(p, p + q; q, i \Omega) = n_{-q} + ig U \frac{1}{\mathcal{N}} \sum_{p'} D_{p', q}(\omega) \Gamma(p', p + q; \omega). \]

\[ D_{p, q}(i \Omega) = \sum_{\omega_n} G(\epsilon_{p - q}, i \omega_n) G(\epsilon_p, i \omega_n + i \Omega_m) \]

\[ G(\epsilon_p, i \Omega) = \frac{1}{iZ^{-1}(\omega + \Omega) - \epsilon_p}. \]

The functions \( D_{p', q}(\omega) \) are related to the polarization functions of Eq.(65,66) of the MT, when frequency is continued
to real values and $p' \sim p_F$. We define

$$\tilde{\Pi}^{1,2}(q, \omega) = -i \nu_0 D_q^{1,2}(\omega) \Gamma_{1,2}(p_F, p_F - q; q, \omega) \quad (E2)$$

where, in place of the $\sum_{p'}$ appearing in Eq.(E1) we multiply by $\nu_0$ after having put $|p'| = p_F$. The resulting functions $D_q^{1,2}$ of Eq.(E2) are redefined as ($\frac{\omega}{Z\nu_F q} < 1$)

$$\nu_0 D_q^1(\omega) = \frac{Z\nu_0}{2\pi} \left[ 1 - i \frac{\omega}{Z\nu_F q} \sqrt{1 - \frac{\omega^2}{(Z\nu_F q)^2}} \right],$$

$$\nu_0 D_q^2(\omega) = \nu_0 \int \frac{d\theta}{2\pi} (1 - \cos \theta) D_{k,q}^2(\omega) \bigg|_{k=p_F}. \quad (E3)$$

In fact, following [18], we consider two ranges of energy values: a low energy one ($\omega < \Omega^*_c$), $(i = 1)$, and an high energy one ($\omega < \Omega^*_c$), $(i = 2)$, with $\nu_0 D_{k,q}^i(q, \omega) = \Pi^i(q, \omega)$.

Limiting ourselves to the FL energy range, $i = 1$, for the moment, Eq.(E2), becomes

$$i \tilde{\Pi}^1(q, \omega) = -i \nu_0 D_{p_F,q}^1(\omega) \Gamma(p_F, p_F - q; q, \omega), \quad (E4)$$

where

$$\nu_0 A = \frac{1}{(2\pi)^2} \frac{2\pi k_F A}{\hbar v_F^c} = \frac{1}{\pi} \frac{k_F A}{\hbar v_F^c}.$$ 

$\nu_0$ is the 2-d density of states at $\epsilon_F$ per unit volume $A$.

Here we are assuming that, in this energy range, $\Gamma(p, p - q; q, i \Omega)$ does not depend on the angle $\theta_{p',q} = \theta_{p'q}$ except for an average of $\sin^2 \theta/2 \sim (q/(2k_F))^2 \sim 1/2$. We have also put $|p|, |p'| = k_F$, so that the only $\Gamma$ dependence is $\Gamma(q, i \Omega)$. This choice, together with that of the onsite interaction $U_c$, provides the reference result we are looking for.

We expect a collective mode of compressibility type embedded in the particle-hole excitation continuum. In the low temperature Fermi liquid limit, the p-h continuum has a boundary of the kind $\min \{ \epsilon_{k+q} - \epsilon_k \} = Z\nu_F q$. We find a collective mode $\omega = Z z\nu_F q$ with $z$ complex and $\Re \{ z \} < 1$ and negative imaginary part which is related to the lifetime of the mode.

Here

$$\Pi^1(q, \omega) = \frac{Z \nu_0}{2\pi} \left[ 1 - i \frac{\omega}{Z\nu_F q} \sqrt{1 - \frac{\omega^2}{(Z\nu_F q)^2}} \right].$$

So that

$$\left\{ \left[ \nu_0 D_{p_F,q}^1(\omega) \right]^{-1} - i g U_c \right\} \left[ i \tilde{\Pi}(q, \omega) \right] = -i n_{-\bar{q}} \quad (E5)$$

This provides the equation for the FL collective excitation mode ($Z \nu_0 = U_c^{-1}$):

$$\left[ 1 - i \frac{z}{\sqrt{1-z^2}} \right]^{-1} - i g = 0, \quad z = \frac{\omega}{Z\nu_F q} < 1$$

$$\Rightarrow 1 - i g = g \frac{z}{\sqrt{1-z^2}}. \quad (E6)$$

The homogeneous equation can be cast in the form: $1/\chi^{(0)} - U = 0$ if

$$i \nu_0 D_{k,q}^1(q, \omega) = i \Pi^{1(0)}(q, \omega) = -i \chi^{(0)} \quad (E7)$$

The contribution to the polarization function from high-energy excitations $\Pi^2(q, i \Omega)$ has a completely local q-
calculating the dc conductivity in metals

In the case of $\nu_0 D^2_{\nu_0, q}$, it was enough to put $p' \to p_F$, but, in the case of $\hat{\Pi}^2_0(\omega)$, it is important to keep a dependence on the scattering angle $\theta_{k, q} = \vec{k}, \vec{q}$ explicitly, and to integrate over it. In fact, we cannot neglect the dependence of the vertex $\Gamma$ appearing in the Bethe-Salpeter Eq.(E1) on the scattering angle $\theta_{k, q} = \vec{k}, \vec{q}$. As it is usual when calculating the dc conductivity in metals, we assume that this angular dependence is 

$$\sim (1 - \cos \theta)$$

in Eq.(E3). The factor $\Gamma$ expresses the growing predominance of forward scattering with declining temperature, which contributes less than wide angle scattering, to the effective "collision rates". A $q$-dependence has been added in $D^2_{k, q}(\omega)$ of Eq.(E3) by including a $k$-dependent correction to the conformal local SYK Green function $G_c$:

Here we use $G_c(i \omega_n)$, extended to include a $k$ dependence

$$G^2_c(\vec{k}, i \omega_n) \approx -\frac{1}{\Sigma^0_c(i \omega_n)} + \frac{\epsilon_{k}}{|\Sigma^0_c(i \omega_n)|^2}.$$  \hspace{1cm} (E8)

Here $\Sigma^0_c(i \omega_n)$ is the self-energy corresponding to the zero order approximation. Eq.(E8) can be viewed as an expansion of the high energy total Green function to lowest order in $\epsilon_k$, or can be obtained, by assuming random hopping between sites, so that, according to Eq.(E2), the propagator is:

$$\nu_0 D^2_{\nu_0}(i \Omega_n) = \left( \frac{1}{\nu} \sum_k \right) \sum_{\omega_n} \frac{1}{\Sigma^0_c(i \omega)} \frac{1}{\Sigma^0_c(i \omega + i\nu)} \left[ 1 + \left( \frac{\epsilon_{k-q}}{\Sigma^0_c(i \omega)} + \frac{\epsilon_{k}}{\Sigma^0_c(i \omega + i\nu)} \right) + \frac{\epsilon_{k-q}}{\Sigma^0_c(i \omega)} \frac{\epsilon_{k}}{\Sigma^0_c(i \omega + i\nu)} \right] \left( \frac{i \nu}{\nu} \right)^{\nu}.$$  \hspace{1cm} (E9)

The $\omega_n$ sum can be transformed into an integral. A lengthy but straightforward calculation provides a function of $k,q, \vec{k} \cdot \vec{q} = k \cos \theta$ and $i \Omega$:

$$f(k, q, \cos \theta; i\Omega) = -\frac{8}{U_c} \ln \frac{U_c}{W} + \frac{\epsilon_{k} \epsilon_{k-q}}{U_c^2} \frac{2}{\Omega} \arctanh \frac{\Omega}{\Omega_c} + \frac{2i}{U_c} \left\{ \frac{(\epsilon_{k-q} - \epsilon_{k})}{W} + \frac{(\epsilon_{k-q} + \epsilon_{k})}{W} \right\} \left( \frac{\nu}{\nu_c} \arctanh \frac{\nu}{\nu_c} \right).$$

We neglect the second and the fourth term (for a p-h pair in a p-h symmetric system is $\epsilon_{k-q} + \epsilon_{k} = 0$), obtaining

$$f(k, q, \cos \theta; i\Omega) \approx -\frac{8}{U_c} \ln \frac{U_c}{W} + \frac{2i}{U_c} \left\{ \frac{(\epsilon_{k-q} - \epsilon_{k})}{W} \right\}.$$

We finally get:

$$\nu_0 D^2_{\nu_0}(\omega) = -\nu_0 W \frac{8}{U_c} \ln \frac{U_c}{W} \left[ 1 + i \frac{1}{8 \ln \frac{U_c}{W}} \frac{\bar{\nu}_F q}{W} \right].$$  \hspace{1cm} (E10)

We now rewrite Eq.(E1) for the vertex as follows ( $i = 1, 2$):

$$\Gamma(p_F, p_F - q; \omega) = n_{-q} + U_c \left[ g_{11} \nu_0 D^1_{\nu_0}(\omega) + g_{12} \nu_0 D^2_{\nu_0}(\omega) \right] \Gamma(p_F, p_F - q; \omega).$$

By posing

$$\tilde{\Pi}^{1,2}(q, \omega) = -i \nu_0 D^1_{\nu_0}(\omega) \Gamma(p_F, p_F - q; \omega),$$

we get a system, which, using $Z \nu_0 U_c = 1$ and $z = \omega/(Z \bar{\nu}_F q)$, takes the form:
The results are given in Fig.1,2 for real and imaginary part of \( \omega / (Z \tilde{\nu} F q) \) vs \( q = U / U_c \).

When \( q \) increases there is a monotonic flattening of \( \Im (\omega_q) \) with \( Z \tilde{\nu} F q / W \) with a saturation at large \( q \), as can be seen by plotting \( d \omega_q / dq \) vs. \( q \). As \( \Re \omega_q \) is strictly linear with \( q \) in a large range of values of \( q \), the plot of Fig.(10) shows the behavior of this derivative.

The polarization function of the coupled system, given by Eq.(E2), satisfies the equation:

\[
\left( \frac{1 - i \sqrt{1 - v_F^2 g_{11}}}{i g_{12}} - i g_{11} \right) - \frac{Z}{\nu \ln \nu} \left( \frac{1 + i \frac{1}{8 \ln \nu} \tilde{\nu} q / W}{i g_{22}} \right) = -i n_{-q} \left( \frac{1}{1} \right)
\]  

(E12)

Given a transferred momentum \( q \), the energy of the corresponding collective excitation makes the determinant of the matrix on the left hand side of Eq.(E13) vanish. Here, \( g_{11} U_c \) is assumed to be the residual interaction within the low energy FL due to the SYK cluster, \( g_{22} U_c \sim J \) parametrizes the interaction within the SYK cluster, while \( g_{12} U_c \) provides the coupling between the two.

To zero order in perturbation, \( \chi^{(0)} = \nu_0 D^1_q (\omega) \), in the limit \( g \to 0 \) the solution is \( z \to 1 \), giving a strictly linear dispersion, \( \omega \propto v_F^* q \).

When the couplings are non vanishing, the mode dispersion keeps being substantially linear, but the physical velocity is renormalized and an imaginary part arises. The real and imaginary part of the function \( z = \omega (v_F^* q) \) are plotted in Fig.s(7,8) as a function of \( g_{11} \) and \( g_{22} \), for \( v_F^* q / W = 0.1 \). Plots are reported for \( U_c / W = 10 \) and \( Z = 0.1 \).

The limitation \( z = \frac{v_F^* q}{v_F^*} < 1 \) implies that we track the collective excitation mode inside the p-h continuum only. In Fig.(8) we plot the real part \( \Re \epsilon \{ z \} \) vs \( g_{11} \), choosing \( g_{22} = 2 \), when the ratio of the couplings between the two systems \( b = g_{12} / g_{11} \) is \( b = 0, 0.4, 0.6, 0.8 \).

The effective velocity of the excitation mode, \( \Re \epsilon \{ z \} \), decreases with respect to the unperturbed value \( v_F^* \) and saturates to about 90% of the unperturbed value when \( g_{12} \) increases. As shown in the inset, the saturation is even faster, when \( g_{11} \) is kept fixed (in our case at the value \( g_{11} = 1 \)) and \( g_{22} > g_{11} \) is increased.

The imaginary part of the energy of the mode \( \Im (\epsilon_q) \) is zero at \( g_{11} = 0 \) and increases mildly, in absolute value, with increasing of \( g_{11} \), as reported in Fig.(8) vs \( g_{11} \) at \( g_{22} = 1.8 \). It is remarkable that, when \( g_{11} > g_{22} \), \( \Im (\epsilon_q) \) vanishes. Simultaneously the slope of the mode increases up to value one, for \( b \to 1 \). This appears in Fig.(8) and, more explicitly, in Fig.(9) which is a plot vs \( b \) with \( g_{11} = g_{22} \) for various values \( g_{11} = 0.8, 1.2, 1.8 \) and \( v_F^* q / W = 0.8 \). When \( g_{12} \) increases in Eq.(E13), a real term \(- \tilde{\nu} F_{12}^* \) grows in the determinant, which reduces \( \Im (\epsilon_q) \). In these conditions, the dispersion tends to the boundary of the particle-hole continuum (\( \Re \epsilon \{ z \} \to 1 \)), while the corresponding imaginary part in Fig.8 vanishes. This feature appears clearly in Fig.s 9 a,b and is the signature of the splitting of a bound state out of the particle-hole continuum with linear dispersion and velocity \( > v_F^* \). We interpret this as an acoustic plasmon which, however, requires strong coupling of the MFL to the SYK cluster, to be tackled even further, far beyond the present perturbative approach. When \( q \) increases, there is a monotonic increase of \( \Im \epsilon_q \) with \( v_F^* q / W \) and a saturation at large \( q \), as can be seen by plotting \( d \omega_q / dq \) vs. \( q \) (see Fig. 10).

To sum up the results of this Appendix, we can conclude that the low energy FL on the lattice appears quite robust with respect to interaction with incoherent local disordered SYK clusters, when only the lowest perturbative order is included and no disorder, in the continuum, \( k \to 0 \), limit. The Fermi surface is still well defined, but the liquid becomes a MFL. The hydrodynamic collective excitation, the would-be acoustic plasmon, is also rather well defined. At strong coupling, in the limit \( U_c \to J \), its dispersion tends to the boundary of the p-h continuum and the imaginary part, which blurs the mode, vanishes. We expect that a well defined acoustic plasmon is on the verge to emerge as a bound state at low energies, splitted off the p-h continuum.

**Appendix F: The superconducting critical temperature at low temperature**

In this Appendix we use an Eliasberg\textsuperscript{45} approach to the superconducting critical temperature, \( T_c \), assuming that pairing is driven by the diffusive excitation modes introduced in Section D. As explained in the main text, \( \Omega^* \) is the energy scale of \( T_c \) and the dependence on the coupling strength turns out to be non BCS-like. We report here the derivation.
FIG. 7: Renormalization of the velocity of the linear dispersion $v_F$ in presence of coupling to high energy correlations. Here we plot the real part of $\omega/(v_F^2 q)$ versus $g_{11}$ at $g_{22} = 1.8$ and $v_F^2 q/W = 0.1$. Here $g_{12} = b g_{11}$ is the coupling strength to the high energy SYK fluctuations ($b = 0, 0.4, 0.6, 0.8$, respectively, with $U_c/W = 10$ and $Z = 0.1$).

FIG. 8: Imaginary part of $\omega/(v_F^2 q)$ versus $g_{11}$ corresponding to the real part of Fig.7 ($g_{12} = b g_{11}$, with $b = 0, 0.6, 0.8$, respectively), at $g_{22} = 1.8$ and $v_F^2 q/W = 0.1$. Note the flattening at zero for $g_{11} > 3.0$, when $b \to 1$.

FIG. 9: Real [a]) and Imaginary [b]) part of $\omega/(v_F^2 q)$ versus $b = g_{12}/g_{11}$ for various values of $g_{ii} \equiv g_{11} = g_{22}$, with $g_{ii} = 0.8, 1.2, 1.8$, respectively, at $v_F^2 q/W = 0.8$. 
FIG. 10: Plot of $\frac{\partial \text{Im} \{\omega\}}{\partial \text{Re} \{\omega\}}$ vs $v_F^* q / W$. As $\text{Re} \{\omega\}$ is strictly linear with $q$ in a large range of values of $q$, the derivative with respect to $\text{Re} \{\omega\}$ can be interpreted as derivative with respect to $q$. Here $g_{11} = g_{22} = 1.8$ and $g_{12} = 0.8 \times g_{11}$ ($Z = 0.1, U_c / W = 10$, as always).

In the mean field Hamiltonian, in the Nambu$^{67}$ representation,

$$H(k) = \sum_{ij} [b_{ij} \hat{c}_{ij} + \Delta_{ij} \hat{c}_{ij}^\dagger + \text{h.c.}] = \sum_k \left( \begin{array}{cc} \epsilon_k^\uparrow c_{-k}^\dagger & \Delta_k \end{array} \right) \left( \begin{array}{c} \xi_k \Delta - \xi_k \end{array} \right) \left( \begin{array}{c} \epsilon_k^\downarrow c_{k} \end{array} \right) \right),$$

(F1)

the one electron Green’s function $\mathcal{G}(p, i \omega_n)$ and the electronic self-energy $\Sigma(p, i \omega_n)$ are $2 \times 2$ matrix defined by the Dyson equation

$$[\mathcal{G}(p, i \omega_n)]^{-1} = [\mathcal{G}_0(p, i \omega_n)]^{-1} - \Sigma(p, i \omega_n),$$

(F2)

$$= i Z^{-1} \omega_n - \xi_p \sigma_3 - \phi(p, i \omega_n) \sigma_1,$$

(F3)

where $\mathcal{G}_0(p, i \omega_n)$ is the one-electron Green’s function for the non interacting system, $\sigma_i$ are Pauli matrices and $\xi_k = \epsilon_k - \mu$. The $\epsilon_k$’s are single particle electron energies and $\mu$ is the chemical potential. We do not include Coulomb electron-electron interaction, so that the self-energy $\Sigma \propto \sigma_1$ is just offdiagonal. The approximation used for the self-energy is$^{68}$:

$$\Sigma(p, i \omega_n) = -\frac{1}{\beta} \sum_{p' \nu'} \sigma_3 G(p', i \omega_n) \sigma_3 |g(p, p')|^2 \mathcal{D}(p - p', i \omega_n - i \omega_n),$$

(F4)

where $g(p, p')$ is the coupling with the bosonic modes and $\mathcal{D}(p - p', i \Omega_n)$ is the response function in imaginary frequency to the bosonic modes. The latter can be represented in terms of its imaginary part, $B(q, i \Omega_n) = -\frac{1}{\pi} \text{Im} \{\mathcal{D}(q, i \Omega_n)\}$ as

$$\mathcal{D}(q, i \omega_n - i \omega_n) = \int_0^\infty d\Omega B(q, \Omega) \left\{ \frac{1}{i \omega_n - i \omega_n - \Omega} - \frac{1}{i \omega_n - i \omega_n + \Omega} \right\},$$

where ($\Omega_q = \hat{D}_Q q^2$) and, in our case

$$B(q, i \Omega) = -\frac{1}{\pi} \left[ \frac{\Omega_q \Omega}{\Omega_q^2 + \Omega_q^2} \right] \mathcal{T}_Q$$

(F5)

To make matches with the usual Eliashberg theory, the isotropic gap model which we consider here provides the
dimensionless coupling function:

$$\alpha^2 F(\Omega) = \frac{1}{N_q} \sum_q \delta(\Omega - \Omega_q) \nu(0) \int dq' |g(k_F, \omega'; q)|^2; \quad (F6)$$

as an integral over the transferred momentum \( q \). We have to integrate over all \( q' s \):

$$\frac{1}{N_q} \sum_q \Rightarrow \frac{\tilde{a}^2}{(2\pi)^2} 2\pi \int dq \frac{d\Omega_q}{d\Omega} = \frac{\tilde{a}^2}{2\pi} \int d\Omega_q, \quad (F6)$$

where \( \tilde{a}^2 \) is the average diffusion area. We obtain

$$\Sigma(p, i \omega_n) = -\frac{1}{\beta} \sum_{q', \nu} \sigma_3 \mathcal{G}(p', i \omega_n) \sigma_3 \int d\Omega_q \frac{\tilde{a}^2}{4\pi} \int d\Omega_q |g_{p, p'}(\Omega_q)|^2 B(\Omega_q, \Omega) \right\} \frac{1}{i \omega_n - \Omega - \Omega} \frac{1}{i \omega_n - \Omega - \Omega} \quad (F7)$$

with

$$[\mathcal{G}(p, \omega)]^{-1} = Z^{-1} \omega - \left\{ \xi_p - i Z^{-1} \left[ |\omega| \frac{\epsilon_F}{\Omega^*} \ln \left( \frac{F}{W} \right) + \frac{\alpha}{Z} \nu_0 |\omega|^2 \ln \frac{Z |\tilde{v}_F| k_F}{|\omega|} \right] \right\} \sigma_3 - \Xi(\omega) \sigma_1 \quad (F7)$$

The mean field \( \Delta(\omega) = Z(\omega) \Xi(\omega) \) has to be determined in the following.

In Eq.(F7) the inverse lifetime of the quasiparticles of Eq.(68) of MT appears. Here \( \alpha \) is a numerical factor of order one.

Note the difference with the usual approach: diffusivity implies Eq.(F5) here, while usually is \( \mathcal{B}(q, \Omega) \sim \frac{\Omega_q}{q^2 + \Omega^2} \) in the Eliashberg approach.

On the other hand, from Eq.(F2): The final form of the selfenergy is:

$$\Sigma(p, \omega) = \nu_0 \int_{-\infty}^{\infty} d\omega' \Re \left\{ \frac{Z^{-1} \omega - \Xi(\omega') \sigma_1}{[\mathcal{P}(\omega')]^{1/2}} \right\} \int_0^{\infty} d\Omega \frac{d\Omega_q}{4\pi} |g_{p, p'}(\Omega_q)|^2 B(\Omega_q, \Omega)$$

$$\times \left[ \frac{f(-\omega')}{\omega - \omega' - \Omega - i 0^+} + \frac{f(\omega')}{\omega - \omega' + \Omega + i 0^+} + \frac{N(\Omega)}{\omega - \omega' - \Omega + i 0^+} + \frac{N(\Omega)}{\omega - \omega' + \Omega + i 0^+} \right],$$

$$\mathcal{P}(\omega) = Z^{-2} \omega^2 + Z^{-2} \left[ |\omega| \frac{\epsilon_F}{\Omega^*} \ln \left( \frac{F}{W} \right) + \frac{\alpha}{Z} \nu_0 |\omega|^2 \ln \frac{Z |\tilde{v}_F| k_F}{|\omega|} \right]^2 - \Xi^2(\omega), \quad (F8)$$

where \( N(\Omega) = [e^{\beta \Omega} - 1]^{-1} \) and \( f(\omega) = [e^{\beta \omega} + 1]^{-1} \) are the Bose and Fermi occupation probabilities. The term in curly brackets arises from \( \Im m \left\{ \nu_0 \int_{-\infty}^{+\infty} d\omega' \sigma_3 \mathcal{G}(p', \omega') \sigma_3 \right\} \), which turns into a real part from the inverse of \([\mathcal{G}(p', \omega')]^{-1}\) given in Eq.(F7). The critical temperature is the one at which \( \Delta \sim 0 \) and can be dropped in the denominator, but the gap equation has to be satisfied.

In all the further calculations we neglect the thermal excitations and drop \( N(\Omega) \). Observing that the integration variable \( \Omega_q \) has the meaning of the diffusive energy (see Eq.(F5)) it is clear that it cannot be integrated at energies above \( \Omega \). We also use the parameter equality \( Z \nu_0 = U_c^{-1} \) and we take \(|g_{k_F \omega}(\Omega_q)|^2 = g^2 \) constant \((|g|^{-1} \sim \text{time} \ (h = 1 \text{ here}) \). We get:

$$\Delta(\omega) \approx \frac{\tilde{a}^2}{D_q} \int_0^{\infty} \frac{d\omega}{|\omega|} \frac{d\Omega_q}{d\Omega} \ln \frac{U_c}{W} \Re \{\Delta(\omega')\} \frac{1}{U_c} \int_0^{U_c} d\Omega \times \int_0^{\Omega} \frac{d\Omega_q}{4\pi} |g_{k_F \omega}(\Omega_q)|^2 \left\{ \frac{1}{\pi} \left[ \Omega_q \Omega \Omega_q + \Omega_q^2 \right] T_q \right\} 2 \left\{ f(-\omega') - f(\omega') \right\} \frac{1}{\Omega + \omega'} \frac{1}{1 - \omega'/\Omega^*}, \quad (F9)$$

We concentrate on \( \omega = 0 \) and we deal with two contributions to \( \Delta(\omega') \) separately, \( \Delta(\omega') = \Delta_a + \Delta_b \), where the first arises from integration over \( 0 < \omega' < \Omega^* \sim W^2/U_c \) and the second from integration over \( \Omega^* < \omega' < U_c \).

In the first case, observing that the range of \( \omega' \) cannot be larger than \( \Omega^* \), but the Fermi function selects \( \omega' \sim 0 \), we neglect \( \omega' \) in the denominators of the curly bracket obtaining:
\[
\Delta_b(0) = \frac{\tilde{a}^2}{D_Q} \int_0^{U_c} d\omega' \frac{\Delta_\omega}{\omega'} \ln \left( \frac{\omega'}{\omega} \right) \left| g |^2 \int_0^{U_c} d\omega \frac{\Delta_\omega}{2 \pi} T_Q \{ f(-\omega') - f(\omega') \} \right.
\]
\[
\approx \frac{\tilde{a}^2}{D_Q} T_Q \frac{\Delta_\omega}{8\pi} \frac{\Delta_\omega}{\omega'} \ln \left( \frac{\omega'}{\omega} \right) \frac{\Delta_\omega}{4U_c} \int_0^{U_c} d\omega \frac{\Delta_\omega}{2 \pi} T_Q \{ f(-\omega') - f(\omega') \}
\]

Now the \( \Delta_\omega \) contribution. We neglect \( \Omega \) in the denominator in the curly bracket of Eq.(F9) and we include the \( \omega^2 \) term of the inverse lifetime, only.

\[
\Delta_b(0) = \frac{\tilde{a}^2}{D_Q} \int_0^{U_c} d\omega' \frac{\Delta_\omega}{2 \nu_0|\omega'|^2} \ln \frac{Z_{\nu_0,\omega'}}{|\omega'|} \frac{1}{U_c} \int_0^{U_c} d\Omega \frac{d\Omega}{4\pi} |g_{k,\omega}(\Omega)|^2
\]
\[
\times \left[ \frac{1}{\pi} \left[ \frac{\Omega_q^2}{\Omega_q^2 + \Omega_q^2} \right] T_Q \right] 2 \left\{ f(-\omega') - f(\omega') \right\}
\]
\[
\approx \frac{\tilde{a}^2}{D_Q} \frac{|g|^2}{4\pi^2} \frac{\Delta_\omega}{4U_c} \int_0^{U_c} d\omega' \frac{\Delta_\omega}{2 \nu_0|\omega'|^2} \ln \frac{Z_{\nu_0,\omega'}}{|\omega'|} \frac{1}{U_c} \int_0^{U_c} 2\Omega d\Omega \frac{1}{\Omega + \omega'}
\]

where we have disregarded the terms \( \propto e^{-\beta\omega'} \) in the last integral. The rest of the derivation can be found in the main text.

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