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Citation	Sachdev, Subir. 2010. Strange metals and the AdS/CFT correspondence. Journal of Statistical Mechanics: Theory and Experiment 2010: P11022.
Published Version	doi://10.1088/1742-5468/2010/11/P11022
Accessed	February 19, 2015 9:14:23 AM EST
Citable Link	http://nrs.harvard.edu/urn-3:HUL.InstRepos:7524883
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Strange metals and the AdS/CFT correspondence

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Abstract. I begin with a review of quantum impurity models in condensed matter physics, in which a localized spin degree of freedom is coupled to an interacting conformal field theory in $d = 2$ spatial dimensions. Their properties are similar to those of supersymmetric generalizations which can be solved by the AdS/CFT correspondence; the low energy limit of the latter models is described by a AdS₂ geometry. Then I turn to Kondo lattice models, which can be described by a mean-field theory obtained by a mapping to a quantum impurity coupled to a self-consistent environment. Such a theory yields a ‘fractionalized Fermi liquid’ phase of conduction electrons coupled to a critical spin liquid state, and is an attractive mean-field theory of strange metals. The recent holographic description of strange metals with a AdS₂ × R² geometry is argued to be related to such mean-field solutions of Kondo lattice models.

Submitted to: *Special Issue of JSTAT*.

Plenary talk at Statphys24, Cairns, Australia, July 2010

1. Introduction

The ‘strange metal’ is a non-zero temperature phase of electrons in solids which appears to be common to most ‘correlated electron’ compounds. These are compounds with transition-metal or rare-earth elements with a crystal structure which often promotes electronic transport primarily within a single two-dimensional layer of atoms. The strange metal is typically found at temperatures (T) above low temperature phases which display antiferromagnetism and superconductivity. It is defined as ‘strange’ because the temperature and frequency dependence of many observables deviate strongly from those expected from conventional Fermi liquid theory: most famously, the resistance increases linearly with T , over a wide range of T values. While the cuprate high temperature superconductors are the most prominent compounds displaying strange metal behavior [1], similar regimes are also found in ruthenium oxides [2], iron pnictides [3], organic metals [4], and heavy-fermion compounds [5].

The AdS/CFT correspondence was originally discovered as a tool for describing strongly-coupled gauge theories [6]. In $D = 4$ spacetime dimensions, Yang Mills gauge theory with $\mathcal{N} = 4$ supersymmetry (abbreviated as SYM4) is characterized by a single dimensionless coupling constant which remains invariant under the renormalization group (RG). Consequently, the theory is conformally invariant for all values of the coupling, and can be viewed as the simplest interacting conformal field theory in $D = 4$ spacetime dimensions: a CFT4. With a $SU(M)$ gauge group, this theory was argued to be equivalent to a string theory on an $AdS_5 \times S^5$ background, where AdS is anti-de Sitter, a symmetric space with constant negative curvature. More usefully, in the $M \rightarrow \infty$ limit for the low energy physics, the string theory can be approximated simply by classical Einstein gravity on AdS_5 . Thus, remarkably, correlations of gauge theories in $D = 4$ can be related to properties of the Einstein gravity in a negatively curved space in $D = 5$. The latter space has an emergent dimension, which can be interpreted as a RG energy scale.

CFTs also arise in condensed matter physics in many different contexts. In Section 2, we will briefly review the CFT3s arising near the quantum critical points of certain quantum antiferromagnets in $d = 2$ spatial dimensions. It was argued in Ref. [7] that such CFT3s could also be usefully analyzed via the AdS/CFT correspondence. In $D = 3$, a supersymmetric analog of the CFT3s in Section 2 is Yang-Mills theory with $\mathcal{N} = 8$ supersymmetry and a $SU(M)$ gauge group (SYM3), and this maps in the $M \rightarrow \infty$ limit to gravity on AdS_4 . Unlike the $D = 4$ case, the coupling constant of SYM3 flows generically to a strong-coupling fixed point, and so there is no free coupling and the low energy physics is conformal. Thus it is always ‘quantum critical’, and it is the first solvable strongly-interacting quantum critical theory in $D = 3$. The condensed matter CFT3s of Section 2 also have strongly interacting quantum critical points, many of whose properties have resisted accurate solution by other available methods. Even though they are not realized as the large M limit of a non-Abelian gauge theory, it was argued [7] that they could be modeled via the AdS/CFT correspondence. In this

context, it should be kept in mind that the operator content of the strongly-coupled fixed point of SYM3 is far removed from the non-Abelian gauge fields in terms of which it was written down at high energy scales; thus its quantum critical physics is no more the physics of non-Abelian gauge fields than that of the models of Section 2. Indeed, it is not unreasonable to view SYM3 at small M as a supersymmetric generalization of the famous classical $D = 3$ Ising model at its critical point. Such applications of the AdS/CFT correspondence to CFTs in condensed matter have been reviewed by the author in other articles [8, 9], and so will not be explored here. In a recent paper [10], we have argued that the AdS/CFT results can yield useful information for the frequency and temperature dependence of transport coefficients which have been studied in experiments in a variety of condensed matter systems.

The CFTs discussed so far, and those in Section 2, are states of quantum matter which are effectively at *zero density*. Their low energy spectrum is relativistic and particle-hole symmetric, similar to that found in *e.g.* pure, undoped graphene. To access states analogous to the strange metal, we have to study quantum matter at non-zero density. We do this by turning on a chemical potential, μ , which couples to a globally conserved charge of the CFT; we choose $\mu = 0$ to correspond to the zero density state. As long as $|\mu| \ll T$, the AdS/CFT correspondence can be extended straightforwardly, and many new results for quantum-critical transport in condensed matter have been obtained by this method [11]. These have also been reviewed elsewhere [8, 9], and will not be discussed further here.

We turn, finally, to states of non-zero density quantum matter at low temperature, which have $T \ll |\mu|$. Here, even for the supersymmetric CFTs at large M , the application of the AdS/CFT correspondence is not immediate. The CFT has scalar fields with exactly flat directions in their potential, and when placed at a non-zero μ at $T = 0$ these flat directions appear to lead to an instability of the theory. Nevertheless, one can presume that the strongly coupled CFT at non-zero μ continues to have a gravitational description in the AdS language. In the absence of a precise derivation of the theory on AdS, the spirit of effective field theory can be used to postulate a phenomenological action on AdS, which then predicts interesting new physics in the doped-CFT at low T . Just such a strategy has been used in a large number of recent studies. Two broad classes of states have been obtained by this method. States in one class display the condensation of a charged [12, 13, 14, 15] or a neutral [16, 17] scalar field (or both [18]); we will not study this class here. States in the second class have no broken symmetries and display evidence of metallic behavior and Fermi surfaces [19, 20, 21, 22, 23, 24, 25, 26, 27, 28, 29, 30]. Interestingly, the Fermi surface quasiparticles have non-Fermi liquid damping, and the resistance can have a linear dependence on T for a particular value of an effective parameter *i.e.* the AdS theory at non-zero μ leads to a holographic description of the strange metal.

However, many key properties of the holographic strange metal so obtained are quite mysterious from a condensed-matter perspective. For a CFTD, the AdS_{D+1} theory at non-zero μ factorizes [22] at low energies to a $\text{AdS}_2 \times R^{D-1}$ geometry, and

this factorization is a key determinant in the unusual properties of the Fermi surface. One immediate consequence is that the singular features in the quasiparticle energy depend primarily on the frequency ω , and this quantum criticality is described by the AdS₂ geometry alone: thus crucial information on the spatial dependence of the strange metal quantum fluctuations appears to be missing. Related to this is the troubling presence of a non-zero entropy density which survives the $T \rightarrow 0$ limit, thus apparently violating the third law of thermodynamics. Finally, the underlying field content leading to this holographic metal phase is not clear.

This paper will review and give additional perspective on a recent proposal [29] connecting Kondo lattice models to the holographic strange metal. We will do this here by adding matter degrees of freedom to the zero density CFT ‘one-at-a-time’. As noted above, we begin in Section 2 by describing how CFTs arise in two-dimensional quantum antiferromagnets. In Section 3 will add a single spin degree of freedom to the CFT: models of this type can be reliably addressed both by traditional condensed matter methods, and in supersymmetric cases by the AdS/CFT correspondence. We will find a close correspondence in the physical properties in the two cases, including an emergence of AdS₂ in the gravitational description. Thus in the single impurity context, a satisfactory physical interpretation of the AdS₂ geometry will be obtained.

Then we will turn to Kondo lattice models and their holographic interpretation in Section 4. In the limit of large dimension, or long-range exchange interactions, such models can be solved by a mapping to a quantum impurity model coupled to a self-consistent environment. The mean-field solution describes a ‘fractionalized Fermi liquid’ (FFL or FL*), with a Fermi surface of conduction electrons coupled to a critical spin liquid. We will argue that such a solution has a close correspondence to the holographic theory with the AdS₂ \times R^{D-1} geometry. This description of the holographic theory leads to simple interpretations of its physical properties. For the Kondo lattice, the mean-field theory of the critical spin liquid does not have collective gauge excitations which are generically expected to be present in realistic spin liquid with finite-range interactions. The holographic theory reaches a similar mean-field but without infinite range interactions: thus it is likely to be amenable to systematic improvements. We hope such improvements will eventually lead to a description of realistic spin liquids and FL* states.

2. Quantum antiferromagnets and CFT3s

We begin with a brief review of the connection between lattice quantum antiferromagnets in $d = 2$ and CFT3s, also reviewed in [9]. The lattice antiferromagnets are described by the Heisenberg exchange Hamiltonian

$$H_J = \sum_{i < j} J_{ij} \hat{S}_i^a \hat{S}_j^a + \dots \quad (1)$$

where $J_{ij} > 0$ is the antiferromagnetic exchange interaction and \hat{S}_i^a ($a = x, y, z$) are $S = 1/2$ spin operators on the sites, i , of a regular lattice: thus they obey the

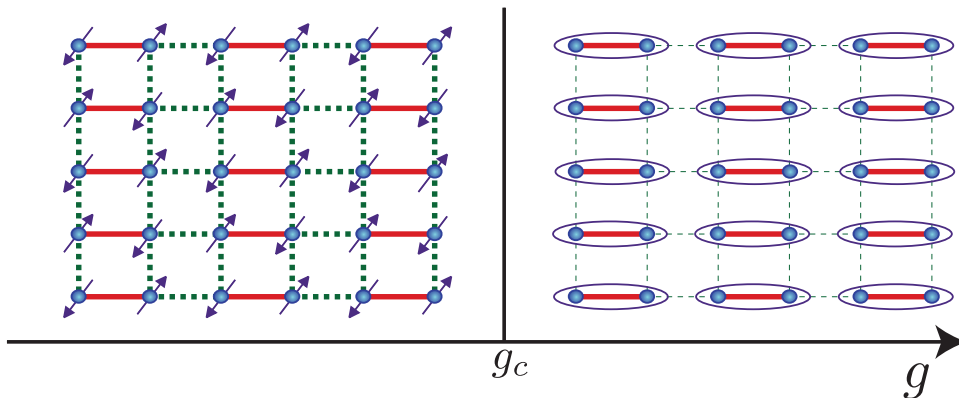


Figure 1. The dimer antiferromagnet. The full red lines represent an exchange interaction J , while the dashed green lines have exchange J/g . The ellipses represent a singlet valence bond of spins $(|\uparrow\downarrow\rangle - |\downarrow\uparrow\rangle)/\sqrt{2}$.

commutation relations

$$[\hat{S}_i^a, \hat{S}_j^b] = i\epsilon_{abc}\delta_{ij}\hat{S}_i^c \quad (2)$$

and $\sum_a (\hat{S}_i^a)^2 = 3/4$ for each i .

Different phases, quantum phase transitions and low energy field theories are obtained depending upon whether there are an even or an odd number of $S = 1/2$ spins per unit cell of the lattice.

2.1. Even number of $S = 1/2$ spins per unit cell

Let us consider first the simpler case of an even number of $S = 1/2$ spins per unit cell. The canonical model is the dimer antiferromagnet, illustrated in Fig. 1. The $S = 1/2$ spins reside on the sites of a square lattice, and have nearest neighbor exchange equal to either J or J/g . Here $g \geq 1$ is a tuning parameter which induces a quantum phase transition in the ground state of this model. At $g = 1$, the model has full square lattice symmetry, and this case is known to have a Néel ground state which breaks spin rotation symmetry. This state has a checkerboard polarization of the spins, just as found in the classical ground state, and as illustrated on the left side of Fig. 1. It can be characterized by a vector order parameter φ^a which measures the staggered spin polarization

$$\varphi^a = \eta_i S_i^a \quad (3)$$

where $\eta_i = \pm 1$ on the two sublattices of the square lattice. In the Néel state we have $\langle \varphi^a \rangle \neq 0$, and we expect that the low energy excitations can be described by long wavelength fluctuations of a field $\varphi^a(r, \tau)$ over space, r , and imaginary time τ . On the other hand, for $g \gg 1$ it is evident from Fig. 1 that the ground state preserves all symmetries of the Hamiltonian: it has total spin $S = 0$ and can be considered to be a product of nearest neighbor singlet valence bonds on the J links. The simplicity of this large g ground state relies crucially on the ‘dimerized’ structure of the Hamiltonian; the fact that there are an even number of $S = 1/2$ spins per unit cell.

It is clear that the $g = 1$ and $g \gg 1$ states are qualitatively distinct, and so there must be a quantum phase transition at a critical $g = g_c$. We can deduce the quantum field theory for this phase transition by using conventional Landau-Ginzburg arguments: we use the order parameter φ^a , and write down the simplest continuum action consistent with the symmetries of the Hamiltonian; this leads to the partition function

$$\begin{aligned} \mathcal{Z} &= \int \mathcal{D}\varphi^a(r, \tau) \exp\left(-\int d^2r d\tau \mathcal{L}_\varphi\right) \\ \mathcal{L}_\varphi &= \frac{1}{2} [(\partial_\tau \varphi^a)^2 + v^2(\nabla \varphi^a)^2 + s(\varphi^a)^2] + \frac{u}{4} [(\varphi^a)^2]^2 \end{aligned} \quad (4)$$

The transition is now tuned by varying $s \sim (g - g_c)$. Notice that this model is identical to the Landau-Ginzburg theory for the thermal phase transition in a $d + 1$ dimensional ferromagnet, because time appears as just another dimension. From this we conclude that the quantum phase transition is described by the famous Wilson-Fisher fixed point of Eq. (4). This was originally discovered by an analysis of the theory in $D = 4 - \epsilon$ spacetime dimensions, using ϵ as an expansion parameter. Since then, extensive numerical and analytical studies have shown that the fixed point is present also in $D = 3$, where it describes a non-trivial CFT3. For the dimer antiferromagnet, very convincing evidence that the quantum criticality is described by the Wilson-Fisher CFT3 is presented in [31].

In experiments, the best studied realization of the dimer antiferromagnet is TlCuCl_3 . In this crystal, the dimers are coupled in all three spatial dimensions, and the transition from the dimerized state to the Néel state can be induced by application of pressure. Neutron scattering experiments by Ruegg and collaborators [32] have clearly observed the transformation in the excitation spectrum across the transition, and these observations are in good quantitative agreement with theory[33].

2.2. Odd number of $S = 1/2$ spins per unit cell

For this case, we can work with the Hamiltonian in Eq. (1), but with the J_{ij} respecting the full space group symmetry of the square lattice. Thus the nearest neighbor J_{ij} must all be equal to each other, unlike the dimer antiferromagnet above. With no further range interactions, the ground state has Néel order, as we discussed in Section 2.1. A variety of routes have been investigated [33] to continuously destroy the Néel order, and we generically represent them as tuning a coupling g in Fig. 2. The phase so-reached was argued [34] to have valence bond solid order (VBS). The VBS state is superficially similar to the dimer singlet state in the right panel of Fig. 1: the spins primarily form valence bonds with near-neighbor sites. However, because of the square lattice symmetry of the Hamiltonian, a columnar arrangement of the valence bonds as in Fig. 1, breaks the square lattice rotation symmetry; there are 4 equivalent columnar states, with the valence bond columns running along different directions. More generally, a VBS state is a spin singlet state, with a non-zero degeneracy due to a spontaneously broken lattice symmetry. A VBS state has been observed in the organic antiferromagnet $\text{EtMe}_3\text{P}[\text{Pd}(\text{dmit})_2]_2$ [35, 36].

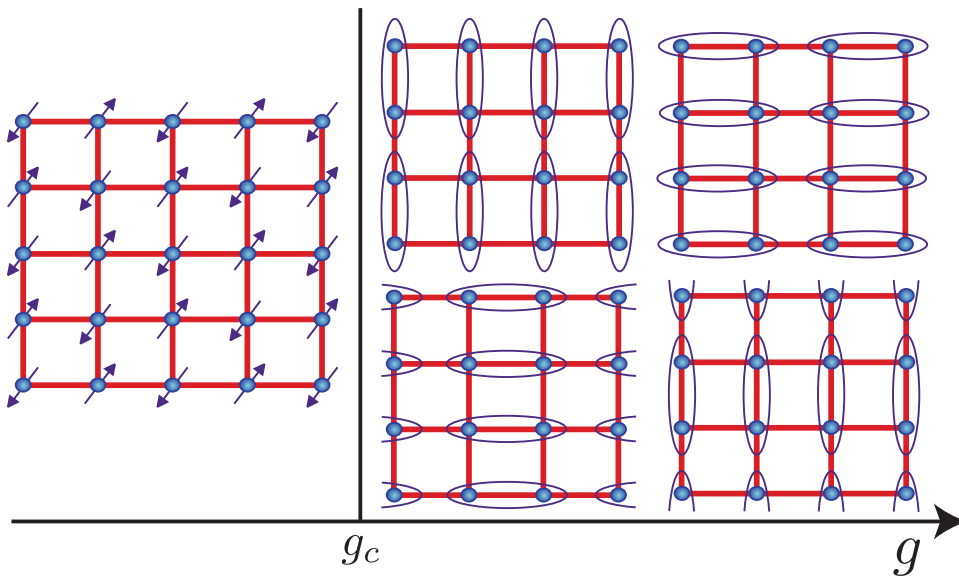


Figure 2. A frustrated square lattice antiferromagnet on the square lattice, with the Hamiltonian preserving the full space group symmetry of the square lattice. The valence bond solid (VBS) state for $g > g_c$ is four-fold degenerate, depending upon the crystallization pattern of the singlet valence bonds.

A direct transition at $g = g_c$ between the Néel and VBS states involves two distinct broken symmetries: spin rotation symmetry, which is broken only in the Néel state for $g < g_c$, and a lattice rotation symmetry, which is broken only in the VBS state for $g > g_c$. The rules of Landau-Ginzburg-Wilson theory imply that there can be no generic second-order transition between such states. It has been argued that a second-order Néel-VBS transition can indeed occur [37], but the critical theory is not expressed directly in terms of either order parameter. It involves a fractionalized bosonic spinor z^α ($\alpha = \uparrow, \downarrow$), and an emergent gauge field A_μ . The key step is to express the vector field φ^a in terms of z^α by

$$\varphi^a = z_\alpha^* (\sigma^a)^\alpha_\beta z^\beta \quad (5)$$

where σ^a are the 2×2 Pauli matrices. Note that this mapping from φ^a to z^α is redundant. We can make a spacetime-dependent change in the phase of the z^α by the field $\theta(r, \tau)$

$$z^\alpha \rightarrow e^{i\theta} z^\alpha \quad (6)$$

and leave φ^a unchanged. All physical properties must therefore also be invariant under Eq. (6), and so the quantum field theory for z^α has a U(1) gauge invariance, much like that found in quantum electrodynamics. The effective action for the z^α therefore requires introduction of an ‘emergent’ U(1) gauge field A_μ (where $\mu = x, \tau$ is a three-component spacetime index). The field A_μ is unrelated the electromagnetic field, but is an internal field which conveniently describes the couplings between the spin excitations of the antiferromagnet. As for Eq. (4), we can write down the quantum field theory for

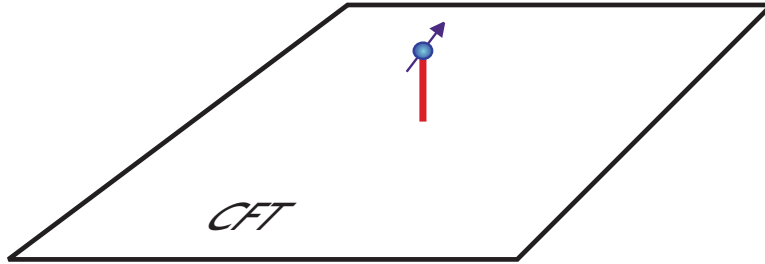


Figure 3. A quantum spin coupled via an exchange interaction to a CFT in 2+1 dimensions.

z^α and A_μ by the constraints of symmetry and gauge invariance, which now yields

$$\mathcal{Z} = \int \mathcal{D}z^\alpha(r, \tau) \mathcal{D}A_\mu(r, \tau) \exp\left(-\int d^2r d\tau \mathcal{L}_z\right)$$

$$\mathcal{L}_z = |(\partial_\mu - iA_\mu)z^\alpha|^2 + s|z^\alpha|^2 + u(|z^\alpha|^2)^2 + \frac{1}{2w^2}(\epsilon_{\mu\nu\lambda}\partial_\nu A_\lambda)^2 \quad (7)$$

For brevity, we have now used a “relativistically” invariant notation, and scaled away the spin-wave velocity v ; the values of the couplings s, u, w are different from, but related to, those in Eq. (4). The Maxwell action for A_μ is generated from short distance z^α fluctuations, and it makes A_μ a dynamical field. This theory has a ‘Higgs’ phase where z^α condenses like the Higgs boson: this we can identify as the Néel state. The ordinary Coulomb phase with z^α gapped appears as a ‘spin liquid’ state with a collective gapless, spinless excitation associated with the A_μ photon. Non-perturbative effects [34] associated with the monopoles in A_μ (not discussed here), show that this spin liquid is ultimately unstable to the appearance of VBS order.

An interesting question now is whether the transition between the Néel and VBS states as described by (7) is a CFT3. The existence of a CFT3 fixed point has been established order-by-order in the $1/N$ expansion, where the spinor index $\alpha = 1 \dots N$. However, the issue remains unsettled for $N = 2$ [38, 39].

For our purposes here, the CFT3s described by Eqs. (4) and (7) are non-supersymmetric analogs of the CFT3 realized by SYM3. Insights gained from the AdS/CFT correspondence are described elsewhere [9, 10].

3. Quantum impurity in a CFT

As we discussed in Section 1, we will move away from the zero density CFTs of Section 3 by adding a single defect localized in space. This will eventually allow us to address the non-zero density case in the following section.

For the quantum antiferromagnets of Section 3, the simplest interesting defect is a single spin \hat{S}^a coupled to the antiferromagnet by an exchange coupling, J , as shown in Fig. 3.

More generally, the configuration of Fig. 3 belongs to a wide class of ‘Kondo’ problems. Usually, the bulk CFT is rather simple: it is a free electron system whose

Fermi surface excitations form an infinite set of CFT2s of free chiral fermions in 1+1 spacetime dimensions; furthermore only a single CFT2 of the free fermions is coupled to the impurity. In such cases, all the quantum correlation effects arise solely in the vicinity of the impurity. In the simplest and most common case of quantum spin coupled to a Fermi surface, there is no interesting quantum-criticality in the low energy limit. There are only two fixed points of the RG, $J = 0$ and $J = \infty$: at the unstable $J = 0$ fixed point, the impurity decouples from the bulk CFT, while at the stable $J = \infty$ fixed point there are only innocuous potential scattering perturbation [40]. However, the situation becomes far more non-trivial if the bulk fermions acquire an additional ‘channel’ or ‘flavor’ index; then, under suitable conditions, a non-trivial stable fixed point is obtained at an intermediate $J = J^*$ [41]. It is the analog of this ‘multi-channel Kondo fixed point’ [42] which we will explore in this section in the more general setting of bulk CFTs with interactions in $D > 2$ spacetime dimensions.

As a first example, let us couple the impurity spin \hat{S}^a to the class of dimer antiferromagnets described by the field theory (4). We can represent the impurity spin by a coherent state path integral over the fluctuations of a unit vector $n^a(\tau)$; then the partition function of the quantum impurity problem becomes

$$\begin{aligned} \mathcal{Z} &= \int \mathcal{D}\varphi^a(r, \tau) \mathcal{D}n^a(\tau) \delta([n^a(\tau)]^2 - 1) \exp\left(-\int d\tau \mathcal{L}_{\text{imp}} - \int d^2r d\tau \mathcal{L}_\varphi\right) \\ \mathcal{L}_{\text{imp}} &= \frac{i}{2} \mathcal{A}^a \frac{dn^a}{d\tau} + J n^a(\tau) \varphi^a(0, \tau) \end{aligned} \quad (8)$$

Here the quantum spin commutation relations in Eq. (2) for the impurity spin are implemented by the Berry phase term in \mathcal{L}_{imp} where \mathcal{A}^a is any function of $n^a(\tau)$ obeying $\epsilon^{abc}(\partial\mathcal{A}^b/\partial n^c) = n^a$. Equivalently, we may represent the impurity spin by a ‘slave’ fermion χ^α , and then the partition function in Eq. (8) can be written as

$$\begin{aligned} \mathcal{Z} &= \int \mathcal{D}\varphi^a(r, \tau) \mathcal{D}\chi^\alpha(\tau) \exp\left(-\int d\tau \mathcal{L}_{\text{imp}} - \int d^2r d\tau \mathcal{L}_\varphi\right) \\ \mathcal{L}_{\text{imp}} &= \chi_\alpha^\dagger \frac{\partial \chi^\alpha}{\partial \tau} + J \chi_\alpha^\dagger [(\sigma^a)_\beta^\alpha \varphi^a(0, \tau)] \chi^\beta. \end{aligned} \quad (9)$$

Actually, the partition function in Eq. (9) has a conserved fermion number $n_\chi = \chi_\alpha^\dagger \chi^\alpha$, and so splits into different sectors labeled by the possible values of $n_\chi = 0, 1, 2$. The mapping to Eq. (8) requires restriction to the sector with $n_\chi = 1$; this constraint can be implemented by a Lagrange multiplier, which we have not written out explicitly.

The quantum impurity problem defined by Eq. (8) or (9) is amenable to a RG analysis using an expansion in $\epsilon = 4 - D$. An extensive theoretical study has been carried out by this method [43, 44, 45], and we now summarize the main results. When the bulk theory is at the Wilson-Fisher CFT3 fixed point, the impurity coupling J flows to a stable fixed point at some $J = J^*$. Some of the characteristics of this fixed point are:

- The correlations of the impurity fermion χ^α and the impurity spin $\hat{S}^a = (1/2)\chi_\alpha^\dagger (\sigma^a)_\beta^\alpha \chi^\beta$ decay with a power-law in time, with non-trivial ‘impurity’ exponents which can be computed order-by-order in ϵ .

- The impurity response to a uniform external field is characterized by an impurity susceptibility which has a Curie form $\chi_{\text{imp}} = \mathcal{C}/T$, where \mathcal{C} is a non-trivial universal number which can be computed in the ϵ expansion. This response is that of an ‘irrational’ free spin, because $\mathcal{C} \neq S(S+1)/3$, with $2S$ an integer.
- There is a finite ground state entropy, S_{imp} , at $T = 0$. This entropy is also ‘irrational’ because $S_{\text{imp}} \neq k_B \ln(\text{an integer})$.

Extensive numerical studies [46, 47, 48, 49] have also been carried out for the above quantum impurity problem, and the results so far are in good agreement with the theoretical expectations.

We also mention here an alternative strong-coupling formulation [50] of the above quantum impurity coupled to the dimer antiferromagnet described by \mathcal{L}_φ . This alternative formulation turns out to be suitable for a determination of the universal critical properties in an expansion in $\epsilon = D - 2$. For this formulation we turn from the ‘soft-spin’ formulation of the bulk critical theory in Eq. (4), to a ‘hard-spin’ formulation in terms of a unit-length field $n^a(r, \tau)$. Then, as is well known, the bulk Wilson-Fisher fixed point is accessed by a $D = 2 + \epsilon$ expansion of the O(3) non-linear σ -model. For the quantum impurity physics, it has been argued that the fixed length limit requires that we send the impurity-bulk coupling J to infinity. In other words, the orientation of the impurity spin \hat{S}^a is fixed to be parallel to that of the bulk field $n^a(r, \tau)$ at $r = 0$. Finally, because the impurity spin orientation is fixed, we only need a spinless fermion χ to account for the presence/absence of the impurity. With this reasoning, we obtain the partition function

$$\begin{aligned}
 \mathcal{Z} &= \int \mathcal{D}n^a(r, \tau) \mathcal{D}\chi(\tau) \delta([n^a(r, \tau)]^2 - 1) \exp\left(-\int d\tau \mathcal{L}_{\text{imp}} - \int d^2r d\tau \mathcal{L}_n\right) \\
 \mathcal{L}_n &= \frac{1}{2g} \left[\left(\frac{\partial n^a}{\partial \tau}\right)^2 + c^2 (\nabla n^a)^2 \right] \\
 \mathcal{L}_{\text{imp}} &= \chi^\dagger \left(\frac{\partial}{\partial \tau} - iA_\tau \right) \chi \quad , \quad A_\tau \equiv \frac{1}{2} \mathcal{A}^a \frac{dn^a(0, \tau)}{d\tau}
 \end{aligned} \tag{10}$$

where \mathcal{A}^a is now a function of $n^a(0, \tau)$, and the gauge potential A_τ is the pullback of \mathcal{A}^a from S^2 . The conserved fermion number can now only take the values $n_\chi = 0, 1$. Here the restriction to $n_\chi = 1$ is trivially implemented because there is only a single fermion state without a fermion spin index: we simply omit χ from the functional integral, while including the Wilson line source term $\exp(i \int d\tau A_\tau)$. The claim [50] is that the $\epsilon = D - 2$ expansion of the partition function in Eq. (10) describes the same universal fixed point as the $\epsilon = 4 - D$ expansion of the partition function in Eq. (9). Notice that the theory (10) has only a single coupling constant g , and this reaches the same fixed point as in the bulk theory. The impurities properties are nevertheless non-trivial and universal, and are entirely a consequence of the Berry phase of the impurity. The theoretical results from this formulation have been successfully compared to numerical studies away from the bulk critical point, within the ordered Néel phase [51, 52].

Next, let us turn to the bulk CFT3 associated with an odd-number of $S = 1/2$ spins per unit cell, the CP^{N-1} model in Eq. (7). Now, there are $S = 1/2$ excitations z^α in the bulk, and so the impurity spin can be more efficiently screened by the environment. To the z^α , the impurity spin appears as a static external $\text{U}(1)$ gauge charge, and so the bulk+impurity theory takes a form similar to Eq. (10). We have [53]:

$$\begin{aligned} \mathcal{Z} &= \int \mathcal{D}z^\alpha(r, \tau) \mathcal{D}A_\mu(r, \tau) \mathcal{D}\chi(\tau) \exp \left(- \int d\tau \mathcal{L}_{\text{imp}} - \int d^2r d\tau \mathcal{L}_z \right) \\ \mathcal{L}_{\text{imp}} &= \chi^\dagger \left(\frac{\partial}{\partial \tau} - iA_\tau(0, \tau) \right) \chi \end{aligned} \quad (11)$$

Like Eq. (10), there is no coupling constant associated with the impurity theory, and so the impurity responses are naturally universal. It should be emphasized that the theory in Eq. (11) is *different* from those in Eqs. (9) and (10): the bulk CFTs correspond to the two cases in Section 2, and so the impurity dynamics is also distinct. The properties of theory in Eq. (11) have been studied in some detail using the $1/N$ expansion [53, 54, 55, 56], and there have also been recent numerical studies of this case [39, 57, 39].

Finally, let us turn to supersymmetric gauge theories, and consider a quantum impurity problem associated with a bulk CFT of SYM4 with the $\text{SU}(M)$ gauge group. Such a problem was considered recently by Kachru, Karch and Yaida [58, 59]. Their impurity was represented by a localized fermion χ^b with $b = 1 \dots M$ a $\text{SU}(M)$ color index. The action for their field theory was

$$\begin{aligned} \mathcal{S} &= \int d^3r d\tau \mathcal{L}_{\text{SYM}} + \int d\tau \mathcal{L}_{\text{imp}} \\ \mathcal{L}_{\text{imp}} &= \chi_b^\dagger \frac{\partial \chi^b}{\partial \tau} + i\chi_b^\dagger \left[(A_\tau(0, \tau))_c^b + v^I (\phi_I(0, \tau))_c^b \right] \chi^c \end{aligned} \quad (12)$$

Here A_μ and ϕ_I are bulk fields of SYM4 which are adjoints under $\text{SU}(M)$, $I = 1 \dots 6$, \mathcal{L}_{SYM} is the Lagrangian of the bulk SYM4 CFT, and v^I is a unit 6-vector determining the specific choice of the quantum impurity. The similarity of Eq. (12) to Eqs. (9), (10), and (11) should now be strikingly evident: in all cases we have an impurity localized fermions, and these are coupled to the bulk CFT by a universal gauge-like coupling. While the present supersymmetric theory has no direct application to condensed matter models, it has the advantage of being solvable by the AdS/CFT correspondence in the limit of $M \rightarrow \infty$. Such a gravitational solution has been presented by Kachru *et al.* [58, 59], who showed that the low energy physics of the quantum impurity is associated with a AdS_2 geometry in the gravity theory (see also Refs [60, 61, 62]). Further, the physical properties of the model in Eq. (12) were found to be qualitatively identical to those listed above for Eq. (9); in particular, the AdS_2 solution also has a non-zero ground state impurity entropy. Thus we may conclude [29] that there is an intimate connection between the quantum impurity models considered in this section, and quantum gravity on AdS_2 .

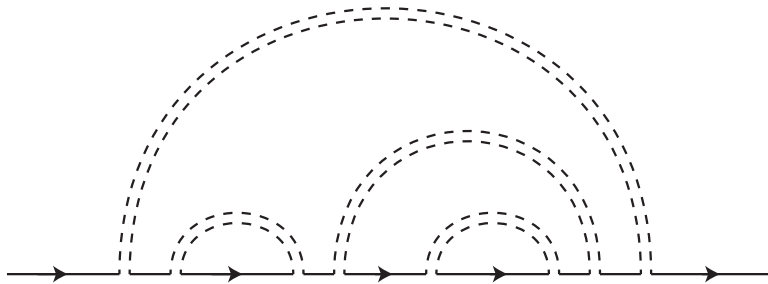


Figure 4. The rainbow graphs. The full line is the fermion, and double-dashed line is the interaction $D(\tau)$. Each line, full or dashed, carries a $SU(N)$ index α .

3.1. Large N solution

This subsection will illustrate the above general concepts on quantum impurities by an explicit solution in a simple limiting case. We will look at a large N limit in which a fermionic field is a vector of N components. It is also possible to set up a large N limit which has the character of a matrix-large N [63], but that is not easily solvable and will not be considered here.

We begin with the theory in Eq. (9), and assume that the bulk field φ^a has Gaussian correlations. This neglects bulk interactions which are ultimately necessary for an accurate description of the critical properties; however, this omission will not be crucial for the calculation discussed below. Ultimately, the justification of non-Gaussian bulk correlations relies on the large dimension or long-range limit of the models to be discussed in Section 4 (note that the bulk Gaussian approximation was not made in the theoretical studies noted above [43, 44, 45, 53, 54, 55, 56]). Anticipating Section 4, we assume that the correlation of φ^a for the impurity physics are fully characterized by the 2-point correlation

$$\langle \varphi^a(0, \tau) \varphi^b(0, \tau') \rangle = \delta^{ab} D(\tau - \tau') \quad (13)$$

Integrating out φ^a from Eq. (9), we obtain a ‘local’ partition function which involves a functional integral over fields that depend only upon τ

$$\begin{aligned} \mathcal{Z} = & \int \mathcal{D}\chi^\alpha(\tau) \mathcal{D}\lambda(\tau) \exp \left(- \int d\tau \left(\chi_\alpha^\dagger \frac{\partial \chi^\alpha}{\partial \tau} + i\lambda(\chi_\alpha^\dagger \chi^\alpha - N/2) \right) \right. \\ & \left. + \frac{2J^2}{N} \int d\tau d\tau' D(\tau - \tau') \chi_\alpha^\dagger(\tau) \chi_\beta^\dagger(\tau') \chi^\beta(\tau) \chi^\alpha(\tau') \right) \end{aligned} \quad (14)$$

where the indices $\alpha, \beta = 1 \dots N = 2$. However, we have written the partition function in a manner so that it can be used for general N , and the limit $N \rightarrow \infty$ is well-defined. Indeed, an explicit solution can be obtained in the large N limit [64], as we now describe.

An examination of the Feynman graph expansion shows that the limit of large N is dominated by [65] the ‘rainbow’ (or ‘non-crossing’) graphs for the fermion Green’s function: see Fig. 4. The summation of these graphs can be written analytically in terms of the following self-consistency conditions on the fermion self energy. As usual,

we define the fermion Green's function G by

$$G(\tau)\delta_\beta^\alpha = -\left\langle \mathcal{T}\chi^\alpha(\tau)\chi_\beta^\dagger(0) \right\rangle, \quad G(\omega_n) = \int_0^{1/T} d\tau e^{i\omega_n\tau} G(\tau) \quad (15)$$

where \mathcal{T} is imaginary-time ordering, and ω_n is a Matsubara frequency. This Green's function is expressed in terms of the self energy by

$$G(\omega_n) = \frac{1}{i\omega_n - \bar{\lambda} - \Sigma(\omega_n)} \quad (16)$$

where $\bar{\lambda}$ is the saddle-point value of $i\lambda$. Then, in the large N limit, it is not difficult to show that the fluctuations of λ about its saddle point can be neglected, and the self energy is given by

$$\Sigma(\tau) = 4J^2 D(\tau)G(\tau). \quad (17)$$

Finally, the fermion number constraint $n_\chi = N/2$, is equivalent to

$$G(\tau \rightarrow 0) = -\frac{\text{sgn}(\tau)}{2}. \quad (18)$$

We are now faced with the mathematical problem of solving the Eqs (16,17,18) for the unknown functions $G(\tau)$ and $\Sigma(\tau)$, and the value of $\bar{\lambda}$. While this is not difficult to do numerically [64] for a general $D(\tau)$, we present here an analytic solution in the limit of low energies for the case of critical correlations in $D(\tau)$. When the bulk theory is a CFT, we expect a power-law decay $D(\tau) \sim \tau^{-\gamma}$, with γ a critical exponent. We generalize this to $T > 0$, with the 'conformal' form

$$D(\tau) = A \left| \frac{\pi T}{\sin(\pi T\tau)} \right|^\gamma, \quad -1/T < \tau < 1/T \quad (19)$$

where A is some real constant. This is the general form of a $T > 0$ correlator at $x = 0$ for a CFT2, and holds also for \mathcal{L}_φ in the upper-critical dimension $D = 4$ which is the only case where the present Gaussian approximation for φ^a correlations is appropriate. Note that Eq. (19) is supposed to be valid at energies well below the ultraviolet cutoff $\sim J$; in other words, for $1/|\tau|, T \ll J$.

We will now show that the solution of Eqs (16,17,18) has the following form at long times [64, 66]

$$G(\tau) = B \text{sgn}(\tau) \left| \frac{\pi T}{\sin(\pi T\tau)} \right|^\rho, \quad -1/T < \tau < 1/T \quad (20)$$

and determine the exact values of B and the exponent ρ . Again Eq. (20) holds only for low energies with $1/|\tau|, T \ll J$. We have used the particle-hole symmetric nature of the constraint in Eq. (18) to conclude that G should be an odd function of τ . Also, note that we are not concerned that Eq. (20) does not obey Eq. (18) as $\tau \rightarrow 0$, because Eq. (20) does not apply in this limit.

We can now perform the Fourier transform of Eqs. (17,19,20) to obtain the low frequency behavior of the Green's function and the self energy:

$$\begin{aligned}
 G(\omega_n) &= [iB\Pi(\rho)] \frac{T^{\rho-1} \Gamma\left(\frac{\rho}{2} + \frac{\omega_n}{2\pi T}\right)}{\Gamma\left(1 - \frac{\rho}{2} + \frac{\omega_n}{2\pi T}\right)} \\
 \Sigma_{\text{sing}}(\omega_n) &= [i4J^2 AB\Pi(\rho + \gamma)] \frac{T^{\rho+\gamma-1} \Gamma\left(\frac{\rho+\gamma}{2} + \frac{\omega_n}{2\pi T}\right)}{\Gamma\left(1 - \frac{\rho+\gamma}{2} + \frac{\omega_n}{2\pi T}\right)}
 \end{aligned} \tag{21}$$

with

$$\Pi(s) \equiv \pi^{s-1} 2^s \cos\left(\frac{\pi s}{2}\right) \Gamma(1-s). \tag{22}$$

We have noted that the contribution to Σ in Eq. (21) is only the singular low frequency term, and an additional cutoff-dependent constant has been omitted.

It now remains to determine if the proposed solution in Eq. (21) obeys the Dyson equation in Eq. (16). We need only obtain agreement in the low frequency limit, where we find that after canceling $\bar{\lambda}$ with the regular part of the self-energy, the Dyson equation reduces for the singular contributions simply to [64]

$$G(\omega_n)\Sigma_{\text{sing}}(\omega_n) = -1. \tag{23}$$

Note that we have assumed that the bare $i\omega_n$ frequency dependence in Eq. (16) is subdominant to the singular contribution from the self energy in the low frequency limit; this requires $\gamma < 2$. Remarkably, we find that the frequency dependent expressions in Eq. (21) can indeed satisfy the constraint in Eq. (23) for *all* ω_n provided we choose the exponent [64, 66]

$$\rho = 1 - \frac{\gamma}{2}, \tag{24}$$

and the prefactor

$$B = [4J^2 A\Pi(1 - \gamma/2)\Pi(1 + \gamma/2)]^{-1/2}. \tag{25}$$

Thus we have the surprising conclusion that the innocuous-looking Eqs (16,17,18) which solve the partition function in Eq. (14) in the large- N limit have a solution which has a conformally-invariant structure at low energies [66, 67]. We then observe that the supersymmetric model in Eq. (12) also has correlators which have a conformal structure, inherited from the AdS₂ geometry [22, 68]. This is then further evidence for the striking connection between the quantum impurity models of this section and the theories of gravity on AdS₂.

It is also useful to collect results for the impurity spin correlation function

$$C(\tau) = \left\langle \hat{S}^a(\tau)\hat{S}^a(0) \right\rangle \tag{26}$$

from the present large N solution. Using the SU(2) relation $\hat{S}^a = (1/2)\chi_\alpha^\dagger(\sigma^a)_\beta^\alpha \chi^\beta$, and evaluating Eq. (26) in the large N limit, we find $C(\tau) = -(N^2/2)G(\tau)G(-\tau)$, and so

$$C(\tau) = \frac{B^2 N^2}{2} \left| \frac{\pi T}{\sin(\pi T\tau)} \right|^h \tag{27}$$

where the exponent [64, 69]

$$h = 2 - \gamma. \tag{28}$$

Of course, in the large N limit we have $h = 2\rho$, but we have introduced an independent exponent h for $C(\tau)$ because we expect that at higher orders in $1/N$ we have $h \neq 2\rho$. On the other hand, it has been argued [44] that the exponent relationship in Eq. (28) is *exact*, and holds to all orders in the $1/N$ expansion for the theory in Eq. (14). This exponent relationship is a consequence of the fact that the Gaussian field φ^a and the spin operator \hat{S}^a are conjugate operators in the second term of \mathcal{L}_{imp} in Eq. (4), and the co-efficient of this term reaches a fixed-point value in the RG.

Finally, this large N computation can also be used to compute the impurity entropy. This requires a somewhat more involved computation [66, 67], and will not be presented here.

4. From a quantum impurity to lattice models

The quantum impurity models discussed in Section 3 now appear to be very well understood: the results of different expansion methods are consistent with each other, and with a variety of numerical studies. And for the supersymmetric impurity models, the AdS/CFT method yields similar results using the geometry of AdS_2 .

In this section, we move beyond impurity models to a variety of lattice models. In condensed matter studies, this is done by using various expansion methods or physical arguments to motivate a specific mean-field decoupling of the quantum lattice model to a preferred ‘impurity’ spin coupled to a bulk ‘environment’. The resulting mean-field theory then has a structure similar to the models considered in Section 3. However, now the ‘environment’ is built out of the same degrees of freedom that yielded the ‘impurity’ spin. This fact leads to an additional self-consistency condition that supplements the solution of the impurity model; it is their combination which then yields the mean-field predictions for the quantum lattice model.

As we will see below, the resulting mean-field theory of the lattice model has strong similarities to the classical gravity theories of AdS_{D+1} at non-zero μ which were outlined in Section 1. In the latter theories, the geometry factorizes to $\text{AdS}_2 \times R^{D-1}$ at low energies; it is this factorization which will be seen to be related to the ‘impurity’ + ‘environment’ factorization of the condensed matter mean-field theories. A notable fact is that the factorization is motivated in the gravity theory from a very different reasoning from that in the condensed matter model. Thus the appearance of similar result in two very different approximations is quite surprising, and indicates a robustness of the theory that should be well worth understanding better.

Returning to the condensed matter perspective, let us describe the mapping from lattice models to self-consistent quantum impurity models. Such a mapping has been carried out for a variety of models [64, 70, 71, 72, 73, 74, 75], which are all versions of

the Kondo lattice Hamiltonian

$$H = H_J + \sum_k \epsilon_k c_{k\alpha}^\dagger c_k^\alpha + \frac{J_K}{2} \sum_i \hat{S}_i^a c_{i\alpha}^\dagger (\sigma^a)_{\beta}^{\alpha} c_i^\beta. \quad (29)$$

Here H_J is a quantum spin model just as in Eq. (1). To these localized spins, we have added mobile conduction electrons: c_k^α is the Fourier transform of the electron operator c_i^α on site i and ϵ_k is the electron dispersion. Finally, J_K is the Kondo exchange coupling between the conduction electrons and the spins.

The most direct mean-field analyses of H appear in lattice models with random infinite-range exchange interactions [64, 74]: the J_{ij} being independent Gaussian random variables with zero mean. Note that the disorder is ‘quenched’ *i.e.* each J_{ij} is independent of time, but is chosen at random from a Gaussian distribution. However, similar mean field equations also arise in the large spatial dimension limit of non-random Kondo lattice models [71, 72, 73, 74, 75].

Such mean field models yield solutions corresponding to the two classes of non-magnetic metallic states expected in Kondo lattice models [79, 80, 81]:

- A Fermi liquid (FL) with a ‘large Fermi surface’, which can be viewed as arising from the RG flow to large J_K . Here the electrons forming the \hat{S}_i^a spins, along with the c_i^α electrons, become part of the Luttinger count which determines the volume enclosed by the Fermi surface.
- A fractionalized Fermi liquid (FFL or FL*) with a ‘small Fermi surface’, in which the effects of J_K can be accounted for perturbatively. Here the \hat{S}^a spins form a spin liquid, while the conduction electrons form a Fermi surface whose volume counts only the density of the c_i^α conduction electrons.

Let us describe the mean-field structure of the FL* phase so obtained [64, 74]. It was found that correlations of the spin liquid sector of this phase are described by a quantum impurity theory which is identical to Eq. (14). However, this theory now has an additional self-consistency condition that the ‘environment’ spins φ^a are the same as the impurity spin \hat{S}^a : thus the two-point correlator of φ^a which appears in Eqs (13) and (14) should be proportional to the two-point correlator of \hat{S}^a in Eq. (26): *i.e.*

$$D(\tau) \propto C(\tau). \quad (30)$$

It was further shown that a solution of this self-consistency relation is only possible if the spectrum is gapless and has a power-law form. Then from Eqs. (19) and (27) we have the exponent relation

$$h = \gamma. \quad (31)$$

Combining this with the exact relation in Eq. (28) for the quantum impurity model, we obtain [64]

$$h = 1. \quad (32)$$

This is the value that corresponds to ‘marginal Fermi liquid’ behavior [76], as we will see shortly. The same value of h is obtained in large dimension solution of non-random

lattice models [73, 75]. The gravity approach does not fix the value of such exponents: they are related to the mass of fermions in AdS_{D+1} and these are free parameters in present analyses.

We have now described a mean-field state, which applies to both random and non-random models, which is a critical spin liquid. It can be viewed as a large dimension analog of the ‘Spin Bose Metal’ [77, 78], in that it can be written as a theory of bosons at non-zero density which do not Bose condense, but form a gapless liquid [64]. This mean-field state has a non-zero ground state entropy density, which descends from the entropy of the quantum impurity problem. It is this critical spin liquid which has been proposed to realize the low energy physics of AdS_{D+1} in a non-zero chemical potential μ .

The main claim of Ref. [29] was that the theory of the holographic metal realizes the FL^* phase, in a situation in which the \hat{S}_i^a spins are in the critical spin liquid state described in Section 3.1. The evidence for this claim so far is the gapless conformal form of the spin and fermion correlations in Section 3.1, the connection with AdS_2 of the impurity models in Section 3, and the non-zero ground state entropy density. Additional evidence comes from the self-energy of the conduction electrons, c_i^α . We can compute the conduction electron self-energy $\Sigma_c(\omega_n)$ in the FL^* phase by perturbation theory in J_K ; at second order in J_K we have the contribution [74]

$$\begin{aligned} \Sigma_c(\omega_n) &\propto J_K^2 T \sum_{\epsilon_n} \sum_k \frac{1}{i(\omega_n + \epsilon_n) - \epsilon_k} C(\epsilon_n) \\ &\propto -i\pi N_0 J_K^2 T \sum_{\epsilon_n} \text{sgn}(\omega_n + \epsilon_n) C(\epsilon_n) \\ &= J_K^2 T^h \Psi(\omega_n/T) \end{aligned} \tag{33}$$

where $\Psi(\omega_n/T)$ is a scaling function, and N_0 is the density of conduction electron states at the Fermi level. Provided $h < 2$, this is a non-Fermi liquid form of the electron self energy. It is also the same result as that obtained for the holographic metal in Refs. [22, 26, 30]. A similar analysis can be done by the present methods of the transport properties [74], and again agreement is obtained with the holographic results [22, 26, 30].

For the exponent $h = 1$ obtained [64] by the self-consistency requirement on lattice models in Eq. (32), the self energy in Eq. (33) has the marginal Fermi liquid form [76].

5. Conclusion

In conclusion, we note that theory of the FL^* phase reviewed here [64, 74], and obtained by the mapping to quantum impurity models, is an attractive candidate for describing strange metal phases. At the semi-phenomenological level, it does provide a satisfactory description of experimental observations. It is indeed quite remarkable and surprising that a similar theory has now appeared from the very different starting point of the AdS/CFT correspondence.

However, the quantum-impurity description of the FL^* phase is not believed to

be complete [79, 80]. The spin liquid constituent of this phase should have emergent gauge excitations (like the A_μ gauge field in Eq. (7)), and these are surely essential for a complete description of spatial correlations. So it would be interesting to find the appropriate gauge fields in the holographic theory. In this context, the recent work of Nickel and Son is notable [82], as they argue that theory of the holographic metal may indeed be missing such emergent gauge fields.

We also note the interesting recent work of Kachru *et al.* [59] showing a transition from a FL* phase to a FL phase using string theory.

Acknowledgements

I am grateful to Sean Hartnoll and Shamit Kachru for very helpful discussions. This research was supported by the National Science Foundation under grant DMR-0757145 and by a MURI grant from AFOSR.

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