ARTICLE OPEN Check for updates Emergence of nodal Bogoliubov quasiparticles across the transition from the pseudogap metal to the *d*-wave superconductor

Maine Christos $(1)^{1}$ and Subir Sachdev $(1)^{1}$

We model the pseudogap state of the hole- and electron-doped cuprates as a metal with hole and/or electron pocket Fermi surfaces. In the absence of long-range antiferromagnetism, such Fermi surfaces violate the Luttinger requirement of enclosing the same area as free electrons at the same density. Using the Ancilla theory of such a pseudogap state, we describe the onset of conventional *d*-wave superconductivity by the condensation of a charge *e* Higgs boson transforming as a fundamental under the emergent SU(2) gauge symmetry of a background π -flux spin liquid. In all cases, we find that the *d*-wave superconductor has gapless Bogoliubov quasiparticles at 4 nodal points on the Brillouin zone diagonals with significant velocity anisotropy, just as in the BCS state. This includes the case of the electron-doped pseudogap metal with only electron pockets centered at wavevectors (π , 0), (0, π), and an electronic gap along the zone diagonals. Remarkably, in this case, too, gapless nodal Bogoliubov quasiparticles emerge within the gap at 4 points along the zone diagonals upon the onset of superconductivity.

npj Quantum Materials (2024)9:4; https://doi.org/10.1038/s41535-023-00608-0

INTRODUCTION

The remarkable phase diagram of the cuprates^{1–4} has inspired an outpouring of theoretical and experimental work to explain their highly exotic phenomenology. Among the most extensively studied phases are the pseudogap metal, a phase characterized by a carrier density that deviates from the expectations required by Luttinger's theorem for a conventional Fermi liquid^{5–7} and *d*-wave superconductivity which sets in at lower temperatures as an instability of the pseudogap phase^{8,9}.

However, despite years of theoretical and experimental progress, a clear understanding of how superconductivity emerges from the experimentally observed pseudogap parent state and its associated small Fermi surface or Fermi arcs remains lacking. This work seeks to provide some basic answers as to what the experimental signatures of the transition from the pseudogap to superconductivity are. Although the pseudogap phase and its associated violation of Luttinger's theorem have been studied most extensively in the hole-doped cuprates, recent photoemission experiments in the electron-doped cuprates have provided evidence for a reconstructed Fermi surface at dopings where long-range antiferromagnetic order is believed to be absent^{10–12}. The pairing in the electron-doped case is also believed to be d-wave¹³. We will, therefore, separately consider both the electron-doped and hole-doped cases in this work.

A number of works^{14–24} have developed a model of the pseudogap metal in which the violation of the Luttinger theorem is associated with zeros of the electron Green's functions. Here, we view these zeros as a signal of the existence of an additional sector of neutral spinon excitations, which are required by non-perturbative extensions of the Luttinger theorem^{25,26}. As we will see below (and has also been argued earlier²⁷), a full treatment of the spinon sector is essential in understanding how the nodal Bogoliubov quasiparticles in the *d*-wave superconductor emerge from the pseudogap metal.

We will employ a theory²⁸ of the pseudogap metal with fermionic spinons coupled to an SU(2) gauge field moving in a background of π -flux²⁹⁻³². The fermionic spinons are coupled to physical electrons, which carry the doping via a charge e boson $B^{30,32,33}$, which transforms under the same gauge SU(2) symmetry as the spinons. In the hole-doped case, due to the presence of the spin liquid, the normal state electron Fermi surface will have pockets associated with hole density *p* rather than the free electron hole-density value $1 - p^{30,32-37}$. When *B* condenses, the gauge symmetry is fully broken, and various symmetry-breaking orders, including *d*-wave superconductivity and charge order, can be inherited by the electrons. Within this approach, superconductivity and charge order are treated on equal footing and can be viewed as low-temperature, competing instabilities of a fractionalized Fermi liquid (FL*) pseudogap phase. (Previous work^{30,32,33} has considered the condensation of such a boson from an incoherent normal state which does not have pocket Fermi surfaces of electrons or holes and with a U(1) staggered flux spin liquid rather than the π -flux spin liquid. The staggered flux spin liquid has a charge e boson whose condensation leads to dwave superconductivity but not the additional possibility of charge order; moreover, it has a trivial monopole instability³⁸, so it is unlikely to have a significant regime of stability).

In this work, we will consider the transition from the pseudogap phase with electron and/or hole pockets and a π -flux spin liquid to a conventional *d*-wave superconductor. We will compute electronic observables in the superconducting phase via the framework of the Ancilla model^{39–45}. While earlier work²⁷ has considered superconductivity as a similar confinement transition from a phenomenological model of the pseudogap Fermi surfaces, the Ancilla model has the benefit of providing a microscopic model for the complete fermion dispersion in the Brillouin zone, which emerges in an approximation of the Hubbard model⁴¹.

The rest of this paper will be organized as follows. In the Section "Ancilla model," we will introduce the Ancilla model, and in the

¹Department of Physics, Harvard University, Cambridge, MA 02138, USA. 쩓email: mainechristos@g.harvard.edu; sachdev@g.harvard.edu

Section "Mean-field theory," its mean-field representation, which we will use to compute various electronic properties of the pseudogap normal state and *d*-wave superconductor.

In Section "Superconductor spectra with hole-doping", we will describe the phenomenology of our theory on the hole-doped side, where the pseudogap normal state is captured by hole-like pockets enclosing a volume associated with hole density p. We will show in the framework of the Ancilla model that the hole pocket Fermi surfaces of the pseudogap undergo a transition first to a *d*-wave superconductor with 12 nodes, and then to 4 nodes as the strength of the superconducting pairing is increased. We will also compute how the Fermi velocity and v_{Δ} of these nodes evolve with the superconducting pairing strength.

In Section "Superconductor spectra with electron-doping," we will turn our focus to the electron-doped side of the cuprate phase diagram. In this case, the normal state Fermi surface will be an FL^{*} state with either (*i*) only electron-like pockets in the anti-nodal region of the Brillouin zone centered at wavevectors (0, π) and (π , 0) or (*ii*) both anti-nodal electron-like pockets and hole-like pockets in the nodal region^{11,46–52}. Perhaps surprisingly, we find that even in the first case where the normal state Fermi surface only exists at the anti-nodal region and any states in the nodal region are fully gapped, a condensed superconducting pairing will immediately lead to the re-emergence of nodes near ($\frac{\pi}{2}, \frac{\pi}{2}$), while the anti-nodal region is gapped out by the pairing. We will also explore how the velocities of these nodes evolve as a function of the *B* condensate in the electron-doped case.

RESULTS

Ancilla model

In this section, we will discuss the model we use to compute the spectral properties of the *d*-wave superconductor. The model is the Ancilla model of ref. ³⁹, which has been shown to have all the ingredients needed to reproduce photo-emission data in both the pseudogap and Fermi liquid regime^{41,42}. A schematic of the model is shown in Fig. 1.

The general idea is to map the low energy physics of the singleband Hubbard-like model of the *c* layer to a model with free electrons on the *c* layer coupled to a bilayer square lattice antiferromagnet of the S_1 and S_2 layers. But we emphasize that the S_1 and S_2 layers are just *ancilla qubits*, i.e., they are useful quantum degrees of freedom employed at intermediate stages to obtain a wavefunction with non-trivial entanglement on the *c* layer



Fig. 1 Schematic of Ancilla model. We show a schematic of the Ancilla model **a** described in Section "Ancilla model". We have colored the *c* electrons blue, the first layer of spins S_1 green, and the second layer of spins S_2 red. The green and red bonds in the first and second layers of spins represent Heisenberg exchange of spins represents antiferromagnetic coupling J_{\perp} , and the dotted lines between the physical *c* electrons and the first layer of spins represent Kondo coupling J_K . The lines connecting the *c* electron sites denote the *c* electrons' hopping. The red, green, and blue coloring corresponding to the third, second, and first layers, respectively, will be kept consistent throughout the paper. The original model is a Hubbard on the *c* layer, and the Hubbard *U* has been canonically transformed away by adding two Ancilla layers of a bilayer antiferromagnet⁴².

alone. There is a passing similarity to "hidden layers" in current models of machine learning wavefunctions⁵³, but it is important that in our case, the hidden layers are quantum, not classical, as that is the key to obeying the Luttinger–Oshikawa constraints on Fermi surface volumes. An explicit example of an ancilla wavefunction for the *FL*^{*} pseudogap metal was presented in ref.³⁹: a product of Slater determinants on the top two and bottom layers was projected onto the physical top *c* layer by taking the overlap with rung singlets on the **S**₁ and **S**₂ layers. In the analytical theories^{39,40,45} of the Ancilla model, the projection is performed by emergent gauge fields. In the following, we shall not consider the influence of the emergent gauge fields as they are hugged in all the phases considered here. Consequently, we maintain that the low-energy dispersions of the fermionic excitations described below will apply also to the single-band Hubbard model.

An earlier work³⁶ formally obtained the same FL* pseudogap metal directly in the single-band model while using the bosonic spinon \mathbb{CP}^1 formulation of the spin liquid. But this single-band formulation does not yield a theory for a transition from FL* to the large doping Fermi liquid, as is possible in the Ancilla model^{39,40}. Moreover, the bosonic spinon spin liquid is dual to the fermionic spinon π -flux spin liquid⁵⁴, and the mapping of the bosonic spinon FL* formulation to the fermionic spinon theory used here becomes evident in the Ancilla model. The fermionic spinons are essential to obtaining the FL* electronic spectrum across the Brillouin zone and not just near the zone diagonals. The fermionic spinons are also key to obtaining 4 nodal points in the *d*-wave superconductor (in contrast to the 8 nodal points in the SC* state obtained by pairing Fermi pockets in the bosonic spinon approach⁵⁵).

As has been discussed in ref. ⁴², we can derive the Ancilla model by an extension of the method used to introduce paramagnons in the theory of correlated metals. We start with single-band Hubbard model

$$H_{U} = -\sum_{i,j} \sum_{\alpha} t_{i,j}^{c} c_{i,\alpha}^{\dagger} c_{j,\alpha} + U \sum_{i} n_{i\uparrow} n_{i\downarrow}$$
(1)

and exactly decouple the Hubbard term by the paramagnon field $\mathbf{\Phi}_i$

$$H_{\boldsymbol{\Phi}} = -\sum_{\boldsymbol{i},\boldsymbol{j}} \sum_{\alpha} t_{\boldsymbol{i},\boldsymbol{j}}^{c} \boldsymbol{c}_{\boldsymbol{i},\alpha}^{\dagger} \boldsymbol{c}_{\boldsymbol{j},\alpha}$$
$$+ \sum_{\boldsymbol{i}} \left[\frac{3}{8U} \boldsymbol{\Phi}_{\boldsymbol{i}}^{2} - \boldsymbol{\Phi}_{\boldsymbol{i}} \cdot \boldsymbol{c}_{\boldsymbol{i},\alpha}^{\dagger} \frac{\boldsymbol{\sigma}^{\alpha\beta}}{2} \boldsymbol{c}_{\boldsymbol{i},\beta} \right].$$
(2)

In the traditional paramagnon method, Φ is treated as a nearly Gaussian field whose correlators are damped by coupling to the Fermi surface of *c*. Here, we identify the Φ paramagnon with the rung-triplet excitation of the S_1 and S_2 layers in Fig. 1; such a triplet excitation is clearly present when J_{\perp} is large. Indeed, the model in Fig. 1 (described explicitly below) can be mapped back to the single-band Hubbard model in (1) via a Schrieffer–Wolff transformation valid for small J_{K}/J_{\perp}^{42} . For other values of J_{K}/J_{\perp} , we need the fluctuations of emergent gauge fields to project out the ancilla layers, but as argued above, we expect such gauge fluctuations to not modify the low energy excitations in the phases considered below.

The Hamiltonian of the Ancilla model in Fig. 1 is:

$$H = -\sum_{ij} \sum_{a} t^{c}_{ij} c^{\dagger}_{i,a} c_{j,a} + H_{c,f_1} + H_{f_1,f_2} + H_{f_1,f_1} + H_{f_2,f_2}$$
(3)

In the above, the two spin flavors of *c* electrons are the physical degrees of freedom that carry the doping and are fixed to have filling:

$$\frac{1}{N}\sum_{i}\sum_{\alpha}c_{i,\alpha}^{\dagger}c_{i,\alpha}=1-p \tag{4}$$

Where in the above *p* denotes the hole doping. H_{c,f_1} denotes a Kondo coupling between the *c* electrons on each side and a layer

of spins S_1 :

$$H_{c,f_{1}} = J_{K} \sum_{i} c_{i,\alpha}^{\dagger} \boldsymbol{\sigma}^{\alpha\beta} c_{i,\beta} \cdot \boldsymbol{S}_{1,i}$$

= $J_{K} \sum_{i} c_{i,\alpha}^{\dagger} \boldsymbol{\sigma}^{\alpha\beta} c_{i,\beta} \cdot f_{1,i,\gamma}^{\dagger} \boldsymbol{\sigma}^{\gamma\delta} f_{1,i,\delta}$ (5)

Where we have chosen to represent the spins of the first layer $S_{1,i}$ with fermionic spinons f_1 , subject to the local constraint:

$$\sum_{a} f_{1,i,a}^{\dagger} f_{1,i,a} = 1$$
 (6)

The term H_{f_1,f_2} describes the antiferromagnetic coupling between the layer of S_1 spins and a second layer of spins labeled S_2 :

$$H_{f_{1},f_{2}} = J_{\perp} \sum_{i} \boldsymbol{S}_{1,i} \cdot \boldsymbol{S}_{2,i}$$

= $J_{\perp} \sum_{i} f_{1,i,\alpha}^{\dagger} \boldsymbol{\sigma}^{\alpha\beta} f_{1,i,\beta} \cdot f_{2,i,\gamma}^{\dagger} \boldsymbol{\sigma}^{\gamma\delta} f_{2,i,\delta}$ (7)

We have introduced a second set of fermionic spinons f_2 to represent the S_2 , which are subject to their own local constraint:

$$\sum_{a} f_{2,i,a}^{\dagger} f_{2,i,a} = 1 \tag{8}$$

The terms H_{f_1,f_1} and H_{f_2,f_2} describe Heisenberg exchange interactions between the spins in the first and second layers, respectively. In order to reproduce the Fermi arcs seen in experiments in the pseudogap phase², the model must contain deconfined fractional degrees of freedom in order to violate Luttinger's theorem^{26,56}. We will therefore take H_{f_2,f_2} to be described by the π -flux spin liquid²⁹

$$H_{f_{2},f_{2}} = -it^{f_{2}} \sum_{\langle ij \rangle} f^{\dagger}_{2,i,a} e_{ij} f_{2,j,a} \,. \tag{9}$$

We have written the saddle point π -flux spin liquid in the second layer of spins in the gauge previously used in^{28,54} where $e_{i,i} = -e_{i,i}, e_{i,i+x} = 1, e_{i,i+y} = (-1)^x$.

Mean-field theory

After a mean field decoupling, the model can be written in the following form:

$$H = \sum_{ij} \sum_{a} \left[-t^{c}_{|i-j|} c^{\dagger}_{i,a} c_{j,a} - t^{f_{1}}_{|i-j|} f^{\dagger}_{1,i,a} f_{1,j,a} \right]$$

$$- it^{f_{2}} \sum_{\langle ij \rangle} \sum_{a} f^{\dagger}_{2,i,a} e_{ij} f_{2,j,a}.$$

$$+ \sum_{i} \sum_{a} \left[\Phi \left(c^{\dagger}_{i,a} f_{1,i,a} + f^{\dagger}_{1,i,a} c_{i,a} \right) + iB_{1,i} f^{\dagger}_{2,i,a} f_{1,i,a} - iB_{2,i} e_{a\beta} f_{2,i,a} f_{1,i,\beta} + h.c. \right]$$
(10)

In the above, the chemical potentials μ_c, μ_{f_1} and μ_{f_2} must be adjusted such that Eqs. (4), (6), and (8) are satisfied. In practice, we set the chemical potentials on an 80×80 momentum space grid and set an error threshold of 0.01 for each filling. We also allow for next, next-next, and next-next-next nearest neighbor terms for the c and f_1 electron dispersions. For the hole-doped system, the hopping parameters in⁴¹ were found to best match photoemission data taken in the pseudogap regime⁵⁷. Therefore, for the hole-doped case, we will take $t_{0,1}^{c} = 0.22 \text{ eV}, t_{1,1}^{c} = -0.034 \text{ eV},$ $t_{2,0}^c = 0.036 \text{ eV}$, and $t_{2,1}^c = -0.007 \text{ eV}$ for the *c* electron dispersion and $t_{0.1}^{f_1} = -0.1$ eV, $t_{1.1}^{f_1} = 0.03$ eV, and $t_{2.0}^{f_1} = 0.01$ eV. The above hoppings were fit to photo-emission data, assuming $\Phi = 0.09$ and p = 0.206. We have not included the pairing terms that will appear in a mean-field decoupling of the first layer Heisenberg interactions, as there should be no pairing in the first layer in the pseudogap phase. In the second Ancilla layer, we have taken $t^{f_2} = 0.14 \text{ eV}$. The f_1 and f_2 spinons are coupled via the twocomponent, complex boson $B_i = \begin{pmatrix} B_{1,i} \\ B_{2,i} \end{pmatrix}$ which is a spin singlet under global SU(2) spin rotations. B_i is viewed as the Higgs field in the context of this theory.

Various possible phases of Eq. (10) have been studied in previous work on the Ancilla model^{39,41,42}. The pseudogap phase corresponds to the case where J_K is much larger than J_{\perp} . In this case, *B* is gapped but Φ is condensed and the gauge symmetry of the spin liquid is unbroken. The ground state, in this case, is described by a fractional Fermi liquid (FL^{*}) state, and it is possible to choose parameters such that the electrons will hybridize with the f_1 spinons and form hole-like pockets with associated hole density *p*, where the spectral weight of the *c* electrons is highest on the front-side pocket closest to the center of the Brillouin zone as in the first row of Fig. 2. A Fermi liquid can be realized in the case where J_{\perp} in Eq. (3) is much larger than J_K , which leads to Φ becoming gapped. In this case, the f_1 and f_2 spinons will form singlets at each site, and the *c* electrons will exhibit a conventional Fermi liquid Fermi surface with hole density 1 - p.

In this work, we will consider starting from a normal state where only Φ is condensed such that the Fermi surface has hole density 1 - p and ask how the electronic spectrum evolves as a *d*-wave superconductor set in when *B* condenses on top of this normal state. As discussed in previous work²⁸ for the case of the *π*-flux spin liquid and in^{30,32,33} for the case of the U(1) staggered flux spin liquid, the different ways in which the two components of *B* condense can break different symmetries corresponding to distinct orders, which may be inherited by the physical electrons if Φ is also condensed. Expanding on the two band minima of the *π*-flux dispersion, we have the following expression for the chargon *B* in terms of B_{a+} and B_{a-} , the continuum degrees of freedom associated with the + and – minima of the *π*-flux mean field dispersion:

$$B_{a}(\mathbf{r}) = -B_{a,+}e^{i\pi(x+y)/2} + (1+\sqrt{2})B_{a,-}e^{i\pi(x-y)/2}$$
for x even
(11)

$$B_{a}(\mathbf{r}) = (1 + \sqrt{2})B_{a,+}e^{i\pi(x+y)/2} - B_{a,-}e^{i\pi(x-y)/2}$$
for x odd
(12)

In the above a is a label that runs over Nambu gauge indices. We will focus on the case where B condenses in such a way that a d-wave pairing is inherited by the physical electrons. In this case, the following continuum order parameter will be condensed:

$$\Delta = \epsilon_{ab} B_{a+} B_{b-} \tag{13}$$

We can then choose

$$B_{a+} = \frac{1}{\sqrt{2}}(-b,b) \quad B_{a-} = \frac{1}{\sqrt{2}}(b,b)$$
 (14)

as a mean-field ansatz for a pairing, which will be inherited by the c electrons, and ask how the electronic observables will evolve when b is nonzero.

Superconductor spectra with hole-doping

In this section, we will discuss some qualitative features of the superconductor, which sets in when *B* condenses in a normal state where $\Phi \neq 0$. We show an example in Fig. 2 of the electronic spectral density over the transition of an FL^{*} normal state to a *d*-wave superconductor as *B* is condensed.

There are several features of the electron spectra which are of particular relevance to experiments. One important question we will address is the number of nodes our theory predicts will appear where *B* is condensed, given the experimental evidence for 4 nodes^{3,4} in the Brillouin zone in the hole-doped superconducting state. We will also study the evolution of the velocities v_F and v_{Δ} as *b* becomes nonzero, as well as discuss the phenomenology of the pairing on the electron-doped side of the phase diagram.



Fig. 2 Spectral density in superconductor for hole-doped normal state with hole pockets. We show the electron spectral density (**a**, **c**, **e**) and band structure along a diagonal cut in the Brillouin zone (**b**, **d**, **f**) for different values of *b* and Φ (see Eq. (14) for the definition of *b*). In the band structure plots, the bands are colored with RGB values where blue = *c* electrons, green = *f*₁ fermions, and red = *f*₂ fermions. We have plotted all bands on the Nambu basis, thus the spectrum is always particle-hole symmetric. When *B* = 0 but $\Phi > 0$ (top row), the electron spectral function shows hole-like Fermi pockets in the nodal region of the Brillouin zone formed from hybridization between the *c* and *f*₁ electrons. When *b* is nonzero and a *d*-wave superconducting order is inherited by the *c* fermions (middle and bottom rows), all of the states at the Fermi level are gapped out except for a node on the front side of the original hole pocket of the parent state. The *c* electron velocity perpendicular to the $k_x = k_y$ cut is shown to increase if Φ is made larger (bottom row). Spectral densities are normalized by their maximum, value A_0 . All spectral densities are computed with a lifetime parameter of 0.005*i*. In practice, when plotting dispersion and spectral functions, we use the gauge of³¹, which is manifestly translationally invariant for the *n*-flux spin liquid dispersion, though we note in the case when *B* is not condensed, the spinon bands are not gauge invariant and do not appear in any physical observable.

The first question we will address is how many nodes there are in the superconducting state. We find that, similar to past studies²⁷, the answer to this question depends on the values of b and Φ . There are two possibilities for our chosen parameters, which are depicted in Fig. 3.

If the particle-hole symmetry breaking in the first layer of spinons is taken to be small, then there is a small but finite window of $0 < b < b_c$ where the spectra show 3 nodes in each quadrant of the Brillouin zone, or 12 nodes total. However, the appearance of a window of *b* with 12 nodes results from the particle-hole asymmetry of the f_1 spinon bands, which was found to be small when the next and next-next nearest neighbor hoppings in the second layer were fit to experiment⁴¹. Thus, this feature persists for a very small window of *b* before two of the three nodes in each quadrant of the Brillouin zone annihilate and we are left with a spectrum with 4 nodes, the scenario which is born out in experiments^{3,4}.

Additional plots of the dispersion throughout the Brillouin zone are shown in Supplementary Figs. 1–3. In the above discussion on the number of nodes, we have assumed $t^{f_2} = .14 \text{ eV}$. While the overall magnitude of t^{f_2} should not change the number of nodes, the sign of t^{f_2} will determine which 2 of the 3 nodes along the diagonal annihilate first and, therefore, qualitatively change the mean-field dispersion. Since the case where $t^{f_2} < 0$ seems not to display the universal behavior discussed above, we consider it separately in Supplementary Fig. 4.

We also study the evolution of the Fermi velocities as b is increased for various values of Φ . Our results are shown in Fig. 4.

There will be two independent velocities, v_F which is defined as the velocity parallel to the $k_x = k_y$ contour in the Brillouin zone, and v_{Δ} , the velocity perpendicular to this contour. Since a region of 12 nodes appears for a relatively small window of *b*, we consider the velocity only in the case of the node that falls along the original *c* electron Fermi surface and has the highest overlap



Fig. 3 Number of nodes in a hole-doped superconductor. We show the evolution of the Fermi surface along a diagonal cut through the nodal region of the Brillouin zone, focusing on the region where the hole pocket is present in the normal state. **a** The Fermi surface formed from the hybridized *c* electrons and f_1 spinons with an isolated Dirac cone from the π -flux spin liquid of the f_2 layer. For very small values of *b*, when the f_2 spinons hybridize with the *c* electrons and f_1 spinons, the Fermi surface has 12 nodes, as shown in (**b**). For larger values of *b*, there are 4 nodes on the Fermi surface, as shown in (**c**).



Fig. 4 Node velocities in superconductor for hole-doped normal state. We show the velocities v_F (a) and v_{Δ} (b) as a function of b for different values of Φ . For small b, v_F takes on the Fermi velocity of the normal state Fermi surface at $k_x = k_y$, while v_{Δ} begins at 0 when B = 0 and increases with finite b as the effective pairing grows. For large b, v_F approaches the Fermi velocity of the un-hybridized c electron bands, while v_{Δ} tends to 0.

with the *c* electrons. We compute v_F and v_{Δ} by discretizing diagonal cuts through a quadrant in the Brillouin zone into 80,000 momentum points and performing a least squares fit on the 500 momentum points nearest the node.

The nodal velocity perpendicular to $k_x = k_y, v_{\Delta}$, begins at zero when b = 0. When *B* is condensed, and *b* is small relative to Φ , the superconducting pairing is inherited by the *c* electrons and as a result, v_{Δ} becomes finite and increases with *b* as the effective pairing gaps out any states which are not on the Brillouin zone diagonal. v_{Δ} will continue to increase until *b* is roughly of the same order as Φ where v_{Δ} attains a maximum. When *b* is sufficiently large relative to Φ, v_{Δ} will begin to decrease as the layer of *c* electrons becomes effectively decoupled from the first and second layers of spinons, which are pushed away from the Fermi level. For large enough *b*, the *c* electrons spectral density will resemble the original Fermi surface of the decoupled *c* electrons and v_{Δ} will tend towards zero as *b* increases.

The nodal velocity along $k_x = k_y$, v_{Fr} begins at a finite value defined by the normal state Fermi velocity and monotonically increases with *b* until it saturates in the limit where $b \gg \Phi$ to the value of the Fermi velocity of the decoupled *c* electron bands at the Fermi surface. The ratio of v_{Δ} to v_F is small for all values of *b* as the Fermi velocity v_F originates mostly from the Fermi velocity of the velocity v_{Δ} is 0 in the normal state and is a higher order effect in Φ and *b*.

Superconductor spectra with electron-doping

In this section, we will discuss the spectra of an FL^{*} to SC transition on the electron-doped side of the phase diagram. The principal difference from the hole-doped case is that we now expect instead of having hole-like pockets near $(\frac{\pi}{2}, \frac{\pi}{2})$, we will have either electron pockets in the anti-nodal region of the Brillouin zone near $(0, \pi)$ and $(\pi, 0)$ as in the first row of Fig. 5 or both electron like pockets at the anti-node and hole-like pockets in the nodal region as in the first row of Fig. 6^{46–52}.

For our computations with a normal state with only electron pockets, we will take the same *c* electron hoppings as on the hole-doped side but change the f_1 spinon hoppings since these were previously obtained by fitting photoemission data taken on hole-doped cuprates⁴¹, and have no established values on the electron-doped side. In the first layer of spinons $t_{1,0}^{f_1} = -0.1 \text{ eV}$ as before, but choose the next nearest neighbor hopping $t_{1,1}^{f_1} = -0.07 \text{ eV}$ and set all other hoppings in the second layer to zero. We keep the spin liquid dispersion hopping $t_2^{f_2} = 0.14 \text{ eV}$. We find whether the normal state has only electron pockets at the anti-node or both electron pockets at the anti-node and hole pockets near $(\frac{\pi}{2}, \frac{\pi}{2})$ depends on the value of Φ , with larger Φ gapping out the hole pockets. For our computations of a normal state with both electron and hole pockets like that of Fig. 6, we take $t_{1,1}^{f_1} = -0.06 \text{ eV}$ and $t_{0,2}^{f_1} = 0.02 \text{ eV}$ and all other parameters the same as above.

M. Christos and S. Sachdev



Fig. 5 Spectral density in superconductor for electron-doped normal state with only electron pockets. Evolution of spectral density of *c* electrons (**a**, **d**), dispersion along a diagonal cut through the Brillouin zone (**b**, **e**), and dispersion along a vertical cut through the $(0, \pi)$ (**c**, **f**) for B = 0 (top row) and B > 0 (bottom row). All plots are computed at *electron doping* p = 0.15. We note the nodes for finite b along the diagonal, which were not previously present in the *c* electron density in the normal state.



Fig. 6 Spectral density in superconductor for electron-doped normal state with both electron and hole pockets. We show the evolution of the electron spectral density (\mathbf{a} , \mathbf{d}) and dispersion when superconductivity sets in for the case of a normal state as *positive* electron doping, which has both a hole-like pocket in the nodal region and an electron-like pocket in the anti-nodal region at electron doping p = 0.15. Dispersions are shown for a cut along the diagonal of the Brillouin zone (\mathbf{b} , \mathbf{e}) and for a cut that connects the anti-node and Brillouin zone center (\mathbf{c} , \mathbf{f}). We note that while the second-row electronic spectral density shows a finite electronic spectral weight in other regions of the Brillouin zone compared to the node, all other points except the node have a finite, albeit sometimes small gap.

We will discuss the number of nodes separately for the two types of Fermi surfaces mentioned above, beginning first with the normal state where the only Fermi surfaces are electron pockets at the anti-node. Naively, it might be expected for this case that the *d*-wave superconductor, which will be inherited by the *c* electrons when *B* condenses, will be fully gapped; however, this is not what we observe. For any finite *b*, the electron pockets of the normal state, which appeared in the anti-nodal region, will become fully gapped, as shown in the rightmost column of Fig. 5, but nodes will re-appear along the diagonal in the nodal region of the Brillouin zone as shown in the central column of Fig. 5. These nodes which at b = 0 were associated with the Dirac points of the π -flux spin liquid will hybridize with the c and f_1 bands but cannot be gapped unless an additional symmetry such as spin rotation symmetry is strongly broken. If the above scenario is excluded, there will always be 4 nodes on the diagonal when B is condensed for a normal state with only electron pockets, assuming a positive spin liquid hopping in the gauge we have chosen. We note that for small b, the normal state Fermi surfaces at the anti-node have a gap, which may be



Fig. 7 Node velocities in superconductor for electron-doped normal state. We show v_F (a) and v_{Δ} (b) for positive electron doping as a function of b for several different values of Φ . In the limit of large b, both velocities show the same asymptotic behavior as in the hole-doped case.

very small, and the node that appears for small b initially has a low c electron spectral weight.

For the case of a normal state that has both electron pockets at the anti-node and hole pockets at the node, we observe the same transition from 4 to 12 nodes as b is increased, as we observed in the hole-doped case.

In all of our analysis on the electron-doped side of the phase diagram, we have assumed $t^{f_2} = 0.14$. However, changing the sign of t^{f_2} will result in qualitatively different behavior in the number of nodes, similar to the hole-doped case as shown in Supplementary Figs. 5 and 6. However, as was shown in Appendix 3 of²⁸, only the former sign of t^{f_2} corresponds to a chargon potential, which favors the continuum superconductor ansatz we have taken here.

We also show how v_F and v_{Λ} of the superconductor nodes on the Brillouin zone diagonal evolve for positive electron doping as a function of *b* and for different values of Φ in Fig. 7 for the choice of normal state with only electron pockets. For this choice of the normal state, we study a narrower range of Φ than in the holedoped case since we wish to choose Φ such that the normal state is gapped at the node. The velocities in the case for which there are additional hole pockets are similar to the behavior shown in Fig. 4. Since there is no Fermi surface observable in the electron spectral density in the nodal region at b = 0, both v_F and v_{Δ} immediately jump to a finite value for finite b. For small b, v_F and v_{Δ} are roughly equal as they are essentially just inherited from the π -flux spin liquid's Dirac points, which are isotropic. As b increases, we see v_F will first slowly increase as band repulsion, which flattens the velocity competes with b, but in the limit $b \gg \Phi$, v_F ultimately returns to the Fermi velocity of the decoupled c electrons as in the hole-doped case. Similar to the behavior of v_{Λ} at large b in the hole-doped case, here v_{Δ} decreases as b increases, ultimately tending towards zero when $b \gg \Phi$.

Excitation energy and quasi-particle residue

We also show the excitation energy and quasi-particle residue of the energetically lowest-lying excitation in the superconducting state, as shown in Fig. 8. We plot both quantities along a contour defined by finding the k_x momentum corresponding to where the energy of the lowest-lying excitation is smallest for a given k_y momentum value. Effectively, this contour is well approximated by choosing momenta along the original, decoupled *c*-electron Fermi surface. We choose the first k_y value plotted to be at the location of the single node, which appears in the superconducting state in either the hole-doped or electron-doped case. We may then contrast the behavior of these quantities in the hole-doped and electron-doped cases. While in the electron-doped case, the excitation energy along the above-specified contour increases monotonically as momentum is varied from the location of the node at $(\frac{\pi}{2}, \frac{\pi}{2})$ to the Brillouin zone edge, the electron-doped case shows non-monotonic behavior as one moves along the original *c*-electron Fermi surface, away from the node. In the electron-doped case, we observe a peak in the excitation energy roughly midway between the node and edge of the Brillouin zone, consistent with the behavior observed in¹². In the electron-doped case, the quasi-particle weight Z_k is mostly flat with a very slight dip between the node and anti-node, whereas in the hole-doped case, the quasi-particle weight shows monotonic behavior as momentum is varied away from the node at $(\frac{\pi}{2}, \frac{\pi}{2})$.

DISCUSSION

In this work, we have studied how various electronic observables evolve when the pseudogap metal transitions to a *d*-wave superconductor in the framework of the Ancilla model³⁹.

For a hole-doped normal state and positive spin liquid hopping in our chosen gauge, as discussed in Section "Superconductor spectra with hole-doping", we initially find a *d*-wave superconductor with 12 nodes, and then a transition to 4 nodes as the pairing strength is increased. When the normal state is chosen to reproduce experimental photo-emission data, the regime of 12 nodes is small, and the generic case for large *b* is a superconducting state with 4 nodes. We also found the velocities v_F and v_{Δ} associated with the surviving nodes differ in scale, with v_{Δ} much smaller than v_F for all values of Φ and *b* we have studied and tending towards 0 in the limit where $b \gg \Phi$. It is, therefore, clear that the velocities are not directly related to the spinon velocities in the π -flux phase, which are isotropic, and this is an important difference from earlier work^{30,32,33}.

We have also separately studied the FL^{*} to superconductor transition for the electron-doped case in Section "Superconductor spectra with electron-doping". In this case, we find that a normal state with both electron-like and hole-like pockets leads to the same transition from 12 to 4 nodes as observed in the hole-doped case as a function of *b*. However, surprisingly, we find that in the case with only electron-like pockets near the Brillouin zone edge, the FL^{*} state immediately transitions to a state with 4 nodes along $k_x = k_y$, as shown in Fig. 5. This feature is unique to the electron-doped side of the phase diagram, is striking in that nodes which are not observable in the electron spectral density in the FL^{*} case immediately reappear for any finite *b*. Unlike in the hole-doped case, v_F and v_{Δ} begin with nearly equal values since the surviving node for small *b* is associated with the spin liquid Dirac point



Fig. 8 Quasi-particle excitation and residue. We show the energy of the nearest quasi-particle excitation to the Fermi level (a) and quasi-particle residue at this excitation energy (b) for the hole-doped case using the normal state shown in Fig. 2. We also show the excitation energy (c) and quasi-particle residue (d) for the electron-doped case where there is only an electron pocket at the anti-node using the normal state shown in Fig. 5. Finally, we also show the excitation energy (e) and quasi-particle residue (f) for the normal state of Fig. 6 where there are both electron and hole pockets.

rather than the *c* electron Fermi surface. This aspect of the electron-doped pairing follows as a direct consequence of the mean-field dispersion of the π -flux spin liquid (though this behavior would be the same had we considered another Dirac spin liquid with the same number of nodes such as the U(1) staggered flux spin liquid state). Thus, for electron-doped super-conductors with only electron pockets in the normal state, it is reasonable to state that nodal Bogoliubov quasiparticles are remnants of Dirac spinons made visible by the onset of superconductivity.

All our analysis was carried out for a pseudogap metal without a long-range antiferromagnetic order. However, the appearance of

antiferromagnetic order at low temperatures within the superconducting phase (as is the case in the electron-doped cuprates) should not invalidate any of our computations, and so we believe our results should continue to apply.

Recent numerical studies^{58,59} of an electron-doped t-J model found robust d-wave superconductivity, and the authors speculated that their d-wave superconductor was fully gapped. From our analysis here, we maintain that a conventional d-wave superconductor, with pairing strength not as large as the Fermi energy, always has 4 nodal points along the zone diagonals. A fully gapped d-wave superconductor requires some additional features and can be reached via the following routes:

- (i) We start from an FL* metal with electron pockets and then pair the electron pockets. At this point, the f_2 Ancilla spin liquid is still 'alive' in the superconducting state, and so such a fully gapped *d*-wave superconductor is an SC* state. Furthermore, the π -flux spin liquid is ultimately unstable⁶⁰, and this implies that a π -flux-SC* state is not stable.
- Starting from a conventional *d*-wave superconductor with 4 nodal points, the onset of strong, co-existing antiferromagnetic order can gap out the nodes when they annihilate in pairs across the magnetic Brillouin zone boundary.
- (iii) Finally, if the pairing interaction becomes as large as the Fermi energy in a conventional *d*-wave superconductor, the four nodal points can meet at the origin (or at (π, π)) and annihilate with each other.

It appears unlikely to us that any of these 3 routes apply to the study in refs. ^{58,59}, and so we believe their superconductor does have 4 nodal points at low temperatures and possibly only electron pockets in the normal state.

In summary, our work has provided testable predictions of what signatures conventional *d*-wave superconductivity will have if it originates from a pseudogap phase containing fractional degrees of freedom described by the π -flux spin liquid. It would be interesting to extend the approach taken in this work to capture other relevant phases in the under-doped cuprates, such as charge order⁶¹. We also note that for the case of superconductivity, the quantities we computed will not necessarily be different among different Dirac spin liquids⁶². It is, therefore, interesting to conceive experimental tests that may be capable of distinguishing between different FL^{*} normal states. We leave these possibilities to future work.

METHODS

Computing spectral densities

All plots are computed from a tight-binding implementation of the model in Eq. (3) in momentum space using the parameters for hoppings and dopings mentioned throughout the text. Chemical potentials in each layer of the Hamiltonian described in Eq. (3) are determined by doping for each set of parameters within an error threshold of 0.01 via the bisection method. An 80×80 grid in momentum space is used to fix the chemical potentials. Spectral densities are computed as:

$$A(\boldsymbol{\omega}, \boldsymbol{k}) = -\frac{1}{\pi} \ln \left[G_{cc}(\boldsymbol{\omega}, \boldsymbol{k}) \right]$$
(15)

All spectral functions are computed on a quarter of the Brillouin zone with a 200×200 grid. A lifetime parameter of 0.005 eV is used when computing spectral densities.

Computing quasiparticle weight

The quasiparticle weight $Z_{\mathbf{k}}$ is computed by first computing the inverse of the greens function $G^{-1}(\omega, \mathbf{k}) = \omega - H(\mathbf{k})$, which is then diagonalized by a unitary transformation $U_{\mathbf{k}}$ such that $U_{\mathbf{k}}^{\dagger}H(\mathbf{k})U_{\mathbf{k}} = G_D^{-1}(\omega, \mathbf{k})$ where $G_D^{-1}(\omega, \mathbf{k})$ is diagonal. For an excitation at positive energy $\Delta_{\mathbf{k}}$, we then compute the quasiparticle residue as:

$$Z_{k} = U_{c\uparrow,a} U_{a,c\uparrow}^{\dagger}$$
(16)

where *a* labels the eigenvector of U_k corresponding to energy – Δ_{kr} , such that the quasiparticle weight is always computed at negative energies.

DATA AVAILABILITY

All of the code which generated the data used during this study is made available at the following publicly accessible repository: https://zenodo.org/records/10076755.

Received: 6 August 2023; Accepted: 23 November 2023; Published online: 05 January 2024

REFERENCES

- Keimer, B., Kivelson, S. A., Norman, M. R., Uchida, S. & Zaanen, J. From quantum matter to high-temperature superconductivity in copper oxides. *Nature* 518, 179–186 (2015).
- Proust, C. & Taillefer, L. The remarkable underlying ground states of cuprate superconductors. Annu. Rev. Condens. Matter Phys. 10, 409–429 (2019).
- Damascelli, A., Hussain, Z. & Shen, Z.-X. Angle-resolved photoemission studies of the cuprate superconductors. *Rev. Mod. Phys.* 75, 473–541 (2003).
- Vishik, I. M. et al. ARPES studies of cuprate Fermiology: superconductivity, pseudogap and quasiparticle dynamics. *Annu. Rev. Condens. Matter Phys.* 12, 105008 (2010).
- Sobota, J. A., He, Y. & Shen, Z.-X. Angle-resolved photoemission studies of quantum materials. *Rev. Mod. Phys.* 93, 025006 (2021).
- Yang, H.-B. et al. Reconstructed Fermi surface of underdoped Bi₂Sr₂CaCu₂O_{8+δ} cuprate superconductors. *Phys. Rev. Lett.* **107**, 047003 (2011).
- Shen, K. M. et al. Nodal quasiparticles and antinodal charge ordering in Ca_{2-x-} Na_xCuO₂Cl₂. Science **307**, 901–904 (2005).
- Tsuei, C. C. & Kirtley, J. R. Pairing symmetry in cuprate superconductors. *Rev. Mod. Phys.* **72**, 969–1016 (2000).
- Ding, H. et al. Spectroscopic evidence for a pseudogap in the normal state of underdoped high-T_c superconductors. *Nature* **382**, 51–54 (1996).
- Helm, T. et al. Magnetic breakdown in the electron-doped cuprate superconductor Nd_{2-x}Ce_xCuO₄: the reconstructed Fermi surface survives in the strongly overdoped regime. *Phys. Rev. Lett.* **105**, 247002 (2010).
- He, J. et al. Fermi surface reconstruction in electron-doped cuprates without antiferromagnetic long-range order. *Proc. Natl Acad. Sci. USA* **116**, 3449–3453 (2019).
- 12. Xu, K.-J. et al. Bogoliubov quasiparticle on the gossamer Fermi surface in electron-doped cuprates. *Nat. Phys.* **19**, 1834–1840 (2023).
- Horio, M. et al. *d*-wave superconducting gap observed in protect-annealed electron-doped cuprate superconductors Pr_{1.3-x}La_{0.7}Ce_xCuO₄. *Phys. Rev. B* **100**, 054517 (2019).
- Dzyaloshinskii, I. Some consequences of the Luttinger theorem: the Luttinger surfaces in non-Fermi liquids and Mott insulators. *Phys. Rev. B* 68, 085113 (2003).
- Stanescu, T. D. & Kotliar, G. Fermi arcs and hidden zeros of the Green function in the pseudogap state. *Phys. Rev. B* 74, 125110 (2006).
- Berthod, C., Giamarchi, T., Biermann, S. & Georges, A. Breakup of the Fermi surface near the Mott transition in low-dimensional systems. *Phys. Rev. Lett.* 97, 136401 (2006).
- 17. Yang, K.-Y., Rice, T. M. & Zhang, F.-C. Phenomenological theory of the pseudogap state. *Phys. Rev. B* **73**, 174501 (2006).
- Sakai, S., Motome, Y. & Imada, M. Evolution of electronic structure of doped Mott insulators: reconstruction of poles and zeros of Green's function. *Phys. Rev. Lett.* **102**, 056404 (2009).
- Robinson, N. J., Johnson, P. D., Rice, T. M. & Tsvelik, A. M. Anomalies in the pseudogap phase of the cuprates: competing ground states and the role of umklapp scattering. *Rep. Progr. Phys.* 82, 126501 (2019).
- Skolimowski, J. & Fabrizio, M. Luttinger's theorem in the presence of Luttinger surfaces. *Phys. Rev. B* 106, 045109 (2022).
- 21. Fabrizio, M. Spin-liquid insulators can be Landau's Fermi liquids. *Phys. Rev. Lett.* **130**, 156702 (2023).
- 22. Wagner, N. et al. Mott insulators with boundary zeros. *Nat. Commun.* 14, 7513 (2023).
- Zhao, J., Mai, P., Bradlyn, B. & Phillips, P. Failure of topological invariants in strongly correlated matter. *Phys. Rev. Lett.* **131**, 106601 (2023).
- Peralta Gavensky, L., Sachdev, S. & Goldman, N. Connecting the many-body Chern number to Luttinger's theorem through Středa's formula. *Phys. Rev. Lett.* 131, 236601 (2023).
- Oshikawa, M. Topological approach to Luttinger's theorem and the Fermi surface of a Kondo lattice. *Phys. Rev. Lett.* 84, 3370 (2000).
- Senthil, T., Vojta, M. & Sachdev, S. Weak magnetism and non-Fermi liquids near heavy-fermion critical points. *Phys. Rev. B* 69, 035111 (2004).
- 27. Chatterjee, S. & Sachdev, S. Fractionalized Fermi liquid with bosonic chargons as a candidate for the pseudogap metal. *Phys. Rev. B* **94**, 205117 (2016).
- Christos, M. et al. A model of *d*-wave superconductivity, antiferromagnetism, and charge order on the square lattice. *Proc. Nat. Acad. Sci. USA* **120**, e2302701120 (2023).
- Affleck, I. & Marston, J. B. Large-*n* limit of the Heisenberg-Hubbard model: implications for high-*T_c* superconductors. *Phys. Rev. B* 37, 3774–3777 (1988).

- 10
- Wen, X.-G. & Lee, P. A. Theory of underdoped cuprates. *Phys. Rev. Lett.* 76, 503–506 (1996).
- 31. Wen, X.-G. Quantum orders and symmetric spin liquids. *Phys. Rev. B* **65**, 165113 (2002).
- Lee, P. A., Nagaosa, N. & Wen, X.-G. Doping a Mott insulator: Physics of hightemperature superconductivity. *Rev. Mod. Phys.* 78, 17 (2006).
- Lee, P. A., Nagaosa, N., Ng, T.-K. & Wen, X.-G. SU(2) formulation of the t-J model: application to underdoped cuprates. *Phys. Rev. B* 57, 6003–6021 (1998).
- Sachdev, S. Quantum phases of the Shraiman-Siggia model. Phys. Rev. B 49, 6770–6778 (1994).
- Kaul, R. K., Kim, Y. B., Sachdev, S. & Senthil, T. Algebraic charge liquids. Nat. Phys. 4, 28–31 (2007).
- Qi, Y. & Sachdev, S. Effective theory of Fermi pockets in fluctuating antiferromagnets. *Phys. Rev. B* 81, 115129 (2010).
- Mei, J.-W., Kawasaki, S., Zheng, G.-Q., Weng, Z.-Y. & Wen, X.-G. Luttinger-volume violating Fermi liquid in the pseudogap phase of the cuprate superconductors. *Phys. Rev. B* 85, 134519 (2012).
- Song, X.-Y., He, Y.-C., Vishwanath, A. & Wang, C. From spinon band topology to the symmetry quantum numbers of monopoles in Dirac spin liquids. *Phys. Rev. X* 10, 011033 (2020).
- Zhang, Y.-H. & Sachdev, S. From the pseudogap metal to the Fermi liquid using ancilla qubits. *Phys. Rev. Res.* 2, 023172 (2020).
- Zhang, Y.-H. & Sachdev, S. Deconfined criticality and ghost Fermi surfaces at the onset of antiferromagnetism in a metal. *Phys. Rev. B* 102, 155124 (2020).
- 41. Mascot, E. et al. Electronic spectra with paramagnon fractionalization in the single-band Hubbard model. *Phys. Rev. B* **105**, 075146 (2022).
- Nikolaenko, A., Tikhanovskaya, M., Sachdev, S. & Zhang, Y.-H. Small to large Fermi surface transition in a single band model, using randomly coupled ancillas. *Phys. Rev. B* **103**, 235138 (2021).
- Nikolaenko, A., von Milczewski, J., Joshi, D. G. & Sachdev, S. Spin density wave, Fermi liquid, and fractionalized phases in a theory of antiferromagnetic metals using paramagnons and bosonic spinons. *Phys. Rev. B* 108, 045123 (2023).
- Zhou, B. & Zhang, Y.-H. Ancilla wavefunctions of Mott insulator and pseudogap metal through quantum teleportation. Preprint at *arXiv* https://doi.org/10.48550/ arXiv.2307.16038 (2023).
- Sachdev, S. Quantum Phases of Matter (Cambridge University Press, Cambridge, UK, 2023).
- Armitage, N. P. et al. Doping dependence of an n-type cuprate superconductor investigated by angle-resolved photoemission spectroscopy. *Phys. Rev. Lett.* 88, 257001 (2002).
- Matsui, H. et al. Evolution of the pseudogap across the magnet-superconductor phase boundary of Nd_{2-x}Ce_xCuO₄. *Phys. Rev. B* **75**, 224514 (2007).
- 48. Song, D. et al. Electron number-based phase diagram of $Pr_{1-x}LaCe_xCuO_{4-\delta}$ and possible absence of disparity between electron- and hole-doped cuprate phase diagrams. *Phys. Rev. Lett.* **118**, 137001 (2017).
- 49. Li, Y. et al. Hole pocket–driven superconductivity and its universal features in the electron-doped cuprates. *Sci. Adv.* **5**, eaap7349 (2019).
- Kartsovnik, M. V. et al. Fermi surface of the electron-doped cuprate superconductor Nd_{2-x}Ce_xCuO₄ probed by high-field magnetotransport. *Annu. Rev. Condens. Matter Phys.* 13, 015001 (2011).
- 51. Breznay, N. P. et al. Interplay of structure and charge order revealed by quantum oscillations in thin films of $Pr_2CuO_{4\pm\delta}$. *Phys. Rev. B* **100**, 235111 (2019).
- Helm, T. et al. Correlation between Fermi surface transformations and superconductivity in the electron-doped high-T_c superconductor Nd_{2-x}Ce_xCuO₄. *Phys. Rev. B* **92**, 094501 (2015).
- Moreno, J. R., Carleo, G., Georges, A. & Stokes, J. Fermionic wave functions from neural-network constrained hidden states. *Proc. Natl Acad. Sci. USA* 119, e2122059119 (2022).
- Wang, C., Nahum, A., Metlitski, M. A., Xu, C. & Senthil, T. Deconfined quantum critical points: symmetries and dualities. *Phys. Rev. X* 7, 031051 (2017).
- 55. Moon, E. G. & Sachdev, S. Underdoped cuprates as fractionalized Fermi liquids: transition to superconductivity. *Phys. Rev. B* **83**, 224508 (2011).
- Senthil, T., Sachdev, S. & Vojta, M. Fractionalized Fermi liquids. *Phys. Rev. Lett.* 90, 216403 (2003).
- He, R.-H. et al. From a single-band metal to a high-temperature superconductor via two thermal phase transitions. *Science* **331**, 1579–1583 (2011).

- Jiang, H.-C. & Kivelson, S. A. High temperature superconductivity in a lightly doped quantum spin liquid. *Phys. Rev. Lett.* **127**, 097002 (2021).
- Jiang, H.-C., Kivelson, S. A. & Lee, D.-H. Superconducting valence bond fluid in lightly doped 8-leg t-J cylinders. *Phys. Rev. B* 108, 054505 (2023).
- Zhou, Z., Hu, L., Zhu, W. & He, Y.-C. The SO(5) deconfined phase transition under the fuzzy sphere microscope: approximate conformal symmetry, pseudo-criticality, and operator spectrum. Preprint at *arXiv* https://doi.org/10.48550/ arXiv.2306.16435 (2023).
- Comin, R. & Damascelli, A. Resonant X-ray scattering studies of charge order in cuprates. Annu. Rev. Condens. Matter Phys. 7, 369–405 (2016).
- Shackleton, H., Thomson, A. & Sachdev, S. Deconfined criticality and a gapless Z₂ spin liquid in the square-lattice antiferromagnet. *Phys. Rev. B* **104**, 045110 (2021).

ACKNOWLEDGEMENTS

We thank Patrick Lee for probing questions on previous work²⁸, Z.-X. Shen, thank you for informing us about ongoing photoemission experiments in the electron-doped cuprates¹². Steve Kivelson, thank you for the discussions on refs. ^{58,59}, and our collaborators Zhu-Xi Luo, Henry Shackleton, Ya-Hui Zhang, and Mathias Scheurer on ref. ²⁸. We also thank Chenyuan Li and Ilya Esterlis for helpful discussions on related work and Alex Nikolaenko for helpful discussions on⁴¹. This research was supported by the U.S. National Science Foundation grant No. DMR-2245246 and by the Simons Foundation (651440, S.S.).

AUTHOR CONTRIBUTIONS

Both authors formulated the research and wrote the paper. M.C. performed the numerical computations.

COMPETING INTERESTS

The authors declare no competing interests.

ADDITIONAL INFORMATION

Supplementary information The online version contains supplementary material available at https://doi.org/10.1038/s41535-023-00608-0.

Correspondence and requests for materials should be addressed to Maine Christos or Subir Sachdev.

Reprints and permission information is available at http://www.nature.com/ reprints

Publisher's note Springer Nature remains neutral with regard to jurisdictional claims in published maps and institutional affiliations.

Open Access This article is licensed under a Creative Commons Attribution 4.0 International License, which permits use, sharing, adaptation, distribution and reproduction in any medium or format, as long as you give appropriate credit to the original author(s) and the source, provide a link to the Creative Commons license, and indicate if changes were made. The images or other third party material in this article are included in the article's Creative Commons license, unless indicated otherwise in a credit line to the material. If material is not included in the article's Creative Commons license and your intended use is not permitted by statutory regulation or exceeds the permitted use, you will need to obtain permission directly from the copyright holder. To view a copy of this license, visit http:// creativecommons.org/licenses/by/4.0/.

© The Author(s) 2024