SCIENTIFIC REPORTS

Received: 22 July 2016 Accepted: 02 February 2017 Published: 22 March 2017

OPEN Entropic Origin of Pseudogap **Physics and a Mott-Slater Transition** in Cuprates

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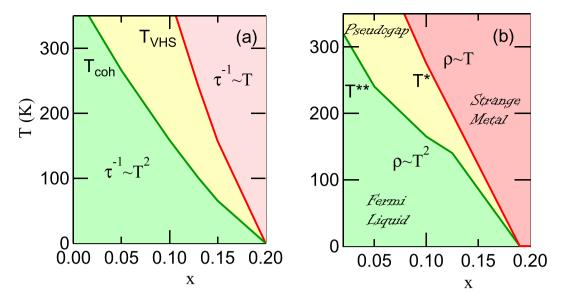
We propose a new approach to understand the origin of the pseudogap in the cuprates, in terms of bosonic entropy. The near-simultaneous softening of a large number of different q-bosons yields an extended range of short-range order, wherein the growth of magnetic correlations with decreasing temperature T is anomalously slow. These entropic effects cause the spectral weight associated with the Van Hove singularity (VHS) to shift rapidly and nearly linearly toward half filling at higher T, consistent with a picture of the VHS driving the pseudogap transition at a temperature $\sim T^*$. As a byproduct, we develop an order-parameter classification scheme that predicts supertransitions between families of order parameters. As one example, we find that by tuning the hopping parameters, it is possible to drive the cuprates across a transition between Mott and Slater physics, where a spin-frustrated state emerges at the crossover.

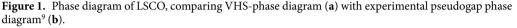
Evidence is growing that the 'pseudogap phase' found in cuprates is home to one or more competing phases, including a variety of stripe, spin-, or charge-density wave (S/CDW) phases¹⁻¹². The CDW phase, in particular, has stimulated considerable interest¹³⁻²². Many of these phases, including superconductivity, seem to appear at temperatures well below the pseudogap temperature T^* , so the exact relation between the pseudogap and these other phases remains elusive. Indeed, the real puzzle is understanding why the pseudogap phase bears so little resemblance to a conventional phase transition. Here we demonstrate that the pseudogap phenomenon arises from strong mode coupling, where a large number of density waves with similar ordering wave-vectors q attempt to soften and condense at the same time. The associated entropy leads to anomalously low transition temperatures and extended ranges of short-range order, characteristic of the pseudogap phase. In an extreme case, this can lead to a condensation bottleneck where long-range order is absent, leaving behind a novel entangled spin-frustrated phase.

To understand how entropy drives short range order, consider the Hubbard model with on-site repulsion U, nearest neighbor hopping energy t, and exchange $J = 4t^2/U$. A simple mean field solution correctly captures an antiferromagnetic ground state with ordering vector $q = (\pi, \pi)$ and gap $2\Delta \sim U$. However, the same calculation yields a Neel temperature $T_N \sim U$, instead of the strong coupling result, $\sim 1/U$. The resolution of this problem is the electronic entropy: repeating the calculation for magnetic order with any other q-vector, including ferromagnetism (q=0) also produces a gap of order U. While the (π, π) state is the ground state, the other states are higher in energy by no more than a factor $\sim J$. Hence, at low T the energy wins out, and the system condenses into its ground state, while at higher temperature entropy wins and the system lowers its free energy by maximizing its entropy, forming a mixture of all these competing phases. Thus, longe-range order is lost at a $T_N \sim l^{23}$. The problem is how to incorporate this entropic effect into a calculation of the phase diagram. Similar problems have arisen in the past. In Overhauser's theory of CDWs in alkali metals²⁴, the q-vectors for all points on a spherical Fermi surface become unstable simultaneously and cannot be handled one at a time. These effects are often referred to as phonon entropy²⁵, or more generally as boson entropy, and can be analyzed via vertex corrections which take proper account of mode coupling^{26,27}. Here we develop a similar theory for the cuprates via a self-consistent renormalization calculation of the vertex corrections. Similar effects are likely to be present in many other families of correlated materials.

The ultimate origin of the bosonic entropy is localization. We provide two examples of this. First, in the above discussion of AFM, large U means single spins tend to localize on individual copper atoms and the effective hopping is renormalized from t- to J-scales, leading to much flatter dispersions, and to small differences in energy

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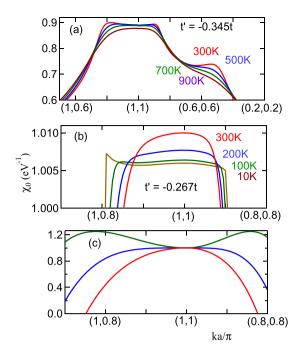
between different configurations of spins. For our second example, we consider a situation where an electron-hole pair –our electronic boson–can form a bound state, an exciton. When excitons are strongly bound, they can condense, creating a form of a CDW or SDW^{28–30}. In the Bose condensation of excitons, the mean-field transition is found to correspond to the temperature at which excitons are formed. However, since the excitons are localized in real space, they are greatly spread in q. When fluctuations are included, the real transition lies at a much lower T, when all excitons condense into the lowest q state.

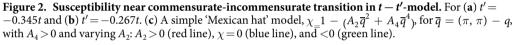
This paper is organized in the following Sections: Key findings, Methods, Results (DFT-Lindhard), Results (Beyond RPA), and Discussion. The Results section has been split into two parts. The first [DFT-Lindhard susceptibility] shows that the Lindhard susceptibility can be separated into two components, one fermionic, associated with Fermi surface nesting, and more important at low *T*, the other bosonic, associated with Van Hove singularity (VHS) nesting, and usually more important at higher *T*, where Pauli blocking becomes unimportant. The second Results section [Beyond RPA susceptibility] shows how the competition between these components can be incorporated into Moriya-Hertz-Millis theory by extending the calculations to realistic band structures. A number of specialized issues are covered in the Supplementary Material.

Key findings

DFT-Lindhard susceptibility. Our study illustrates the transformative role of the density functional theory (DFT) in many-body perturbation theory (MBPT)³¹, and provides a deeper understanding of the bare (Lindhard) susceptibility χ_0 [Methods Section]. Textbook calculations of χ_0 replace all band structure effects by a simple electron gas with a parabolic band. When realistic band structures are included, the Lindhard function becomes a nesting map³², showing all the weighted nesting vectors of the various Fermi surfaces – in particular, a scaled map of each Fermi surface, $q = 2k_F$ (where k_F is the Fermi wave vector), associated with intraband nesting [Results Section (DFT-Lindhard)]. Indeed, we find that χ_0 can serve as a basis for classifying phase transitions, analogous to the spectrum generating algebras of nuclear physics [Supplemental Materials Section I]. We find that doping-temperature phase space can be divided up into fluctuation maps, or basins of attraction, where the character of the dominant nesting vector changes. These fluctuation maps lead to the idea of equivalence classes, where materials with very different sets of hopping parameters nevertheless display topologically equivalent fluctuation maps. In particular, we find that LSCO lies in a different equivalence class from other cuprates. From each equivalence class, we can define a reference family having a minimal number of hopping parameters, that can be used to interpolate between different cuprates. In particular, we find that all cuprates lie on a single line in hopping parameter space, and that this tuning allows us to understand that LSCO is more strongly correlated than other cuprates. However, the susceptibility differs in one key way from most other properties of a Fermi liquid. While many properties are controlled entirely by states near the Fermi level, the susceptibility can also contain important contributions from states deep in the Fermi sea, here referred to as bulk contributions. In the cuprates, we find that χ_0 contains not only sharp features associated with the Fermi surface nesting map, but also contains a smoothly varying background, peaking at $q = (\pi, \pi)$ and giving rise to the near- (π, π) -plateau in the susceptibility. This background is associated with the VHS peak, which lies a distance $T_{VHS} = (E_F - E_{VHS})/k_B$ below the Fermi level, Fig. 1(a). Since it peaks away from the Fermi level, the corresponding features have a strong temperature dependence, associated with the smearing of the Fermi functions that enter in the expression for χ_0 , an effect referred to as Pauli blocking.

Role of entropy and VHS nesting in pseudogap physics. McMillan's phononic entropy referred to a competition wherein phonons with many different *q*-vectors try to soften at the same time²⁵, preventing any one





phonon from condensing until very low T and giving rise to an extended range of short-range order. We find that similar strong entropic effects arise from electronic bosons (electron-hole pairs), predominantly associated with the VHS. In turn, the entropic effects modify VHS physics, causing the main spectral weight associated with the VHS to shift rapidly towards half filling as T increases, not following the evolution of the density of states (DOS). This anomalous shift suggests that the VHS may play a large role in pseudogap physics, Fig. 1⁹. The key anomaly in pseudogap physics lies not below the pseudogap temperature but above T, Fig. 1(b), where the resistivity in the strange metal phase varies linearly with T, $\rho \sim \overline{T}$. The pseudogap temperature T^* indicates the onset of deviations from linear resistivity, while below a lower temperature T^{**} the resistivity becomes purely quadratic, suggestive of a Fermi liquid regime. While the strange metal phase is suggestive of the quantum critical phase in the Heisenberg quantum phase transition³³, there is no obvious reason why the resistivity of the Heisenberg model should be linear in T, nor any natural candidate for either the pseudogap (since the Heisenberg model has a well-defined ordered phase), or for the transition at T^{**}. In contrast, these features arise naturally in our VHS model, Fig. 1(a). For $T > T_{VHS}$ Pauli blocking is completely ineffective, leaving the expected VHS scattering rate $\tau^{-1} \sim T$. Below T_{VHS} this blocking of the VHS starts to turn on, while coherent Fermi surface features appear, until below a coherent-incoherent crossover, $T < T_{coh}$, the VHS scattering is suppressed, leaving behind a Fermi liquid $\tau^{-1} \sim T^2$. These results are consistent with experimental findings that the pseudogaps in Bi2201 and Bi2212 terminate near the VHS doping x_{VHS}^{34-36} .

Mode coupling and order parameter competition. The bosonic entropy can be incorporated into the DFT-Lindhard function by modifying Moriya's mode-coupling approach³⁷ to include realistic band-structures via a susceptibility DOS [Results Section (Beyond RPA)]. We use this technique to calculate the growth of the correlation length in LSCO, finding that the Mermin-Wagner theorem is satisfied leading to a broad regime of short-range order. We also find a temperature range where the correlation length does not grow at all. By tuning the hopping parameters, we find that this is associated with a 'supertransition' between two families of order parameters. This is particularly interesting as it involves competition between (π, π) Mott antiferromagnetic order driven by VHS fluctuations and a $(\pi, \pi - \delta)$ Slater SDW driven by conventional Fermi surface nesting. We are able to define an *order parameter DOS* that diverges at the transition: Fig. 2 shows the susceptibility near the transition for two different values of t'/t. At the transition, the susceptibility at (π, π) changes from a maximum to a local minimum, and the susceptibility is quite flat across the entire (π, π) -plateau. This is especially true in frame (b), which is tuned close to the T = 0 transition (note change of scale). It is convenient to introduce a susceptibility density-of-states (DOS) to quantify the number of competing states. This behavior can be approximately captured in a simple 'Mexican hat' model, Fig. 2(c). A similar analysis for Bi2201 finds a transition between $(\pi, \pi - \delta)$ and antinodal nesting order, consistent with the SDW-to-CDW transition found in most cuprates.

Methods

Our calculation is a form of many-body perturbation theory (MBPT) based on Hedin's scheme³¹. The scheme involves four elements: electrons are described by Green's functions *G* with DFT-based dispersions renormalized by a self-energy Σ ; electronic bosons [electron-hole pairs] are described by a spectral weight [susceptibility] χ

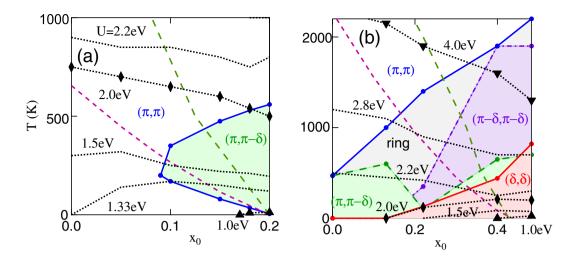


Figure 3. Mean-field phase diagrams of cuprates. DFT-based models of LSCO (a) and Bi2201 (b). Black dots indicate mean-field phase boundaries for different *U* values, including U = 4 (down triangles), 2 (diamonds), or 1 eV (up triangles). Here x_0 is the doping at T = 0 (plots are at constant E_F), but for qualitative purposes we can assume $x \simeq x_0$.

renormalized by vertex corrections Γ . Neglecting vertex corrections, the self-energy can be calculated as a convolution of *G* and $W = U^2 \chi$ (GW approximation). This approach has been used to solve the energy gap problem in semiconductors, where the Γ correction leads to excitons via the solution of a Bethe-Salpeter equation^{38,39}, and in extending DMFT calculations to incorporate more correlations [e.g., DMFT + GW, etc.]. Our approach here is to extend our previous GW calculations [quasiparticle-GW or QPGW⁴⁰] to approximately include vertex corrections.

In QPGW we introduce an auxiliary function $G_Z = Z/(\omega - \epsilon_k^{QP})$, where the dressed, or QP dispersion is $\epsilon_k^{QP} = Z \epsilon_k^{DFT}$, and ϵ_k^{DFT} is the bare, or DFT dispersion. G_Z behaves like the Green's function of a Landau-type QP – a free electron with renormalized parameters that describes the low-energy dressed electronic excitations. However, this is a non-Fermi liquid type QP, since the frequency-integral of $Im(G_Z)$ is Z and not unity. That is, the Z-QP describes only the coherent part of the electronic dispersion, and is not in a 1:1 correspondence with the original electrons. The importance of such a correction can be readily demonstrated. Since a Z-QP has only the weight Z of a regular electron, the susceptibility [a convolution of two Gs] is weaker by a factor of Z² than an ordinary bare susceptibility. To match this effect in the Stoner criterion requires introducing an effective $U_{eff} = ZU$. In contrast, MBPT calculations in semiconductors typically set the GW-corrected Green's function to $G_{GW}^{-1} = \omega - \epsilon_k^{QP}$, where ϵ_k^{QP} is the average GW-renormalized dispersion⁴¹, thereby missing the reduced spectral weight of the low-energy, coherent part of the band.

In this paper the DFT-Lindhard susceptibility χ_0 is calculated at intermediate coupling level, using Z-corrected first-principles dispersions to describe the coherent part of the quasiparticle-GW dressed carriers⁴⁰. We calculate χ_0 based on the dressed dispersion $\epsilon_Z = Z\epsilon_0$, assuming a doping-independent Z = 0.5. We use a single-band model of the cuprates, working in a purely magnetic sector, with the Hubbard U controlling all fluctuations; there is a competition in this model between near-nodal and antinodal nesting (ANN) which mimics the SDW-CDW competition in cuprates, sharing the same nesting vectors^{22,32}.

Results: DFT-Lindhard susceptibility

Sorting the susceptibility peaks: basins of attraction for cuprates. Fluctuation maps. We define fluctuation maps in x-T space as maps of domains ('basins of attraction') of classes of q-vectors at which $\chi_0(q, 0)$ is maximum. Figure 3 displays fluctation maps for $La_{2-x}Sr_xCuO_4$ (LSCO) (a) and $Bi_2Sr_2CuO_{6+x}$ (Bi2201) (b), using the full DFT dispersion. We fix an energy scale by taking t = 0.4195 eV, appropriate for LSCO⁴². Most of the colored regions are associated with FS nesting, while the white region is mainly associated with (π, π) VHS nesting, as will be discussed below. This figure confirms our earlier finding that LSCO, Fig. 3(a), has a very different phase diagram from most other cuprates, here represented by Bi2201, Fig. 3(b)³². In the phase labeled 'ring', the susceptibility is approximately constant along a ring of *q*-vectors surrounding (π, π) . We note in particular that antinodal nesting (red-shaded region) emerges naturally in most cuprates at higher doping, but is absent in LSCO. Shown also in Fig. 3 are T_{VHS} (dark red-dashed line) and the shift of the DOS-peak (thin green-dashed line), and a series of mean-field phase boundaries, based on a Stoner criterion,

$$U\chi_0(q,\,\omega=0) = 1.$$
 (1)

There is a critical value of $U = U_0$ below which there is no phase transition at x = 0. For $U < U_0$, the phase boundary is dome-shaped, but with a peak shifted away from the T = 0 value of x_{VHS} to a value close to the line $x_{VHS}(T)$ (red-dashed line). For $U > U_0$ the phase boundary rapidly evolves to a strong coupling form where the highest T transition occurs near half-filling. Consistent with this, we will show that VHS phenomena are dominated by entropy effects, and thus scale with the (π, π) susceptibility and not with the peak DOS. In all cases, at sufficiently high *T*, FS nesting is washed out and the dominant instability is at $q = (\pi, \pi)$. In comparing these phase boundaries to experiments in cuprates, it should be kept in mind that the dotted curves refer to constant *U*, whereas *U* is likely to be strongly screened near a VHS⁴³.

Reference families. The above fluctuation maps allow us to introduce equivalence classes of models for cuprates, and in particular to determine simple reference families that can be used to search parameter space beyond the physical cuprates. To understand differences between various cuprates, and to gauge their proximity to the pure Hubbard model, it is necessary to add a third axis to the usual *T*- and doping-phase diagrams. Here we find reference families for which differences between physical cuprates' band structures can be parametrized by a single hopping parameter.

By a Wannier-type downfolding of DFT results, each cuprate can be modelled by a single CuO_2 band crossing the Fermi level, so their parameter space consists of hopping parameters, traditionally t, t', t'', In fact, since t sets an energy scale, the band dispersion is determined by the ratios t'/t, t''/t, To reduce the number of hopping parameters, we introduce a notion of equivalence, wherein a set of materials are equivalent if they have the same fluctuation map. From this equivalence, the DFT parameters can be mapped onto an equivalent *reference family* depending only on two parameters, namely t'/t and t''/t. We find that when the physical cuprates are approximated by their reference families they lie close to a particular cut in this 2D space.

Figure 4 compares the DFT-derived fluctuation maps with similar maps calculated for a variety of simpler dispersions. The left-hand frames (a), (c), (e), refer to LSCO, the right-hand frames (b), (d), (f), to Bi2201. The DFT results [middle row, frames (c), (d)] are compared to two reference families, (*i*) a minimal cut defined by t'' = 0 [top row, (a), (b)] or (*ii*) the Pavarini-Andersen cut⁴⁴, t'' = -t'/2 [bottom row, (e), (f)]. We find that reference family (*ii*) semi-quantitatively reproduces the DFT derived fluctuation maps for both cuprates. On the other hand, setting t'' = 0 works well for LSCO, Fig. 4(a), but fails for Bi2201 Fig. 4(b). We further find that a reasonable choice of t'/t is the one which matches the Van Hove singularity (VHS) doping of the full DFT dispersion. We note that both reference families have the additional desirable feature that they evolve from the state with t' = 0– i.e., the original Hubbard model.

An advantage of using the reference families is that this lets us systematically vary parameter space to study the evolution from one cuprate to another, or to extrapolate the parameter space to better understand the origin of various phases. Figure 5(a) shows the cuts in t' - t''-space associated with reference families (i) (t'' = 0, red line) and (ii) (t'' = -t'/2, blue line). The dotted lines, Fig. 5(a), indicate ground-state transitions at x = 0, from (π , π) to (π , $\pi - \delta$) (black dots) and from (π , $\pi - \delta$) to (δ , δ) order (red dots). The black dotted line is the Mott-Slater line discussed in the Results Section (Beyond RPA). The densities-of-states (DOS) shown for the two families in Fig. 5(b,c) provide insight into the origin of the crossovers. Note that the use of these reference cuts demonstrates the clear evolution of the VHS between 2D and 1D behavior.

For reference family (*i*), Fig. 5(*c*), the DOS reveals a crossover associated with a change from 2D to 1D nesting. This is seen by the DOS at t' = -t/2 which shows the characteristic $1/\sqrt{E}$ divergence expected for a 1D system. To explore the origin of the 1D-behavior we consider the electronic dispersion for reference family (*i*) of the form $E = -2t(\cos(k_x a) + \cos(k_y a)) - 4t'\cos(k_x a)\cos(k_y a)$. If either k_x or $k_y = 0$ and t' = -t/2 then E = -2t. Therefore the FS at the VHS energy reduces to just the *x* and *y* axes. This shows that the strong 1D nature of the corresponding FS leads to a strong divergence in the DOS at the VHS.

Near t' = -t/2, the DOS has two peaks, one at the saddle-point VHS found in 2D materials, the other at the leading edge, as found in 1D systems, Fig. 5(c) (violet line). For t' < -0.45t (green-black), the leading edge is larger, suggestive of weakly-coupled chains, while for t' > -0.28t (green dotted arrow) the leading edge peak turns into a local minimum, Fig. 5(c) inset, at the green dotted arrow. This quasi-1D behavior is suggestive of the nematic phase found in cuprates⁴⁵. Indeed, an earlier VHS renormalization group calculation found that charge order transitions splitting the degeneracy of the VHSs at $(0, \pi)$ and $(\pi, 0)$ [which would now be called a nematic transition] are only possible if $t'/t < -0.276^{43}$. Remarkably, t' = -0.28t approximately coincides with the commensurate-incommensurate transition found in Fig. 2(b) to lead to a severe ξ depression.

For family (*ii*) a non-zero t'' modifies the dispersion such that it breaks up the 1 dimensionality of the FS and weakens the DOS divergence, as seen in Fig. 5(b). However, a residual 1 dimensionality remains in that there is conventional FS nesting in the ANN region, which is responsible for the CDW. For the red, orange, yellow, and green curves we find a resemblance to that of the DOS for t'' = 0. At the light blue curve there is a topological transition where a pocket centered at $(\pi, 0)$ splits off from the rest of the Fermi surface. For larger t', one now sees two features in the DOS, a peak followed by a step at lower energies. The peak is the topological transition where the pockets split off, and the step is where the $(\pi, 0)$ pocket disappears. The inset illustrates this progression showing the FS at the DOS peak energy for various t'.

Origins of structure in the susceptibility: competition between FS and VHS nesting. Separating *FS maps and broad background*. In order to understand the structure in the fluctuation maps we need to analyze the susceptibilities more deeply. The susceptibility contains sharp features which are a map of the Fermi surface and which quantify the strength of the Fermi surface nesting. In addition, the susceptibility contains a smooth, intense background feature which peaks at (π, π) . This peak plays an important role in the cuprates, producing large entropic effects. This peak can shift the balance of the FS nesting to *q*-vectors closer to the peaks, and in special cases, it can lead to commensurate nesting away from the FS nesting vector at exactly (π, π) . Moreover, as *T* increases, coherent FS features are washed out, leaving behind only the commensurate bulk contribution. Here we determine that this background can be considered to be a bosonic VHS (b-VHS), in that the electronic excitation comes predominantly from the region of the VHS. It is analogous to the Van Hove excitons found in optical spectra⁴⁶, but present in the intraband susceptibility. The evolution of this feature contains hints as to the origin of the pseudogap.

Consider the expression for the $\omega = 0$ bare susceptibility that enters into the Stoner criterion,

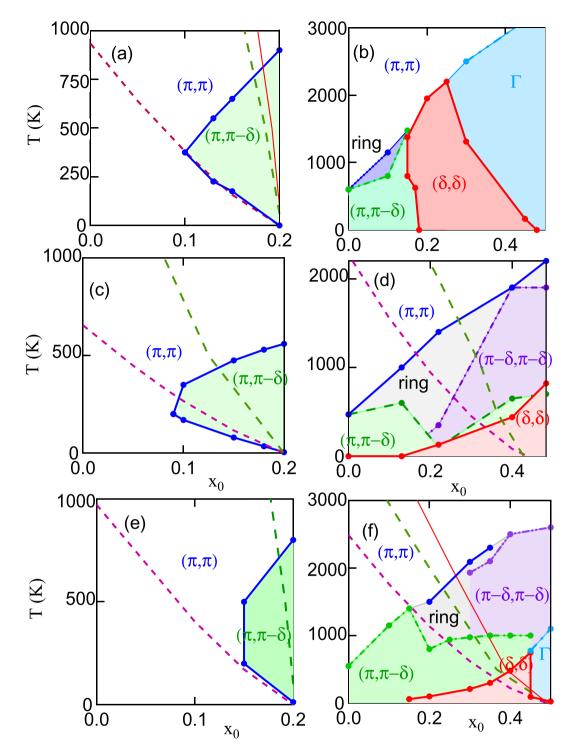
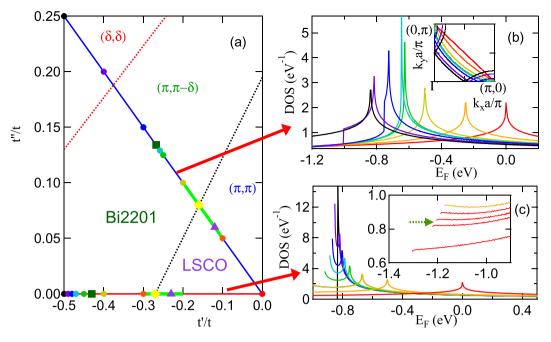
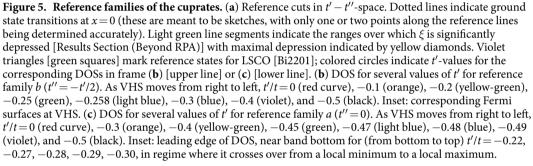


Figure 4. Fluctuation fingerprints of cuprates. Maps of *q*-vectors of largest susceptibility for DFT-based models of LSCO (**c**) and Bi2201 (**d**) compared to t - t' - t' reference phase a [t''=0] with t' = -0.23t (**a**) or -0.43t (**b**), and reference phase b [t''=-0.5t'] with t' = -0.12t (**e**) or -0.258t (**f**). Recall, from Fig. 3, that x_0 is the doping at T = 0, but we assume $x \simeq x_0$.

$$\chi_0(q) = \sum_k \frac{\Delta f_{k,q}}{\epsilon_k - \epsilon_{k+q}},\tag{2}$$

where $\Delta f_{k,q}$ is the Pauli blocking factor





$$\Delta f_{k,q} = f(\epsilon_{k+q}) - f(\epsilon_k). \tag{3}$$

The denominator is smallest if $\epsilon_k = \epsilon_{k+q}$, but in this case the numerator vanishes at T = 0 unless ϵ_k is at the Fermi level. Even in this case, the term contributes little weight to the susceptibility integral unless the *q*-shifted FSs are tangential, which is generally satisfied if $q = 2k_F$, giving rise to a folded map of the FS – the ridge of FS nesting³².

The Fermi surface nesting features in the susceptibility map for Bi₂Sr₂CuO₆ (Bi2201), at x = 0.13 and $\omega = 5$ meV, are illustrated in Fig. 6, with χ'_0 given in frame (a), and the corresponding map of χ''_0 in frame (b). Due to the energy δ -function, χ''_0 contains only the FS contribution⁴⁷, except an extra feature at Γ associated with the DOS. In Fig. 6(d), we plot χ'_0 (red line) and χ''_0 (blue line) along a cut $\Gamma \to (\pi, 0) \to (\pi, \pi) \to \Gamma$ in momentum space. It can be seen that, away from Γ , the peaks in both components fall at the same *q*-values, and the relative weights of the peaks are comparable. Moreover, the anti-nodal nesting peak near Γ in χ''_0 is stronger than the near- (π, π) peak. However, χ'_0 is largest on the near- (π, π) peaks, since the FS contribution is riding on the large background contribution to the susceptibility. Here we demonstrate that VHS nesting plays a dominant role in creating this background, thereby providing a deeper understanding of the VHS vs FS nesting competition.

Two approaches are taken to eliminate the FS nesting features to see the background susceptibility by itself. First, raising *T* (violet long-dashed line in Fig. 6(d)) washes out the sharp FS features first, while leaving the background features behind. Secondly, in frame (c), we plot the difference between χ'_0 and a suitably scaled quantity proportional to χ''_0 ; this is also plotted as the green curve in frame (d). Since χ'_0 and χ''_0 are Kramers-Krönig transforms, they have different lineshapes, but since the FS map is nearly one-dimensional, a large part of it is still cancelled in taking the difference. From frame (d) it can be seen that both approaches lead to similar results for the bulk susceptibility, and that this contribution provides the dominant contribution to virtually the whole (π , π) plateau structure. Here the background contribution arises as follows. Even when $\epsilon_k \neq \epsilon_{k+q}$, there will be a finite contribution to the susceptibility, Eq. 2, reduced by the factor $1/(\epsilon_k - \epsilon_{k+q})$, as long as the two energies lie on opposite sides of the Fermi level (for T=0). This contribution will be especially significant near a VHS, which has a large spectral weight.

Background peak as VHS nesting. Here we demonstrate the connection between the bulk contribution to χ_0 and VHS nesting by deconvolving the bare susceptibility, Eq. 2, into its various ϵ_k components. To understand this procedure, recall that the VHS gives rise to susceptibility features near Γ (DOS) and (π , π). The latter are typically

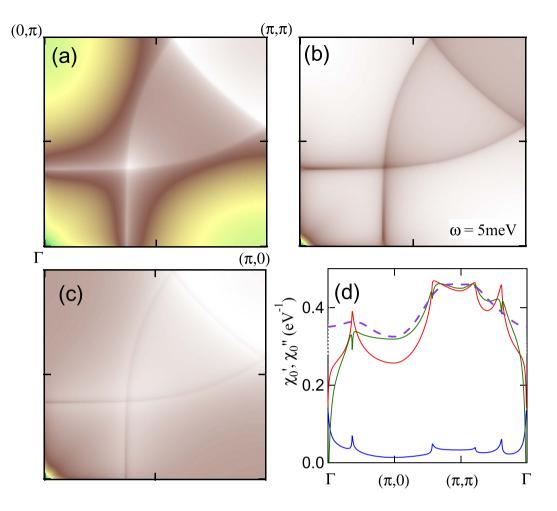


Figure 6. Extracting the bulk χ_0 contribution. (a) Susceptibility χ'_0 of Bi2201 for full DFT dispersion and x = 0.13, plotted in first Brillouin zone at T = 0, $\omega = 5$ meV. Color scale from minimum (blue) to maximum (white). (b) As in (a), but for χ''_0 . (c) Difference plot of (a) and the scaled version of (b). (d) Cuts of susceptibility from earlier frames plotted along high symmetry lines, with red curve representing data in frame (a), blue curve for frame (b), and green curve for frame (c). Shown also is the curve corresponding to frame (a), but at T = 1000 K (violet long-dashed line) scaled by 1.4 and shifted by -0.045 eV^{-1} to match other curves near (π, π) .

more intense. Thus, to confirm that the background susceptibility is associated with the VHS, we analyze Eq. 2 for χ_0 , assuming that the electron is associated with ϵ_k and the hole with ϵ_{k+q} . We must demonstrate that, when $q \sim (\pi, \pi)$, the dominant contribution to χ_0 has *k* near the VHS at $(\pi, 0)$ [or $(0, \pi)$] and ϵ_k near E_{VHS} .

To demonstrate this we divide the Brillouin zone into four patches, around the Γ , (π , 0), (0, π), and (π , π) points, and calculate separately the susceptibility due to each patch, rewriting Eq. 2 as

$$\chi_0(q) = \sum_{i=1,4} \sum_{\epsilon_k} \chi_{0,i}(q, \epsilon_k), \tag{4}$$

where *i* runs over the four quadrants Γ , $(\pi, 0)$, $(0, \pi)$, and (π, π) . In LSCO, the plateau is predominantly composed of equal contributions from the $(0, \pi)$ and $(\pi, 0)$ patches, as expected for a VHS. We then explore the distribution of values of ϵ_k which contribute to the susceptibility in the $(\pi, 0)$ patch at x = 0, Fig. 7. There is a clear crossover from a dominant near-FS contribution $(\epsilon_k = 0)$ at low *T* to a dominant contribution at high-*T* which lies along a flat (q-independent) line centered at energy $k_B T_{VHS} = E_F - E_{VHS}$. As we vary doping, this flat line shifts towards the Fermi level following T_{VHS} , red-dashed line in Figs 4(a) and 8.

This crossover is associated with Pauli blocking, i.e. the factor Δf in Eq. 5. Since at low *T* the Fermi function is a step, Δf blocks the scattering between states far below E_F . As *T* increases *f* gets smeared out, so that features away from the Fermi surface can contribute to the susceptibility. Since the VHS has a substantial peak, the high-*T* changes are due to the unveiling of this feature. Note that at low *T*, Fig. 7(a–c), there is still scattering from one VHS, but to states near E_F , and hence contributing intensity at $q \neq (\pi, \pi)$. Correspondingly, the weight near (π, π) comes from states in the tail of the VHS extending to E_F .

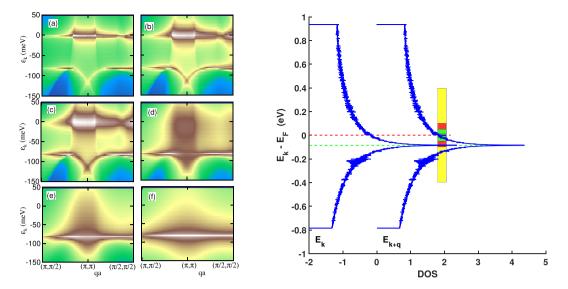


Figure 7. Coherent-incoherent crossover at $\mathbf{x} = 0$: role of VHS nesting. Deconvolution of the susceptibility, χ_0' , for *k*-states in the $(\pi, 0)$ -patch at several temperatures *T*. Contribution to the susceptibility of $\chi_{0,2}(q, \epsilon_k)$, Eq. 4, is plotted along *q*-cuts near (π, π) with the color scale ranging from white for a large contribution to blue for negligible contribution for: T = 0 (a), 25 K (b), 50 K (c), 250 K (d), 500 K (e), and 2000 K (f). Frame (g) illustrates scattering from the VHS at E_k (green dotted line) to states centered at the Fermi level at E_{k+q} . The colored region shows where Pauli blocking is ineffective at T = 25 (blue bar), 50 (black bar), 250 (green bar), 500 (red bar), and 2000 K (yellow bar).

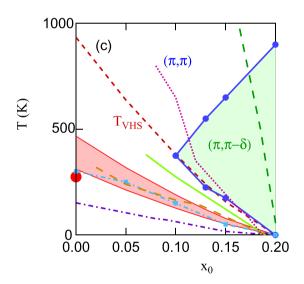


Figure 8. LSCO phase diagram, showing several VHS-related crossovers and regions where the susceptibility peaks at different *q*-vectors. Dominant fluctuations are at commensurate (π, π) (white shaded region) or incommensurate $(\pi, \pi - \delta)$ (green shaded region); crossovers are T_{VHS} (red short-dashed line), DOS peak (green long-dashed line), coherent-incoherent crossover (pink shaded region), T_{γ} (violet dot-dashed curve), and the position of the (π, π) peak vs doping (light blue dot-dashed line). Experimental pseudogaps from ref. 9 (yellow-green solid line and brown dashed line) and ref. 51 (magenta dotted line). Recall, from Fig. 3, that x_0 is the doping at T=0, but we assume $x \simeq x_0$.

While Δf enhances features associated with the VHS as *T* increases, it has the opposite effect on features near the Fermi level. For energy ϵ exactly at E_F , $f(\epsilon)$ is exactly 1/2 at any *T*, so $\Delta f = 0$ if both states are at E_F . As *T* increases, *f* smooths out, so that if $|\epsilon_k|$, $|\epsilon_{k+q}|$ are both $< k_BT$, then $f_k \approx f_{k+q} \approx 1/2$, and therefore, $\Delta f \approx 0$. This explains why the crossover from FS nesting to VHS nesting in Fig. 7 occurs in such a narrow *T* range (pink shaded region in Fig. 8). Passing between Fig. 7(c,d) the Fermi surface contribution goes through a coherent - incoherent crossover – note that the FS nesting contribution is fully incoherent by 500 K.

In Fig. 8 we superpose this crossover, plotted as a pink shaded region, on the fluctuation map for LSCO, with all calculations for the minimal reference family (t'' = 0). This incoherent-coherent crossover scales with T_{VHS} , with boundary lines given by $T_{VHS}/2$ and $T_{VHS}/3$. The same crossover can be seen directly in the susceptibility,

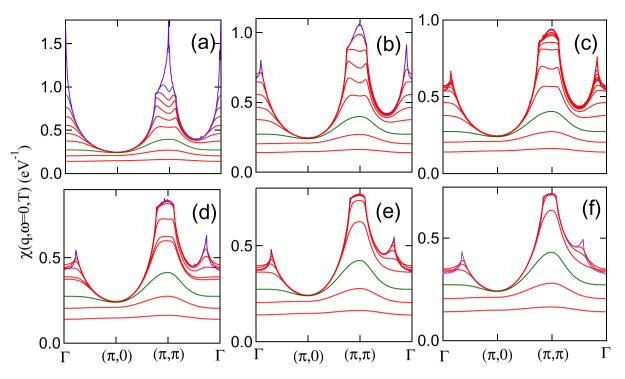


Figure 9. Coherent-incoherent crossover at $\mathbf{x} = \mathbf{0}$ in $\chi_0(T)$. Lindhart susceptibility $\chi'(q,\omega=0)$ of LSCO, plotted along high-symmetry lines at 6 different dopings: (a) x = 0.20 (~VHS), (b) x = 0.18, (c) x = 0.15, (d) x = 0.10, (e) x = 0.05, and (f) x = 0, at (from top to bottom) T = 0, 10, 25, 50, 100, 200, 400, 1000, 2000, and 4000 K, except that in frames (d) and (f) the traces at T = 10 and 25 K are omitted. The T = 1000 K susceptibility is plotted as a green line.

Fig. 9, as the coherent, FS-related features present at low *T* are washed out, leaving only a featureless (incoherent) (π, π) plateau at high *T*.

These effects of Pauli blocking can explain several additional features of Fig. 8. For instance, hole doping causes the (π, π) -plateau to expand in momentum space, leading to a weakening of VHS nesting. Thus, in Bi2201, the transition between (π, π) -nesting and some form of FS nesting shifts to higher *T* with increasing *x*, as seen in Fig. 4(b,d,f). A similar effect is found in LSCO, Fig. 4(a,c,e), for the high-*T* termination of the incommensurate $(\pi, \pi - \delta)$ order (green shaded region). However, in LSCO there is also an anomalous reentrant transition, where (π, π) order reappears below a lower domain boundary (blue solid line). To understand this unusual reentrance, consider first what happens to χ_0 when $E_F = E_{VHS}$. Fig. 9(a). While the VHS peak is clearly present at T = 0 (blue line), finite temperature does not lead to simple thermal broadening but actually splits the peak. For small finite *T* the very sharp VHS at E_F is blocked. As T increases, the blocking spreads over a wider range of energy, leading to the splitting of the VHS feature. Thus, raising *T* causes a blocking first of the sharp VHS peak to produce the lower edge of the domain boundary. Shifting to $x < x_{VHS}$ means that a higher *T* is necessary to block the sharp VHS peak. Note that the lower branch of the boundary closely follows T_{VHS} , especially for the t'' = 0 reference family. This is also why the conventional VHS theory that relies on the sharp VHS peak only holds at relatively low *T* as discussed in conjunction with Fig. 3.

Pauli-blocking and the bosonic-VHS. If we think of χ_0 as the propagator of different electron-hole pairs, then the pair associated with $q = Q \equiv (\pi, \pi)$ has many similarities with the excitonic-VHS seen in optical studies⁴⁶, and we will refer to it as the bosonic-VHS, or b-VHS for short. We clarify the role of Pauli blocking by describing its effect on the b-VHS in detail, using the t'' = 0 reference family for simplicity. To understand the relationship between the b-VHS and entropy effects, we need to know how Pauli blocking "hides" the states associated with the VHS at low *T*, and how they recover as *T* increases. If the renormalized dispersion of a single electron is ϵ_k with wave vector *k*, then an e-h pair at wave vector *q* has a dispersion $\omega_q(k) = \epsilon_{k+q} - \epsilon_k = -2\epsilon_{q-}(k)$, where $\epsilon_{q\pm}(k) = (\epsilon_k \pm \epsilon_{k+q})/2$, and a Pauli blocking factor $\Delta f_{k,q} = f(\epsilon_{k+q}) - f(\epsilon_k)$. Then the corresponding pair DOS is $D_q(\omega) = \sum_k \Delta f_{k,q} \delta(\omega - \omega_q(k)) = \chi_0''(q, \omega)$. For LSCO, the dominant pairs are those at q = Q, the AF nesting vector. The associated dispersion $\omega_Q(\pi, \pi)$, plotted in Fig. 10(a), resembles the electronic dispersion ϵ_k , but with an important distinction: it depends only on $\epsilon_{Q-}(k)$, whereas all of the hopping terms that shift the electronic VHS away from half-filling (t', t'') are contained in $\epsilon_{Q+}(k)$, i.e. the b-VHS is pinned at $\omega = 0$. Since χ_0'' is an odd function of $\omega, \chi_0''(Q, \omega = 0) = 0$. However, while the b-VHS is pinned at $\omega = 0$, Fig. 10(a), its weight vanishes at T = 0, due to the Pauli-blocking factor, $\Delta f_{k,Q} = 0$ near $k = (\pi, 0)$ at T = 0, Fig. 10(b). Finite *T* restores weight, optimally near 1000 K, although Δf always vanishes exactly at $(\pi, 0)$. Hence the bosonic entropy only turns on at finite *T*. Figure 7(g) illustrates how the Pauli window opens with increasing *T*.

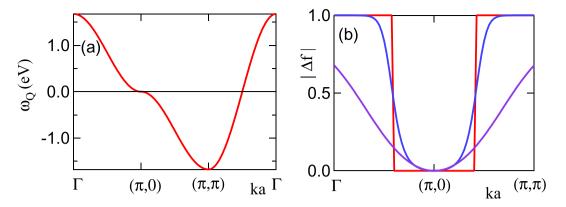


Figure 10. Coherent-incoherent crossover at the VHS. (a) Pair dispersion, ω_Q , as a function of k, and (b) the corresponding weight Δf , at T = 0 K (red), 500 K (blue), and 2000 K (violet) for LSCO (x = 0).

As *T* increases, the Fermi functions spread out, so the FS contribution decreases whereas the bulk contribution grows as Δf becomes unblocked. For the VHS, we expect full unblocking at a temperature near $T \sim T_{VHS}$ (red short-dashed line in Fig. 8). Since *f* varies exponentially with *T*, the unblocking will appear at a temperature $T < T_{VHS}$, which may vary according to the property studied. This is similar to a Schottky anomaly, where a feature at energy Δ above the ground state produces a peak in the heat capacity at a temperature $T_m \sim \Delta/2.4$. The low-*T* fadeout of the b-VHS is controlled by the Pauli blocking factor,

$$\Delta f_{k,Q} = \frac{\sinh x_-}{\cosh x_- + \cosh x_+},\tag{5}$$

where $x_{-} = \epsilon_{Q,-}(k)/k_BT$, $x_{+} = (\epsilon_{Q,+}(k) - E_F)/k_BT$. When the electronic VHS is at the Fermi level, $x_{+} = 0$ and $\Delta f_{k,Q} = tanhx_{-}/2$, while at lower doping the b-VHS is exponentially suppressed, cosh $x_{+} \sim exp(T_{VHS}/T)/2$, which equals 1 when $T \sim T_{VHS}/ln(2)$.

Entropy and a revised VHS scenario. Here we will demonstrate that this b-VHS peak produces large entropic effects. To show this, we calculate several characteristic VHS-dominated properties – the Sommerfeld coefficient of the heat capacity and the susceptibilities at Γ and (π, π) . The entropy is calculated from the standard expression

$$S = -k_B \sum_{k} [f_k \ln(f_k) + (1 - f_k) \ln(1 - f_k)],$$
(6)

with $\gamma = dS/dT$, Fig. 11. The sum is over the electronic dispersion of LSCO, assuming a paramagnetic phase. At T=0, γ is proportional to the DOS, and hence diverges at x_{VHS} , Fig. 11. However, as *T* increases, the peak does not remain near x_{VHS} , but rapidly shifts toward x=0. We plot the peak position T_{γ} both in the inset of Fig. 11 and as a violet dot-dashed line in Fig. 8. Clearly, since γ is a measure of entropy, it is sensitive to sources of entropy away from the Fermi level, when *T* is high enough to reduce the Pauli blocking. It has recently been noted that there should be an excess of low-*T* entropy in the vicinity of a quantum phase transition⁴⁸.

Similar behavior is found for χ_0 . The VHS has logarithmically diverging features in the susceptibility at two different *q*-vectors, an inter-VHS contribution at (π, π) , and an intra-VHS contribution at $\Gamma^{49,50}$. While both peaks cross the Fermi level at x_{VHS} at T = 0, they evolve differently with temperature. This is because the feature at Γ represents the peak in the DOS, and hence it is dominated by near-FS physics. As a result it shifts only weakly with *T*, green long-dashed line in Fig. 8. In contrast, the inter-VHS (π, π) peak is a bulk contribution and hence is sensitive to Pauli blocking, as discussed above. This leads to the rapid thermal evolution of the (π, π) -susceptibility peak all the way to x = 0 (light blue dot-dashed line), scaling with both T_{γ} and T_{VHS} . Fig. 8. However, this behavior seems to be dominated by a separate aspect of the VHS physics. As discussed in ref. 32, the susceptibility has an approximate electron-hole symmetry, with similar spectra for a nearly empty $(x \sim 1)$ or a nearly full $(x \sim -1)$ band. The t' and t'' hopping parameters break this symmetry, and as a result, the low-T susceptibility is roughly symmetrical about x_{VHS} , where it has its largest peak. Here it seems that as T increases, the electron-hole symmetry is restored when T > t'. Since $T_{VHS} \sim t'$, one expects a similar scaling of T vs x for $x \ge 0$. But when x < 0 T_{VHS} continues to increase, whereas the peak in $\chi_0(q = (\pi, \pi))$ stays at x = 0, indicative of a restored electron-hole symmetry (see Supplementary Material Fig. 2).

These entropic effects have profound consequences for VHS physics that were overlooked in earlier mean-field studies^{49,50}. Most earlier studies of cuprate antiferromagnetism used a model valid only near the FS, or at low-*T*, overlooking Pauli unblocking of the VHS feature at large *T*. As a consequence, they predicted that χ_0 would show a maximum for *x* near x_{VHS} , inconsistent with experiment. Figure 3(a) shows revised RPA phase diagrams for several values of *U*, based on the DFT-Lindhard susceptibility. For small enough U ($U < \sim 1.3 \text{ eV}$), the earlier results are recovered, but when *U* is larger a new behavior is found, with a maximum transition temperature at half-filling, decreasing with doping as if approaching a quantum critical point. This now starts to resemble the pseudogap phase diagram of the cuprates. The comparison of experimental pseudogap lines^{9,51} in Fig. 8 will be discussed in the Discussion Section.

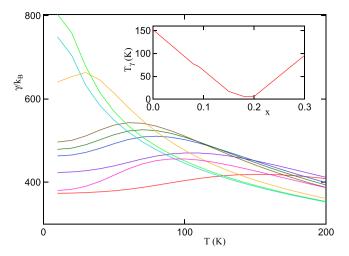


Figure 11. Sommerfeld constant calculated at several dopings. x = 0 (red curve), 0.05 (violet), 0.08 (blue), 0.09 (green), 0.10 (brown), 0.15 (orange), 0.18 (light green), 0.20 (light blue), and 0.30 (magenta). Inset: T_{γ} vs doping.

Finally, when McMillan²⁵ introduced the idea of bosonic entropy, he meant that competition between different phonon instabilities (phonons with different q-vectors) slowed down the tendency to CDW ordering, leading to broad *T*-ranges with only short-range order. In the following section we will show that a similar competition exists for electronic bosons, and that mode coupling allows us to quantify these effects.

Results: Beyond RPA susceptibility

Mode coupling formalism. Above, we have established the outlines of a classification scheme for phase transitions, analogous to the spectrum-generating algebras (SGAs) of nuclear physics. We summarize this development in Supplementary Material Section I.A, and compare it to SGAs in Supplementary Material Section I.C. In the present section we show how competition between condensing phases affects the transition temperature, and how to quantify the number of competing modes. A few striking examples of this competition will be discussed.

From χ_0 we calculate an interacting χ of modified random-phase approximation (RPA) form,

$$\chi(\mathbf{q}, i\omega_n) = \frac{\chi_0(\mathbf{q}, i\omega_n)}{1 + \lambda - U\chi_0(\mathbf{q}, i\omega_n)},\tag{7}$$

where λ is a mode-coupling parameter which must be calculated self-consistently. The imaginary part of χ can be thought of as the density-of-states (DOS) of electronic bosons, electron-hole (e-h) pairs, which may become excitons or excitonic resonances when a long-range Coulomb interaction is turned on. However, static ($\omega = 0$) instabilities depend on a Stoner criterion, and hence on

$$\chi_0'(q,\,\omega=0) = \sum_k \frac{f(\epsilon_{k+q}) - f(\epsilon_k)}{\epsilon_k - \epsilon_{k+q}} = 2 \int_0^\infty \frac{d\omega'}{\pi} \frac{\chi_0''(q,\,\omega')}{\omega'}.$$
(8)

The self-consistent parameter λ is found from a Matsubara sum of the susceptibility

$$\lambda = \frac{A_0 T}{N} \sum_{\mathbf{q}, i\omega_n} \chi(\mathbf{q}, i\omega_n), \tag{9}$$

where *N* is the number of *q*-points, and the summation in Eq. 9 can be transformed:

$$\frac{T}{N}\sum_{\mathbf{q},i\omega_n}\chi(\mathbf{q},i\omega_n) = \int \frac{d^2qa^2}{4\pi^2} \int_0^\infty \frac{d\omega}{\pi} coth\left(\frac{\omega}{2T}\right)\chi''(q,\omega+i\delta)$$
$$\simeq \hat{\lambda}_1 + \hat{\lambda}_2 T, \tag{10}$$

with *a* the in-plane lattice constant,

$$\hat{\lambda}_1 = \int \frac{d^2 q a^2}{4\pi^2} \int_0^\infty \frac{d\omega}{\pi} \chi''(q,\,\omega),\tag{11}$$

$$\hat{\lambda}_2 = \int \frac{d^2 q a^2}{4\pi^2} \chi'(\mathbf{q}, \ 0).$$
(12)

The term $\hat{\lambda}_1$ introduces a small, nonsingular correction²³ to Eq. 9, which we neglect.

In order to explore the role of suseptibility plateaus for realistic band dispersions, the self-consistency equation is evaluated numerically. Note that since χ_0^{-1} has dimensions of energy, a plot of $\chi_0^{-1}(q, 0)$ resembles a dispersion map. Hence we can define a SDOS:

$$\int \frac{d^2 q a^2}{4\pi^2} = \int N_-(X_-) dX_-,$$
(13)

where $X_{-} = \chi_0^{-1}$, is the variable of integration, and we introduce a corresponding susceptibility density of states (SDOS) N_{-} . Eq. 9 then becomes

$$\lambda = \Gamma T A_0 \int_{U_c}^{\infty} dX_- \frac{N_-(X_-)}{X_- - U_c + \delta},\tag{14}$$

with $\delta = U_c - U_{sp}$, $U_c = 1/\chi_0(\mathbf{Q}_0, 0)$. In Section II of Supplementary Materials, we show that Eq. 14 is closely related to excitonic Bose condensation, with λ proportional to the effective number of bosons. Once N_- has been calculated, Eq. 14 can be evaluated numerically. The singular part of the integral is treated analytically, and the remainder numerically, with u and U approximated as constants, $u = 0.8eV^{-1}$, U = 2 eV. It is convenient to fix the SDOS at some temperature T', and then solve Eq. 14 for $T(\xi_{th}, T')$, with self-consistency requiring $T(\xi_{th}, T') = T'$.

Figure 12 illustrates how integrating over the inverse susceptibility, Fig. 12(a), leads to the SDOS, Fig. 12(b), at x = 0 where the near- (π, π) plateau is dominant. Here we use hopping parameters appropriate to the DFT dispersion of Bi2201, t = 419.5, t' = -108.2 and t'' = 54.1 meV (Subsection Mode coupling in Bi2201). This yields a doping phase diagram that is qualitatively similar to that of most cuprates, except LSCO. By comparing Fig. 12(a,b), one can see how features in χ_0^{-1} translate into features in N_- . Thus, the intense, flat-topped peak in N_- at small values of χ_0^{-1} represents the near- (π, π) plateau. Its broad leading edge (smaller χ_0^{-1}) is controlled by anisotropy of the plateau edge between $(\pi, \pi - \delta)$ and $(\pi - \delta, \pi - \delta)$, while the sharp trailing edge corresponds to the local maximum of χ_0^{-1} at (π, π) . For the ANN peak, its leading edge scarcely leaves any feature in N_- , but a local maximum translates into a large peak in N_- . Finite temperature, T = 100K (blue line), rounds off the cusp in χ^{-1} . Fig. 12(c), leading to a step in N_- Fig. 12(d), but otherwise has little effect.

 χ_0^{-1} , Fig. 12(c), leading to a step in N_- , Fig. 12(d), but otherwise has little effect. Instead of such a self-consistent calculation, many treatments of phase transitions, including path-integral approaches, spin-fermion models, and Hertz-Millis^{52,53} quantum critical theories, approximate the susceptibility denominator by a *T*-independent Ornstein-Zernicke(OZ)⁵⁴ form,

$$\chi(\mathbf{q},\,\omega) \sim \frac{1}{\overline{q}^2 + \xi^{-2} + i(\omega/\omega_c)^z},\tag{15}$$

in terms of various deviations from the critical point (in \bar{q} , ω , and a 'tuning parameter' which is proportional to ξ^{-2} , where ξ is the correlation length). Here z is a dynamic exponent and $\bar{\mathbf{q}} = \mathbf{q} - \mathbf{Q}_0$ is the difference between the actual wave number \mathbf{q} and \mathbf{Q}_0 , the wave number where the susceptibility has a peak, $\chi_0(\mathbf{Q}_0, 0) = max_q[\chi'_0(\mathbf{q}, 0)]$ If the OZ parameters are assumed to be T- and doping-independent^{23,37,55,56}, the resulting physics becomes quite simple. For 2D materials, the Mermin-Wagner (MW) theorem⁵⁷ is satisfied, and the RPA transition at T_{mf} turns into a pseudogap onset at $T^* \sim T_{mf}$ with a crossover to long-range order when interlayer coupling is strong enough – in short, not much changes from the mean-field results.

When the OZ form of χ is assumed, $\chi_0(q) = \chi_0(Q) - A_2 \overline{q}^2$, then N_- becomes a constant, which we denote by N_a , and

$$\hat{\lambda}_2 = \frac{a^2}{4\pi} \int \frac{d\bar{q}^2}{\delta + B\bar{q}^2} = N_a \int \frac{dX_-}{\delta + (X_- - U_c)},\tag{16}$$

with $B = A_2 U_c^2$. Thus

$$\xi_{th}^{-2} = \delta/B \tag{17}$$

and $N_a = a^2/4\pi B$. Eq. 14 leads to long range order at T = 0 only, with correlation length $\xi_{th}/a = 1/\sqrt{4\pi N_a \delta}$ given by²³

$$\xi_{th}q_c = e^{T_2/T},$$
 (18)

where $T_2 = \pi A_2 \lambda / 6 \Gamma u a^2$, A_i is the coefficient of q^i , and q_c is a wave number cutoff.

The present $G_Z - W_Z - \Gamma_Z$ model is the simplest model which captures the essential physics of the pseudogap. For a more quantitative comparison with experiment, two additional problems must be solved. First, an extension to fully self-consistent $G - W - \Gamma$ may be needed to see how the susceptibility responds to the gap-opening. Secondly, the term in Eq. 15 proportional to ω^z must be included to describe quantum fluctuations. Since we are primarily interested in the opening of the pseudogap at higher temperatures, we ignored it in our analysis.

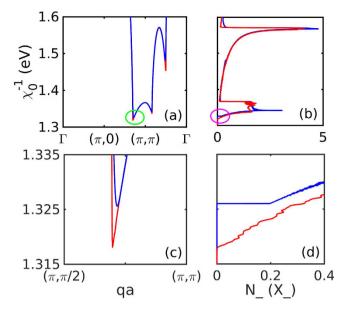


Figure 12. Origin of SDOS features. (a) Inverse susceptibility χ_0^{-1} (**q**) for Bi2201 at doping x = 0. Note that $U_c = min(\chi_0'^{-1})$ (circled region). Curves are at T = 10 (red) and 100 K (blue). (**b**) Corresponding susceptibility density of states (SDOS) $N_-(X_-)$, plotted horizontally, corresponding to the susceptibility of (**a**). (**c**,**d**) Blowups of circled regions in (**a**,**b**).

Strong mode coupling leads to an extended range of short-range order. Thus, mode coupling modifies the Stoner criterion to

$$U\chi_0(q, 0) = 1 + \lambda, \tag{19}$$

where λ satisfies Eq. 14. In the OZ approximation, this becomes

$$\lambda \propto \int_0^{q_c} \frac{q dq}{q^2 + \xi^{-2}} \sim \ln(\xi q_c).$$
⁽²⁰⁾

This OZ approximation misses the essential and strongly *T*-dependent mode-coupling physics, and we develop an alternative formulation of the theory. In particular, the inverse curvature of the susceptibility $1/A_2$ can diverge due to a competition between conventional Fermi-surface nesting and Van Hove singularity (VHS) nesting, and this parameter plays a special role in the theory. In particular, we can get a semiquantitative picture of the correlation length changes by simply accounting for the *T* dependence of $A_2(T)$. We define the order-parameter DOS (OPDOS) as the threshold SDOS,

$$N_{OP}(T) = N_{-}(X_{-} = U_{c}^{+}) = a^{2}/4\pi A_{2}(T)U_{c}^{2},$$
(21)

and use this in the OZ formalism, Eqs 16–18. While this neglects the fact that q_c can also have a strong *T*-dependence, and the contribution away from threshold is not negligible, the simplification should work when $\xi q_c \gg 1$. We believe that this OPDOS quantifies the number of competing modes and provides an essential link to the SGA theory. The connection between Eq. 14 and entropy is further discussed in Supplementary Material Section III.

Figure 13 illustrates the profound effects that strong mode coupling has in LSCO, as well as the complete inability of the OZ approximation to capture this physics. The SDOS, Fig. 13(a), contains VHS-like features characteristic of conventional DOSs. However, the singular behavior of Eq. 20 involves only features near the threshold, $X_- \sim U_c$, which evolve strongly with *T*, see inset to Fig. 13(a). For T > 0, the threshold behavior is always a step at $X_- = U_c$, indicative of a parabolic peak in χ_0 with curvature inversely proportional to the step height, Eq. 21. For analyzing critical phenomena, we define a threshold correlation length ξ_{th} (red solid lines in Fig. 13(b)) in terms of the initial curvature of χ_0 at \mathbf{Q}_0 , Eq. 17. We then evaluate ξ_{th} using the full Eq. 14, but the results can be interpreted in terms of an OZ form with a strongly *T*-dependent step height $N_{OP}(T)$. However, we caution that the experimentally measured $\xi_{1/2}$ is typically taken from the susceptibility half-width. This $\xi_{1/2}$ is generally larger than ξ_{th} due to the fast fall-off of χ near the edge of the (π, π) -plateau. Thus, for the example of Fig. 13(c), we calculated a $\xi_{1/2}$ (filled blue circles) from the full half-width of the renormalized χ , and found that typically $\xi_{1/2} \sim 2\xi_{th}$, where the solid red line in Fig. 13(c) plots ξ_{th} vs 1/T; the values of $\xi_{1/2}$ are in good agreement with experiment (green dot-dot-dashed line)⁵⁸. We note that while the OZ form is used in quantum critical theory^{35,36}, it is explicitly stated that it is to be used only in a limited T-range, and only in the absence of FS nesting³⁶, both of which are violated here.

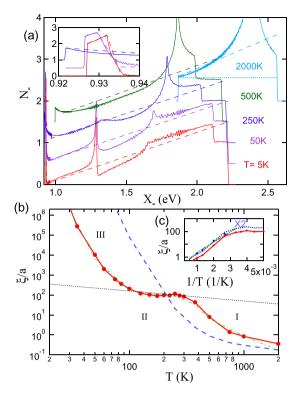


Figure 13. Structures in LSCO correlation length. (a) SDOS for undoped LSCO at several temperatures; light-blue dotted line gives OZ form of SDOS. (b) Corresponding temperature dependence of correlation length ξ_{th} . For x = 0, the red solid line with filled dots is ξ_{th} calculated from the solid lines in (a), while the thin green dot-dashed line is based on the dashed lines in (a), showing that ξ_{th} is relatively insensitive to the structures away from threshold. In contrast, the blue dashed line is based on the shape of the SDOS at T = 2000 K, but shifted and renormalized to match the SDOS at lower *T*, illustrating sensitivity to the leading edge structure. The black dotted line illustrates the scaling $\xi_{th} \propto T^{-1/2}$. (c) Calculated ξ_{th} reploted for x = 0 (red solid line with filled circles) compared with $\xi_{1/2}$ (blue filled circles) and with experiment (green dot-dashed line)⁵⁸; the blue dotted line is twice ξ_{th} .

Figure 13(b) illustrates how strong mode coupling slows down the correlation length divergence. Undoped LSCO (red solid line) shows two regions I and III of exponential growth of ξ_{th} with decreasing *T*, separated by an anomalous region II where ξ_{th} actually decreases with decreasing *T*. While the model captures the MW-like divergence at low *T* (region III), our main interest is in the higher-*T* behavior. In the high-*T* limit (region I), the leading-edge parabolic curvature is quite small, and if it were *T*-independent, as in the OZ approximation, the growth in ξ_{th} would follow the blue dashed line (Eq. 18), but thermal broadening causes the curvature to decrease with increasing *T*, leading to the faster growth of the red solid line. The origin of the anomalous region II for undoped LSCO, Fig. 13(b), can be readily understood from Fig. 13(a), where for T < 1500 K there is excess SDOS weight near U_c , leading to a strong peak (inset) as $T \rightarrow 0$. This feature represents the development of the (π, π) -plateau. In Supplementary Material Section III we present a simple calculation of ξ_{th} , showing how a range of power-law behavior, $\xi_{th} \sim 1/T^{1/2}$ (dotted line in Fig. 13(b)), could arise. While this power-law behavior captures the average *T* dependence of the anomaly, it cannot reproduce the nonmonotonic variation of the correlation length.

Exploring parameter space. The reference families can also be used to interpolate between physical cuprates. Here we use interpolation to understand the origin of the condensation bottleneck found in region II. We find that it is caused by *proximity* to a much stronger transition, a crossover between Mott and Slater physics, and that LSCO is more correlated than other cuprates, lying close to this crossover. Remarkably, this conclusion echoes an early neutron scattering study, which claimed that LSCO is close to an instability that is off of the physical parameter plane⁵⁹.

We study two important cuts in t'/t - t''/t space, a minimal cut (t''=0) and the Pavarini-Andersen [PA] cut (t''=-t'/2) – the latter seems to best capture the physics of the cuprates. By tuning t' we unveil the origin of the condensation bottleneck as a localization-delocalization crossover tied to the crossover from (π, π) - to FS-nesting. As a byproduct, we gain insight into why LSCO is so different from other cuprates, and how cuprates evolve from the pure Hubbard limit (t'=0).

Figure 14(a) shows the *T*-evolution of ξ_{th} for several values of t' along the minimal cut, including the data of Fig. 13(b). For t'/t > -0.17, the system is characterized by commensurate (π, π) order with an effective Neel temperature $T_N \sim 1000$ K [the correlation length grows so rapidly that interlayer correlations will drive a transition to full 3D order]. Similarly, for t'/t < -0.345 there is incommensurate $(\pi, \pi - \delta)$ order with T_N about a factor of 10

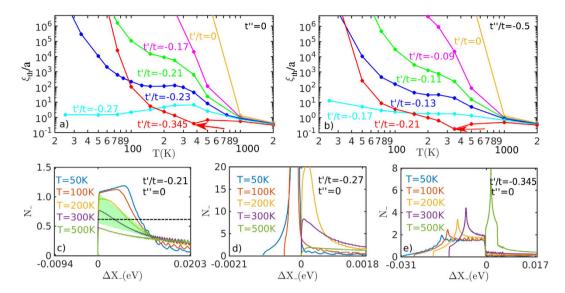


Figure 14. t' dependence of correlation bottleneck. (a) ξ_{th} for the minimal reference family at several values of t'/t [see legend] and t'' = 0. (b) ξ_{th} for the PA reference family at several values of t'/t [see legend] and t'' = -t'/2. (c-e) SDOS at t'' = 0 and three values of t'/t = -0.21 (b) [as in Fig. 13(b)], -0.27 (c), and -0.35 (d). All curves are shifted to line up the (π, π) data at $\Delta X_{-} = X_{-} - X_{-}(\pi, \pi) = 0$. Arrows in (a,b) show dip in ξ_{th} associated with the commensurate-incommensurate transition.

smaller. But for intermediate t'/t the *commensurate-incommensurate transition* is highly anomalous, with correlation length orders of magnitude smaller than expected. Figure 14(b) shows that a similar evolution follows along the PA cut in parameter space. The reason for this anomalous behavior can be seen by looking at the leading edge SDOS in the crossover regime, shown for three values of t' along the minimal cut in Fig. 14(c–e). For ease in viewing, these curves have been shifted to line up the SDOS at (π, π) at all *T*. It is seen that the anomalous collapse of ξ_{th} is associated with a rapid growth of the step height $N_{OP} \sim 1/A_2$, culminating in a near-divergence at $t'_c = -0.27t$, where the leading edge curvature A_2 goes to zero, Fig. 2(b). Note that ξ_{th} drops by 9 orders of magnitude at T = 200 K when t'/t changes from -0.17 to -0.27, then grows by a similar amount at 100 K when t'/t changes from -0.27 to -0.345. This transition may be hard to detect experimentally, since over this same range, the $\xi_{1/2}$ will be frozen at the value corresponding to the half-width of the (π, π) plateau. The green shaded region in Fig. 14(c) shows that the range of the anomalous growth (II) of $\xi_{1/2}$ in Fig. 13(b) coincides with the range of rapid growth of the SDOS leading edge; such behavior is absent if a *T*-independent OZ form (black dashed line) is assumed.

From the relationship between ξ_{th} and the step height A_2 , Eq. 18, we see that a diverging A_2 will cause $\xi_{th} \rightarrow 0$. The divergence arises when the susceptibility at (π, π) crosses over from a maximum to a local minimum. In the latter case, the maximum intensity of $\chi(q, 0)$ is spread along a 'ring' in *q*-space surrounding (π, π) , so that the inverse susceptibility resembles a 'Mexican hat', Fig. 2(c). Approximating the χ^{-1} -dispersion by a Mexican hat form $\chi_0^{-1} - A_2 q^2 + A_4 q^4$ leads to a threshold divergence $N_- \sim (\chi^{-1} - U_c)^{-1/2}$ as A_2 passes through zero, similar to Fig. 14(d) or Fig. 15(a). Note that the SDOS-divergence resembles a 1D VHS in the conventional DOS, even though here the 1-D direction is the radial direction away from (π, π) . The divergence of χ along a ring can be thought of as a 2D analog of Overhauser's effect²⁴. Similar effects are found in Bi2201, next Subsection.

The strength of the anomaly in ξ_{th} at $t'_c = -0.27t$ suggests that it is not an ordinary quantum critical point. Indeed, from Fig. 3 it can be seen to be a 'supertransition' between two domains of attraction, one commensurate at (π, π) , the other incommensurate at $(\pi, \pi - \delta)$ (compare Fig. 3(a,b) at x = 0). Such a supertransition is known from spectrum-generating algebras (SGAs) to be highly anomalous. The diverging OPDOS at the supertransition, also characteristic of SGAs, indicates that many different *q*-vectors compete simultaneously, frustrating the divergence of any particular mode. This can be clearly seen in Fig. 2(b), where the susceptibility is nearly flat. This is the electronic analog of McMillan's phonon entropy: if many phonons are simultaneously excited, the transition is suppressed to anomalously low temperatures. At this transition the *ground state is infinitely degenerate*, representing an anomalous spin frustrated state arising in the absence of disorder, with the frustration arising from strong mode coupling, although whether it represents a spin liquid or spin glass requires further analysis. Note also that when the commensurate transition ξ_{th} is strongly suppressed, $\xi_{th} < a$, arrows in Fig. 14(a,b).

For LSCO, the anomaly in Fig. 13(b) (large red dot in Fig. 8(c)) seems to fall near the coherent-incoherent crossover (pink shaded region), consistent with an early neutron scattering result which indicated that doped LSCO is close to a magnetic QCP⁵⁹. This proximity to a novel disorder-free spin-glass or spin-liquid QCP may play a role in stripe physics. Notably, in LSCO, commensurate (π , π) order disappears rapidly by ~2% doping, being replaced by incommensurate magnetic fluctuations and low-*T* spin-glass effects. Moreover, the commensurate-incommensurate transition involves a highly disordered regime separating two well-ordered phases (Fig. 14(a,b)). Since ordering tends to lower the free energy of the electronic system, the disordered

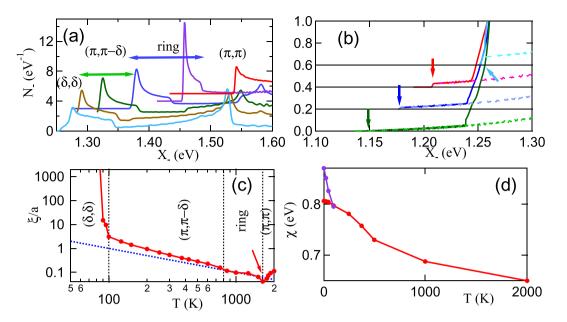


Figure 15. Structures in Bi2201 correlation lengths. Temperature dependence of: (a) SDOS N_{-} for t - t' - t'' reference state of Bi2201, x = 0.20. Datasets are offset for clarity; from highest to lowest, T = 2000 (red), 1000 (violet), 500 (blue), 375 (green), 250 (brown), and 100 K (light blue). (b) Continuation of SDOS in (a) to lower *T*, showing evolution of the leading edge. From top to bottom: T = 100 (light blue), 50 (red), 25 (blue), and 5 K (green). For each curve, a partial SDOS near the ANN peak is superposed to show the SDOS associated with the ANN peak whose onset is marked by arrows. (c) Correlation length vs *T*. Blue dotted line represents $\xi_{1/2} \sim T^{-1}$. Arrow shows dip in $\xi_{1/2}$ associated with commensurate-incommensurate transition. (d) Susceptibility of $\chi(q)$ peak vs *T* for *q* near the (π, π) -plateau (red line) or for ANN nesting (blue line).

regime represents a state of high free energy, and doping across this region can lead to a regime of [nanoscale] phase separation (NPS), which in LSCO is manifest as the stripe $phase^{60}$. It should be noted that also in Cr the commensurate-incommensurate AF transition is first order⁶¹. The very different situation in most other cuprates is discussed in the next Subsection, where it is shown that high-*T* magnetic fluctuations give way at low *T* to a fragile CDW order, consistent with experiment.

Finally, we note that the t' of the Mott-Slater transition is also the point at which the DOS at the bottom of the band changes from a local minimum to a local maximum (green dotted arrow in the inset to Fig. 5(c) – see also Fig. 2), and close to the point at which a nematic transition [splitting the degeneracy of the VHSs at $(0, \pi)$ and $(\pi, 0)$] becomes possible⁴³.

Mode coupling in Bi2201. We noted earlier that the OZ form should not be used in quantum critical theory when FS nesting is present⁵³. The problem can however be addressed by using SDOS. The issue is that at $T = 0 \chi_0$ is nonanalytic, with cusp-like peaks, Fig. 12, so the leading *q*-dependence is not parabolic. Indeed, since the dominant instabilities tend to be associated with double-nesting, the peak shape can be quite anisotropic, $\sim |\overline{q}|$ in some directions and $\sim |\overline{q}|^{1/2}$ in others³². There is a simple relation between the \overline{q} -dependence of the peak susceptibility and the form of $N_-(X_-)$ at threshold. If the SDOS varies as a power law, $N_- = \partial \overline{q}^2 / \partial X_- \sim X_-^p$, then at threshold, $(X_- - U_c) \sim \overline{q}^{2/(p+1)} + \xi_{th}^{-2/(p+1)}$. For a parabolic onset $(X_- - U_c) \sim \overline{q}^2 + \xi_{th}^{-2}$, or a step in N_- at threshold (p=0 or $N_- = const > 0$ for $X_- > U_c$). Instead, for Bi2201 there are cusps in χ_0 near $(\pi, \pi - \delta)$ and an antinodal cusp at (δ, δ) . For both cusps, a smooth power law is found, with $p \sim 2$, or $(X_- - U_c) \sim |\overline{q}|^{2/3} + \xi_{th}^{-2/3}$ at threshold. However, this leads to new problems. When a term $|\overline{q}^p|$ for $p \le 1$ is substituted into Eq. 14, the integral converges at the value of the Alexand for the case can be seen from Fig. 12: at any

However, this leads to new problems. When a term $|\bar{q}^p|$ for $p \leq 1$ is substituted into Eq. 14, the integral converges at threshold, so the Mermin-Wagner effect appears to be absent. That this is not the case can be seen from Fig. 12: at any finite *T*, the cusp is rounded off, and χ_0 varies parabolically, which means A_2 is finite. The cusps still produce a strong effect in that A_2 has an exceptionally large *T*-dependence, $A_2 \rightarrow \infty$ as $T \rightarrow 0$. This leads to an extra strong divergence of ξ_{th} Eq. 18, at $T \rightarrow 0$. But, as we increase *T* from zero, this also means that the corresponding ξ_{th} decreases rapidly, leading to a very fragile phase. We have separately confirmed that the order is also sensitive to ω and disorder.

In Bi2201 and most other cuprates, a larger ratio |t'/t| expands the (π, π) -plateau, leading to a rich phase diagram, Fig. 3(b). In this case the b-VHS is dominant only at high *T*, while FS-related susceptibility features cause a crossover of the (π, π) susceptibility from a maximum at high *T* to a local minimum at lower *T*. This leads to a different kind of strong mode coupling at intermediate *T*, best exemplified by the ring phase, Fig. 4(d,f). In this phase, the leading-edge SDOS has a 'Mexican hat'-like divergence (previous Subsection), Fig. 15(a) near 2000 K, leading to a slow growth in $\xi_{th} \sim T^{-1}$, Fig. 15(c), quite similar to that observed in doped YBCO^{62,63}. Near the onset of ring order, T = 1625 K, the correlation length has a nonmonotonic *T*-dependence, similar to that seen in LSCO, Fig. 14(a,b). While many pseudogap-like features are similar to LSCO, Bi2201 has some novel properties associated with the ANN instability, which can become a CDW when charge order is included in the model. Since the ANN susceptibility peak has a much larger intrinsic curvature, mode-coupling effects are weak. Hence the ANN susceptibility peak grows rapidly at low temperature, and as doping increases, it can surpass the (π, π) -peak, thereby sharply cutting off the growth of near- (π, π) fluctuations. This is illustrated in Fig. 15 for x = 0.20. Notably, since this peak is not associated with strong mode coupling it has a weak SDOS with a very strong *T*-dependence, Fig. 15(b,d), crossing the AFM peak near T = 100 K. The large subsequent growth of the ANN correlation length would ensure the rapid establishment of a true long-range 3D ANN order. However, corresponding to the large T-dependence of χ_{ANN} , there is a similar strong disorder dependence – particularly in the presence of short-range near- (π, π) order, which should cause the ANN correlation length to saturate.

Discussion

Cuprate pseudogap. In Figs 1 and 8, and Results Section (DFT-Lindhard), we discussed a revised VHS scenario in which a number of VHS related features evolve nearly linearly in doping and scale with T_{VHS} . We noted that this scaling resembles the pseudogap crossover $T^*(x)$. Here we provide additional evidence for this identification. Note that T_{VHS} is typically close to the measured T^* in most cuprates, including YBCO and Hg-cuprates⁹ (yellow-green line in Fig. 8) and LSCO (magenta dotted line)⁵¹. Furthermore, the pseudogaps in Bi2201 and Bi2212 terminate near the point $T_{VHS} \rightarrow 0$, i.e., at the conventional VHS^{34-36,64,65}.

The phase diagram of Fig. 8 has a number of further consequences for cuprate physics. First, the VHS onset near T_{VHS} combined with the coherent - incoherent crossover (pink shaded region in Fig. 8) can explain the anomalous transport properties found near the pseudogap, Fig. 1. This parallels transport, where for $T > T^*$, the resistivity ρ varies linearly with T^9 , a behavior expected near a VHS⁶⁶. For lower T, the resistivity is mixed, but $\rho \sim T^2$, as expected for a coherent Fermi liquid below a $T_{coh} < T_{VHS}$. The brown long-dashed line in Fig. 8 represents the corresponding T_{coh} found in YBCO and Hg-cuprates⁹, in excellent agreement with the coherent-incoherent crossover.

We note further that the compressibility should have a peak similar to γ that extrapolates to the VHS doping as $T \rightarrow 0$, consistent with our Fig. 8. In a related cluster-DFT calculation⁶⁷, the pseudogap is associated with a Widom line, related to a peak in the compressibility. This is further discussed in Supplementary Material Section IV. Notably, the Widom line terminates in a first-order phase transition near the VHS. The similarity of this phase diagram to the excitonic liquid-gas transition found in semiconductors⁶⁸ suggests that the b-VHS may be best described as a VHS-exciton^{69,70}.

This picture bears a resemblance to the Barzykin-Pines model of the cuprate pseudogap⁷¹, identifying T_{VHS} and the coherent - incoherent crossover with T^* and $T_{coh} < T^*/3$ in their model. Since their model is related to Kondo lattice physics⁷², the question arises as to whether a similar mode-coupling calculation in heavy-fermion compounds could lead to a similar anomalous entanglement at the f-electron incoherent-to-coherent transition.

Classifying order parameters. Competition between order parameters also arises in a totally different field, nuclear physics, where order parameters are formally classified in terms of a spectrum generating algebra (SGA) (Supplementary Material Section I.C). Order parameters are grouped into families (Lie subalgebras), and an order parameter density-of-states (OPDOS) is defined. In addition to conventional phase transitions, there are 'supertransitions' (called excited-state quantum critical points) where the system is driven across a boundary between two families and the OPDOS diverges. As a byproduct of this work, we develop a similar classification of order parameters for condensed matter systems, where similar phenomena can be observed (Supplementary Material Section I.A).

Strong coupling physics. The condensation bottleneck that we identified in Fig. 14 appears to be a previously unexpected *transition between Mott and Slater physics*, shedding light on the controversies between Anderson and Scalapino^{73,74}. On the Slater-side of the line (larger |t'|), the cuprate physics is dominated by FS nesting with Neel temperatures (strictly, $\xi_{th}/a \sim 1000$) near 100 K, while on the Mott side, the Fermi surface plays no role, and the magnetic transition is at $T \sim 1000$ K, and always lies at (π, π) due to VHS nesting. Thus, the phenomenological strong-coupling Yang-Rice-Zhang model⁷⁵ can only work on the Mott side of the transition, since it requires (π, π) nesting. However, at the crossover, the correlation length remains nanoscale, leading to strong mixing of the two competing phases. Among the physical cuprates, all fall into the Slater domain, except for LSCO, which lies close to the Mott phase. Notably, LSCO shows the strongest signs of nanoscale phase separation in the form of stripe physics [Results Section (Beyond RVB)]. This result is consistent with a recent first-principles calculation of the Hubbard U, which finds LSCO more strongly correlated than other cuprates⁷⁶.

We recall that our self-energy formalism⁴⁰ is able to reproduce most spectral features of the cuprates in terms of competition between d-wave superconductivity, a (π , π) ordered AFM phase, and an ANN CDW. We reproduce not only the photoemission dispersions, limited to the lower Hubbard bands, but optical and x-ray spectra, which depend sensitively on the Mott gap. In a related 3-band model, we reproduced the Zhang-Rice result that the first doped holes are predominantly of oxygen character, with overall dispersions at half-filling agreeing with subsequent DMFT results⁴⁰. While our earlier calculations predicted long-range (π , π) AFM order to persist to too high *T*, this is corrected in the present calculations which have only short range order and reproduce the upper and lower Hubbard band dispersions with broadening ~1/ ξ_{th}^{23} . Moreover, the calculated $\xi_{1/2}$ values (filled blue dots in Fig. 13(c)) agree with experiment in LSCO (green dot-dot-dashed line)⁵⁸. The experimental data are shown only above 300 K, since at lower *T* interlayer coupling drives a transition to long-range order. In the Heisenberg model, $\xi_{1/2}$ is a function only of *T/J*, and these data have been used to adduce the value of the exchange *J*. Hence, our mode coupling calculation successfully reproduces the crossover from the *U*-scale to the *J*-scale physics, as was found earlier for electron-doped cuprates²³.

For more insight into the strong coupling limit, we note that in the two-particle self-consistent approach⁷⁷, λ is determined by a sum rule involving double occupancy, which is fixed by assuming

$$U_{sp}/U = \langle n_{\uparrow}n_{\downarrow} \rangle/(n/2)^2$$

which leads to a saturation of U_{sp} as $U \to \infty$ (or $\langle n_{\uparrow} n_{\downarrow} \rangle \sim 1/U \sim J$). In our calculation, this saturation arises naturally, since U_{sp} can never exceed U_c .

VHS and excitons. While we have demonstrated the important role of the VHS in pseudogap physics, we have not addressed the deeper question of the physical significance of the VHS. We suggest that the significance is primarily topological: the saddle-point VHS is associated with the crossover of a Fermi surface from electron-like to hole-like, and hence represents the point nearest to electron-hole symmetry. This explains its strong role in transport, Fig. 1. While electron-electron scattering is strong, it usually plays only a small role in transport, since if an electron-hole collision has a large effect on current⁷⁸.

Now strong electron-hole scattering is indicative of a tendency to exciton formation, and we believe our results can most easily be understood in terms of a VHS-driven excitonic instability. Thus, the large entropic effects are related to a broad spread of the electrons in *k*-space, consistent with a localization of the real-space electron-hole separation in an exciton. A two-band excitonic insulator transition is analogous to a superconducting BCS-BEC transition; we suggest that in our one-band case this becomes the Mott-Slater transition we have found, with the Mott (BEC) phase dominated by preformed excitons. It is often stated that excitonic effects are unimportant in metals due to screening. However, we are studying the Hubbard model, where excitons could be present near half-filling, and their evolution with doping could contribute to a metal-insulator transition.

Conclusions

The present paper has provided significant progress toward DFT corrected MBPT. (1) We have demonstrated the significant role of bosonic entropy in driving short-range order. Our computational scheme incorporates the DFT-Lindhard function, and should capture strong correlation effects, at least for $T_N \sim 1/U$. (2) We have shown that the DFT-Lindhard function can be used to classify phase transitions, and may yield a condensed matter analog of SGAs. (3) We have shown that entropy effects in a one-band model involve a non-FS contribution associated with the VHS. In turn, these entropic effects modify the VHS physics, causing the spectral weight associated with the VHS [in the heat capacity and susceptibility] to shift rapidly towards half-filling as *T* increases, instead of following the evolution of the (q=0) DOS. This large shift suggests that the role of the VHS in pseudogap physics needs to be reexamined.

Finally, we comment on the limitations of the present mode-coupling framework. In particular, one would like to go beyond the present results to full inclusion of vertex corrections and *excitonic physics*. Another issue is the *evolution of the Fermi surface in the regime of short-range order*. As the FS shrinks from a large to a small size, the susceptibility will also change, and drive changes in the fluctuation maps and the associated properties discussed in this study such as the appearance of a second phase out of a regime of short-range order of a primary phase. With a DFT-corrected Lindhard function, one should be able to obtain insight into these issues.

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Acknowledgements

This work is supported by the US Department of Energy, Office of Science, Basic Energy Sciences grant number DE-FG02-07ER46352, and benefited from Northeastern University's Advanced Scientific Computation Center (ASCC) and the allocation of supercomputer time at NERSC through grant number DE-AC02-05CH11231. Part of this work was done while RSM was on sabbatical at the Advanced Light Source in Berkeley and the Los Alamos National Laboratory, where he benefitted from numerous discussions.

Author Contributions

R.S.M., I.G.B., P.M., C.L. and A.B. contributed to the research reported in this study and the writing of the manuscript.

Additional Information

Supplementary information accompanies this paper at http://www.nature.com/srep

Competing Interests: The authors declare no competing financial interests.

How to cite this article: Markiewicz, R. S. *et al.* Entropic Origin of Pseudogap Physics and a Mott-Slater Transition in Cuprates. *Sci. Rep.* **7**, 44008; doi: 10.1038/srep44008 (2017).

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