Precursor Pairing Correlations and Pseudogaps

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Abstract

I begin by briefly reviewing various experimental results on the pseudo-gap phenomena in underdoped cuprates. I argue that, taken together, all of these lead to a picture of singlet pairing above \( T_c \). I then explore the idea that the pseudogap is a normal state precursor of the superconducting gap due to local, dynamic pairing correlations in a state without long range phase coherence. Early work on simple model systems which exhibit pseudogap anomalies in the normal state of 2D superconductors in a low density, small pair size regime is reviewed and critically re-examined in view of more recent developments. I also describe recent studies of how the underlying d-wave superconducting ground state affects the anisotropy of the pseudogap and the destruction of the Fermi surface.

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I. Introduction

The deviations from Fermi liquid theory (FLT) in the normal state of high \( T_c \) superconductors are now well accepted as experimental facts, even if there is no theoretical consensus about their origin. These anomalies were first established from transport studies in the optimally doped materials, i.e., those with the highest \( T_c \) within a given family. Angle-resolved photoemission spectroscopy (ARPES) then showed that there is a well-defined Fermi surface even though there are no sharp quasiparticle excitations in the normal state. It has recently become clear that the underdoped cuprates exhibit even more remarkable deviations from FLT than the optimally doped materials: not only are the quasiparticles not defined, but even the Fermi surface becomes ill-defined due to the opening of a pseudogap above \( T_c \).

In these lectures, I will first describe the experimental evidence for the pseudogap phenomena as obtained by various probes: NMR, optics, thermodynamics, \( \mu \)SR, photoemission and tunneling. I will then argue that these remarkable observations point to the existence of singlet pairing correlations above the superconducting \( T_c \). I will next describe calculations on model systems that capture some of the physics of the pseudogap materials. The pseudogap is argued to be a natural consequence of local, dynamic pairing correlations in the normal state of low density, small pair-size superconductors.

II. Pseudogap Experiments

I will briefly review experiments on the underdoped cuprates in the pseudogap regime
regime above $T_c$. The word “pseudogap” is used to mean a large suppression of low frequency spectral weight (in contrast to a “hard” gap, which signifies strictly zero spectral weight).

**Spin Dynamics:** Some of the earliest evidence for the suppression of low frequency spectral weight above $T_c$ in underdoped materials came from NMR experiments \[1\]. The spin susceptibility $\chi(T)$, deduced from the Knight shift, is found to be strongly $T$-dependent, decreasing by a factor of three as the temperature is reduced from 300K to $T_c = 60K$ in underdoped YBCO, and then decreases smoothly through $T_c$. At the same time the relaxation rate $1/T_1T$ on the $^{17}$O and $^{89}$Y nuclei, (where the form factors filter out the antiferromagnetic contribution) was also found to be strongly $T$-dependent, with $1/T_1T \sim \chi(T)$. Note that in conventional metals both these quantities are $T$-independent; Pauli spin susceptibility and Korringa relaxation $1/T_1T \sim \chi^2$. For the $^{65}$Cu nuclei the relaxation rate $1/T_1T$ shows a non-monotonic $T$-dependence. With decreasing $T$, it first increases, presumably reflecting a build up of antiferromagnetic correlations, but eventually decreases below 150K, showing that a pseudogap is opening up even for the excitations probed by the Cu nuclei.

All of these anomalies, are collectively called “spin gap” behaviour. I emphasize that, even more so than the spin susceptibility, it is the rapid drop in $1/T_1T$ with decreasing $T$ which directly gives evidence for a pseudogap since $1/T_1T = \lim_{\omega \to 0} \sum_q F(q) \text{Im} \chi(q, \omega)/\omega$ directly probes the low frequency spectral weight in the spin excitation spectrum. (Here $F$ is the form factor which is different for different nuclei).

**Charge Dynamics:** Early studies of the in-plane conductivity $\sigma_{ab}(\omega)$ of YBCO as a function of doping reveal several important features \[2\]. First, the inferred scattering rate $1/\tau$ increases with underdoping, suggesting that at high temperatures the electronic excitations in underdoped materials are even less sharply defined than at optimal doping. Second, there is a considerable suppression of $\sigma_{ab}$ at finite frequencies below 500 cm$^{-1}$ which begins at temperatures much above $T_c$; the onset temperature for this gap-like feature increases with underdoping. Remarkably $\sigma_{ab}(\omega = 250 \text{cm}^{-1}; T)$ has the same qualitative $T$-dependence as $1/T_1T$ for $^{17}$O (see Fig. 4 of Rotter et al. \[2\]). Thus the charge response appears to show the same suppression of spectral weight as the spin response.

However, there is one crucial difference in the low frequency limit of the charge response. Recent optical data at lower frequencies show the appearance of a narrow “Drude-like” peak which shows directly that there are low lying charge excitations \[3\], unlike in the spin channel. This is consistent with earlier d.c. resistivity studies \[4\] in underdoped systems which see a characteristic deviation below the linear resistivity, implying that $\sigma_{ab}(\omega = 0; T)$ increases slightly as $T$ decreases unlike the higher frequency pseudogap behavior noted above. We shall return later to the question of these low lying charge excitations in the a-b plane.

There are many aspects of the c-axis transport which are puzzling, some of which appear to be directly linked to the opening up of a pseudogap. The c-axis optical conductivity \[3,5\] in underdoped YBCO clearly shows loss of spectral weight below 500 cm$^{-1}$ whose onset temperature is much above $T_c$. In contrast to the in-plane conductivity, the pseudogap $\sigma_c(\omega; T)$ appears to directly extrapolate to the inverse of the d.c. resistivity (i.e., there is no narrow peak at low frequencies). One of the puzzling features is the manner in which the optical conductivity sum rule is satisfied: since no weight is building up near $\omega = 0$ (as in the ab plane), it must be transferred to above the gap, but this is not apparent even
up to very high frequencies. Finally, it is found that, at fixed $\omega$ in the pseudogap region, \( \sigma_c(\omega; T) \) tracks the $T$-dependence of the spin susceptibility, once again showing that the spin and charge degrees of freedom are both getting gapped out, and that the c-axis transport is probing the pseudogap in the a-b plane excitation spectrum.

**Specific Heat:** The specific heat experiments of Loram and coworkers on YBCO as a function of doping have been discussed in detail by W. Y. Liang at this school, so I will be very brief. Specific heat probes all degrees of freedom including spin and charge excitations. Once the background subtractions for the lattice contribution are made, the electronic contribution gives direct evidence for a pseudogap [6] developing above $T_c$ with increasing underdoping. In the underdoped samples, the phase transition looks less and less mean-field like, and furthermore, loss of entropy begins to occur at temperatures much higher than $T_c$. The $T$-dependence of the entropy has been argued to be indicative of a degenerate normal state and not consistent with the thermodynamics of a gas of preformed bosons (which, by definition, would be a non-degenerate system above its $T_c$). A direct measurement of the electronic chemical potential of the underdoped systems, along the lines of [7], would be very useful in establishing the quantum degeneracy of the system.

**\( \mu \)SR:** Next we turn to the \( \mu \)SR measurements of the low temperature penetration depth \( \lambda_L(0) \), which is proportional to the $T = 0$ superfluid density $n_s(0)/m^\ast$, by Uemura and coworkers, the first of which appeared as early as 1989 [8]. A plot of $T_c$ versus $1/\lambda_L^2(0)$ for a large variety of materials shows the remarkable scaling $T_c \sim n_s(0)/m^\ast$ for all underdoped cuprates. This clearly suggests that transition is not of the standard BCS mean-field variety. The connection of this important observation by Uemura with pseudogap physics will be discussed later in Section VII.

**Photoemission:** Recent angle-resolved photoemission (ARPES) studies [9–11] on Bi-2212 compounds, have made very important contributions to our understanding of the pseudogap phenomenon. Prior to this work there was no clear understanding of many issues: Does the pseudogap only affect the spin degrees of freedom, or also the charge channel? (Although the optical conductivity data is quite unambiguous in this regard, nevertheless it was only the spin gap which tended to be emphasized in the literature). Does the pseudogap have any k-dependence? What is the connection, if any, of the pseudogap, above $T_c$ with the SC gap below $T_c$? The ARPES data provide answers to all of these questions.

The ARPES studies have found a highly anisotropic suppression of spectral weight that persists far above $T_c$. Near the $(\pi, 0)$ point the pseudogap is the largest and it closes only at a much larger temperature $T^*$, which correlates well with the crossover scale determined from other measurements. Along the zone diagonal the pseudogap vanishes. Remarkably, the pseudogap has the same magnitude (or scale) as the SC gap at each $k$ point (the maximum gap in optimal Bi-2212 is around 35 meV), and it evolves smoothly through $T_c$, with essentially the same $k$-dependence just above $T_c$ as the SC gap below $T_c$. The only difference in the spectroscopic data going through $T_c$ is that there are well defined quasiparticle peaks in the SC state but not above $T_c$. More recently it has been found [11] that both the pseudogap (above $T_c$) and the SC gap (below $T_c$) are “tied to” the Fermi surface that exists above $T^*$ when all gaps are closed. In other words there is an “underlying” Fermi surface which appears to have been gapped out in the pseudogap state. For a detailed review of these results, see the companion set of lectures on ARPES given at this school [12].

Very recently ARPES studies [13] have observed a pseudogap in Bi-2201, which correlates
well with the SC gap in that material (which is around 10 meV as opposed to the 35 meV scale in Bi-2212). This observation is particularly significant because Bi-2201 has one Cu-O layer per unit cell, and thus it shows that bilayers are not essential to the pseudogap phenomena.

**Tunneling:** Finally we conclude this Section by mentioning some very recent STM work by the Geneva group [14] on Bi-2212, in which a pseudogap is observed above $T_c$ in tunneling spectroscopy and found to scale with SC gap. Further, the tunneling spectra show smooth evolution through $T_c$ from the SC state to the pseudogap state.

### III. Survey of Scenarios

The experiments on underdoped cuprates described above give very clear evidence for a SC transition that looks quite different from the BCS mean field transition, and also for a “normal” state above $T_c$, with a pseudogap, that looks quite different from an ordinary Fermi liquid.

A common line of reasoning is that a theoretical understanding of the non-Fermi liquid normal state must come first because only then can one study the superconducting instability in that state. In these lectures, I will take a very different point of view and invert the question: given that the ground state is a superconductor made up of pairs of fermions, must the corresponding normal state necessarily be a Fermi liquid? As I will show the answer turns out to be “No”, and, in fact, under very general conditions the normal state has pseudogap anomalies arising from pairing correlations in a state without phase coherence [15–17]. In this scenario the pseudogap is a precursor to the SC gap. The ARPES data, which establish the connection between the SC gap below $T_c$ and the pseudogap above $T_c$, are a vindication of this approach.

But before turning to the main theme of these lectures, let me briefly discuss some of the alternative approaches to understanding the pseudogap; it is an impossible task to survey the whole literature. One general class of approaches is what I call “non-pairing” scenarios in which the pseudogap has no connection with the d-wave SC gap. In my view, it is very unnatural for non-pairing theories to obtain a pseudogap scale and anisotropy which is closely tied to the SC gap, and also to obtain the observed smooth evolution through $T_c$. Nevertheless, some of these ideas, particularly those related to antiferromagnetic correlations [18, 19], may be relevant to spectral weight suppression on the energy scale of $J \sim 100$ meV or more (for which there is some experimental evidence in ARPES), which is much larger than the SC gap scale (35 meV in optimal Bi-2212).

Another approach involves “spin-charge separation” and the pseudogap is associated with pairing of hypothetical $S = 1/2$, charge neutral fermions called spinons, rather than of real electrons. This approach has its origin in the early resonating valence bond (RVB) ideas of Baskaran, Zou and Anderson [20] which already contained the idea of singlet pairing without phase coherence. It was further developed by Kotliar and by Fukuyama [21], and particularly by Lee and Wen [22]. The “breakup” of physical electrons into spinons and holons (charged, spinless bosons) and their interaction with gauge fields has been central to the more recent developments. The hope is to understand not only the pseudogap phase, but the entire phase diagram from Mott insulator to the overdoped regime. It is not very clear, at least to me, if the holons and spinons have any physical reality (in a trivial sense they do not, since these are gauge-dependent objects) or if they are simply a more convenient basis
in which to do calculations (actually the problem turns out to be strongly coupled even in this basis).

It thus seems useful to ask if the intuitively appealing, original RVB picture of phase-incoherent singlet pairs can be studied in a simpler setting, with the more modest goal of understanding the pseudogap regime, rather than the full phase diagram. It is in this sense that the results discussed below tie in with the RVB idea.

I close this Section by pointing out some other proposals that have something in common with the pairing correlation ideas to be described next. The work of Emery and Kivelson [23] on classical phase fluctuations determining the underdoped $T_c$, and its relation to the present approach, will be discussed in detail in Section VII. In an early paper Doniach and Inui discussed the possibility of the insulator going into a state with incoherent pairs upon doping [24]. There are also several proposals involving mixed boson-fermion models, some of which [25] will be discussed elsewhere in the proceedings of this School, and others [26] will be mentioned later in Section VIII.

IV. Pairing Correlations above $T_c$

One of the reasons why the pseudogap experiments described in Sec. II seem surprising to us is that we are conditioned to think about superconductors in terms of the BCS mean field theory. There are two aspects to the BCS description: (a) pairing, which leads to a gap in the spectrum, and (b) phase coherence, or macroscopic occupation of the $q = 0$ pair state, which leads to the superflow properties. It is a very special feature of BCS theory, and of the conventional metallic superconductors so well described by this theory, that both of these effects happen together at the same temperature $T_c$: as soon as the pairs form they automatically condense. One only needs to look at other phase transitions in nature, say metallic magnetism, to see that this is an exception rather than the rule. In magnets the formation of the moments, the analogue of (a), and their ordering, the analogue of (b), occur at widely separated temperature scales: the Stoner temperature and the Curie temperature respectively.

The small parameter that justifies the BCS mean field approximation for conventional superconductors is $T_c/E_f \ll 1$ or $1/k_F \xi_0 \ll 1$: each pair contains a large number of other pairs within it. Given the low carrier concentration, and the small pair size in the cuprates, it is hardly surprising that one has to re-examine the validity of this approximation. In order to go beyond BCS theory, into a regime where there is no small parameter it is useful to consider the problem of the crossover from BCS pairing to Bose-Einstein condensation of tightly bound pairs. I will be very brief on this topic and refer the interested reader to ref. [27]. I emphasize that I am not claiming that by varying any experimental parameter in the high $T_c$ cuprates one is traversing such a crossover. The crossover picture allows us to bracket an interesting intermediate coupling problem ($\xi_0 \sim k_F$) between two known, and rather different, limiting cases (BCS and Bosons).

The simplest lattice model within which this problem can be studied is the attractive Hubbard model where the pair-size can be tuned by varying the strength of the on-site attraction $|U|$ relative to the nearest neighbour hopping $t$:

$$H = -t \sum_{i,j;\sigma} c_{i\sigma}^\dagger c_{j\sigma} + \text{h.c.} - |U| \sum_i n_{i\uparrow} n_{i\downarrow} + \mu \sum_{i;\sigma} n_{i\sigma}. \quad (1)$$
The chemical potential $\mu$ is used to fix the average density $\langle n \rangle$.

In Fig. 1 we show a schematic phase diagram of this model (away from half-filling so that the ground state is a superconductor). The weak coupling $|U|/t \ll 1$ limit is very well described by BCS theory. The pair size $\xi_0 \gg a$ (the lattice spacing) and upon increasing $T$ the destruction of pairs and of phase coherence occurs at the same $T_c \sim t \exp(-t/|U|)$. The normal state is a Fermi liquid. In the opposite extreme where $|U|/t \gg 1$ the ground state is a condensate of hard core lattice bosons. Upon heating, phase coherence is lost at $T_c \sim t^2/|U|$ (determined by the effective hopping for the bosons) and the state just above $T_c$ is a normal Bose liquid of tightly bound pairs. It is only at a much higher temperature $T^* \sim |U|$ that the pairs dissociate. Our main interest here will be the intermediate coupling regime, particularly above $T_c$.

A closely related continuum model has been studied in some detail in refs. [16,28] where a functional approach was used to study the global phase diagram at all values of the temperature and coupling. It was found that a mean field approximation is able to capture the physics of the SC ground state, and excitations, including collective modes, all the way from the BCS to the composite boson limit at $T = 0$ [28]. However, such an approximation is completely inadequate to describe even the qualitative physics of the normal state as one moves away from the weak-coupling BCS limit. There is a crossover temperature $T^*$ above which the system can be described in terms of free fermions, but below which pairing correlations become important, and a simple mean field description fails. With increasing coupling, or decreasing pair-size, $T^*$ and the transition temperature $T_c$ separate from each other and a proper treatment of fluctuations about the trivial saddle point is necessary to understand the temperature region between $T^*$ and $T_c$ [16,27]. For sufficiently strong coupling, in any dimension, $T^* \gg T_c$ and it is dynamic (frequency dependent) fluctuations which lead to this separation of scales. In 2D this separation between scales persists to much weaker coupling for two reasons: classical phase fluctuations, and the fact that arbitrarily weak attraction leads to the formation of independent bound states [29].

The functional integral approach described above led to a very useful interpolation scheme that gave reliable answers in both the BCS and Bose limits. However, it was least reliable in the intermediate coupling normal state where there is no small parameter to control the calculation. This, however, is the most interesting regime, since it is here that we expect to see characteristic deviations from Fermi liquid theory (FLT) as it crosses over from a weak coupling Fermi liquid limit to a strong coupling normal Bose liquid limit. The question then is whether there is a broad intermediate coupling regime in between these two limiting cases, especially in 2D, where the normal state is a degenerate Fermi system and yet exhibits non-Fermi liquid correlations for $T_c < T < T^*$? And, if so, what are the characteristic deviations from FLT due to pairing correlations? We will use quantum Monte Carlo simulations to address this question in the next Section.

V. Pseudogap in s-wave Models

I now describe results on the anomalous normal state properties of the 2D attractive Hubbard model which have been obtained from extensive quantum Monte Carlo (QMC) simulations [15,17,30]. The finite temperature determinantal QMC, developed by Scalapino and co-workers [31], begins by writing down the coherent-state path integral for the lattice Fermi system with the imaginary time direction discretized. The (quartic) interaction term
is decoupled using a discrete Hubbard-Stratonovich decoupling, using Ising variables as the bosonic fields \[32\]. The Fermi fields are then integrated out leading to an effective bosonic field theory on a discrete space-(imaginary) time lattice. The functional integral over the bosonic (Ising) fields is then performed using Monte Carlo techniques, and various fermionic correlation functions are calculated.

This QMC method is uniquely suited to the problem at hand for a variety of reasons. First, it is an inherently non-perturbative, finite temperature technique. Second, it works best in the intermediate coupling regime (in weak coupling, the pair-size is larger than the finite size systems one can study, and in strong coupling there are large Trotter errors introduced by the imaginary time discretization and also the very slow exploration of phase space). Finally, the attractive Hubbard model does not have a “sign problem” at any filling, unlike most other interesting fermionic systems, thus making it possible to obtain very accurate results. This comes about due to a decoupling scheme introduced by Hirsch \[32\] such that the weight factors in the resulting bosonic theory are positive definite. There are, however, two limitations of all QMC calculations that we must face up to: finite system size and imaginary time. We have studied the properties of interest on lattices of typical size \(8 \times 8\), and in many cases up to \(16 \times 16\), to ensure that our results are not finite size artifacts. The size limitation is not very severe because we are interested in normal state physics, rather than that of the phase transition (at which the correlation length diverges). As regards analytic continuation from imaginary time data to real frequencies, this is a real limitation. In our own work we have developed some tricks for looking at some correlation functions, like the density of states at the chemical potential, and the NMR relaxation rate that allows us to finesse this problem under certain circumstances; see \[15,17\]. But there are no useful results available on many interesting quantities such as the d.c. resistivity. More recently there have been some results on spectral functions and on density of states obtained by using the maximum entropy method to do the analytic continuation \[33\].

The normal state anomalies described below are in fact seen over a broad range of parameters. Here I focus on results at a density \(\langle n \rangle = 0.5\), at a moderate coupling strength of \(|U| = 4t\). For this choice, we have the rough estimate \(k_F \xi_0 \sim 3\). All statistical error bars not explicitly shown in the figures are of the size of the symbols. From the decay of the SC order parameter correlation function we estimate \(T_c \approx 0.1t\); from the spin susceptibility we will deduce below that \(T^* \sim 1t\).

The uniform, static spin susceptibility \(\chi_s\) as a function of \(T\) is shown in Fig. 2. In a Fermi liquid \(\chi_s\) should be \(T\)-independent (Pauli susceptibility) at temperatures smaller than the Fermi energy. In contrast, we find a strongly \(T\)-dependent result with \(d\chi_s/dT > 0\) below a temperature scale \(T^* \approx 1t\) (in fact, we may use this to define the crossover temperature \(T^*\)). The tendency toward strong singlet pair fluctuations in the normal state leads to \(d\chi_s/dT > 0\). This behavior is easy to understand at large \(|U|\) where tightly bound singlet pairs do not contribute to \(\chi_s\) unless they are ionized. What is remarkable is that such correlations persist down to \(|U| = 4t\) where one has a degenerate Fermi system (as shown below).

I next turn to the single-particle density of states (DOS) \(N(\omega)\) where \(\omega\) is measured from \(\mu\). To obtain this quantity from QMC requires analytic continuation which is usually a highly nontrivial problem. We have devised a simple method \[17\] to estimate \(N(0)\) for temperatures much smaller than the frequency scale on which \(N(\omega)\) shows significant variation. Thus in
the absence of a “low frequency scale” we are able to extract \( N(0) \) from the local Greens function at imaginary time \( \tau = \beta/2 \). The resulting \( N(0) \) is plotted in Fig. 2 as a function of temperature. We see that with decreasing \( T \) a pseudo-gap develops at the chemical potential in the normal state, i.e., the DOS is depleted. To emphasize this \( T \)-dependence of the DOS we use the notation \( N_T(0) \). It is interesting to note that \( \chi_s(T) = N_T(0) \) (to within the errors inherent in extracting the latter). While this equality looks like that in Fermi liquid theory (FLT), note that both quantities are strongly \( T \)-dependent, in marked contrast to FLT.

We have studied the NMR relaxation rate \( 1/T_1 \) (using an analytic continuation method similar in spirit to the one described above for the DOS) and found \([13]\) that, unlike the Korringa behavior expected of usual metals, \( 1/T_1 T \) is strongly \( T \)-dependent and \( 1/T_1 T \sim \chi_s(T) \) as seen in Fig. 3. This behavior of \( 1/T_1 T \) is characteristic of a spin gap as discussed in connection with the experiments in Sec. II. Since there is no non-trivial momentum space structure in this model, we cannot discuss the differences between relaxation rates of various nuclei, which depends on a proper description of short range antiferromagnetic correlations. Our discussion only sheds light on those aspects of the data which are not affected by these AFM correlations.

To understand the origin of \( 1/T_1 T \sim \chi_s(T) \) we study \( K(q) = \lim_{\omega \to 0} \text{Im} \chi(q,\omega)/\omega \). We find \([17]\) that \( K(q) \) is of the form \( N_T(0)/\Gamma(q) \) where \( \Gamma \sim v_F q \) is essentially the same as that for the noninteracting system. Thus \( K(q) \) is suppressed more or less uniformly in \( q \)-space with decreasing temperature and its \( T \)-dependence is governed by the pseudo-gap in the DOS. Using \( 1/T_1 T = \sum_q K(q) \), the above form of \( K \) gives a natural explanation of \( 1/T_1 T \sim \chi_s(T) \).

Does the charge channel also exhibit this pseudo-gap? To answer this question we analyze the compressibility \( dn/d\mu \) obtained by numerically differentiating the average density \( \langle n \rangle \) measured as a function of \( \mu \). (We found that \( dn/d\mu \) shows much larger finite size effects than, e.g., \( \chi_s \), which forced us to work on large lattices \((L = 16)\). We see from Fig. 4 that, in sharp contrast to the one-particle DOS, \( dn/d\mu \) is very weakly \( T \)-dependent, i.e., there is no pseudo gap in the charge channel. While this \( T \)-independence superficially looks like FLT, we argue below that in fact it arises due to collective excitations.

It is perhaps worth noting that a simple RPA analysis (particle-hole bubbles) fails to account for the observed behavior in either the spin or charge channels (except for the general trend that attractive interactions lead to a decrease in \( \chi_s \) and an increase in \( dn/d\mu \)). For the susceptibility \( \chi_{\text{rpa}} = N^0/(1 + |U|N^0) \), where \( N^0(0) \) is the noninteracting DOS, RPA fails to account for the \( T \) dependence. On the other hand \( dn/d\mu_{\text{rpa}} = 2N^0(0)/[1 - |U|N^0(0)] \) predicts an entirely spurious instability (phase separation) at \( |U|N_0(0) = 1 \). In fact, pairs form at large \( |U| \) and their residual interactions are repulsive, so that the compressibility of the system remains finite for all \( |U| \). We have calculated \([17]\) the compressibility at \( T = 0 \) within a mean field (Hartree-Fock-Bogoliubov) approach which is capable of dealing with the \( T = 0 \) BCS-boson crossover. We find that it has the same order-of-magnitude as the normal state MC result (note that we do not expect \( dn/d\mu \) to change dramatically as \( T \) goes through \( T_c \)). This lends further support to the argument that collective excitations (singlet pair fluctuations) contribute to the charge channel.

Note that we find qualitative differences between the spin and charge response functions. Quantitative differences are well known even in a Landau FL where \( \chi_s \) and \( dn/d\mu \) are renormalized by different FL parameters. What we see here is much stronger: as a result of
strong interactions, $\chi_s$ and $dn/d\mu$ acquire qualitatively different $T$-dependences. The spin response is dominated by single-particle excitations ($S = 1/2, Q = 1$); this is $T$-dependent because the DOS for these excitations is depleted by collective excitations, i.e. pairs with $S = 0, Q = 2$. The collective excitations, on the other hand, directly contribute to the charge channel. At this point it may be useful to recall the discussion of the transport experiments in Sec. II, where clear evidence for low lying charge excitations was found, in spite of the fact that the single-particle excitations were (pseudo)gapped.

The momentum distribution function $n(k)$ is shown in Fig. 5. Note the rapid variation with $k$, the remnant of a “Fermi surface” though it is broadened by the temperature and by interactions through the pseudogap effect. This, together with the fact that the chemical potential $\mu \gg T$ (see ref. [15]), clearly shows that we are dealing with a highly degenerate Fermi system (and not with a system of preformed bosons!). Thus it makes sense to talk of deviations from Fermi liquid behaviour when discussing the above results. We thus see that the correlations in this regime show very interesting behavior which is intermediate between a Fermi liquid and a Bose liquid.

The DOS pseudogap in the normal state of the attractive Hubbard model has also been verified by self-consistent $T$-matrix calculations [34] and by analytic continuation of Monte Carlo data using maximum entropy methods [33]. The latter study also sees evidence for the precursor of the Bogoliubov-like dispersion of excitations associated with the pseudogap above $T_c$. There is an interesting recent paper [35] which introduces analytically tractable approximations leading to a DOS pseudogap.

VI. Commonly Asked Questions:

Before proceeding further, it may be useful to pause and answer several questions which are commonly asked of the author.

(1) **Does the pseudogap arise from “superconducting fluctuations”?** The answer is no, at least not in the conventional sense that the phrase “SC fluctuations” has been used in the literature. The conventional usage follows Aslamazov and Larkin [36] where it is used to describe small, but singular, corrections to weak coupling BCS theory in the vicinity of $T_c$. The choice of diagrams (the four shown in Fig. 1 of [36]) is controlled by a leading order perturbation theory in the small parameter $\max(T_c/E_f, 1/E_f \tau) \ll 1$, and these lead to significant effects in a small range of temperatures $(T-T_c)/T_c \ll 1$, where the SC correlation length is diverging.

In contrast, the pseudogap anomalies are *not* small corrections to a normal state background; these are large effects over a large temperature range (persisting up to temperatures which are several times $T_c$) and thus lie beyond the standard perturbative SC fluctuation approach. Nevertheless, the standard approach does contain within it the first hint of the pseudogap anomalies, as shown by Varlamov and coworkers [37], who have emphasized the importance of the “DOS” diagram: where the SC fluctuation propagator acts as a self energy dressing the electron propagator. A very similar approach has been taken by Rainer and coworkers (in largely unpublished work) who study systematic perturbative corrections to Fermi liquid theory [38].

(2) **Is two dimensionality and Kosterlitz-Thouless physics sufficient to give rise to the pseudogap?** The answer is no, by itself it is not sufficient. To see this most clearly consider a weak coupling BCS superconductor. Let us denote the BCS mean field
transition temperature by $T_c^0$, and the superfluid stiffness $D_s(T = 0) \simeq \hbar^2 n_s(0)/m^*$. In the weak coupling limit $T_c^0/E_F \ll 1$ which means that $D_s(0) \gg T_c^0$. We use the universal jump condition of Kosterlitz-Thouless (KT) theory $\hbar^2 n_s(T_c^-)/m^* = 2T_c/\pi$ to determine the actual (KT) $T_c$. To make a rough estimate, it suffices to use the mean field form $D_s(T) \simeq D_s(0)(T_c^0 - T)/T_c^0$ in this condition. We thus find that $T_c = T_c^0[1 - O(T_c^0/E_F)]$; (a more accurate theory gives a multiplicative logarithm in the correction term [39]). Since the KT $T_c$ and the mean field $T_c^0$ are so close to each other, there is no significant temperature range over which pseudogap anomalies can exist in the weak coupling limit even in 2D. This clearly points to the importance of being in a regime outside the weak coupling BCS limit.

(3) **Does the pseudogap arise from “preformed bosons”?** Again the answer is no, insofar as the experiments or the calculations presented above are concerned. Certainly if all electrons were tightly bound into “preformed boson” pairs, this would lead to a gap in the one-electron spectrum. But in this case the normal state of the system would be a completely non-degenerate system (since $T_c$ in Bose systems corresponds to the onset of degeneracy: thermal deBroglie wavelength becoming comparable to the interparticle spacing). To the extent that thermodynamic and spectroscopic studies, described above, find the underdoped cuprates to be (consistent with) a degenerate Fermi systems, a bosonic description is untenable. Another important point to mention here is one made by Leggett long ago, and described in some detail in [40]: in a large class of models the observation of nodes (zeros) of the SC gap is by itself sufficient to rule out the “bosonic” regime.

The upshot of all this is that one is neither in the BCS limit where small deviations from weak coupling theory can be understood in terms of the conventional theory SC fluctuations, nor is one in a preformed boson limit, where the state above $T_c$ is a normal Bose liquid, but rather in an intermediate coupling regime.

(4) **Why is a simple model like the attractive Hubbard model relevant for the problem at hand?** The first point to make is that the 2D attractive Hubbard model is not a microscopic model for the high $T_c$ cuprates; this cannot be overemphasized. It just happens to be the simplest lattice model in which: (a) one can study an intermediate coupling superconductor with $\xi_0 \sim k_F$; (b) one can see a clear separation between $T^*$, the crossover scale below which effects of pairing correlations become significant, and the transition temperature $T_c$ at which phase coherence is established; and (c) one can reliably calculate at least some experimentally interesting properties and establish pseudogap anomalies in the temperature range $T_c < T < T^*$.

The simplicity of this model certainly helps in that a new physical situation is being studied, and also it focuses attention on the important ingredients: low density, short coherence length, low dimensionality. Any other more realistic model which has a pseudogap arising from pairing correlations above $T_c$ will have to analyzed in an analogous manner.

At the same time, we must also remember the limitations of a simple model. An obvious one is that the attractive Hubbard model has s-wave pairing. The study of similar models but with a d-wave ground state is comparatively recent, and will be discussed in Section VIII. Another simplifying feature is that we are not capturing some aspects of the strong correlation problem. As a concrete example, note that in the simple models considered here, the electronic density $n$ determines both the size of the underlying Fermi sea, $k_F$, as well as the zero temperature superfluid stiffness $n_s(0)/m^*$ (the importance of which will be made
clear in next Section). In the underdoped cuprates, which are doped Mott insulators with \( x \) doped holes per Cu, the size of the Fermi sea may be large, scaling like \( (1 + x) \), (satisfying Luttinger’s theorem), and yet the superfluid density may be small \( n_s(0)/m^* \sim x \).

**VII. Phase Fluctuations and \( T_c \)**

In all the calculations reviewed above we looked at the properties of the system *coming down in temperature* from above \( T^* \) to below it, and asked how pairing correlations without long range order (LRO) affect normal state properties. It is very useful to ask the complementary question: what are the excitations about the SC ground state that destroy the SC LRO as the temperature is increased, and thus determine \( T_c \)?

An answer to this question was given recently by Emery and Kivelson (EK) in a very nice paper [23]. I will first describe their argument, clarify its relationship to the calculations described in Section V, and finally discuss its relevance to d-wave systems. Recall that in BCS mean field theory, the excitations responsible for loss of order are the (Bogoliubov) quasiparticles coming from thermally broken pairs, which lead to a collapse of the self-consistent gap at \( T^0 \). EK argued, in a model-independent manner, that in a system with small superfluid stiffness \( n_s(T = 0)/m^* \) phase fluctuations are soft and define a phase disordering temperature. For a layered system this is given by \( T^\theta_c \sim \hbar^2 n_s(T = 0)d/m^* \), where \( d \) is the layer spacing. (This is the analogue of the statement that in a magnetic system with well defined moments, the disordering temperature scale is set by the exchange coupling \( J \)). Further, the actual \( T_c \) is given by the smaller of \( T^\theta_c \) and \( T^0 \). In the underdoped materials, which are lightly doped Mott insulators with a low superfluid stiffness, \( T_c = T^\theta_c \sim n_s(T = 0)/m^* \). This also accounts for the Uemura scaling [8].

EK go on to emphasize that the coherence length or pair size \( \xi_0 \) does not enter their estimate, and hence their result describes a completely different physical situation from that discussed in Section V. I believe that this is not correct, for their calculation is based on the implicit assumption that the gap is the largest energy scale in the problem: \( \Delta_0 > T^\theta_c \) (or else BCS gap collapse would determine the phase transition). Writing \( \Delta_0 \sim \hbar v_F/\xi_0 \) and \( n_s(T = 0) \sim k_F^2/d \), the above inequality yields \( k_F \xi_0 < 1 \), which is *the same regime as the one we have been discussing*. Thus by working in the pair size comparable to interparticle spacing regime in simple model systems, we were simulating one of the key relevant conditions in a lightly doped Mott insulator where the superfluid stiffness is much smaller than the SC energy gap. (See also remarks at the end of Section VI).

For a system with an isotropic s-wave gap, phase fluctuations (in the sense of excitations of an XY model) are clearly the most important excitations not included in a simple mean field description. However, in a d-wave system with nodes in the gap the situation is even more interesting. As first pointed out by Lee and Wen [41] in a phenomenological analysis (which is quite independent of their microscopic ideas on spin-charge separation [22]), the SC LRO of the ground state is most effectively lost due to the excitations around the point nodes. These excitations lead to the well-known linear \( T \) reduction in the superfluid stiffness (this much is true even for a d-wave BCS SC). But since \( D_s(T = 0) \) is small to begin with (and this is the non-BCS part of the argument), this linear decrease by itself is sufficient to drive the system normal, without affecting the gap! This appears to lead to an even lower \( T_c \) estimate than that obtained from phase fluctuations alone, and it too naturally leads to the Uemura scaling.
Whether the $T$-dependence of the $n_s(T)$ is dominated by phase fluctuations themselves or by excitations near the nodes, the main conclusion is that, in the regime of interest, it is the vanishing of the superfluid density that controls the transition and determines $T_c$, and not the collapse of the gap (which is the main driving force in BCS). The normal state above $T_c$ must then still have a remnant of the superconducting gap, which is precisely the pseudogap scenario discussed in Section V. I next turn to the very important issue of how those results are modified when the ground state is a d-wave superconductor.

VIII. Pseudogaps in d-wave models

At this point it would be appropriate to discuss the d-wave analogue of the results described in Section V. However, there is no simple known model with a d-wave SC ground state on which reliable quantum Monte Carlo calculations can be done. (The main problem is the fermion sign problem in quantum Monte Carlo simulations). Thus a different approach is required. In an ongoing collaboration with Engelbrecht, Nazarenko and Dagotto, we have used self-consistent, conserving approximations. All of the results described below are based on ref. [42]. Let us consider a simple tight binding model on a 2D square lattice with dispersion $\epsilon_k = -2t(\cos k_x + \cos k_y)$, ($t = 1$), described by

$$H = \sum_{k,\sigma}(\epsilon_k - \mu)c^\dagger_{k\sigma}c_{k\sigma} + \frac{1}{N} \sum_{kk'q} V_{k,k'}c^\dagger_{k\uparrow}c^\dagger_{q-k\downarrow}c_{q-k'\downarrow}c_{k'\uparrow}$$  \hspace{1cm} (2)

where the separable potential $V_{k,k'} = -|U_d|(\cos k_x - \cos k_y)(\cos k'_x - \cos k'_y)$, is chosen to produce a d-wave SC ground state. The goal is to study the normal state properties of this model in a regime where $|U_d|$ is comparable to the bandwidth, and the system is presumably dominated by pairing correlations.

We use the self-consistent $T$-matrix method in which we study the renormalization of the Green’s functions by the exchange of pairing fluctuations. (This method has been used earlier by several authors to study the normal state of s-wave superconductors [34,43]). This requires a numerical solution of the following set of integral equations. The p-p channel vertex $\Gamma$ is given (symbolically) by

$$\Gamma = I - IGG\Gamma$$  \hspace{1cm} (3)

where $I$ is the p-p irreducible vertex and $G$ is the fully renormalized Green’s function

$$G = [G_0^{-1} - \Sigma]^{-1}$$  \hspace{1cm} (4)

defined in terms of the self-energy which satisfies

$$\Sigma = VGG\Gamma$$  \hspace{1cm} (5)

where $V$ is the bare two-body potential. These three equations are formally exact, and to make progress we need to specify an approximation for $I$; we make the simplest choice $I \simeq V$. In addition to being fully self-consistent, this approximation is also conserving in the sense of Baym-Kadanoff.

It is important to remember, however, that there is no small parameter which controls this calculation, and there is an on-going debate in the literature about the “best” set
of diagrams to retain [44]. This can really only be settled by comparing one’s favourite approximation with a more exact calculation such as quantum Monte Carlo, if it exists. Even though the details of the results may well depend upon the precise approximation scheme, I will emphasize those qualitative aspects which I believe to be robust on physical grounds. Also, some of the issues that arise in interpreting the results are clearly much more general, and go beyond the specific model or approximations being discussed here.

The main quantity of interest is the one-electron Greens function and the associated spectral function \( A(k, \omega) = -(1/\pi) \text{Im}G(k, \omega + i0^+) \), which is closely related to the ARPES intensity [45]. (I will not discuss here results on response functions like the spin susceptibility, for lack of space). From the spectral function we will deduce interesting results about pseudogaps and about the “Fermi surface”. The question of defining a “Fermi surface” at finite temperatures, especially in the absence of sharp quasiparticles, is one of general interest in the field of high \( T_c \) superconductors. (See the lectures on ARPES by the author and Campuzano at this summer school for more discussion on this issue [12]). We may identify points \( k^* \) in the Brillouin zone at which the spectral functions disperse through the chemical potential (\( \omega = 0 \)), i.e., \( A(k^*, \omega) \) has a dominant peak at \( \omega = 0 \). The “locus of gapless excitations” \( \{ k^* \} \) then generalizes the notion of a “Fermi surface” (FS). For a strongly correlated system at finite temperatures, this locus need not possess all of the properties that we associate with the usual \( T = 0 \) definition of a FS in a Fermi liquid. For instance, there is no guarantee that its shape or the volume enclosed by it is \( T \)-independent, and we indeed find this in our numerics. One may question this whole concept, but experimentally [12] it has been found in the optimally doped cuprates that a similar definition yields a “Fermi surface” which has many nice features. It is \( T \)-independent, it forms a continuous contour in the repeated zone scheme which is quite similar to what band theory predicts, and there is a rapid variation of the momentum distribution as one crosses this locus.

I will now discuss the results for \( U_d = -8 \) and \( n = 0.5 \) (quarter-filling) shown in Figs. 6 – 8; qualitatively similar results were found for other parameter sets. For details of the calculations, see ref. [42]. All the results are in the non-superconducting state above \( T_c \) as ascertained from the pair susceptibility, and in a degenerate regime, as seen from \( \mu \gg T \) and studies of the momentum distribution \( n(k) \).

The form of the interaction \( V_{k,k'} \) is such that electronic states near the zone diagonal \((0,0) \rightarrow (\pi,\pi)\) are unaffected by the interaction, while those near the zone corners \((\pi,0)\) are very strongly affected. In fact, we find that near \((\pi,0)\) the spectral functions acquire very large widths and a quasi-particle description is not valid at any temperature. Nevertheless, at high temperatures, there is an identifiable dispersion of the broad \( A(k,\omega) \) peaks (see Fig. 6) from which we can determine a “Fermi surface” \( \{ k^* \} \), even in the absence of well-defined quasiparticles, as discussed above. It forms a continuous contour in the repeated zone scheme, although, unlike the experiments [12], it is not \( T \)-independent in our calculations.

As the temperature is lowered below a scale \( T^* \approx 1 \) there is a qualitative change in the spectral functions; see Fig. 7. First the lineshape: the spectral weight of the one broad feature at high \( T \) is now redistributed into a multiple peak structure. We can get further insight into the lineshape by studying the real and imaginary parts of the self-energy \( \Sigma \) as functions of \( \omega \); for details see [42]. Second, the dominant peaks of \( A(k,\omega) \) now exhibit very anomalous dispersion. As \( k \) varies from \((0,0)\) to \((\pi,0)\), the spectral peak approaches \( \omega = 0 \) but never crosses it, either “bouncing back” towards negative \( \omega \) or “stalling” (depending on
This is a clear signature of a Fermi surface crossing being destroyed due to the opening up of a gap, even though there is no long-range order (more appropriately, algebraic order for the 2D model under study). This is very similar to the pseudogap seen by ARPES experiments on underdoped cuprates in the temperature regime $T_c < T < T^*$. In the spirit of the ARPES experiments we estimate the pseudogap magnitude by making scans through $k$-space and noting the position of the spectral function peak which is farthest to the right, i.e. at the largest frequency below zero. We then plot the spectral function pseudogap $\Delta_{ps}$ as a function of $\theta = \arctan(k_y/k_x)$ in Fig. 8 at two temperatures $T = 0.2t$ and $T = 0.75t$.

The first point to note in Fig. 8 is the strong anisotropy of the pseudogap, which is suppressed to zero in an arc about the diagonal. The extent of the nodal arc region, and the magnitude of the maximum gap at low $T$ are both sensitive functions of the choice of parameters as seen in the inset to this Figure. The second important point to note is the $T$-dependence of the pseudogap anisotropy which shows that small gaps are destroyed at lower temperatures compared with larger gaps. This leads to Fermi arcs (of nodes) which expand with $T$, as seen from Fig. 8. At sufficiently high temperatures, when the ordinary dispersion of the spectral functions has been restored, these disconnected arcs must merge to form the continuous Fermi contour above $T^*$.

All of these effects are rather remarkable. First consider the $T$-dependent anisotropy of the pseudogap. Contrast this with gap anisotropy in the broken symmetry state of a d-wave BCS superconductor, where the self-consistent gap equation leads to the solution $\Delta_k(T) = \Delta_0(T)(\cos k_x - \cos k_y)$. Thus the temperature and $k$ dependences factorize, and there is no variation of the anisotropy with $T$; gaps at different $k$’s vanish at the same temperature $T_c$ defined by $\Delta_0(T_c) = 0$. The behaviour of the pseudogap is qualitatively different in that small gaps collapse at lower temperature.

Viewed as a function of decreasing temperature, one obtains the picture of a continuous Fermi surface contour being destroyed in patches as the pseudogap begins to develop starting at the $(\pi,0)$ point and then proceeding towards the diagonal directions. This qualitative prediction of our calculations has been verified in very recent ARPES experiments by the Argonne-UIC group [46]. A schematic representation of the the Fermi Surface destroyed in patches is shown in Fig. 9. Although the novel $T$-dependence discussed above was first described in [42], the general picture shares some qualitative similarities with a phenomenological model proposed in ref. [26], which consists of bosonic pairs around $(\pi,0)$ and fermions in the rest of the zone interacting via a $k$-dependent coupling.

IX. Open Questions

There are many open questions even within the precursor pairing picture of the pseudogap. I will briefly list them here. The reason why the answers are not known, even for the simple models considered, is the lack of reliable, controlled calculations in the parameter regime of interest. Thus one is faced with one of the main dilemmas of the whole field of high $T_c$ theory. On the one hand, one needs to simplify models so that they become more tractable, on the other hand, one probably needs to incorporate more realistic features of the materials in order to better confront experiments.

Transport in the pseudogap regime is a very important open problem for the pairing scenarios. Qualitatively there are two competing effects that need to be considered. The
opening up of a pseudogap in the one-particle excitation spectrum implies that one-particle contribution to the d.c. resistivity $\rho$ will show an insulating upturn: increasing $\rho$ as $T$ decreases. However, we must also include the collective contribution to transport from the pairs, which will show up as a “Drude-like” peak in the the conductivity (a broadened version of the SC state delta function). This contribution, by itself, will show a marked down-turn in the d.c. resistivity with decreasing temperature (paraconductivity). The net effect of both these contributions need to be considered, and it is clear that sufficiently close to $T_c$ the pair-channel will dominate. (Interestingly, a similar interplay between the depleted one-particle contribution and the enhanced pair contribution was seen in the QMC calculation of the compressibility in Sec. V).

As described in Sec II, experimentally the d.c. resistivity shows only a distinct, but relatively modest (on the scale of other pseudogap effects) decrease away from linearity as $T$ decreases. It is often suggested that the spin-charge separation theories, briefly discussed in Section III, are able to naturally account for this. However, there are no quantitative calculations of charge transport by bosons interacting with gauge fields, and the arguments within that framework are also, at the present time, qualitative.

A second important point about transport in the real materials is that it is dominated by regions of $k$-space near the zone diagonal where the electronic states are much more dispersive (large Fermi velocity) than those near $(\pi,0)$; see [12] and references therein. This is also where the pseudogap effects are smallest. These strong $k$-dependences of the the electronic structure, the gaps and the interactions are likely to be crucial for a proper understanding of transport (and other) experiments. In particular, the opening up of the pseudogap starting near the $(\pi,0)$ point at $T^*$, and the way in which it progressively affects larger regions of the Fermi surface with decreasing $T$, as described in Section VIII, will be an important input to detailed transport calculations, as well for understanding the crossovers in other correlation functions.

The effects of magnetic fields in the pseudogap regime, diamagnetism, and magneto-transport are all questions for further study. The disconnected Fermi arcs of Section VIII may have profound consequences in the presence of magnetic fields.

Much remains to be understood about c-axis transport. As far as the pseudogap in c-axis optical conductivity is concerned one might ask why there are no low lying charge fluctuations like those seen in a-b plane transport? The qualitative reason is that tunneling of incoherent pairs along the c-axis is exponentially suppressed relative to the one-particle contribution (which is itself small). Hence transport across planes is dominated by the one-particle contribution which upon the opening of a pseudogap leads to an upturn in the c-axis resistivity with decreasing temperature.

Another very important question which I have not touched on in these lectures is: why are the quasiparticles as seen by ARPES destroyed above $T_c$? In the pseudogap regime, there is a suppression of low-lying particle-hole excitations, but nevertheless an electron can scatter off the pairs. As soon as the pairs become phase coherent below $T_c$ this scattering mechanism freezes out. Detailed calculations along these lines are needed to see if this mechanism is able to account for the experimental observations.

There is an important question on which more experimental input would be welcome. There seems to be some difference in the behaviour of the underdoped bi-layer materials compared with those which have a single CuO plane per unit cell (in particular the 214
LaSrCuO). There are suggestions that the spin-gap is entirely due to bilayer effects [47,48]. The recent ARPES observation of a pseudogap in the one-layer Bi-2201 [13] casts doubt on whether bilayers are necessary for this phenomenon. All of the models described in these lectures are single-plane models. Nevertheless, more studies of one-layer cuprates may shed light on whether or not there is an important distinction between one and two layer systems, especially in so far as the pseudogap is concerned.

Finally, direct experimental probes of pairing correlations above $T_c$ would be very important. Possible avenues for research include characteristic signatures in low frequency collective transport, and precursors of the Josephson effect between a pseudogap metal and a superconductor [49].

X. Conclusions

The main message of these lectures is that, in a sense, the pseudogap anomalies are not very mysterious. Although such deviations from Fermi liquid behaviour above $T_c$ are without precedent, nevertheless, on theoretical grounds they follow as soon as one has a superconducting ground state at $T = 0$ with a small superfluid stiffness and a large SC energy gap. In the simple models discussed above this was accomplished by being in a regime of low density and small pair size. Many open questions remain, including some questions of principle, as discussed in the preceding Section. Nevertheless, the pseudogap anomalies, by themselves, do not appear to require consideration of novel ground states or new types of excitations. Whether the non-Fermi-liquid anomalies above $T^*$, and in particular, those at optimal doping, (which were not discussed in this lecture) force us to consider such new states, or these are associated with even more complicated finite temperature crossovers, is not clear at the moment.

The key open questions are: why do the underdoped cuprates have such a small superfluid density – which is surely related to their being doped Mott insulators – and such a large effective pairing interaction? Quantitative comparison with the experiments must, therefore, await a controlled calculation based on a microscopic model which describes how a Mott insulator upon doping goes into a short coherence-length d-wave superconductor whose normal state is dominated by pairing correlations.

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FIG. 1. Schematic phase diagram of the attractive Hubbard model illustrating the crossover from a BCS superconductor with a Fermi liquid normal state (in the weak coupling $|U|/t \ll 1$ limit) to a condensate of bosons with a normal Bose liquid state above $T_c$ (for $|U|/t \gg 1$). $T_c$ is the transition temperature at which phase coherence is established and $T^*$ is a crossover scale below which pairing correlations become manifest in the normal state.
FIG. 2. The one-particle density of states $N_T(0)$ at the chemical potential (full triangles) and the spin susceptibility $\chi_s(T)$ (open squares) as a function of temperature. These results were obtained from Quantum Monte Carlo simulations [15,17] of the 2D attractive Hubbard model for $|U|/t = 4$, $\langle n \rangle = 0.5$ and $L^2 = 8^2$.

FIG. 3. Parametric plot of NMR relaxation rate $1/T_1 T$ vs spin susceptibility $\chi_s$ of the 2D attractive Hubbard model for $|U|/t = 4$, $\langle n \rangle = 0.5$ with temperature as the implicit variable. (Adapted from ref. [15]). The plot shows that the two quantities track each other: $1/T_1 T \sim \chi_s(T)$. Note that for a Fermi liquid all the data would collapse to a single point.
FIG. 4. The compressibility $dn/d\mu$ for $\langle n \rangle = 0.5$, $U = 0$ (full line) and $|U| = 4$ (open circles), as a function of $T$ obtained on a $16^2$ lattice. The $T = 0$ non-interacting and the $T = 0$ mean field result for $|U| = 4$ are also shown. (From ref. [17]).

FIG. 5. The momentum distribution function $n(k)$ as a function of $k$ for $|U| = 4t$, $\langle n \rangle = 0.5$, and $T = 0.25t$. (From ref. [17]).
FIG. 6. Self-consistent $T$-matrix results for the spectral function $A(k, \omega)$ for a sequence of momenta $k = (x\pi/32, 0)$ with $x = \{0, 6, 9, 12, 15, 18, 21, 24, 27, 32\}$ for $U_d = -8$ and $n = 0.5$ at a high temperature $T = 2.0$ ($T > T^* \simeq 1$). Inset: Brillouin zone with points indicating the momenta $k$ and solid curve showing the $T = 0$ non-interacting Fermi surface. (From ref. [42]).

FIG. 7. Spectral functions $A(k, \omega)$ for the same momenta as in Fig. 6: $k = (x\pi/32, 0)$ with $x = \{0, 6, 9, 12, 15, 18, 21, 24, 27, 32\}$ for $U_d = -8$ and $n = 0.5$, but at a temperature $T = 0.2$ below $T^* \simeq 1$. Notice the changes in lineshape and anomalous dispersion compared with high temperature results of Fig. 6. (From ref. [42]).
momemtum at which the pseudogap opens up and the Fermi surface is destroyed.

FIG. 9. Schematic representation of the results of ref. [42] showing that the Fermi surface

\[ T=0.75 \]

\[ T=0.20 \]

\[ \theta = \theta' = 0, \phi' = \phi = 5. \text{ (From ref. [42].)} \]

Inset: The angle dependence of \( \Delta \) for various densities \( \omega \) and temperatures \( \beta \).