Paired electron pockets in the hole-doped cuprates

Talk online: sachdev.physics.harvard.edu



Paired electron pockets in the hole-doped cuprates



Hole dynamics in an antiferromagnet across a deconfined quantum critical point, R. K. Kaul, A. Kolezhuk, M. Levin, S. Sachdev, and T. Senthil, *Physical Review* B **75**, 235122 (2007).

Algebraic charge liquids and the underdoped cuprates, R. K. Kaul, Y. B. Kim, S. Sachdev, and T. Senthil, *Nature Physics* **4**, 28 (2008).

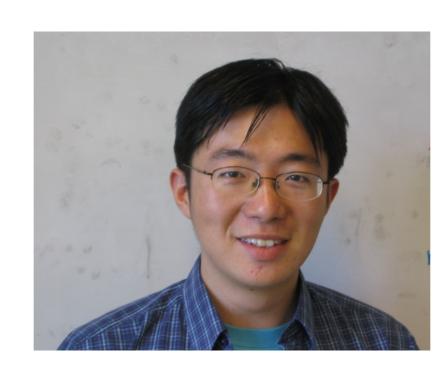
Destruction of Neel order in the cuprates by electron doping, R. K. Kaul, M. Metlitksi, S. Sachdev, and C. Xu, *Physical Review B* **78**, 045110 (2008).

Paired electron pockets in the underdoped cuprates, V. Galitski and S. Sachdev, arXiv:0901.0005





Ribhu Kaul UCSB



Victor Galitski Maryland

Cenke Xu Harvard

Outline

I. Nodal-anti-nodal dichotomy in the cuprates Survey of recent experiments

2. Spin density wave theory of normal metal From a "large" Fermi surface to electron and hole pockets

3. Algebraic charge liquids Pairing by gauge forces, d-wave superconductivity, and the nodal-anti-nodal dichotomy

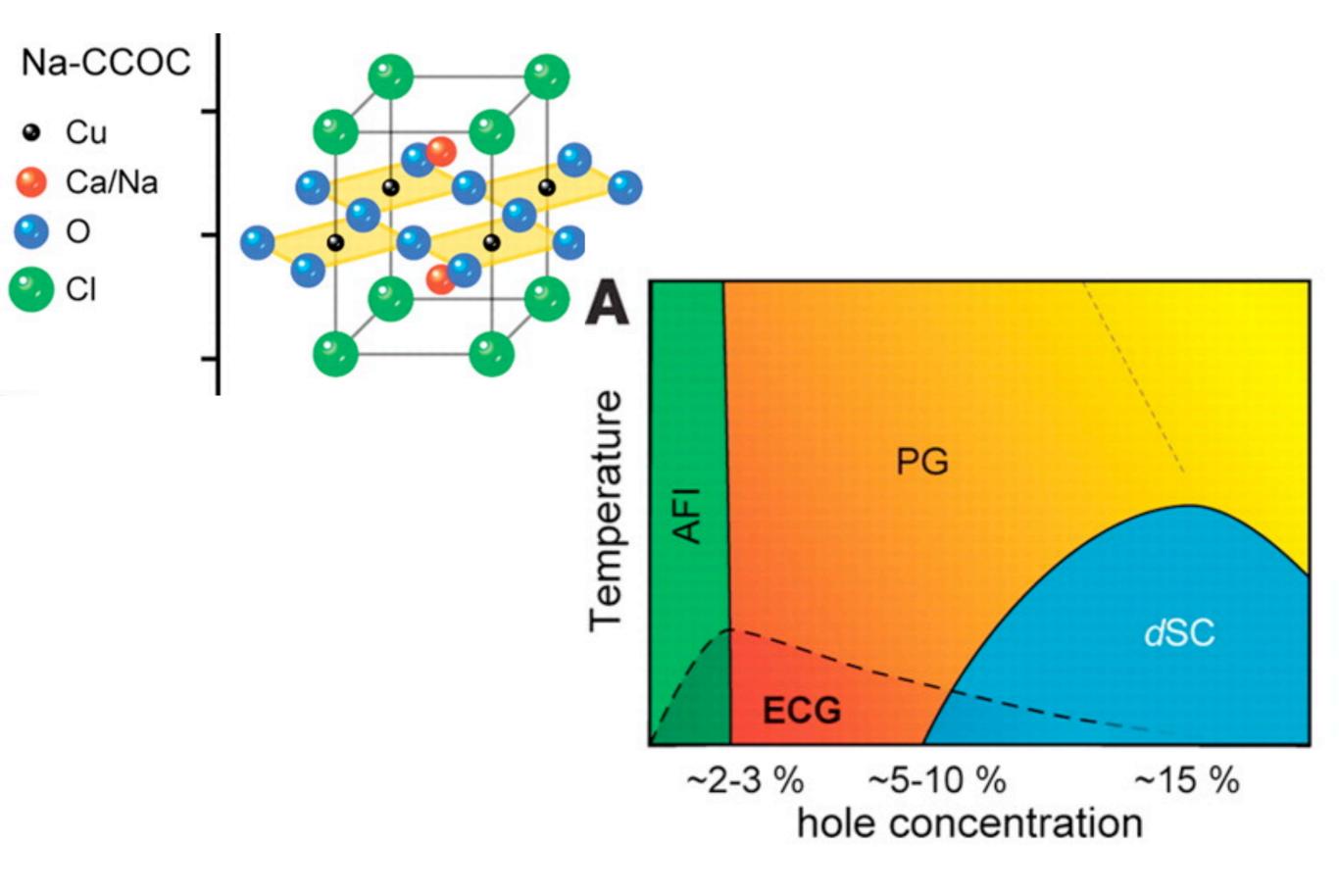
Outline

I. Nodal-anti-nodal dichotomy in the cuprates Survey of recent experiments

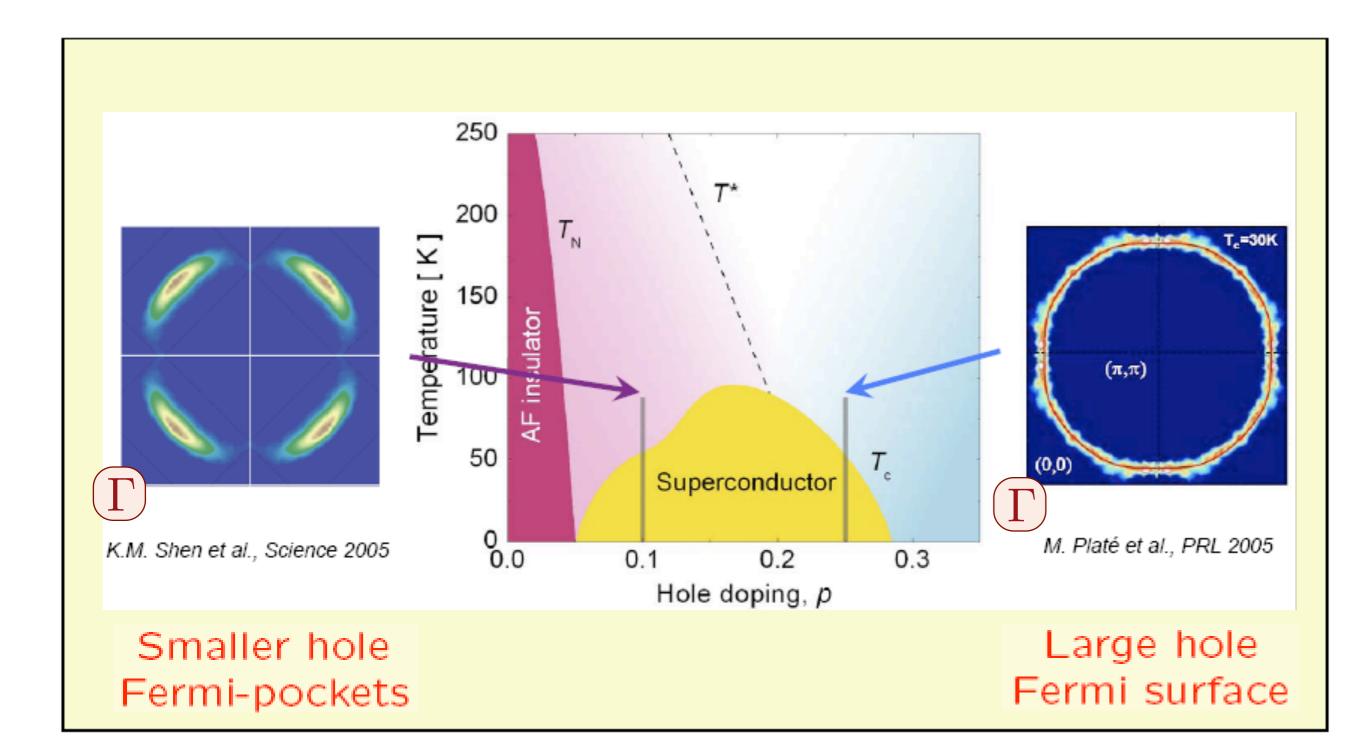
2. Spin density wave theory of normal metal From a "large" Fermi surface to electron and hole pockets

3. Algebraic charge liquids Pairing by gauge forces, d-wave superconductivity, and the nodal-anti-nodal dichotomy

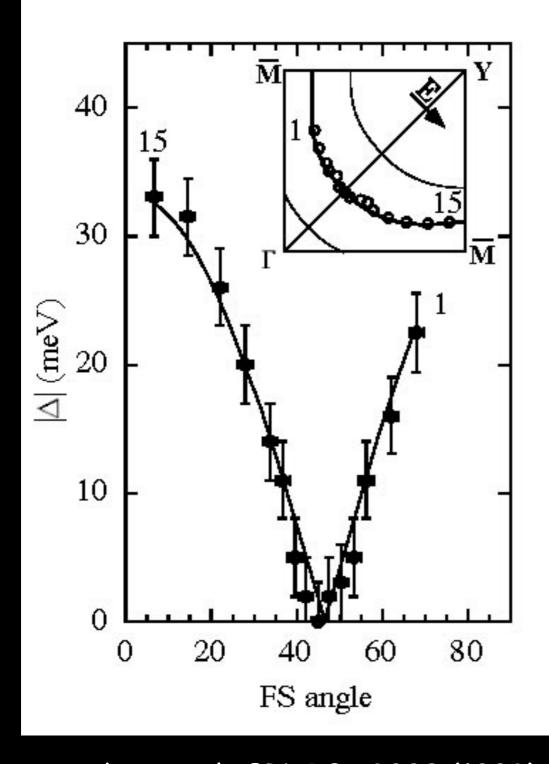
The cuprate superconductors



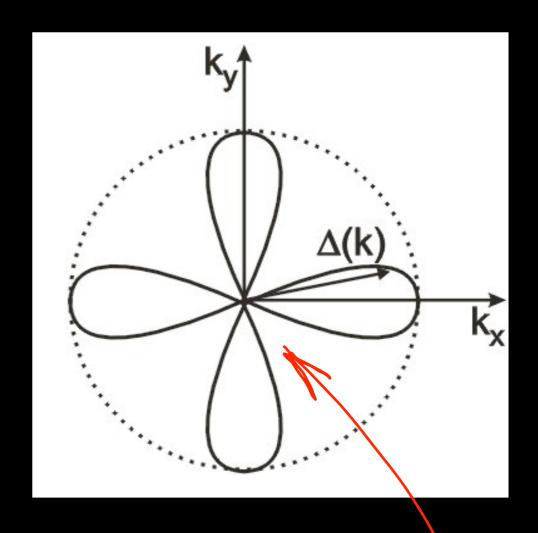
Evolution of the (ARPES) Fermi surface on the cuprate phase diagram



Overdoped SC State: Momentum-dependent Pair Energy Gap $\Delta(\vec{k})$

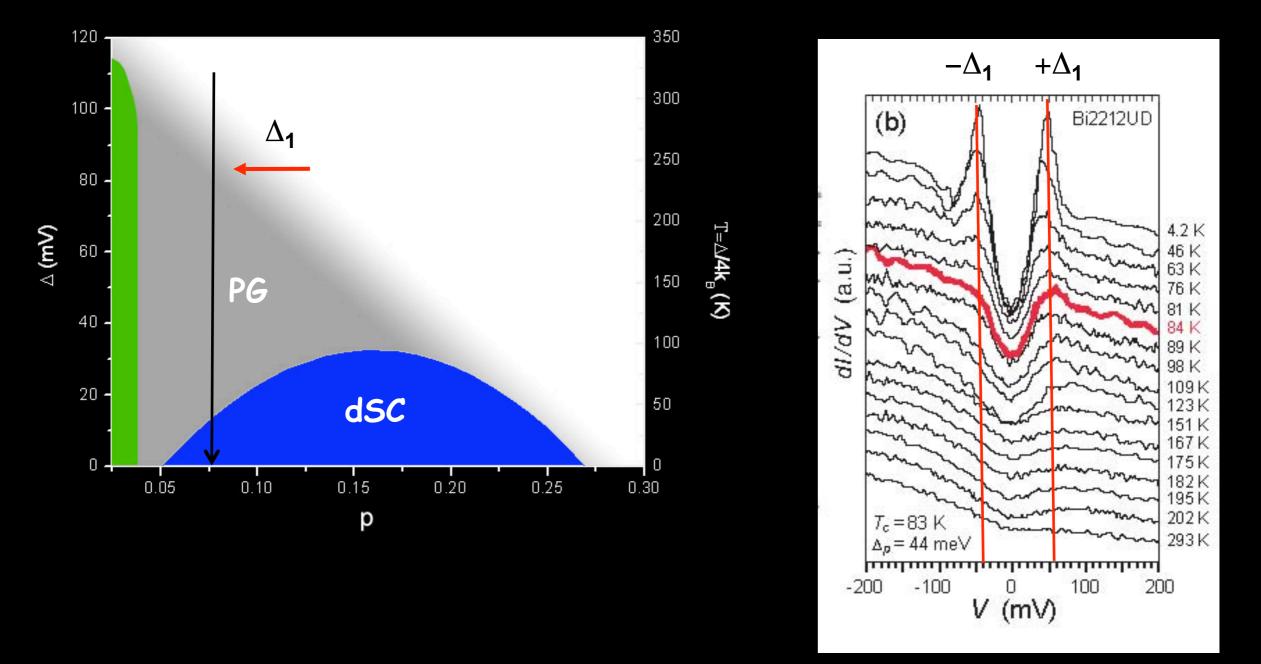


Shen et alPRL 70, 3999 (1993)Ding et alPRB 549678 (1996)Mesot et alPRL 83840 (1999)



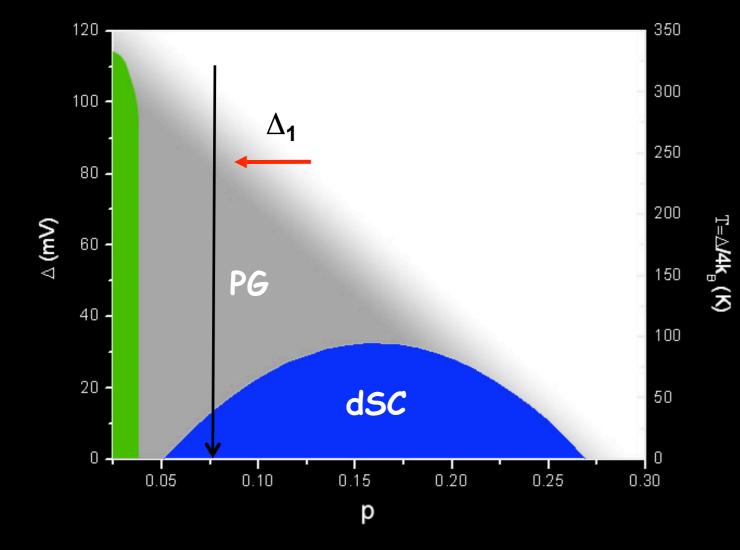
The SC energy gap $\Delta(k)$ has four nodes.

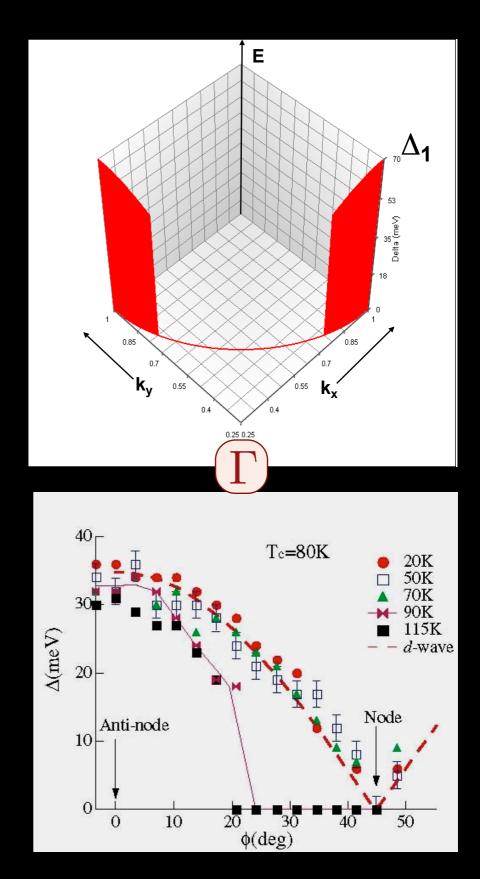
Pseudogap: Temperature-independent energy gap exists T>>T_c



Ch. Renner et al, PRL 80, 149 (1998) Ø. Fischer et al, RMP 79, 353 (2007)

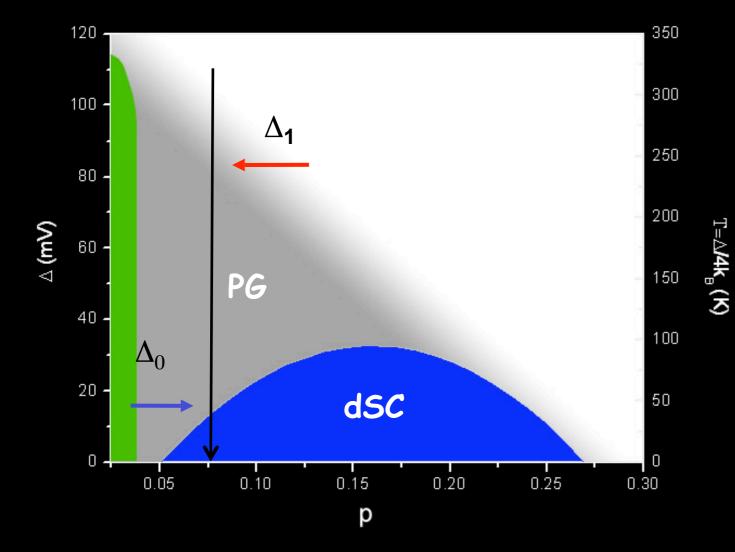
Pseudogap: Temperature-independent energy gap near k~(π ,0)

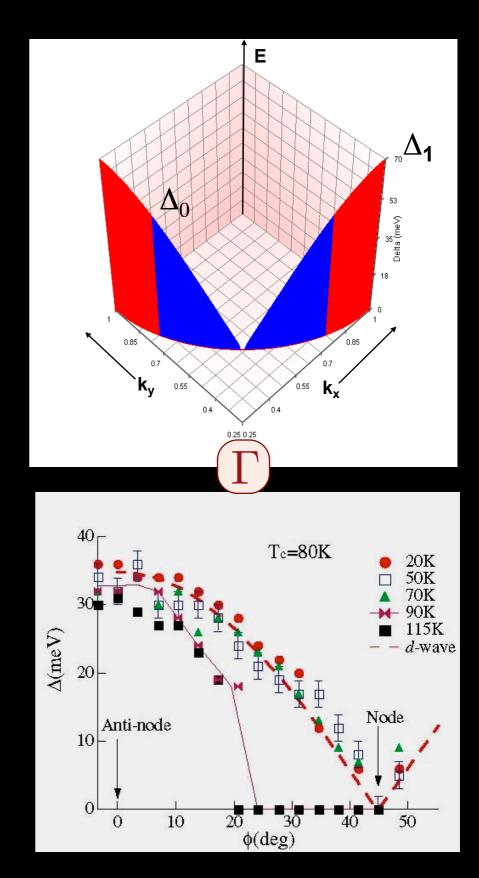




Loeser et al, Science 273 325 (1996) Ding et al, Nature 382 51, (1996) Norman et al, Nature 392, 157 (1998) Shen et al Science 307, 902 (2005) Kanigel et al, Nature Physics 2,447 (2006) Tanaka et al, Science 314, 1912 (2006)

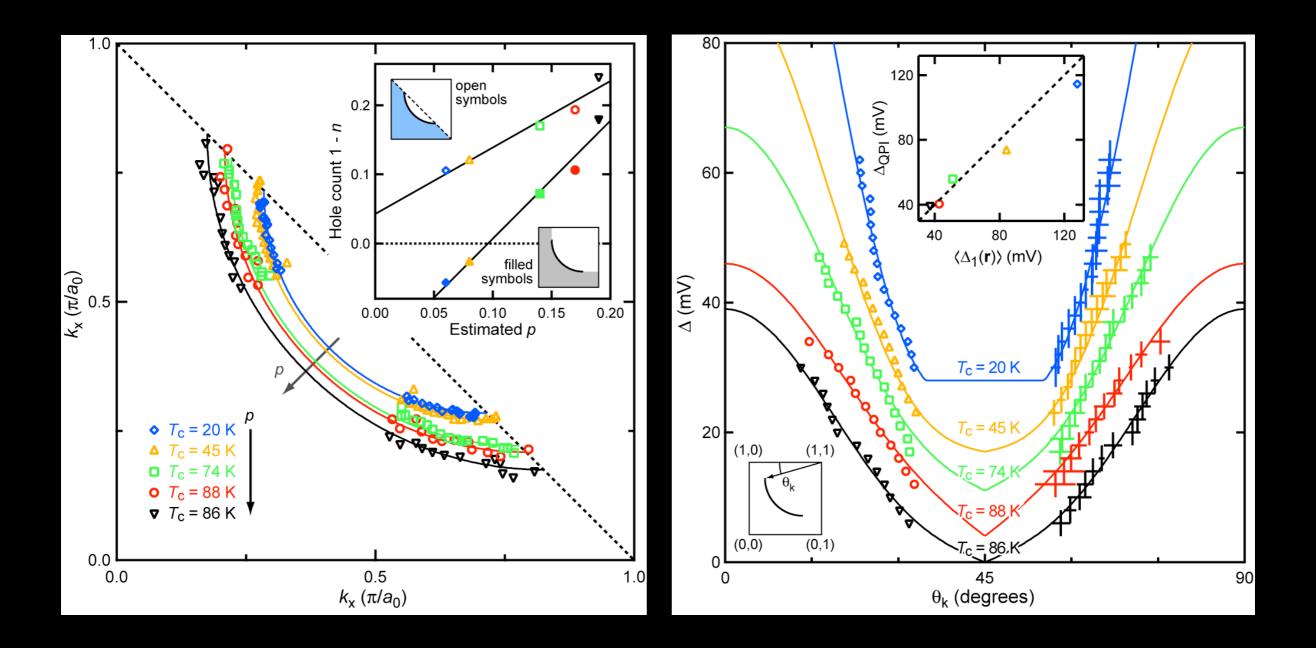
Pseudogap: Temperature-dependent energy gap near node



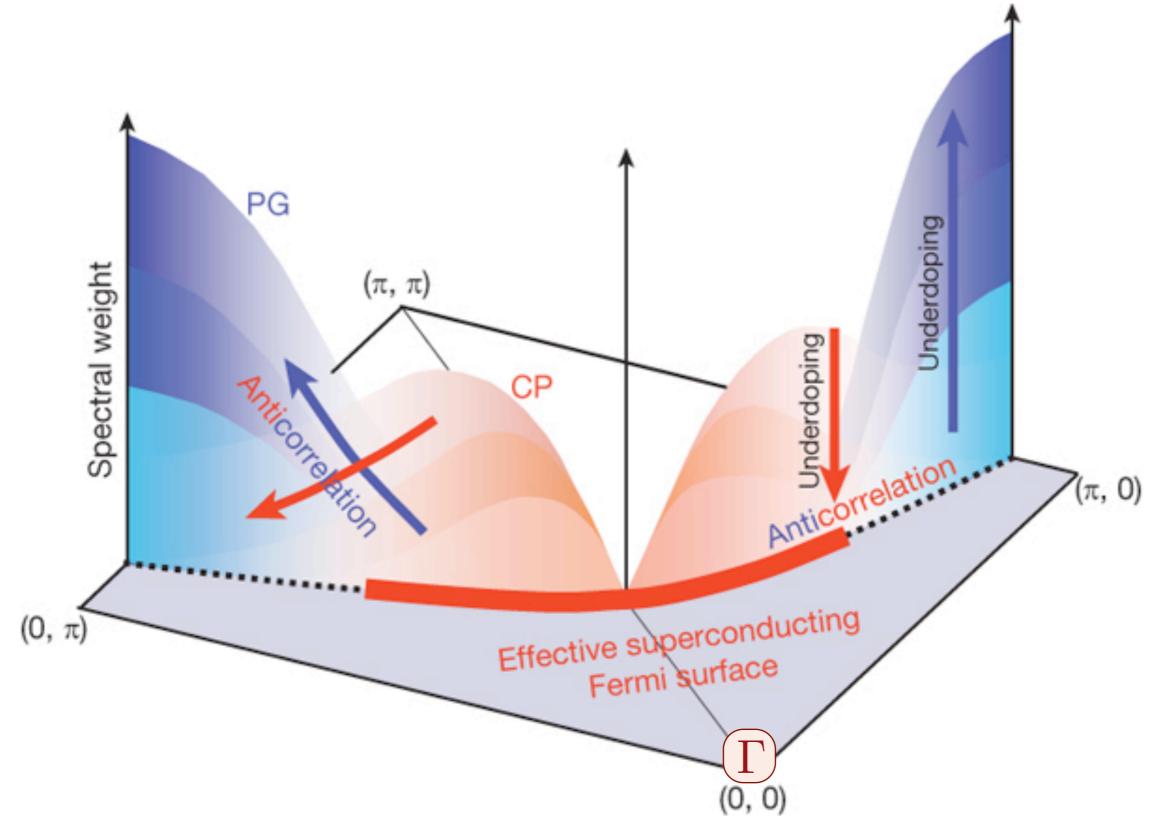


Loeser et al, Science 273 325 (1996) Ding et al, Nature 382 51, (1996) Norman et al, Nature 392, 157 (1998) Shen et al Science 307, 902 (2005) Kanigel et al, Nature Physics 2,447 (2006) Tanaka et al, Science 314, 1912 (2006)

Development of Fermi arc with underdoping



Y. Kohsaka et al., Nature 454, 1072, (2008)



Competition between the pseudogap and superconductivity in the high-Tc copper oxides

T. Kondo, R. Khasanov, T. Takeuchi, J. Schmalian, A. Kaminski, Nature 457, 296 (2009)

Nodal-anti-nodal dichotomy in the underdoped cuprates

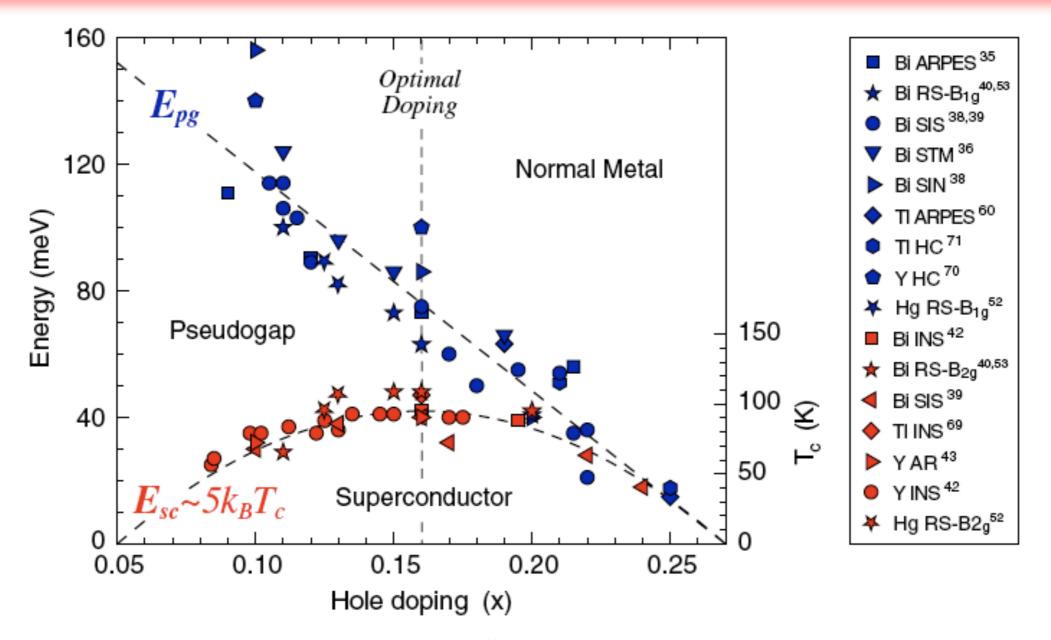
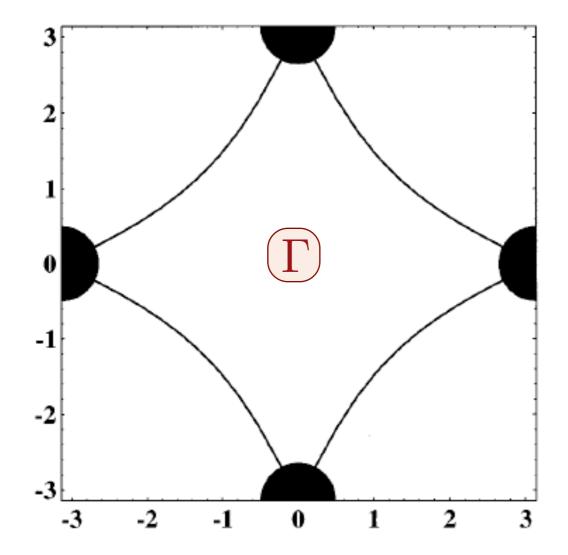


Figure 2. Pseudogap ($E_{pg} = 2\Delta_{pg}$) and superconducting ($E_{sc} \sim 5k_BT_c$) energy scales for a number of HTSCs with $T_c^{max} \sim 95$ K (Bi2212, Y123, T12201 and Hg1201). The datapoints were obtained, as a function of hole doping x, by angle-resolved photoemission spectroscopy (ARPES), tunneling (STM, SIN, SIS), Andreev reflection (AR), Raman scattering (RS) and heat conductivity (HC). On the same plot we are also including the energy Ω_r of the magnetic resonance mode measured by inelastic neutron scattering (INS), which we identify with E_{sc} because of the striking quantitative correspondence as a function of T_c . The data fall on two universal curves given by $E_{pg} = E_{pg}^{max} (0.27 - x)/0.22$ and $E_{sc} = E_{sc}^{max} [1 - 82.6(0.16 - x)^2]$, with $E_{pg}^{max} = E_{pg}(x = 0.05) = 152 \pm 8$ meV and $E_{sc}^{max} = E_{sc}(x = 0.16) = 42 \pm 2$ meV (the statistical errors refer to the fit of the selected datapoints; however, the spread of all available data would be more appropriately described by ± 20 and ± 10 meV, respectively).

S. Hufner, M.A. Hossain, A. Damascelli, and G.A. Sawatzky, Rep. Prog. Phys. 71, 062501 (2008)

Superconductivity in a system with preformed pairs



$$\begin{split} H &= \sum_{q} \varepsilon b_{q}^{\dagger} b_{q} + \sum_{p,q}' V_{p,q} (b_{q}^{\dagger} c_{p\uparrow} c_{q-p\downarrow} + \text{H.c.}) \\ &+ \sum_{p} \xi_{p} c_{p,\sigma}^{\dagger} c_{p,\sigma}; \\ &V_{p,q} = V a^{2} (p_{x}^{2} - p_{y}^{2}) \end{split}$$

FIG. 1. Sketch of the Fermi line and region of the momentum space where pseudogap pairs is formed. The Fermi line shown here was obtained in the tight binding model with diagonal hopping t' = -0.3t; it is similar to the Fermi line observed in the underdoped Bi₂Sr₂CaCu₂O_{8+ δ} (Ref. 5). The shaded disks denote the part of the momentum space where a pseudogap was observed in the experiment. We shall assume that the fermions in these regions are paired into the bosons.

V. B. Geshkenbein, L. B. Ioffe, and A. I. Larkin, Phys. Rev. B 55, 3173 (1997).

Superconductivity in a system with preformed pairs

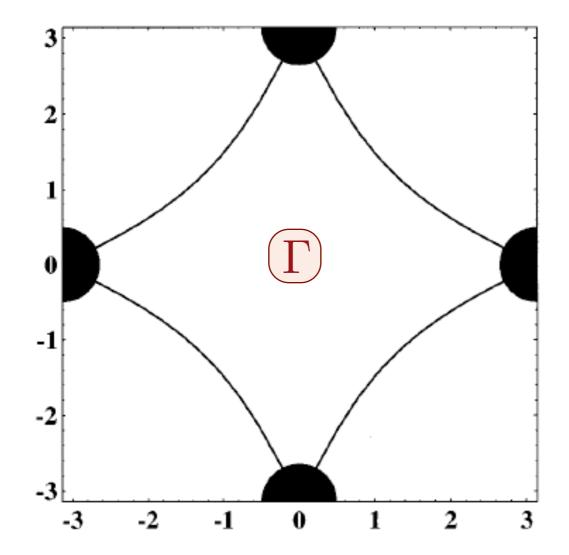


FIG. 1. Sketch of the Fermi line and region of the momentum space where pseudogap pairs is formed. The Fermi line shown here was obtained in the tight binding model with diagonal hopping t' = -0.3t; it is similar to the Fermi line observed in the underdoped Bi₂Sr₂CaCu₂O_{8+ δ} (Ref. 5). The shaded disks denote the part of the momentum space where a pseudogap was observed in the experiment. We shall assume that the fermions in these regions are paired into the bosons.

$$\begin{split} H &= \sum_{q} \varepsilon b_{q}^{\dagger} b_{q} + \sum_{p,q}' V_{p,q} (b_{q}^{\dagger} c_{p\uparrow} c_{q-p\downarrow} + \text{H.c.}) \\ &+ \sum_{p} \xi_{p} c_{p,\sigma}^{\dagger} c_{p,\sigma}; \\ V_{p,q} &= V a^{2} (p_{x}^{2} - p_{y}^{2}) \end{split}$$

Attractive phenomenological model, but theoretical and microscopic basis is unclear

V. B. Geshkenbein, L. B. Ioffe, and A. I. Larkin, Phys. Rev. B 55, 3173 (1997).

Outline

I. Nodal-anti-nodal dichotomy in the cuprates Survey of recent experiments

2. Spin density wave theory of normal metal From a "large" Fermi surface to electron and hole pockets

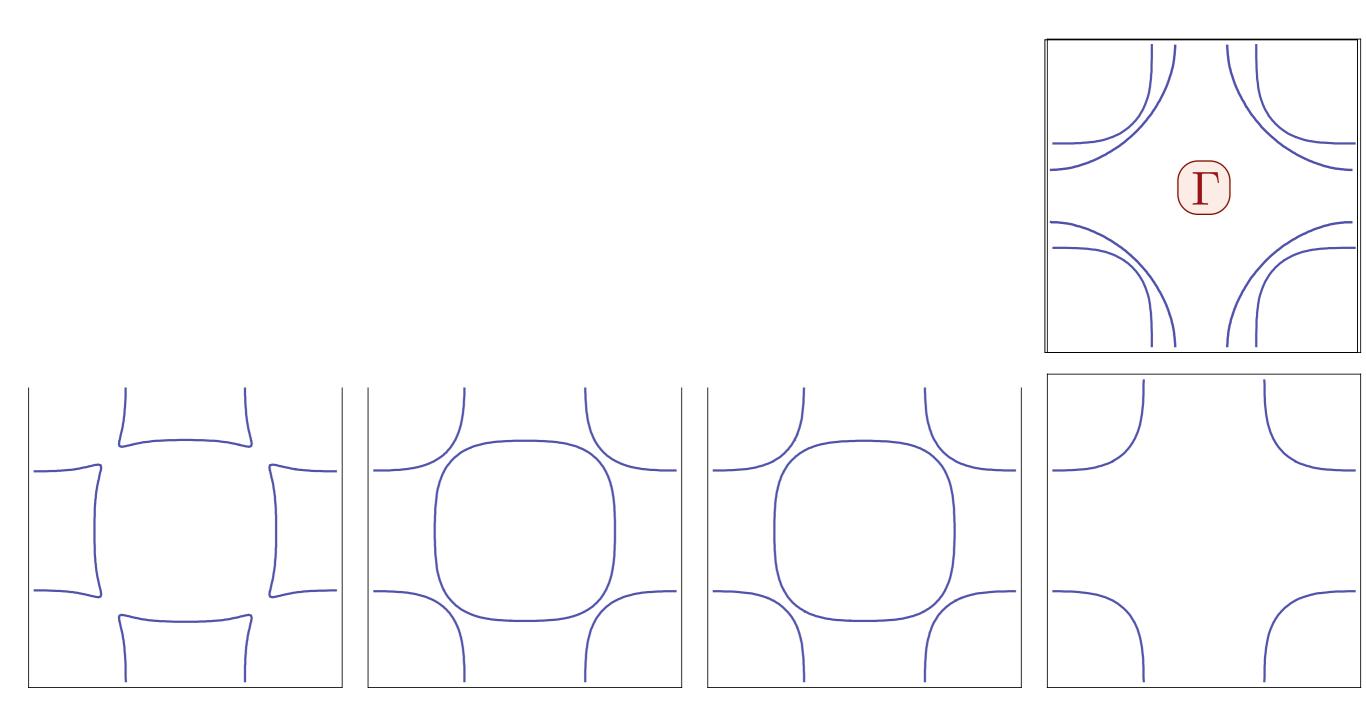
3. Algebraic charge liquids Pairing by gauge forces, d-wave superconductivity, and the nodal-anti-nodal dichotomy

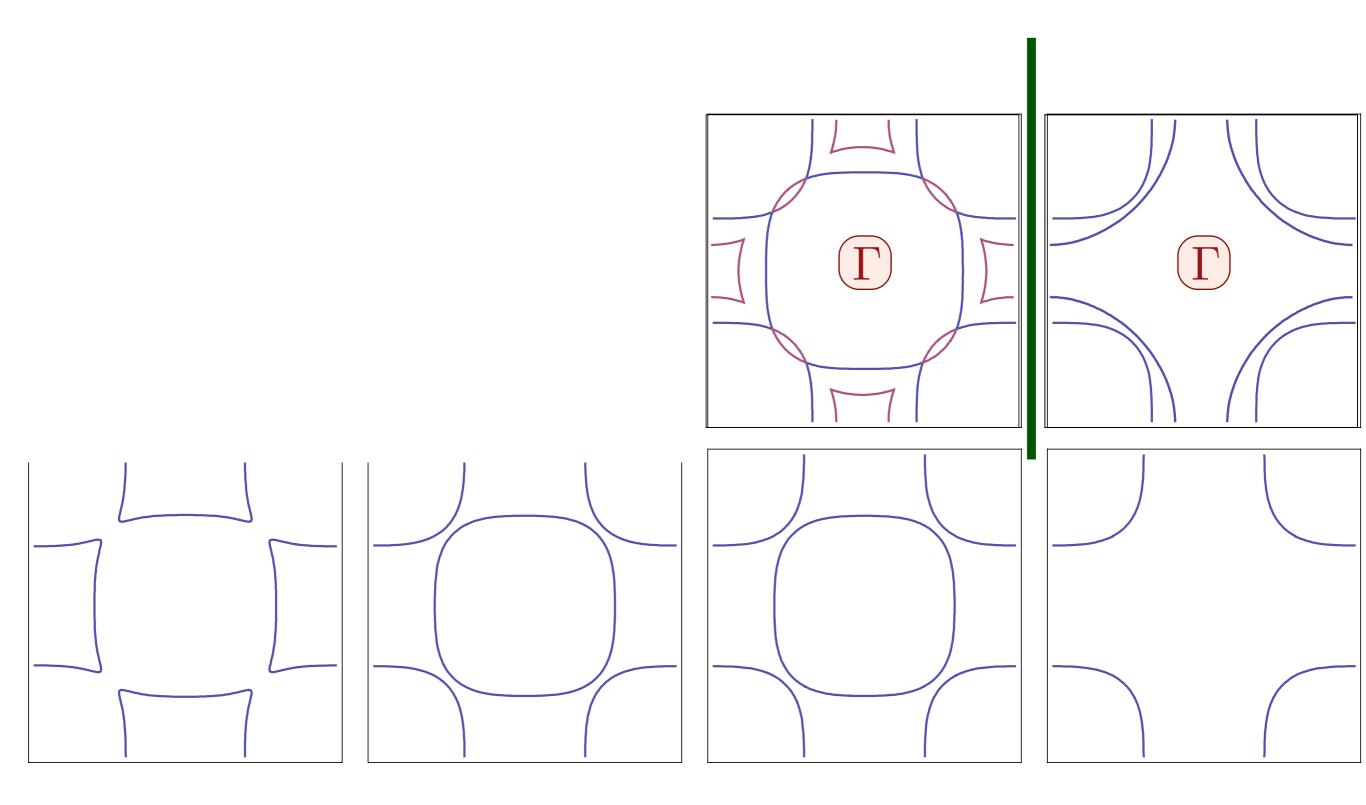
Outline

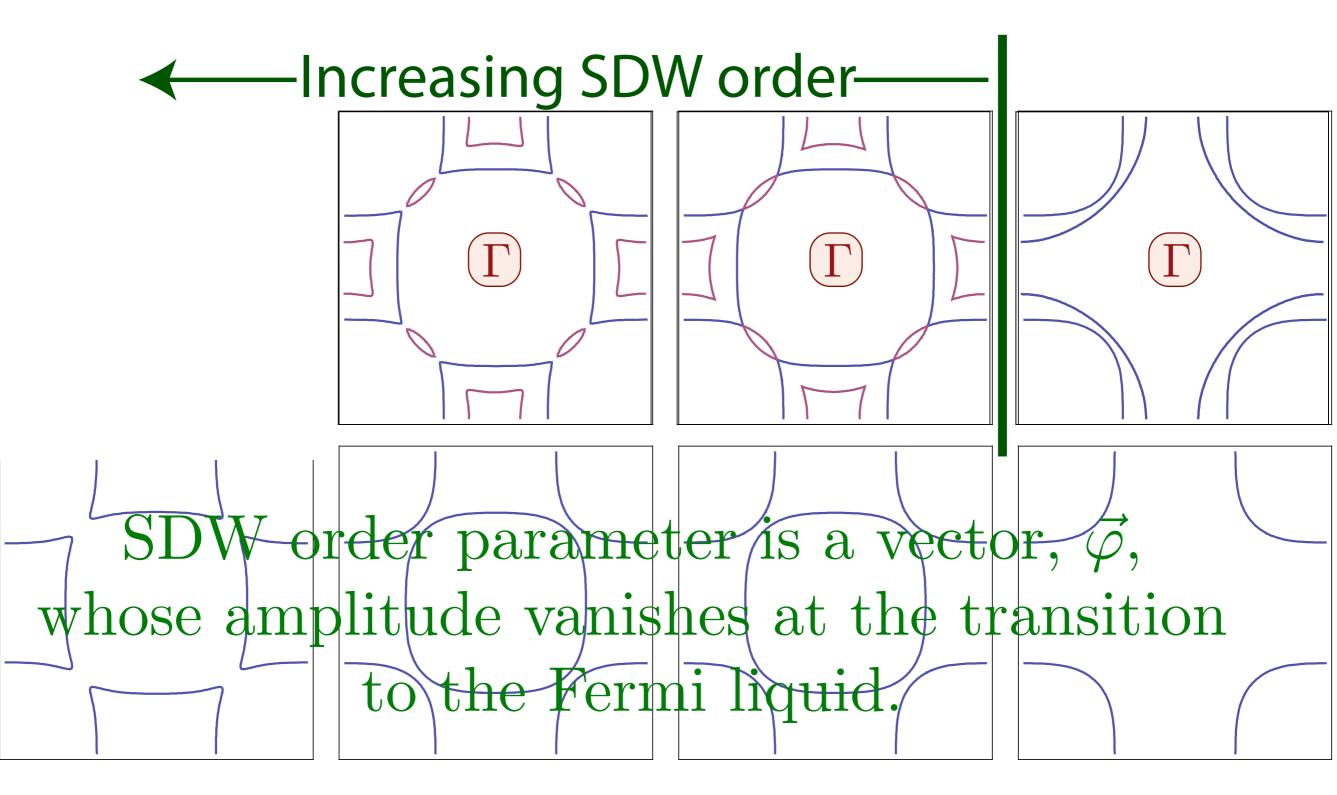
I. Nodal-anti-nodal dichotomy in the cuprates Survey of recent experiments

2. Spin density wave theory of normal metal From a "large" Fermi surface to electron and hole pockets

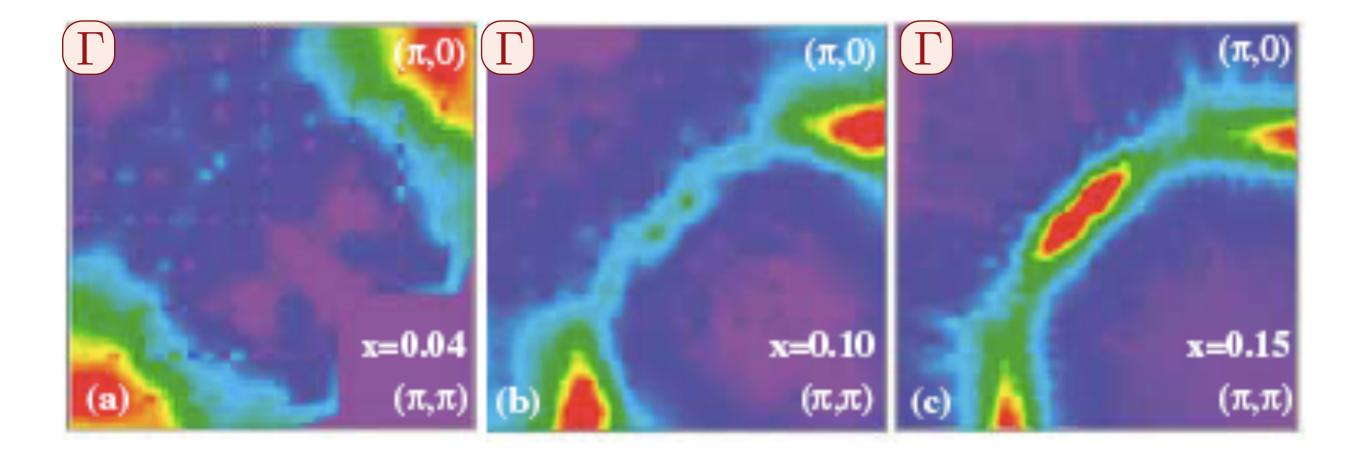
3. Algebraic charge liquids Pairing by gauge forces, d-wave superconductivity, and the nodal-anti-nodal dichotomy



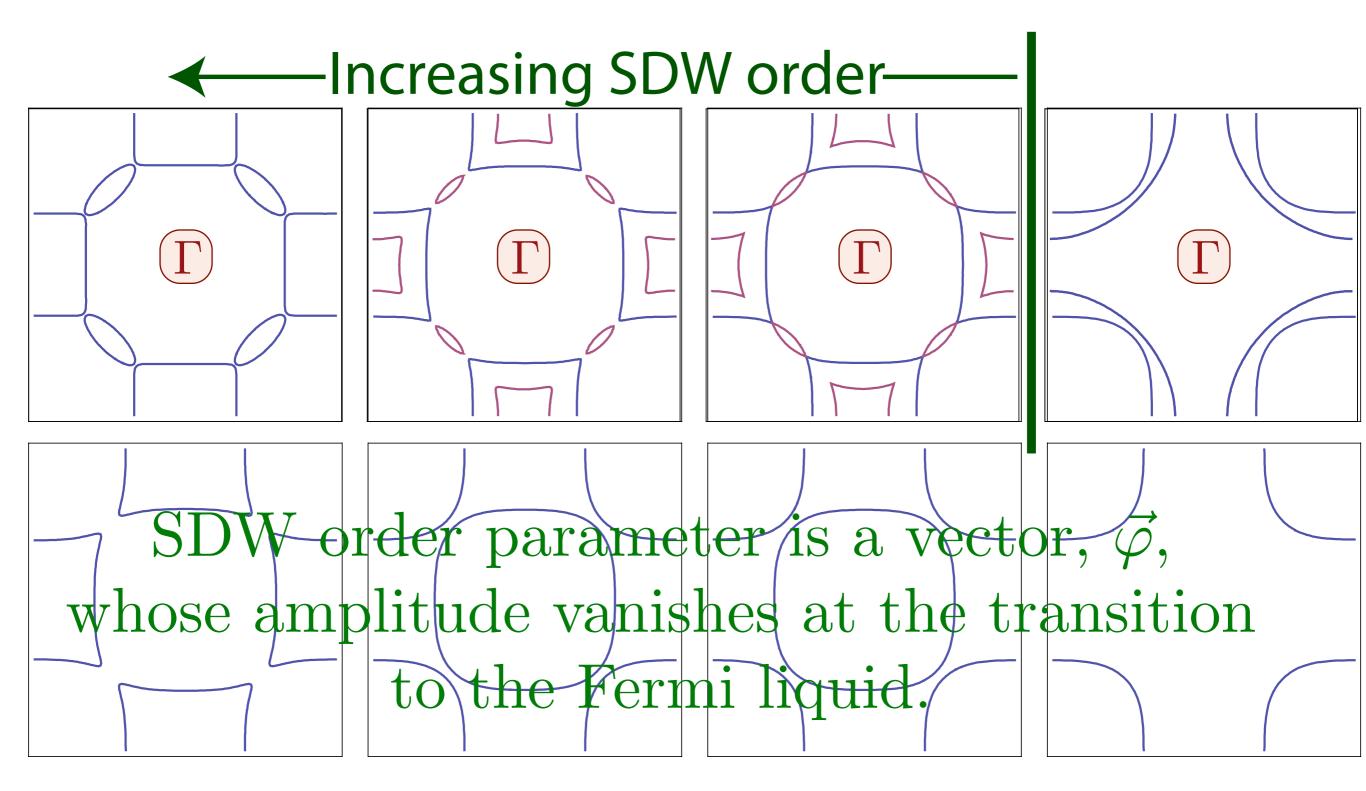


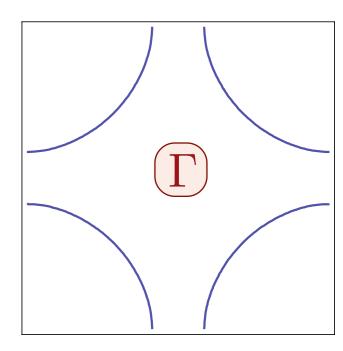


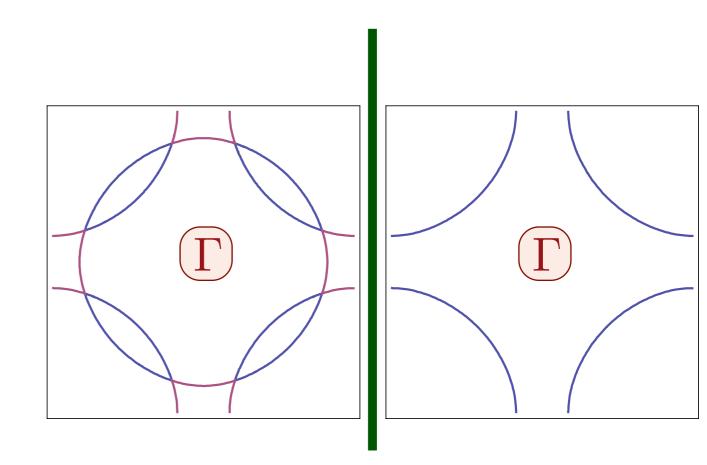
Photoemission in NCCO (electron-doped)

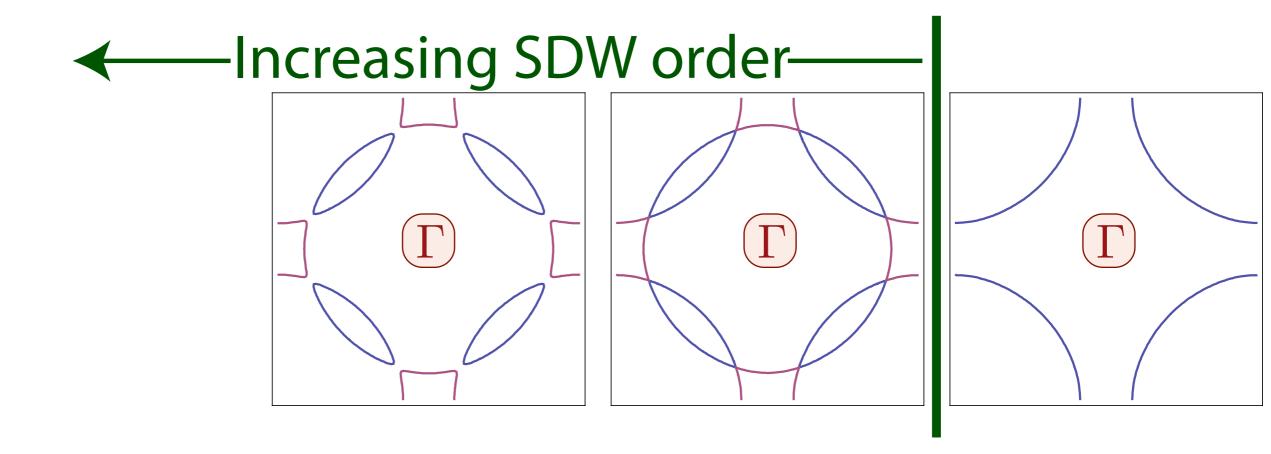


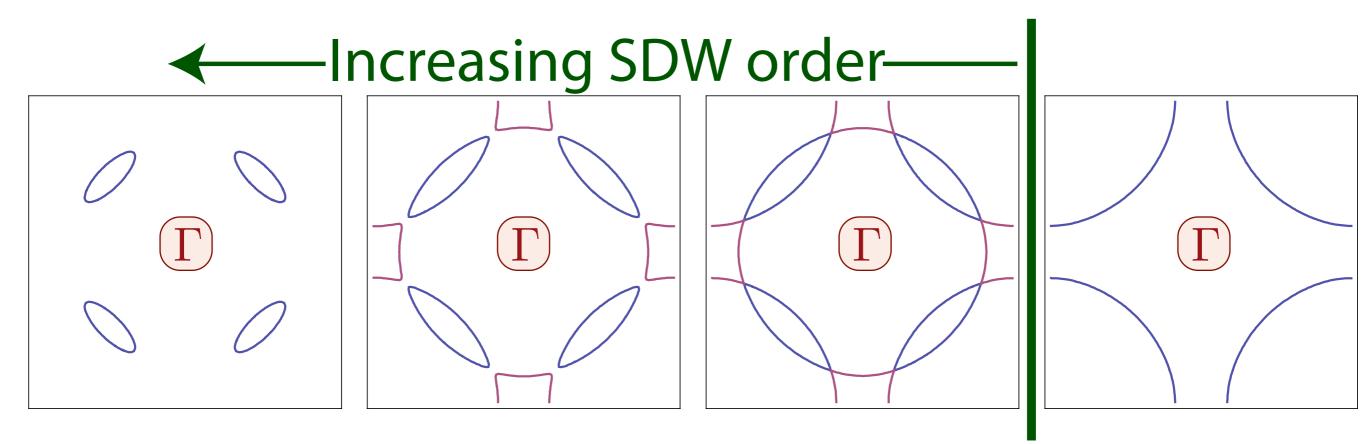
N. P. Armitage et al., Phys. Rev. Lett. 88, 257001 (2002).



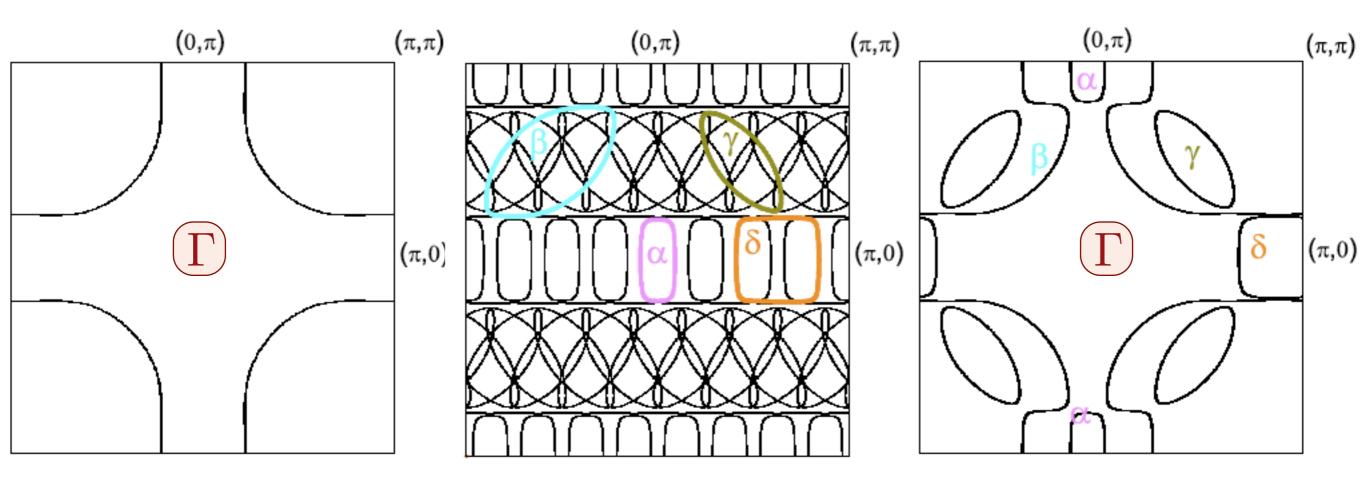








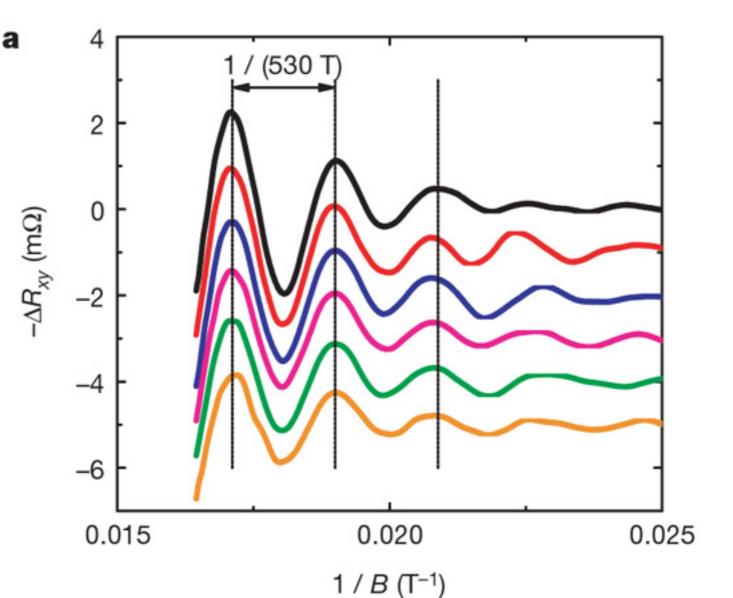
SDW order parameter is a vector, $\vec{\varphi}$, whose amplitude vanishes at the transition to the Fermi liquid.



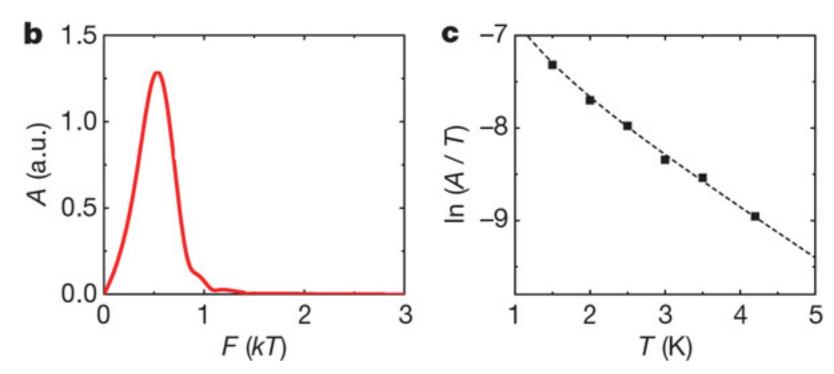
Incommensurate order in YBa₂Cu₃O_{6+x}

N. Harrison, arXiv:0902.2741.

Quantum oscillations and the Fermi surface in an underdoped high T_c superconductor (ortho-II ordered YBa₂Cu₃O_{6.5}). The period corresponds to a carrier density ≈ 0.076 .



N. Doiron-Leyraud, C. Proust, D. LeBoeuf, J. Levallois, J.-B. Bonnemaison, R. Liang, D. A. Bonn, W. N. Hardy, and L. Taillefer, *Nature* **447**, 565 (2007)



Electron pockets in the Fermi surface of hole-doped high-T_c superconductors

David LeBoeuf¹, Nicolas Doiron-Leyraud¹, Julien Levallois², R. Daou¹, J.-B. Bonnemaison¹, N. E. Hussey³, L. Balicas⁴, B. J. Ramshaw⁵, Ruixing Liang^{5,6}, D. A. Bonn^{5,6}, W. N. Hardy^{5,6}, S. Adachi⁷, Cyril Proust² & Louis Taillefer^{1,6}

а b С 20 Y123-II Y123-VIII Y124 1.0 0 10 В, 0.5 R_H (mm³ C⁻¹) 0 -10 90 K -100.0 80 K 70 K 60 K 60 K 30 K -20 50 K -20 35 K 50 K 3 K -0.540 K 40 K 4.2 K 40 K 34 K 50 K 30 K -3060 K 27 K 20 K 10 -1.080 K 15 K 10 K 20 K -30 4.2 K 12 K 20 K — 100 K -401.5 K 150 K B -1.520 30 40 50 10 30 40 50 60 10 0 20 30 40 50 60 20 0 10 0 B (T) B (T) B (T)

Nature 450, 533 (2007)

Outline

I. Nodal-anti-nodal dichotomy in the cuprates Survey of recent experiments

2. Spin density wave theory of normal metal From a "large" Fermi surface to electron and hole pockets

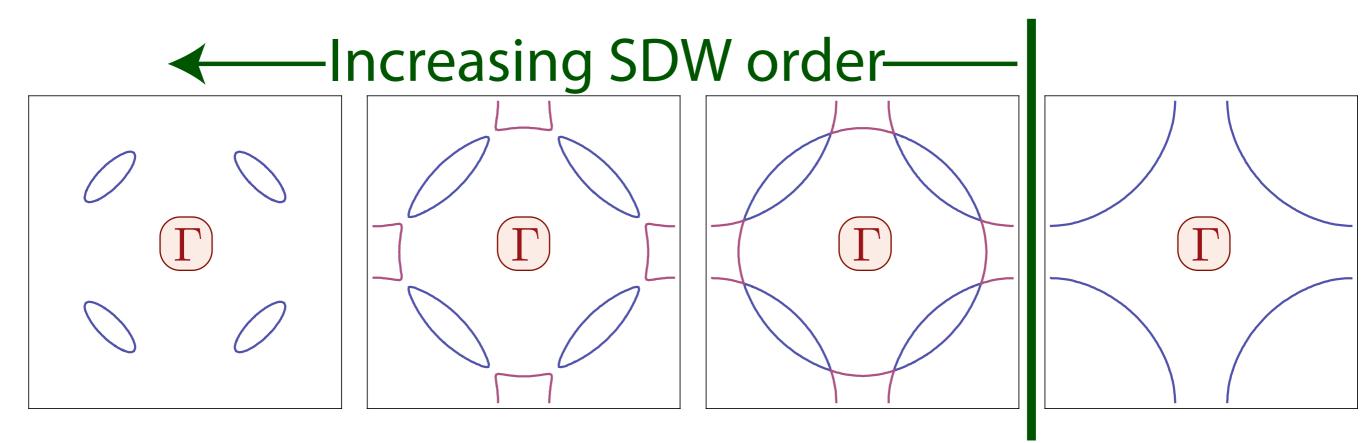
3. Algebraic charge liquids Pairing by gauge forces, d-wave superconductivity, and the nodal-anti-nodal dichotomy

Outline

I. Nodal-anti-nodal dichotomy in the cuprates Survey of recent experiments

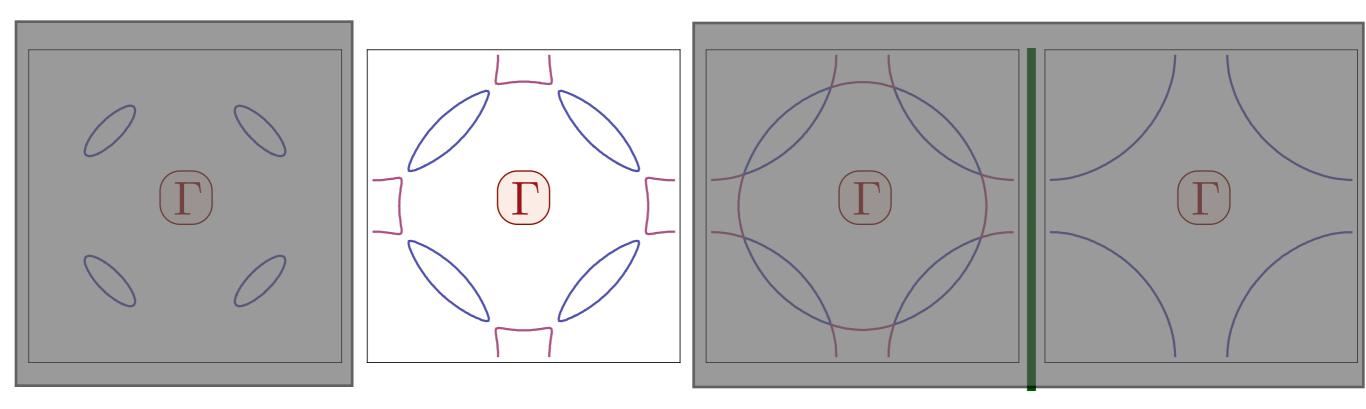
2. Spin density wave theory of normal metal From a "large" Fermi surface to electron and hole pockets

3. Algebraic charge liquids Pairing by gauge forces, d-wave superconductivity, and the nodal-anti-nodal dichotomy



SDW order parameter is a vector, $\vec{\varphi}$, whose amplitude vanishes at the transition to the Fermi liquid.

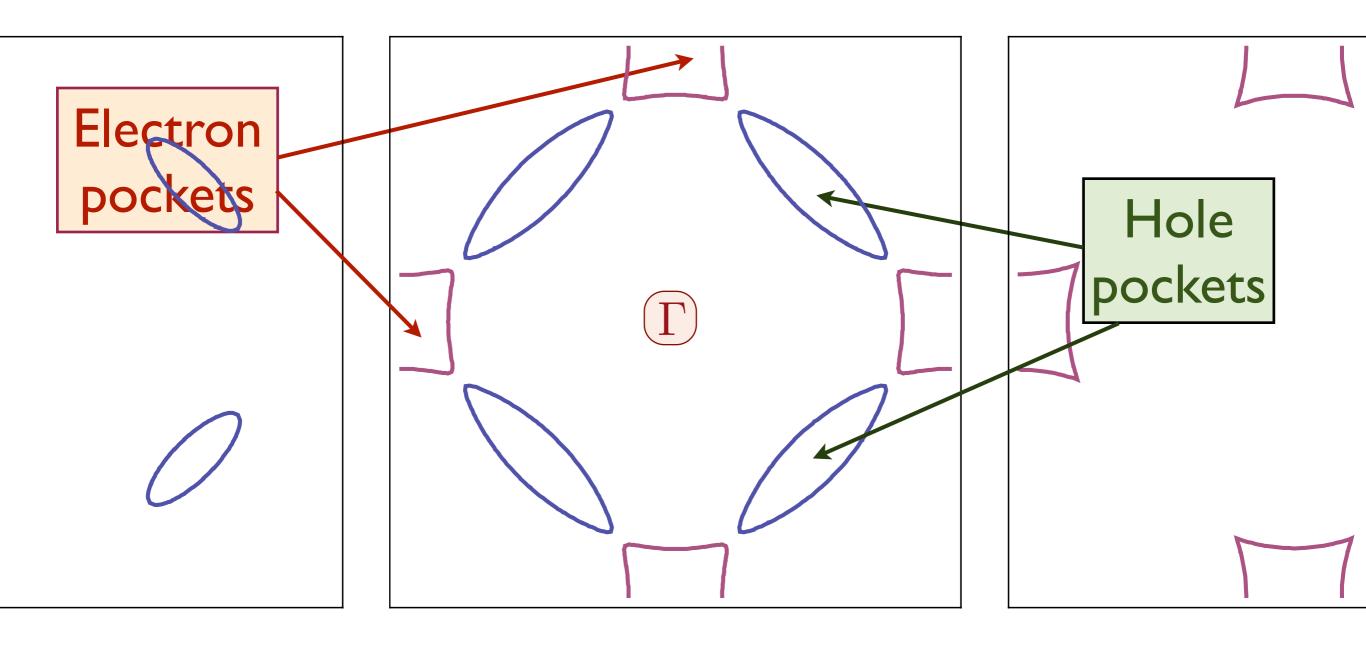
Fermi pockets in hole-doped cuprates

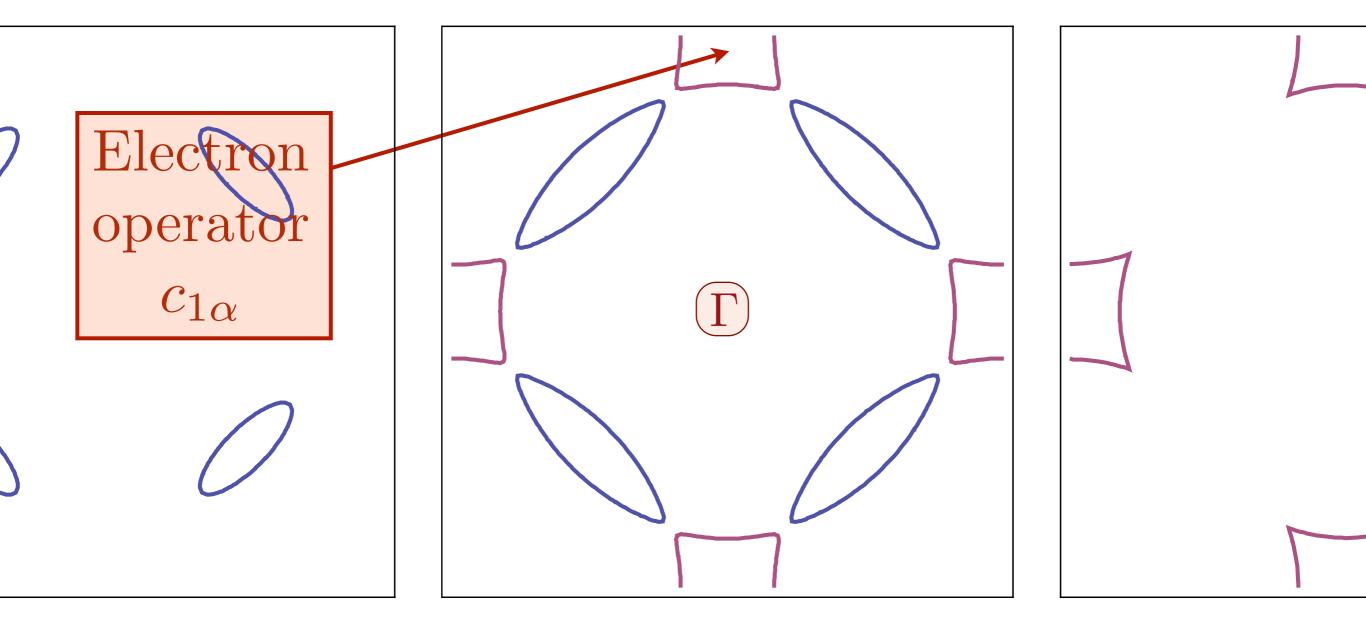


Begin with SDW ordered state, and focus on fluctuations in the *orientation* of $\vec{\varphi}$, by using a unit-length bosonic spinor z_{α}

$$\vec{\varphi} = z_{\alpha}^* \vec{\sigma}_{\alpha\beta} z_{\beta}$$

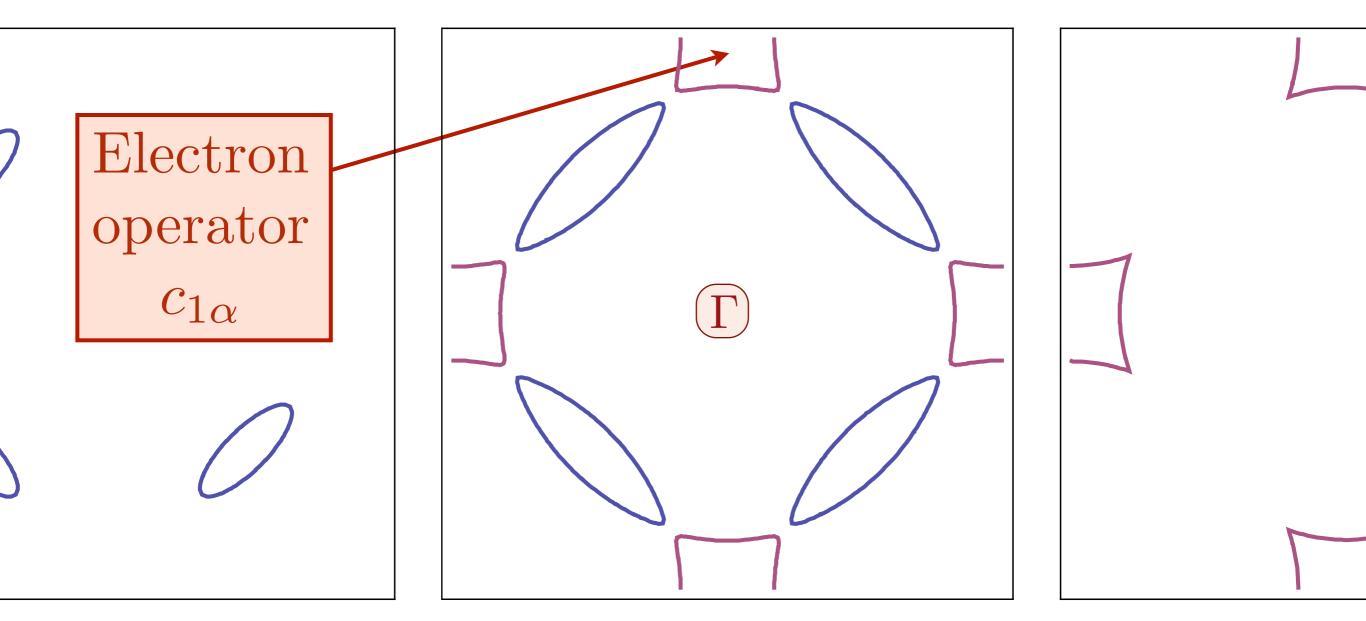
Charge carriers in the lightly-doped cuprates with Neel order





For a uniform SDW order with $\vec{\varphi} \propto (0, 0, 1)$, write

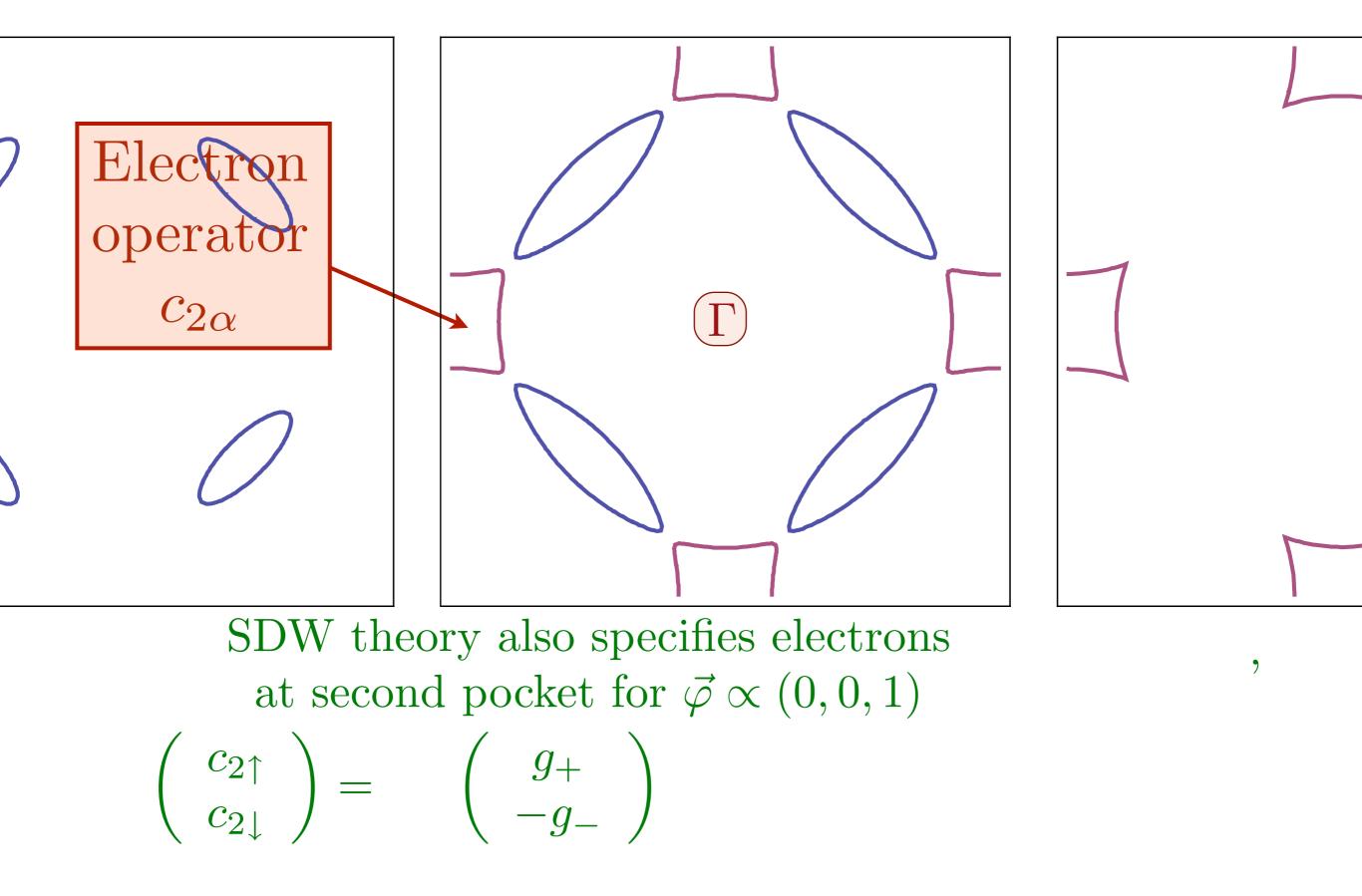
$$\left(\begin{array}{c}c_{1\uparrow}\\c_{1\downarrow}\end{array}\right) = \left(\begin{array}{c}g_{+}\\g_{-}\end{array}\right)$$

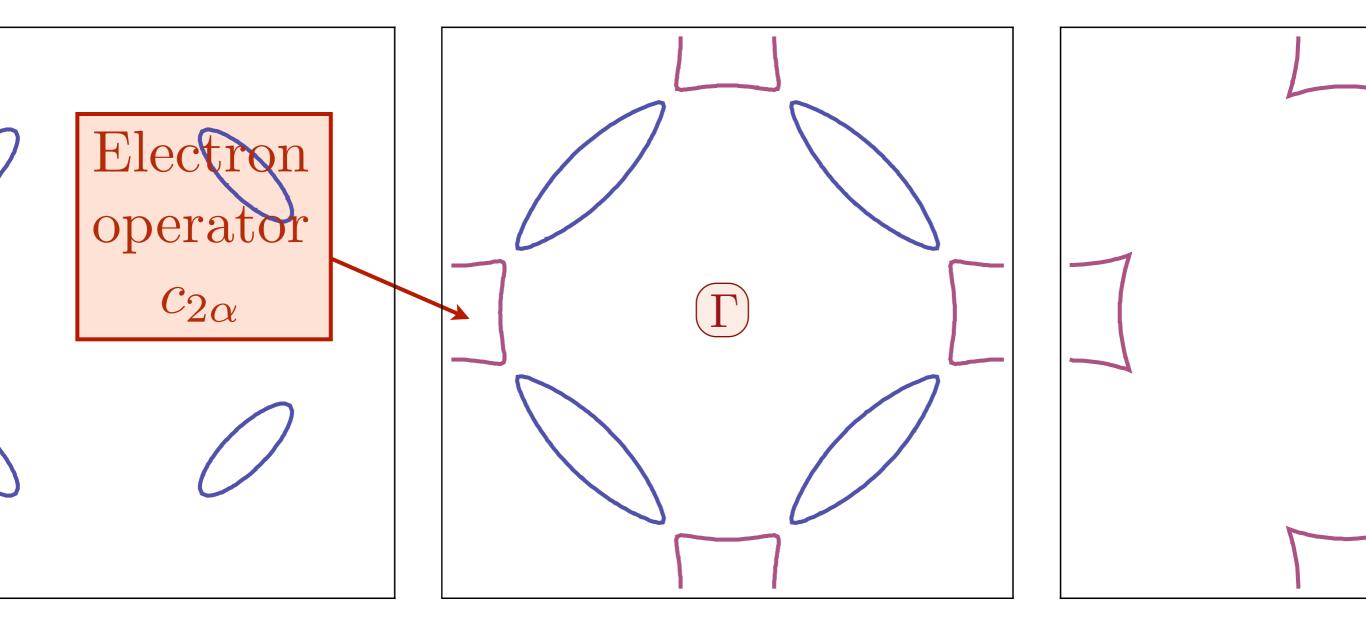


For a spacetime dependent SDW order, $\vec{\varphi} = z_{\alpha}^* \vec{\sigma}_{\alpha\beta} z_{\beta}$,

$$\begin{pmatrix} c_{1\uparrow} \\ c_{1\downarrow} \end{pmatrix} = \mathcal{R}_z \begin{pmatrix} g_+ \\ g_- \end{pmatrix} ; \quad \mathcal{R}_z \equiv \begin{pmatrix} z_{\uparrow} & -z_{\downarrow}^* \\ z_{\downarrow} & z_{\uparrow}^* \end{pmatrix}.$$

So g_{\pm} are the "up/down" electron operators in a rotating reference frame defined by the local SDW order





For a spacetime dependent SDW order, $\vec{\varphi} = z_{\alpha}^* \vec{\sigma}_{\alpha\beta} z_{\beta}$,

$$\begin{pmatrix} c_{2\uparrow} \\ c_{2\downarrow} \end{pmatrix} = \mathcal{R}_z \begin{pmatrix} g_+ \\ -g_- \end{pmatrix} ; \quad \mathcal{R}_z \equiv \begin{pmatrix} z_{\uparrow} & -z_{\downarrow}^* \\ z_{\downarrow} & z_{\uparrow}^* \end{pmatrix}.$$

Same SU(2) matrix also rotates electrons in second pocket.

Low energy theory for spinless, charge -e fermions g_{\pm} , and spinful, charge 0 bosons z_{α} :

$$\mathcal{L} = \mathcal{L}_z + \mathcal{L}_g$$

$$\mathcal{L}_z = \frac{1}{t} \Big[|(\partial_\tau - iA_\tau) z_\alpha|^2 + c^2 |\nabla - i\mathbf{A}) z_\alpha|^2 \Big]$$

+ Berry phases of monopoles in A_μ .

CP¹ field theory for z_{α} and an emergent U(1) gauge field A_{μ} . Coupling t tunes the strength of SDW orientation fluctuations. Low energy theory for spinless, charge -e fermions g_{\pm} , and spinful, charge 0 bosons z_{α} :

$$\mathcal{L} = \mathcal{L}_z + \mathcal{L}_g$$

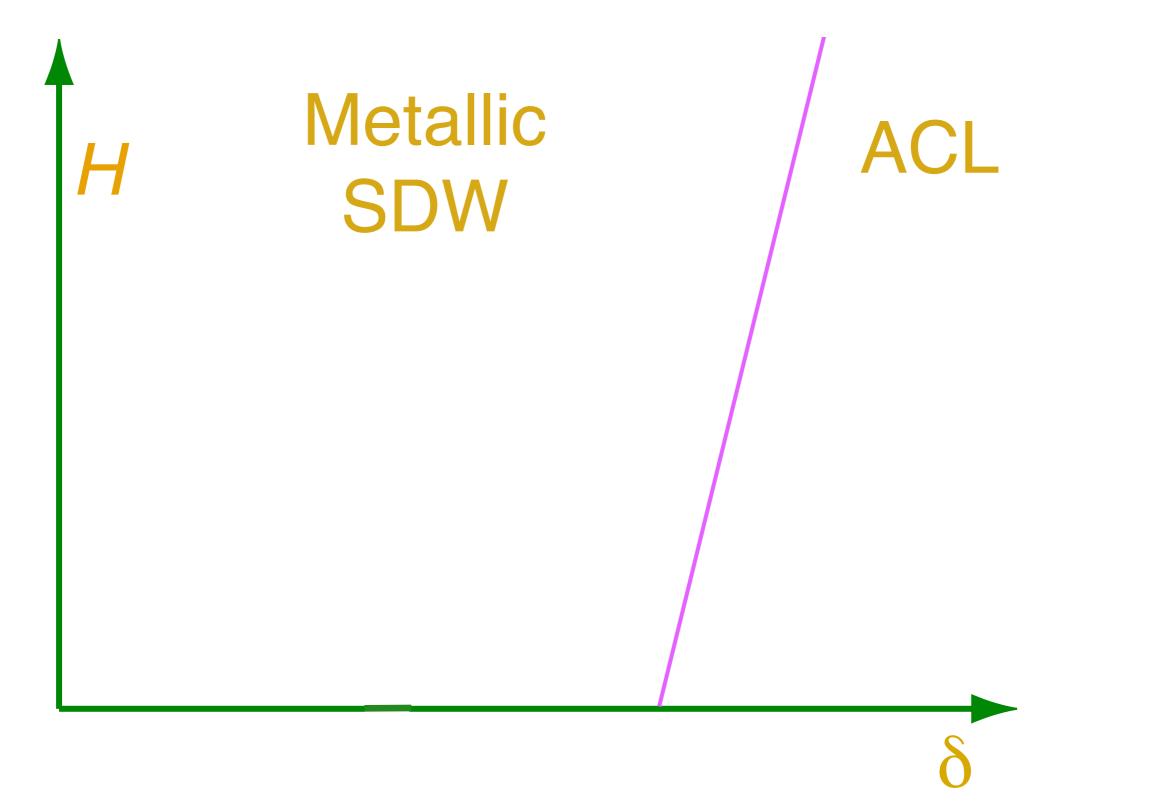
$$\mathcal{L}_z = \frac{1}{t} \Big[|(\partial_\tau - iA_\tau) z_\alpha|^2 + c^2 |\nabla - i\mathbf{A}) z_\alpha|^2 \Big]$$

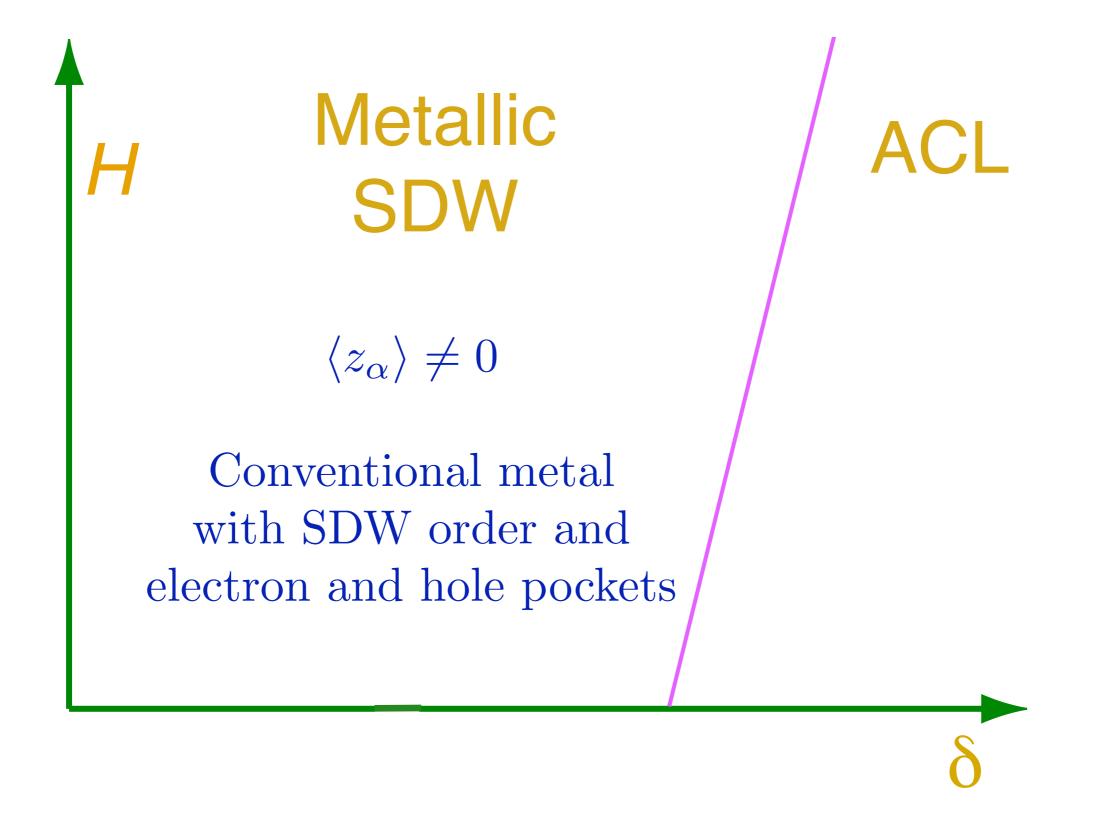
+ Berry phases of monopoles in A_μ .

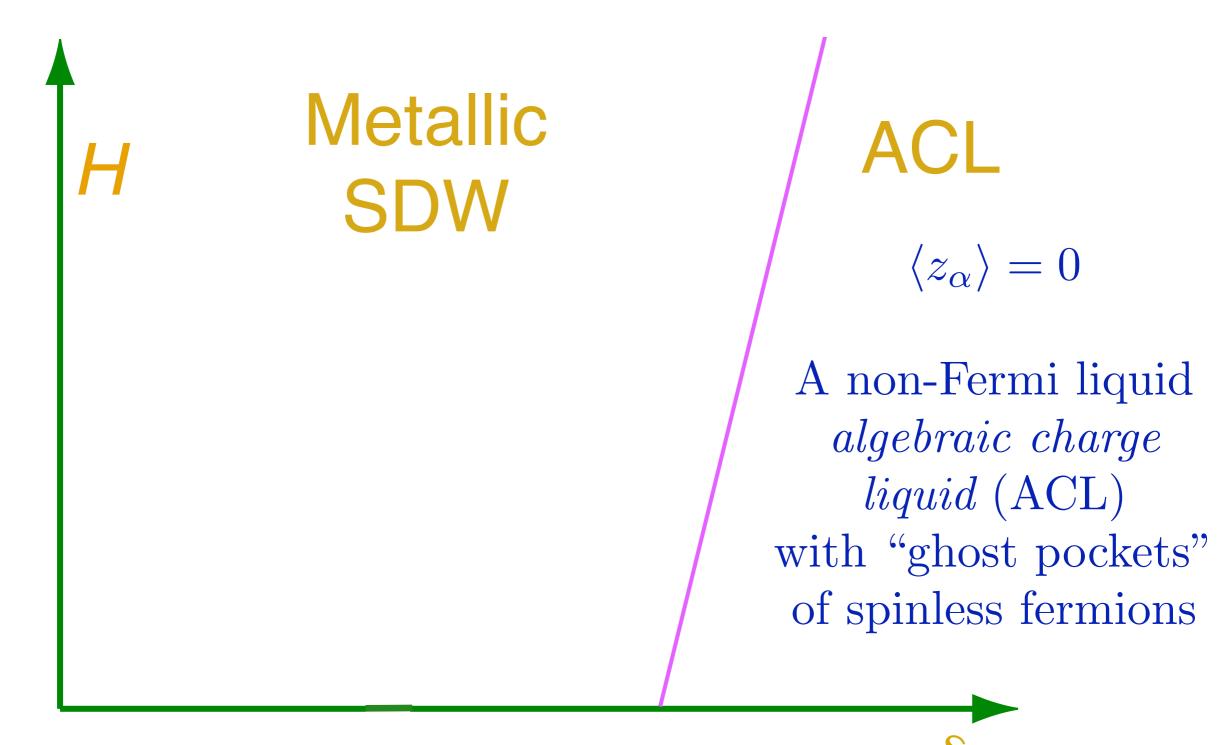
CP¹ field theory for z_{α} and an emergent U(1) gauge field A_{μ} . Coupling t tunes the strength of SDW orientation fluctuations.

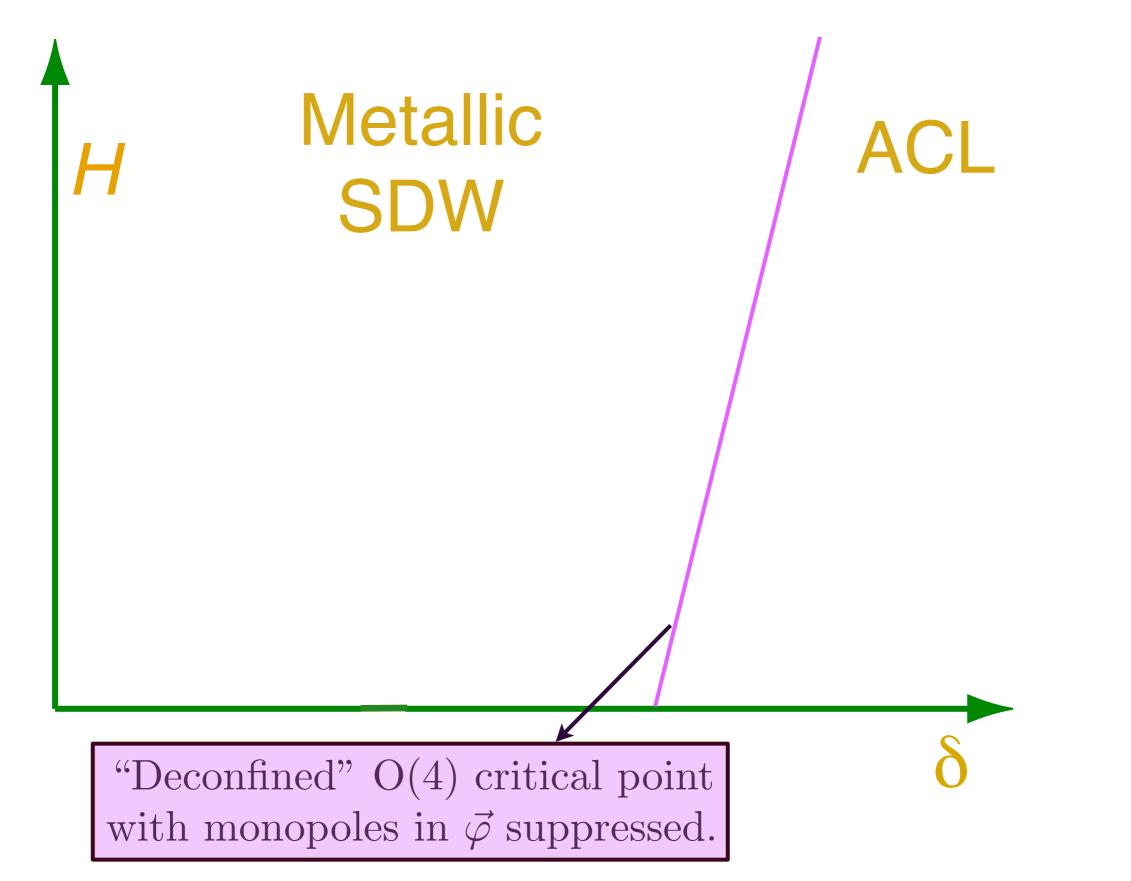
$$\mathcal{L}_g = g_+^{\dagger} \left[(\partial_{\tau} - iA_{\tau}) - \frac{1}{2m^*} (\nabla - i\mathbf{A})^2 - \mu \right] g_+ + g_-^{\dagger} \left[(\partial_{\tau} + iA_{\tau}) - \frac{1}{2m^*} (\nabla + i\mathbf{A})^2 - \mu \right] g_-$$

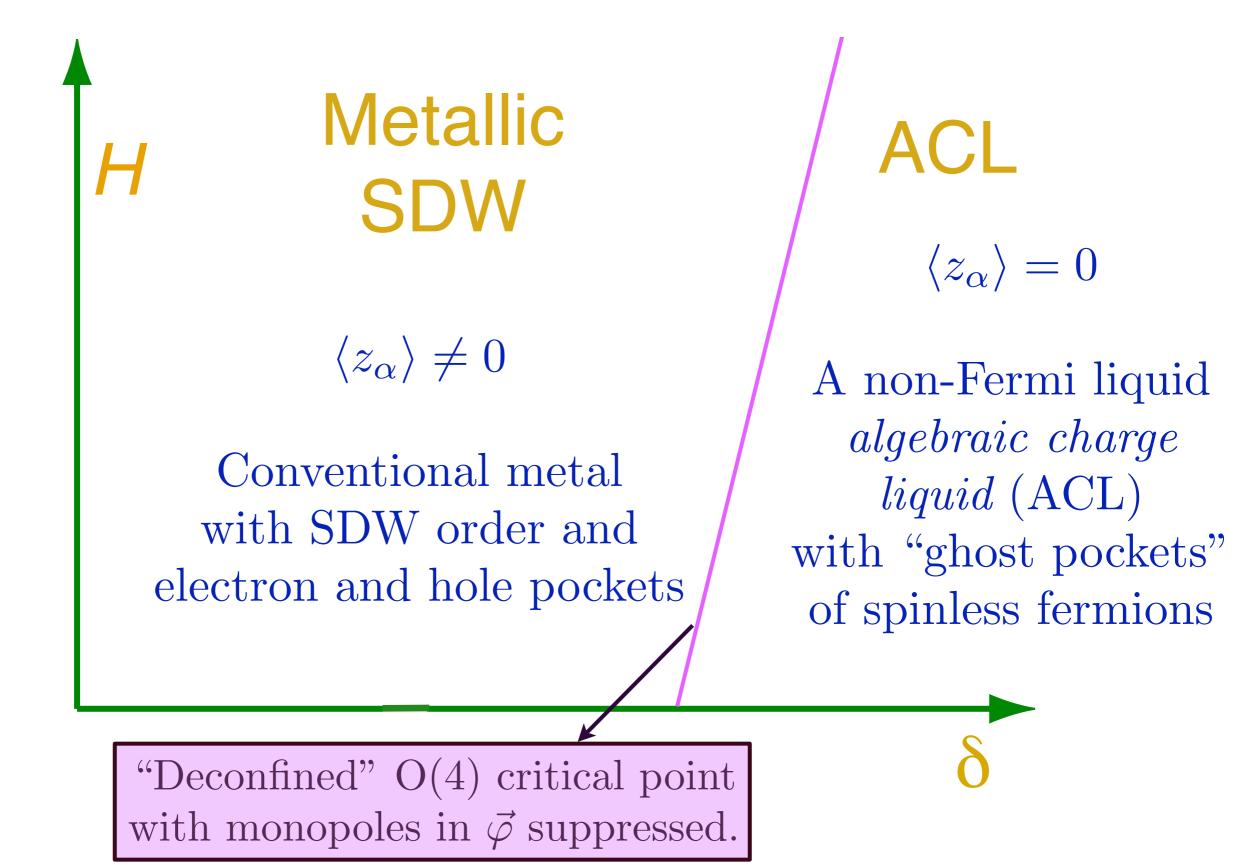
Two Fermi surfaces coupled to the emergent U(1) gauge field A_{μ} with opposite charges











Strong pairing of the g_{\pm} electron pockets

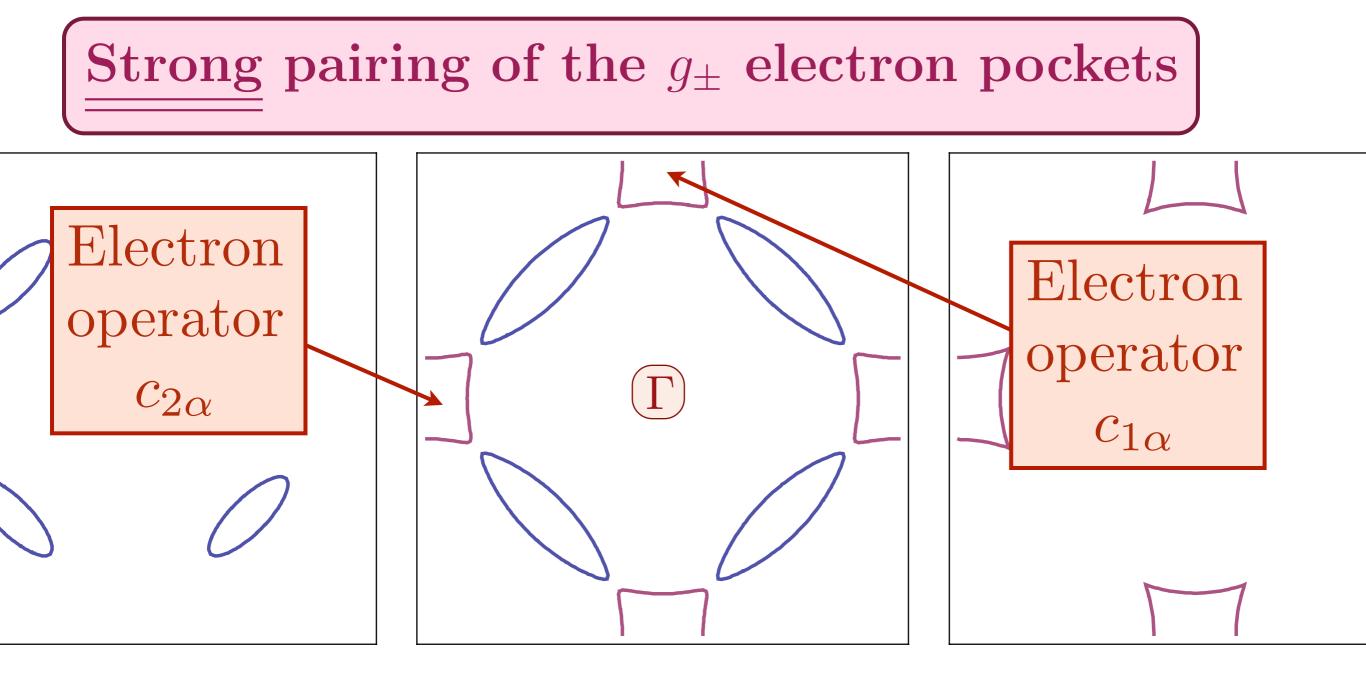
Problem is similar to double layer quantum Hall systems at total filling fraction ν = 1. At large layer spacing we have 2 composite fermion Fermi surfaces each at filling fraction ν = 1/2. At small layer spacing, there is a paired state formed by attractive interaction mediated by antisymmetric gauge field.

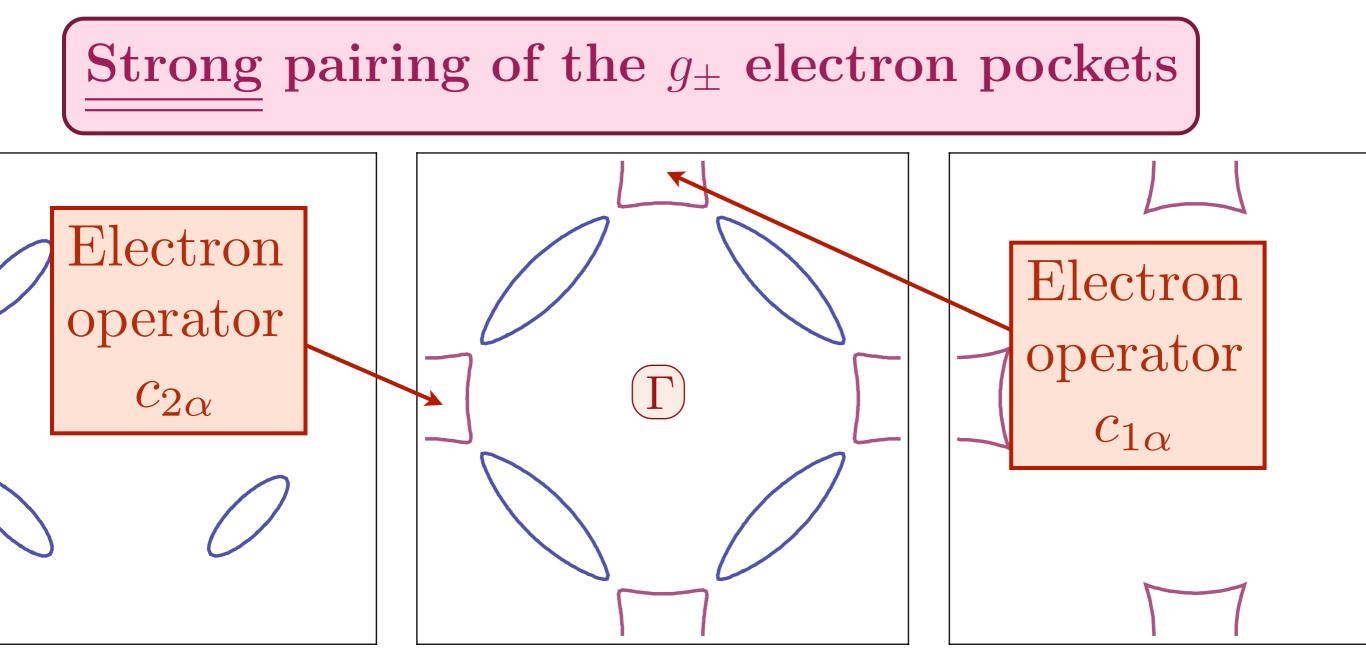
N. E. Bonesteel, I.A. McDonald, and C. Nayak, *Phys. Rev. Lett.* **77**, 3009 (1996). I. Ussishkin and A. Stern, *Phys. Rev. Lett.* **81**, 3932 (1998).

Strong pairing of the g_{\pm} electron pockets

- Problem is similar to double layer quantum Hall systems at total filling fraction ν = 1. At large layer spacing we have 2 composite fermion Fermi surfaces each at filling fraction ν = 1/2. At small layer spacing, there is a paired state formed by attractive interaction mediated by antisymmetric gauge field.
- Gauge forces lead to a *s*-wave paired state with a T_c of order the Fermi energy of the pockets. Inelastic scattering from low energy gauge modes lead to very singular g_{\pm} self energy, but is *not* pair-breaking.

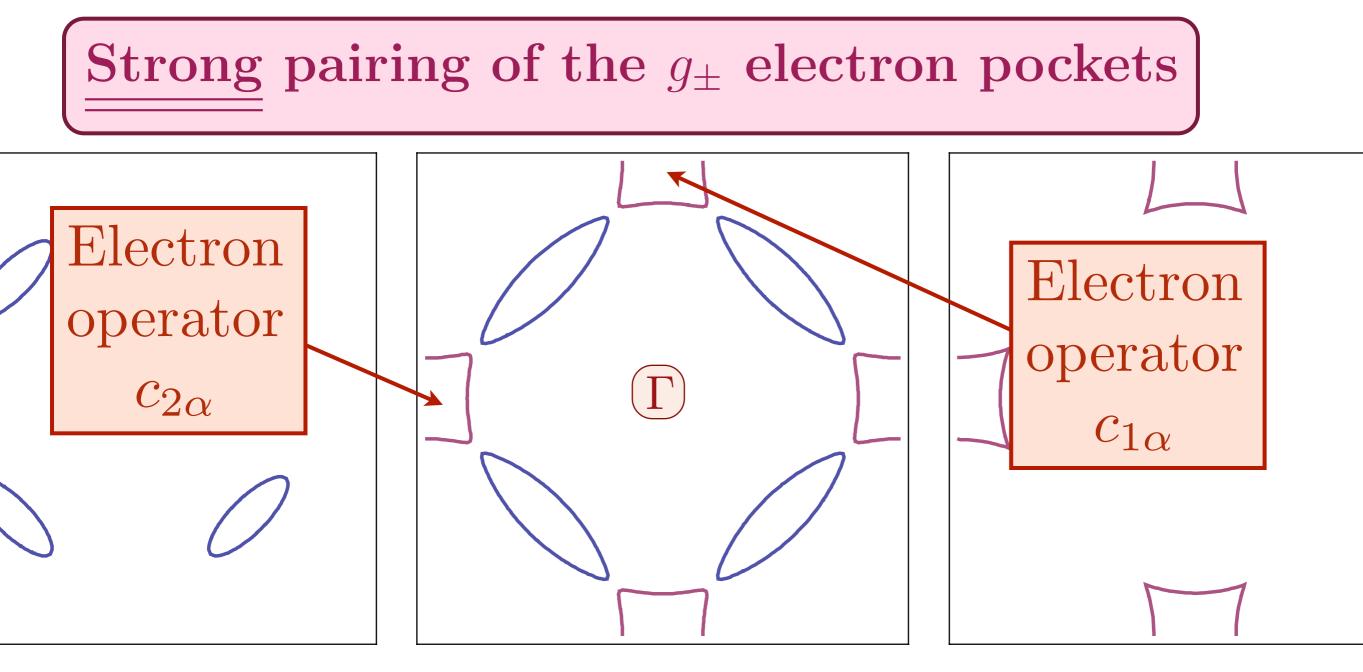
$$\langle g_+g_-\rangle = \Delta$$





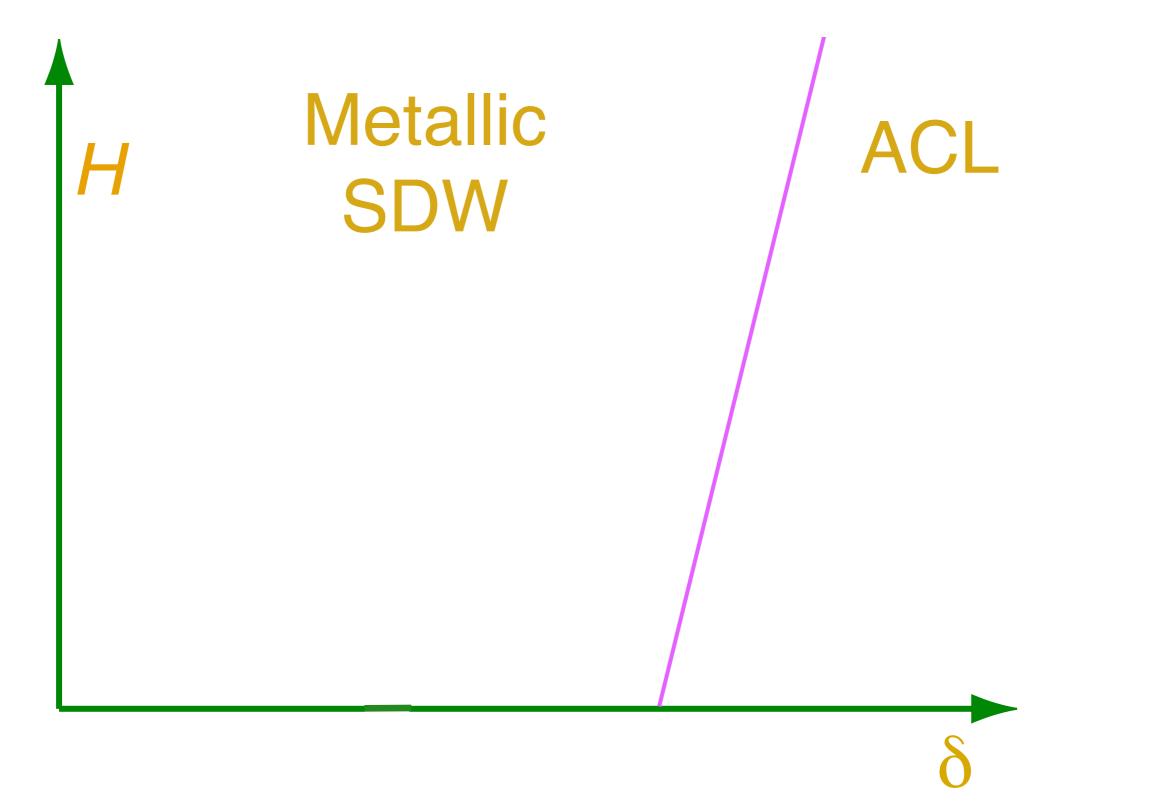
• Transforming back to the physical fermions:

$$\begin{pmatrix} c_{1\uparrow} \\ c_{1\downarrow} \end{pmatrix} = \begin{pmatrix} z_{\uparrow} & -z_{\downarrow}^{*} \\ z_{\downarrow} & z_{\uparrow}^{*} \end{pmatrix} \begin{pmatrix} g_{+} \\ g_{-} \end{pmatrix} ; \quad \begin{pmatrix} c_{2\uparrow} \\ c_{2\downarrow} \end{pmatrix} = \begin{pmatrix} z_{\uparrow} & -z_{\downarrow}^{*} \\ z_{\downarrow} & z_{\uparrow}^{*} \end{pmatrix} \begin{pmatrix} g_{+} \\ -g_{-} \end{pmatrix},$$
we find: $\langle c_{1\uparrow}c_{1\downarrow} \rangle = -\langle c_{2\uparrow}c_{2\downarrow} \rangle \sim \langle |z_{\uparrow}|^{2} + |z_{\downarrow}|^{2} \rangle \langle g_{+}g_{-} \rangle;$

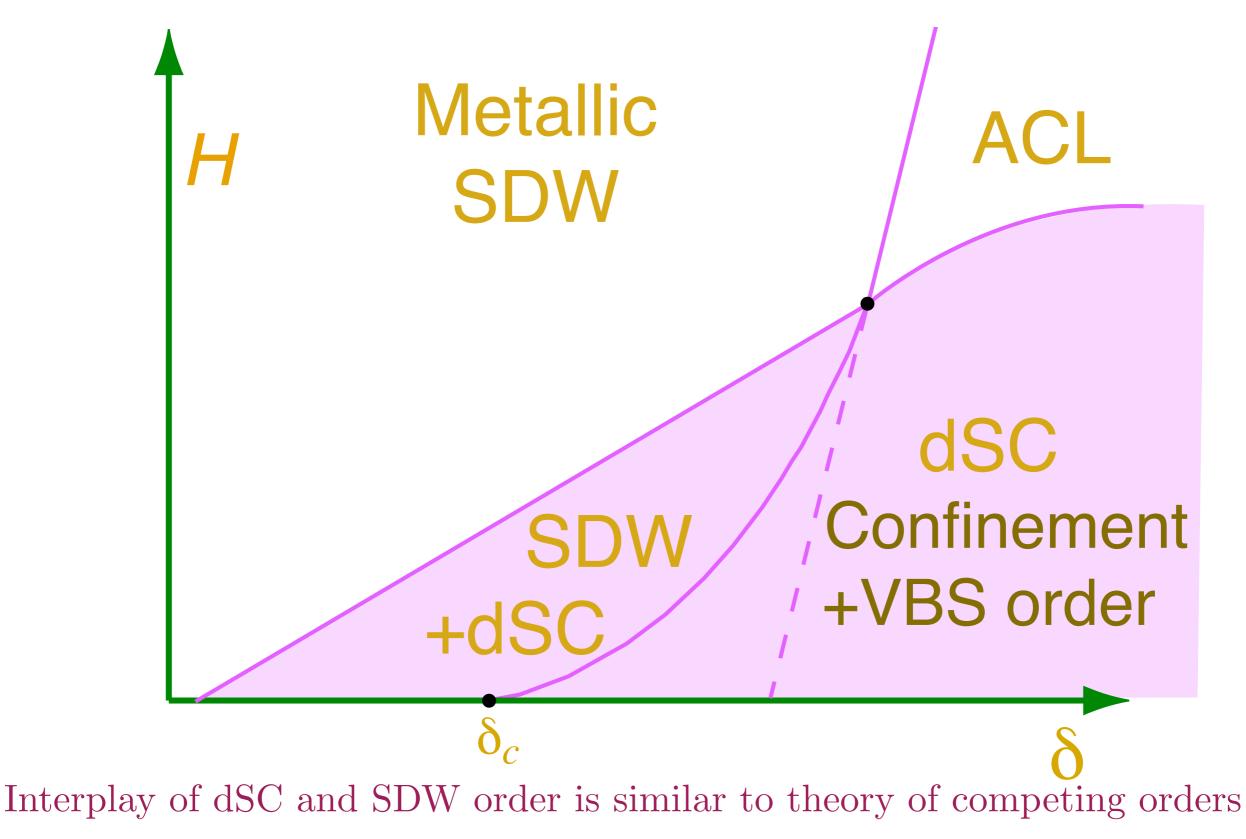


• Transforming back to the physical fermions:

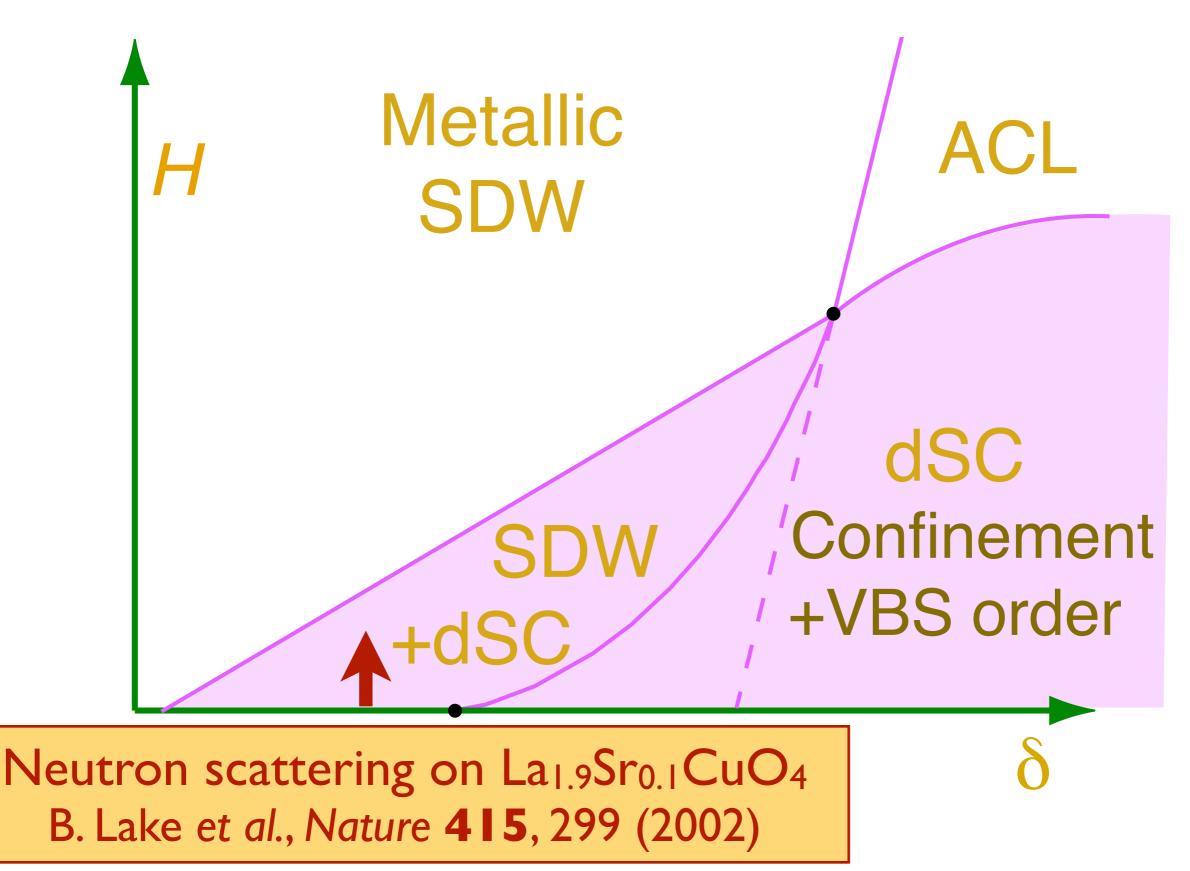
$$\begin{pmatrix} c_{1\uparrow} \\ c_{1\downarrow} \end{pmatrix} = \begin{pmatrix} z_{\uparrow} & -z_{\downarrow}^{*} \\ z_{\downarrow} & z_{\uparrow}^{*} \end{pmatrix} \begin{pmatrix} g_{+} \\ g_{-} \end{pmatrix} ; \quad \begin{pmatrix} c_{2\uparrow} \\ c_{2\downarrow} \end{pmatrix} = \begin{pmatrix} z_{\uparrow} & -z_{\downarrow}^{*} \\ z_{\downarrow} & z_{\uparrow}^{*} \end{pmatrix} \begin{pmatrix} g_{+} \\ -g_{-} \end{pmatrix},$$
we find: $\langle c_{1\uparrow}c_{1\downarrow}\rangle = -\langle c_{2\uparrow}c_{2\downarrow}\rangle \sim \langle |z_{\uparrow}|^{2} + |z_{\downarrow}|^{2}\rangle \langle g_{+}g_{-}\rangle;$
i.e. the pairing signature for the electrons is *d*-wave.

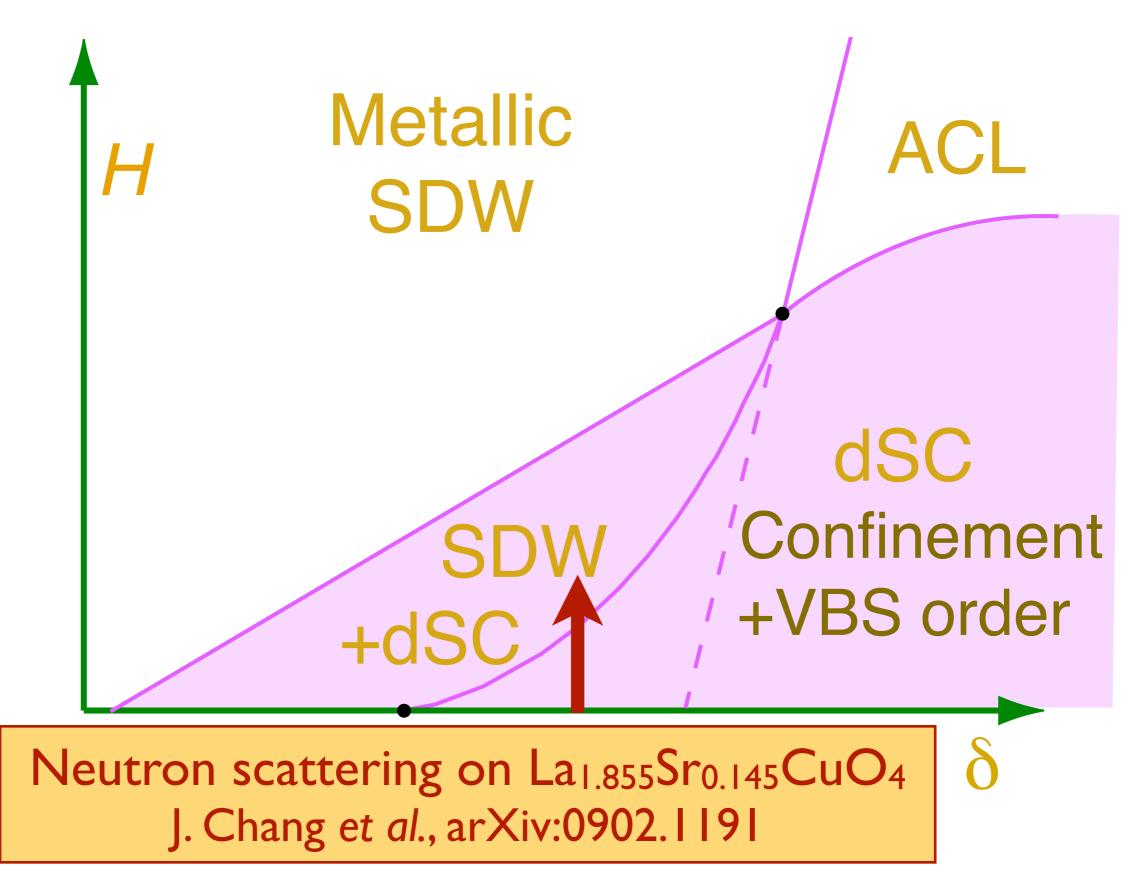


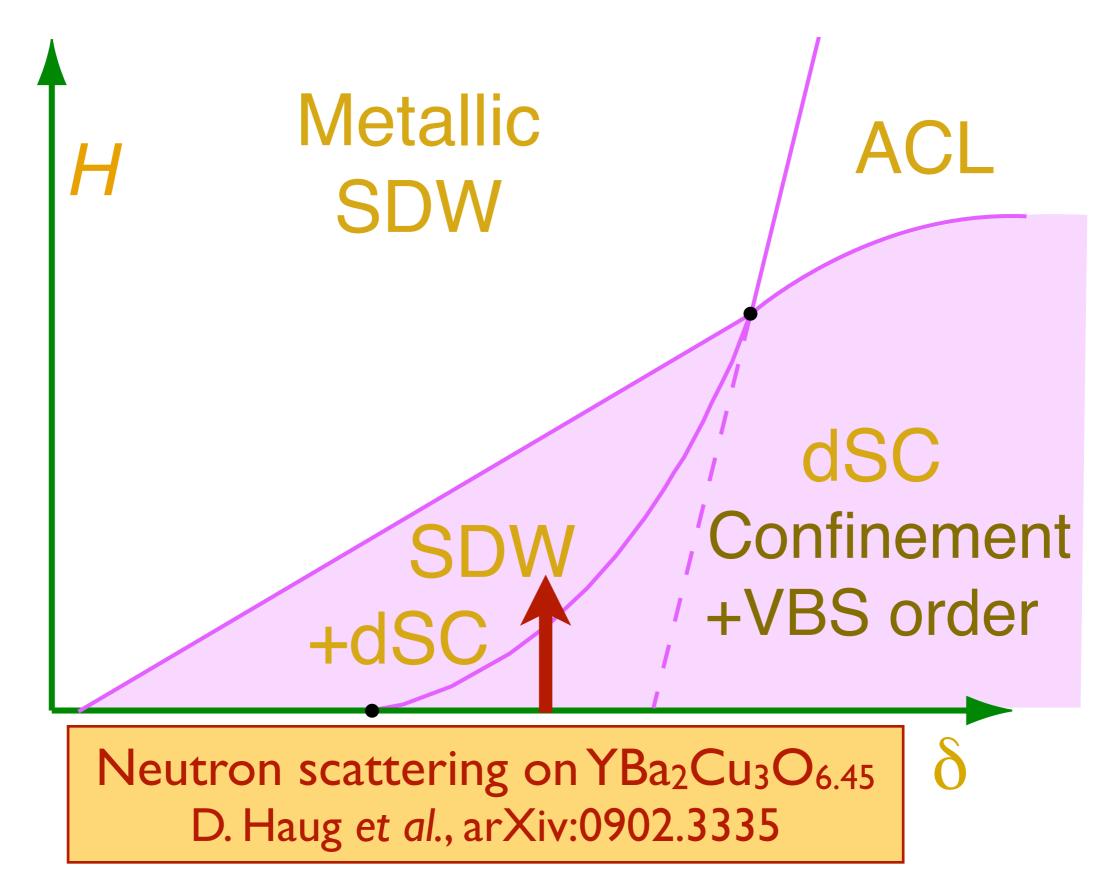
as a function of hole density $\delta \sim t$ and magnetic field H.



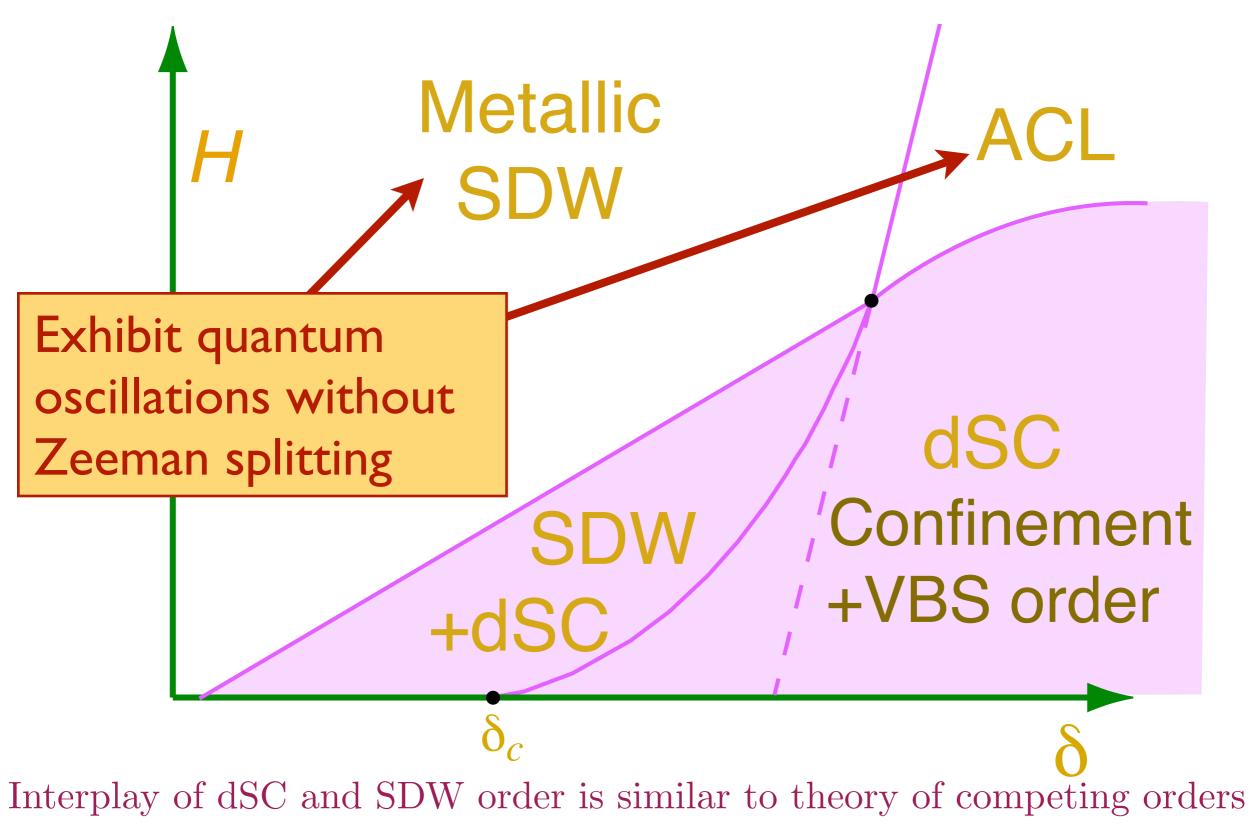
E. Demler, S. Sachdev and Y. Zhang, *Phys. Rev. Lett.* 87, 067202 (2001).







as a function of hole density $\delta \sim t$ and magnetic field H.



E. Demler, S. Sachdev and Y. Zhang, *Phys. Rev. Lett.* 87, 067202 (2001).

Simple numerical estimates

- Oscillation experiments provide the following information:
 - 1. Effective mass: $m^* \sim (1-3)m_e$
 - 2. Area of the Fermi surface: $A_{\rm e} \sim 3\% A_{\rm BZ}$
 - 3. Fermi temperature: $T_{\rm F} \sim (m_{\rm e}/m^*) 700 \,{\rm K}$
- The pairing critical field for the e-pocket:

 $H_{p2}(0) \sim H_{onset\,of\,oscillations} \sim 50\,T$

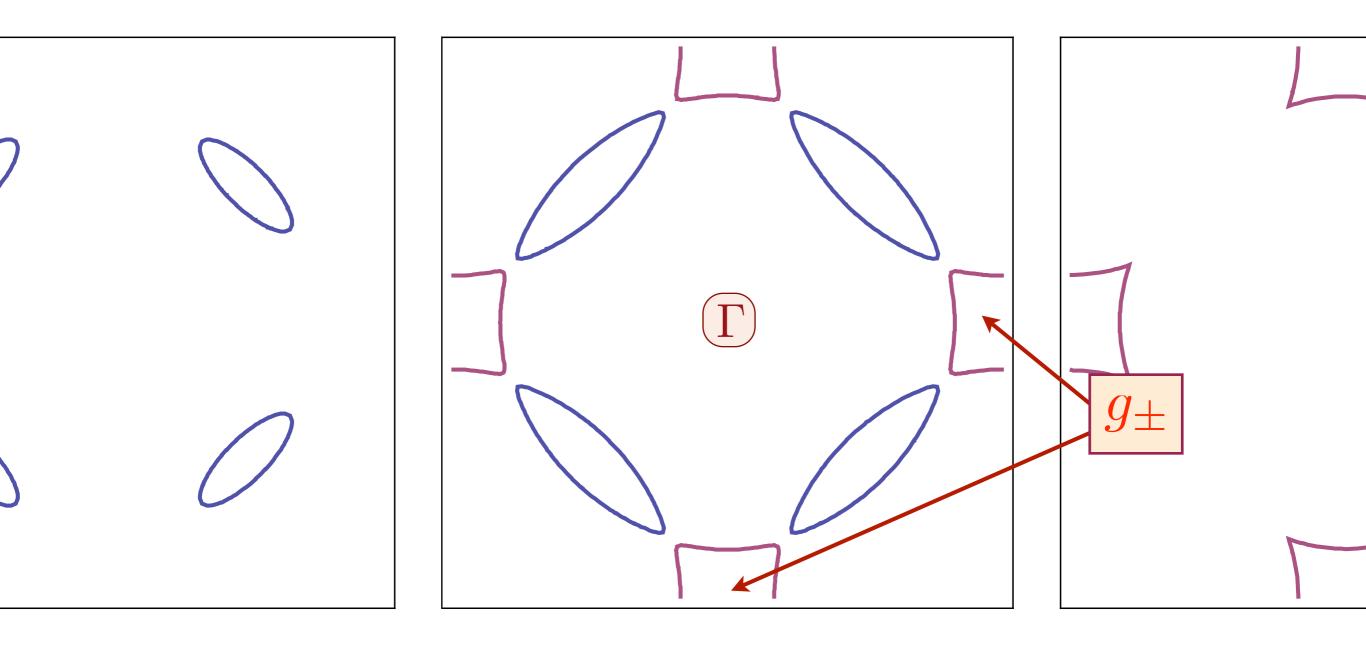
 For numerical estimates, we can use the BCS formula connecting *H*_{p2}(0) and the pairing temperature:

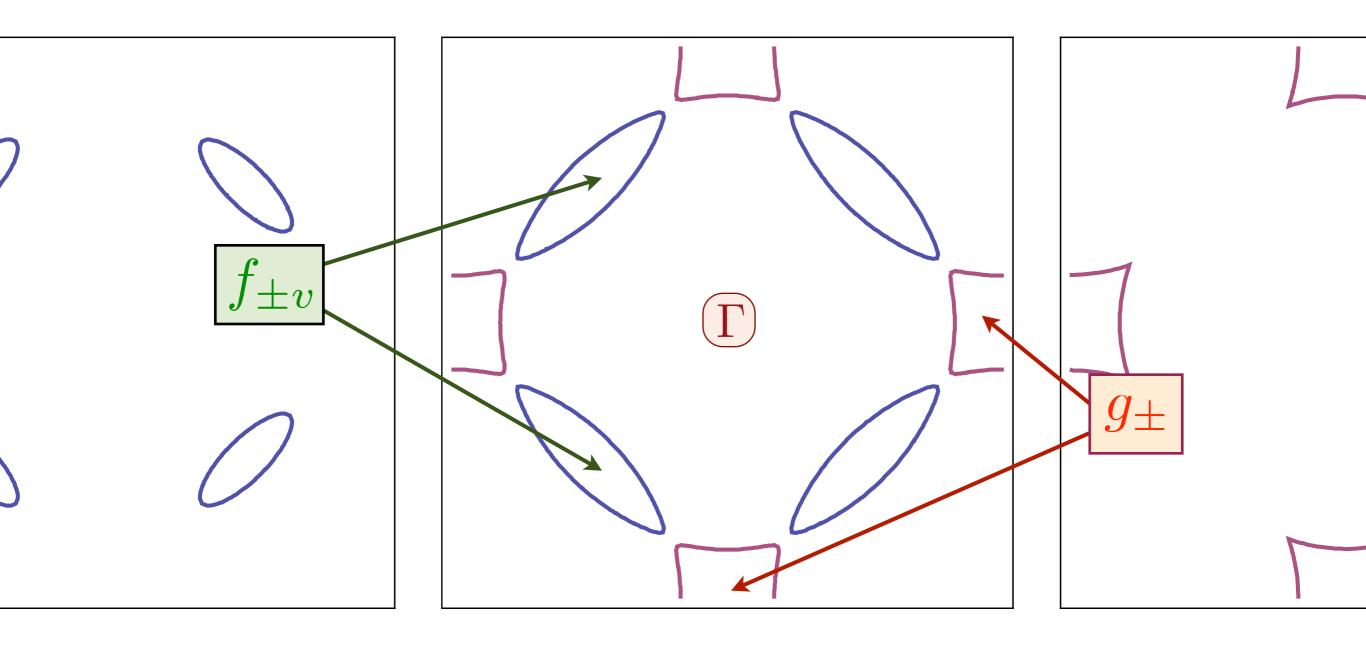
$$\frac{eH_{p2}(0)}{m^*c} = \frac{\pi^2 k_{\mathsf{B}} T_{\mathsf{p0}}^2}{\gamma_E \hbar T_{\mathsf{F}}}$$

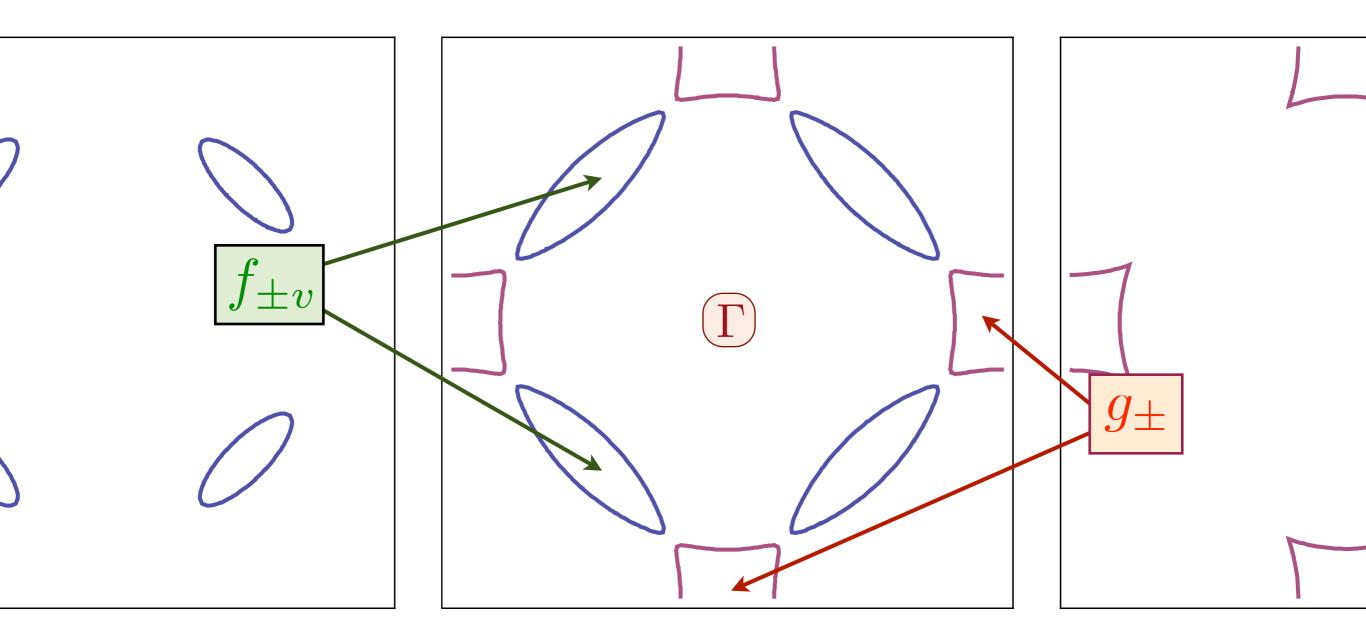
This leads to the pairing temperature of order

 $T_{\rm p0} \ {\rm K} \approx \sqrt{\frac{0.24 m_e}{m^*}} [T_{\rm F} \ {\rm K}] [H_{\rm p2}(0) \ {\rm T}] \sim 100 \ {\rm K} \ {\rm or} \ {\rm 2\Delta} \sim 400 \ {\rm K}$

Strong e-pocket pairing removes Fermi surface signatures from *H*=0 photoemission experiments







Low energy theory for spinless, charge +e fermions $f_{\pm v}$:

$$\mathcal{L}_{f} = \sum_{v=1,2} \left\{ f_{+v}^{\dagger} \left[(\partial_{\tau} - iA_{\tau}) - \frac{1}{2m^{*}} (\nabla - i\mathbf{A})^{2} - \mu \right] f_{+v} \right. \\ \left. + f_{-v}^{\dagger} \left[(\partial_{\tau} + iA_{\tau}) - \frac{1}{2m^{*}} (\nabla + i\mathbf{A})^{2} - \mu \right] f_{-v} \right\}$$

<u>Weak</u> pairing of the f_{\pm} hole pockets

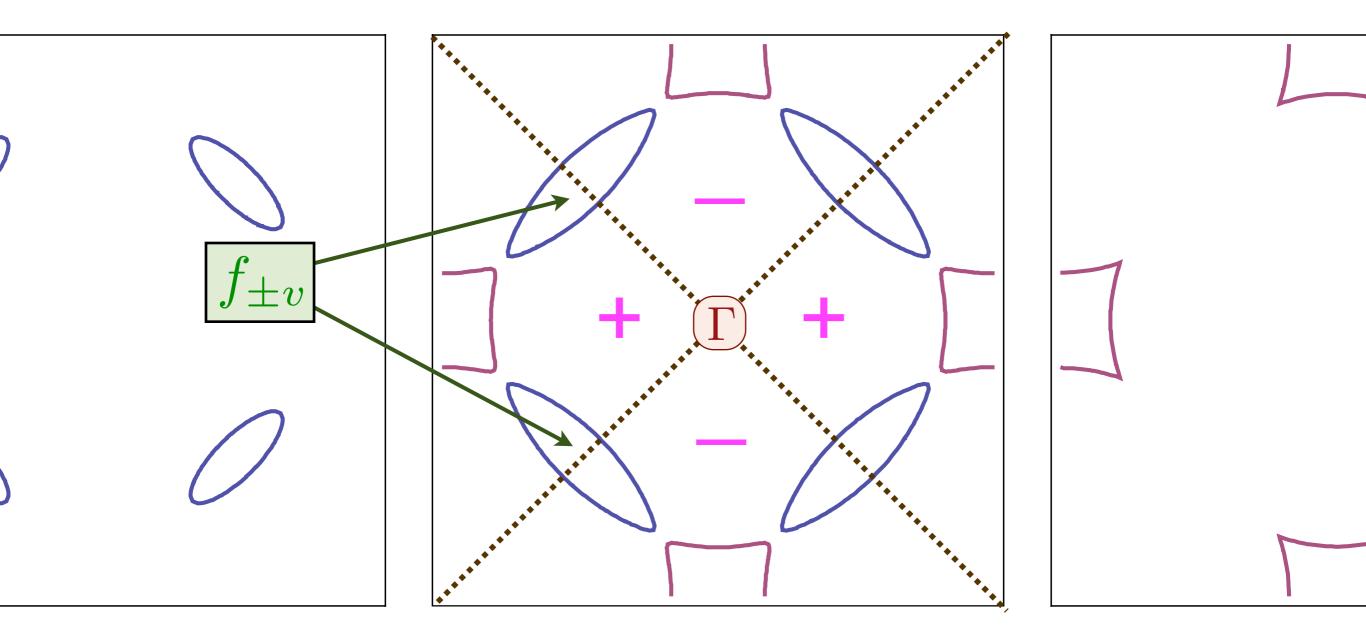
$$\mathcal{L}_{\text{Josephson}} = iJ \left[g_{+}g_{-} \right] \left[f_{+1} \stackrel{\leftrightarrow}{\partial}_{x} f_{-1} - f_{+1} \stackrel{\leftrightarrow}{\partial}_{y} f_{-1} + f_{+2} \stackrel{\leftrightarrow}{\partial}_{x} f_{-2} + f_{+2} \stackrel{\leftrightarrow}{\partial}_{y} f_{-2} \right] + \text{H.c.}$$

V. B. Geshkenbein, L. B. Ioffe, and A. I. Larkin, Phys. Rev. B 55, 3173 (1997).

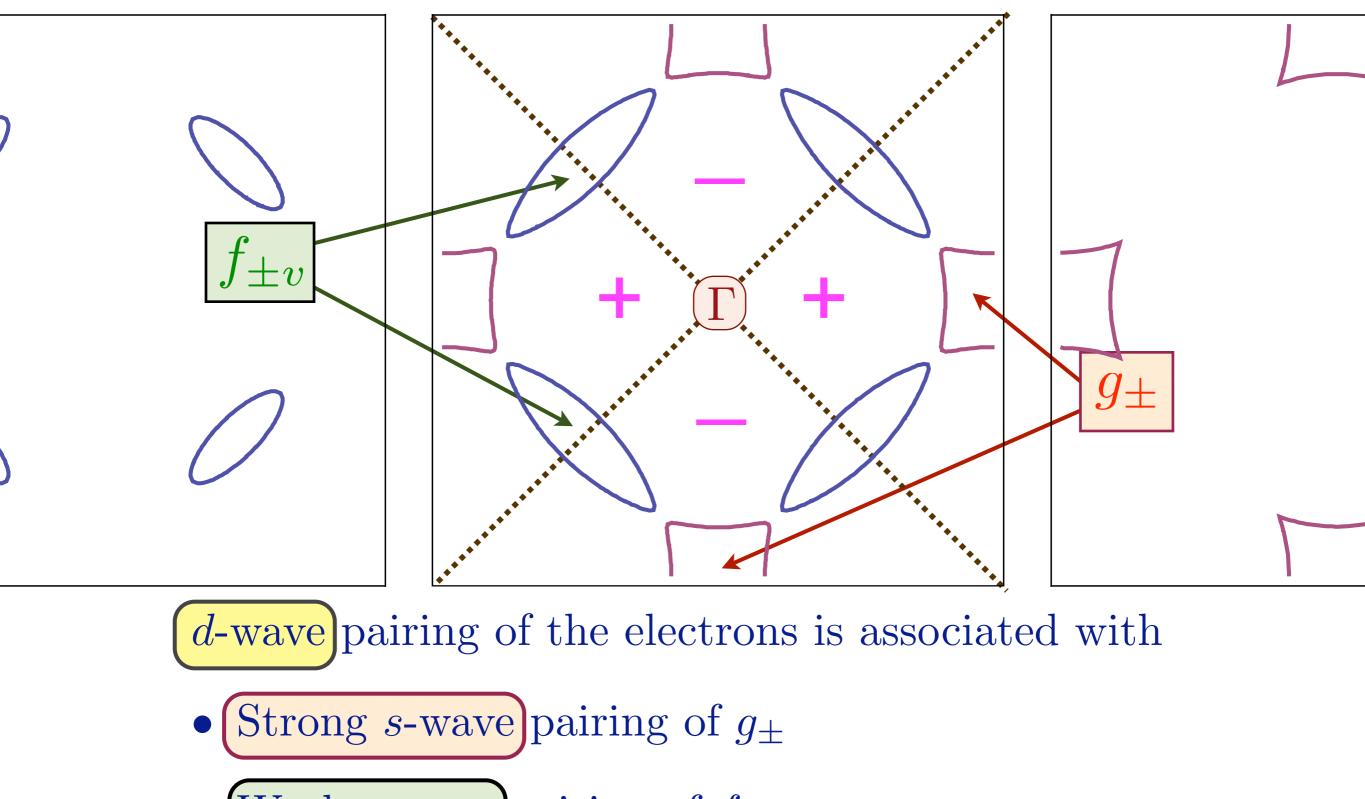
Proximity Josephson coupling J to g_{\pm} fermions leads to p-wave pairing of the $f_{\pm v}$ fermions. The A_{μ} gauge forces are pair-breaking, and so the pairing is weak.

$$\langle f_{+1}(\mathbf{k})f_{-1}(-\mathbf{k})\rangle \sim (k_x - k_y)J\langle g_+g_-\rangle; \langle f_{+2}(\mathbf{k})f_{-2}(-\mathbf{k})\rangle \sim (k_x + k_y)J\langle g_+g_-\rangle; \langle f_{+1}(\mathbf{k})f_{-2}(-\mathbf{k})\rangle = 0,$$

<u>Weak</u> pairing of the f_{\pm} hole pockets



 $\langle f_{+1}(\mathbf{k})f_{-1}(-\mathbf{k})\rangle \sim (k_x - k_y)J\langle g_+g_-\rangle;$ $\langle f_{+2}(\mathbf{k})f_{-2}(-\mathbf{k})\rangle \sim (k_x + k_y)J\langle g_+g_-\rangle;$ $\langle f_{+1}(\mathbf{k})f_{-2}(-\mathbf{k})\rangle = 0,$



• Weak *p*-wave pairing of $f_{\pm v}$.

<u>Conclusions</u>

* Non-Landau-Ginzburg theory for loss of spin density wave order in a metal

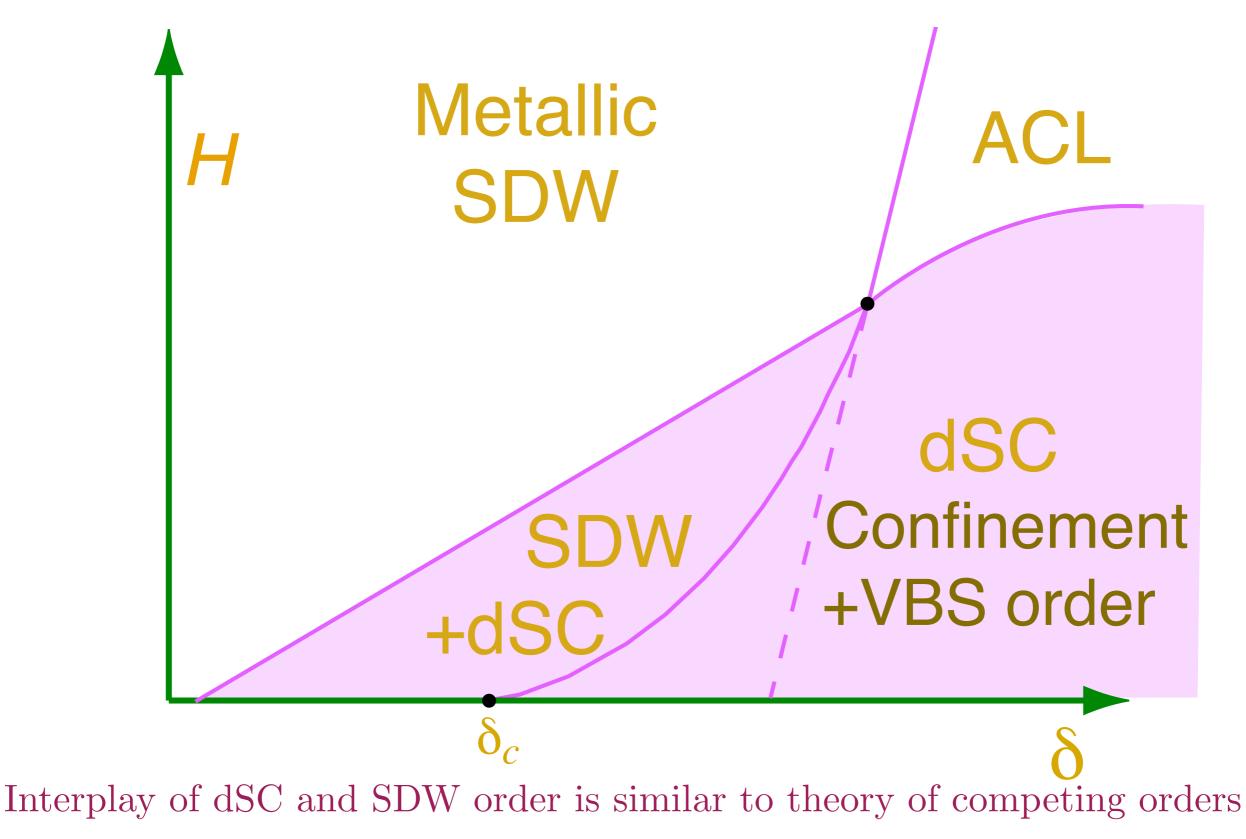
Conclusions

- * Non-Landau-Ginzburg theory for loss of spin density wave order in a metal
- * Natural route to *d*-wave pairing with strong pairing at the antinodes and weak pairing at the nodes

Conclusions

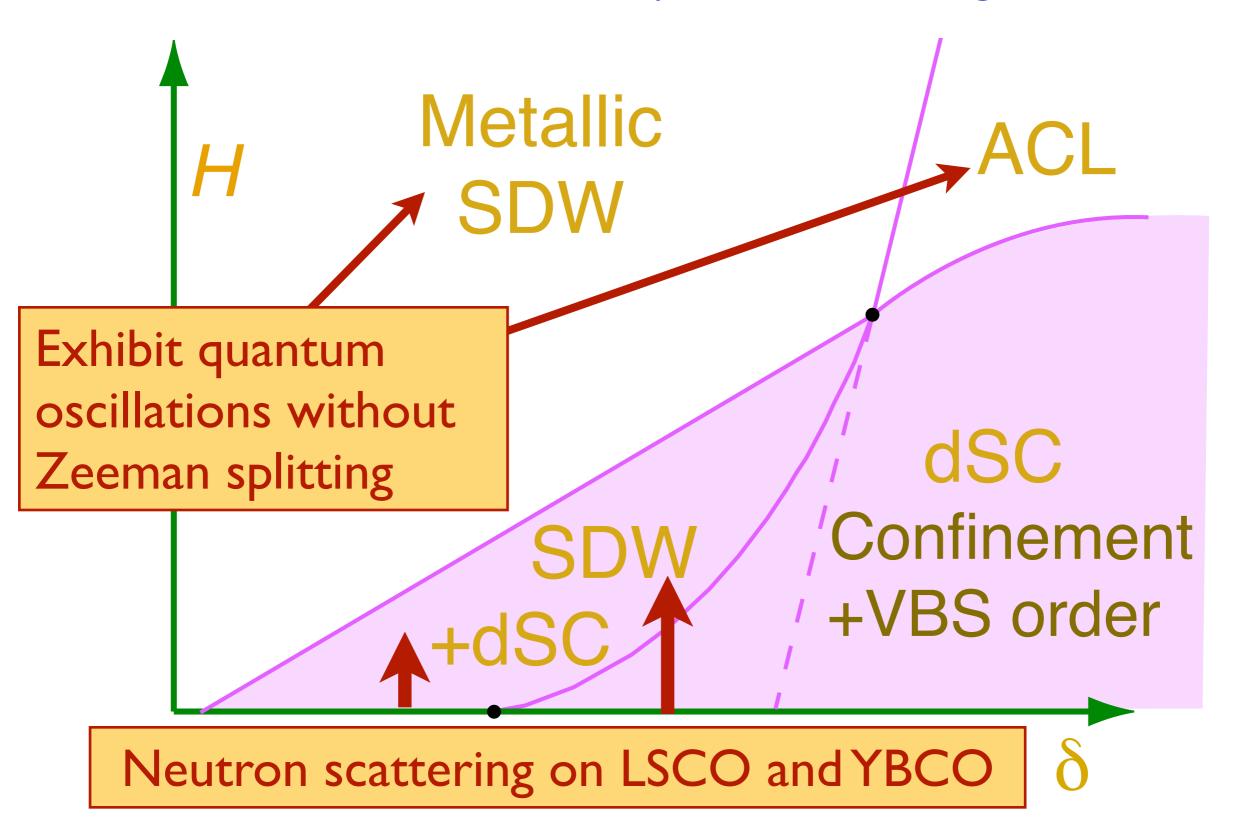
- * Non-Landau-Ginzburg theory for loss of spin density wave order in a metal
- * Natural route to *d*-wave pairing with strong pairing at the antinodes and weak pairing at the nodes
- New metallic state, the ACL with "ghost" electron and hole pockets, is a useful starting point for building field-doping phase diagram

as a function of hole density $\delta \sim t$ and magnetic field H.



E. Demler, S. Sachdev and Y. Zhang, *Phys. Rev. Lett.* 87, 067202 (2001).

as a function of hole density $\delta \sim t$ and magnetic field H.



E. Demler, S. Sachdev and Y. Zhang, *Phys. Rev. Lett.* 87, 067202 (2001).

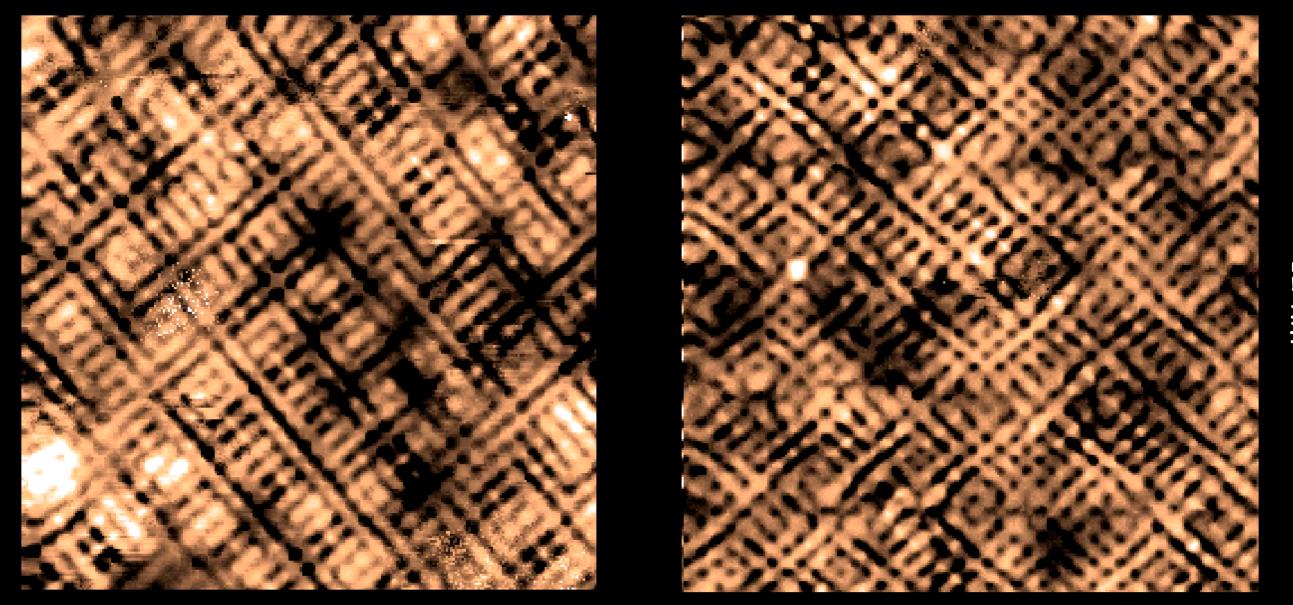
Conclusions

- * Non-Landau-Ginzburg theory for loss of spin density wave order in a metal
- * Natural route to *d*-wave pairing with strong pairing at the antinodes and weak pairing at the nodes
- New metallic state, the ACL with "ghost" electron and hole pockets, is a useful starting point for building field-doping phase diagram

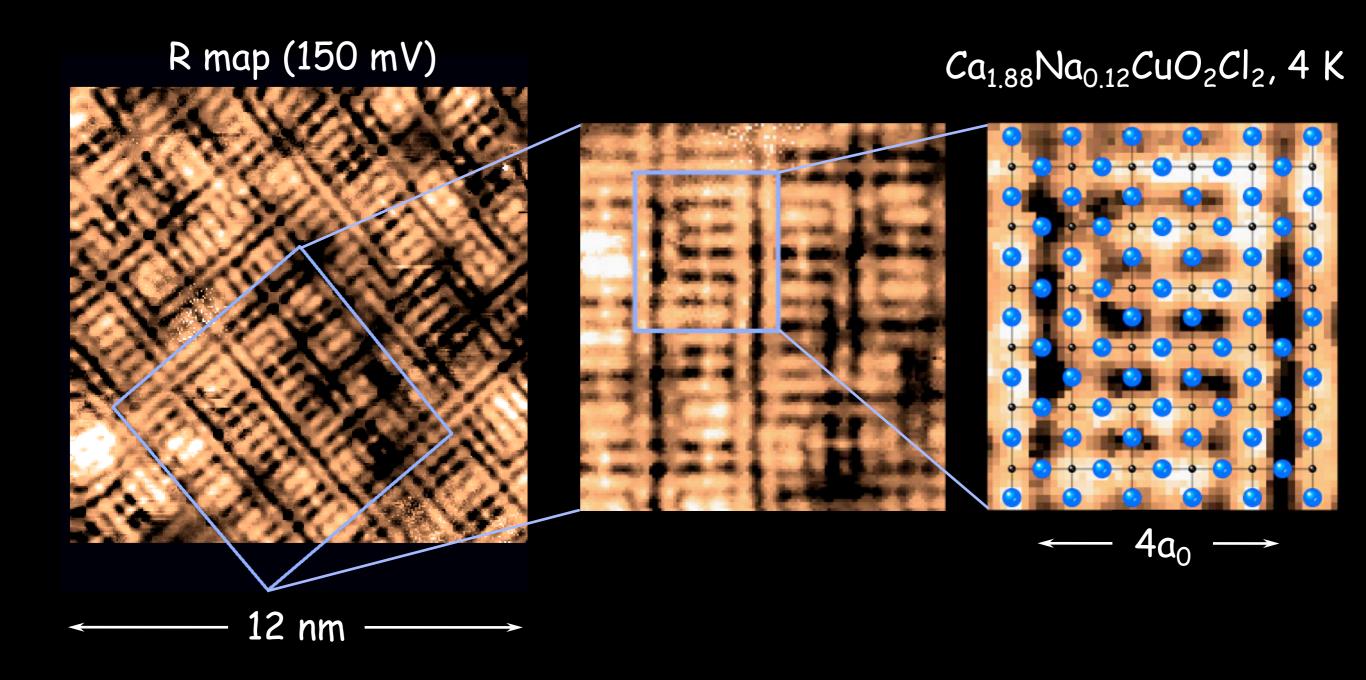
Conclusions

- * Non-Landau-Ginzburg theory for loss of spin density wave order in a metal
- * Natural route to *d*-wave pairing with strong pairing at the antinodes and weak pairing at the nodes
- New metallic state, the ACL with "ghost" electron and hole pockets, is a useful starting point for building field-doping phase diagram
- * Paired electron pockets are expected to lead to valence-bond-solid modulations at low temperature

Tunneling Asymmetry (TA)-map at E=150meV $Ca_{1.90}Na_{0.10}CuO_2Cl_2$ $Bi_{2.2}Sr_{1.8}Ca_{0.8}Dy_{0.2}Cu_2O_y$

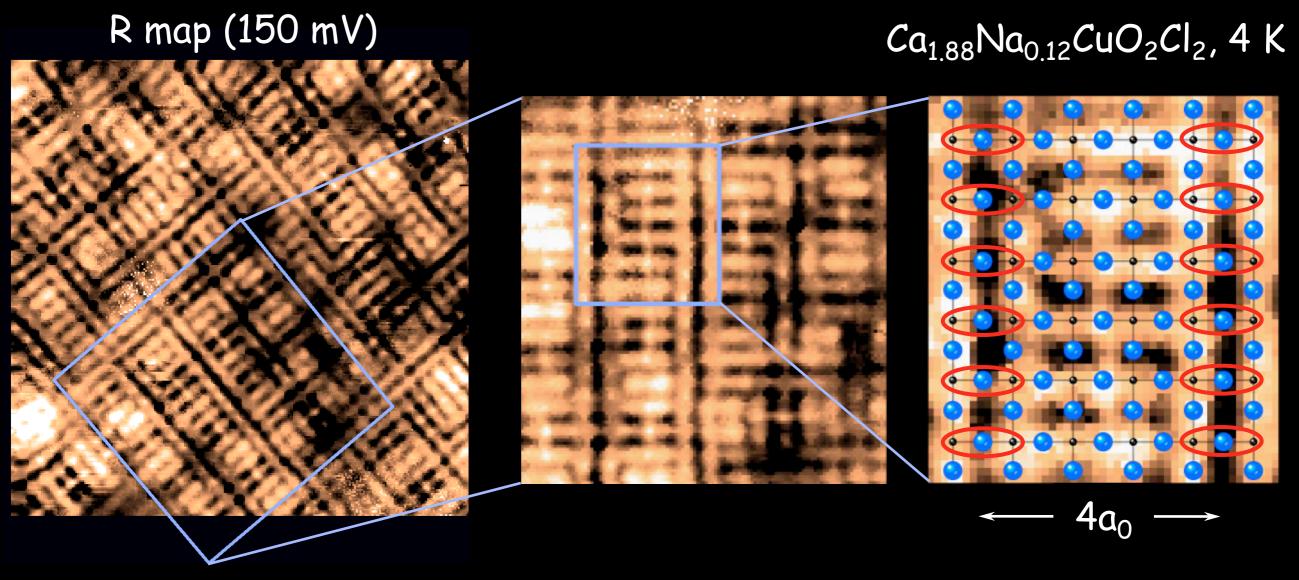


Indistinguishable bond-centered TA contrast with disperse 4a₀-wide nanodomains Y. Kohsaka et al. Science 315, 1380 (2007) TA Contrast is at oxygen site (Cu-O-Cu bond-centered)



Y. Kohsaka et al. Science 315, 1380 (2007)

TA Contrast is at oxygen site (Cu-O-Cu bond-centered)



12 nm ——

Evidence for a predicted valence bond supersolid

S. Sachdev and N. Read, *Int. J. Mod. Phys.* B **5**, 219 (1991). M.Vojta and S. Sachdev, *Phys. Rev. Lett.* **83**, 3916 (1999).

Conclusions

- * Non-Landau-Ginzburg theory for loss of spin density wave order in a metal
- * Natural route to *d*-wave pairing with strong pairing at the antinodes and weak pairing at the nodes
- New metallic state, the ACL with "ghost" electron and hole pockets, is a useful starting point for building field-doping phase diagram
- * Paired electron pockets are expected to lead to valence-bond-solid modulations at low temperature

Conclusions

- * Non-Landau-Ginzburg theory for loss of spin density wave order in a metal
- * Natural route to *d*-wave pairing with strong pairing at the antinodes and weak pairing at the nodes
- New metallic state, the ACL with "ghost" electron and hole pockets, is a useful starting point for building field-doping phase diagram
- * Paired electron pockets are expected to lead to valence-bond-solid modulations at low temperature
- Needed: theory for transition to "large" Fermi surface at higher doping