

On the Condensed Matter Analog of Baryon Chiral Perturbation Theory

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In collaboration with C. Brügger, F. Kämpfer, M. Moser, M. Pepe and U.-J. Wiese

Facultad de Ciencias, UCOL

October 10, 2008

Outline

- 1 Motivation
 - Condensed matter analog of baryon chiral perturbation theory
 - Effective field theory for magnons
- 2 Hole-Doping
 - Construction of effective field theory
- 3 Electron-Doping
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- 4 Spiral Phases
 - Uniform Background Field
 - Dispersion Relations of Indoped Fermions
 - Spiral Phases for Holes
 - Homogeneous Phase for Electrons
- 5 Conclusions

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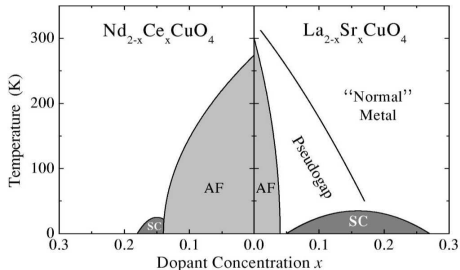
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Motivation: High- T_c superconductivity in cuprates

1986: Bednorz and Müller discover high- T_c superconductivity by doping copper oxide compounds (cuprates):



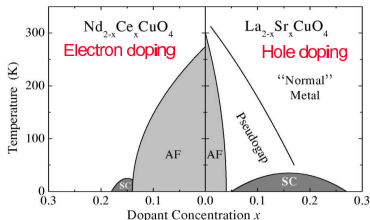
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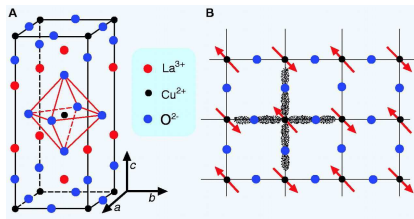
- SC results from doping antiferromagnetic insulators
- Doping possible with both electrons and holes
- Electron-Hole asymmetry

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Phase diagram of cuprates:



Crystal structure:



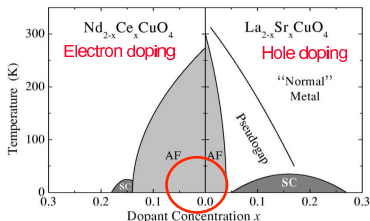
Damascelli, Hussain, and Shen,
 Rev. Mod. Phys. 75 (2003) 473

Orenstein and Millis, Science 288 (2000) 468

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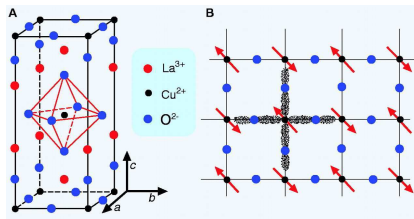
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- SC results from doping antiferromagnetic insulators
- Doping possible with both electrons and holes
- Common structure: CuO_2 layers separated by spacer layers
- Concentrate on antiferromagnetic region: low doping, low T

Microscopic description: The Hubbard model

The **Hubbard Hamiltonian** defined on a square lattice:

$$\begin{aligned}
 H = & -t \sum_{x,i} (c_{x\uparrow}^\dagger c_{x+\hat{i}\uparrow} + c_{x+\hat{i}\uparrow}^\dagger c_{x\uparrow} + c_{x\downarrow}^\dagger c_{x+\hat{i}\downarrow} + c_{x+\hat{i}\downarrow}^\dagger c_{x\downarrow}) \\
 & + U \sum_x c_{x\uparrow}^\dagger c_{x\uparrow} c_{x\downarrow}^\dagger c_{x\downarrow} - \mu \sum_x (c_{x\uparrow}^\dagger c_{x\uparrow} + c_{x\downarrow}^\dagger c_{x\downarrow} - 1)
 \end{aligned}$$

- Parameters:

t : Hopping parameter (nearest neighbors)

U : On-site Coulomb repulsion

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- Parameters:
 - t : Hopping parameter (nearest neighbors)
 - U : On-site Coulomb repulsion
 - μ : Chemical potential for fermion number
- Minimal model for cuprates: contains the relevant physics
- Away from half-filling: Hamiltonian virtually unsolvable from first principles (Neither analytically nor numerically)

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- Symmetries:

$SU(2)_s$: Global spin rotation

$U(1)_Q$: Fermion number conservation

D_i : Displacement by one lattice spacing

O : 90 degrees rotation

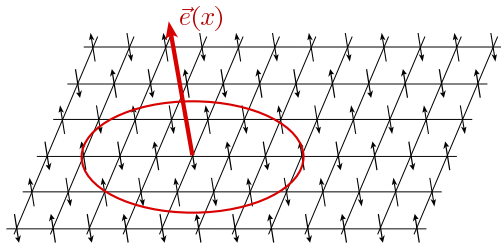
R : Reflection on a lattice axis

T : Time reversal

Antiferromagnetism: Near half-filling (1 fermion per site)

Near half-filling:

- **Antiferromagnetic** alignment of spins is preferred
- Spontaneous symmetry breaking: $SU(2)_S \rightarrow U(1)_S$
- Goldstone's theorem: 2 massless excitations \implies **2 magnons**

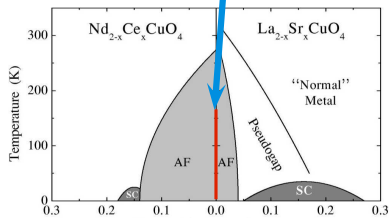


Systematic effective field theory description

	Antiferromagnets	QCD
Spont. symm. breaking	$SU(2)_s \longrightarrow U(1)_s$	$SU(2)_L \otimes SU(2)_R \rightarrow SU(2)_{L=R}$
GB physics	Magnon perturbation theory	Chiral perturbation theory
GB + matter physics	Effective theory presented here	Baryon chiral perturbation theory

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Pure magnon sector: Magnon perturbation theory

Spontaneous global $SU(2)_s \rightarrow U(1)_s$ spin symmetry breaking:

- 2 Goldstone bosons (magnons) described by

$$\vec{e}(x) = (e_1(x), e_2(x), e_3(x)) \in S^2 = SU(2)_s / U(1)_s$$

with $x = (x_1, x_2, t)$

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- Low-energy magnon physics described by nonlinear σ -model

$$\mathcal{L} = \frac{\rho_s}{2} (\partial_i \vec{e} \cdot \partial_i \vec{e} + \frac{1}{c^2} \partial_t \vec{e} \cdot \partial_t \vec{e}) + \dots$$

ρ_s : spin stiffness c : spin wave velocity

Chakravarty, Halperin, and Nelson, PRB 39 (1989) 2344
 Hasenfratz and Niedermayer, Phys. Lett. B268 (1991) 231

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State of affairs: Fermionic sector of effective theory

Earlier attempts by: Shraiman and Siggia, Wen, Shankar, ...

General agreement:

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No agreement on low-energy effective Lagrangian for fermions:

- Conflicting realizations of fermion fields
- Non-unique structure of terms in Lagrangians

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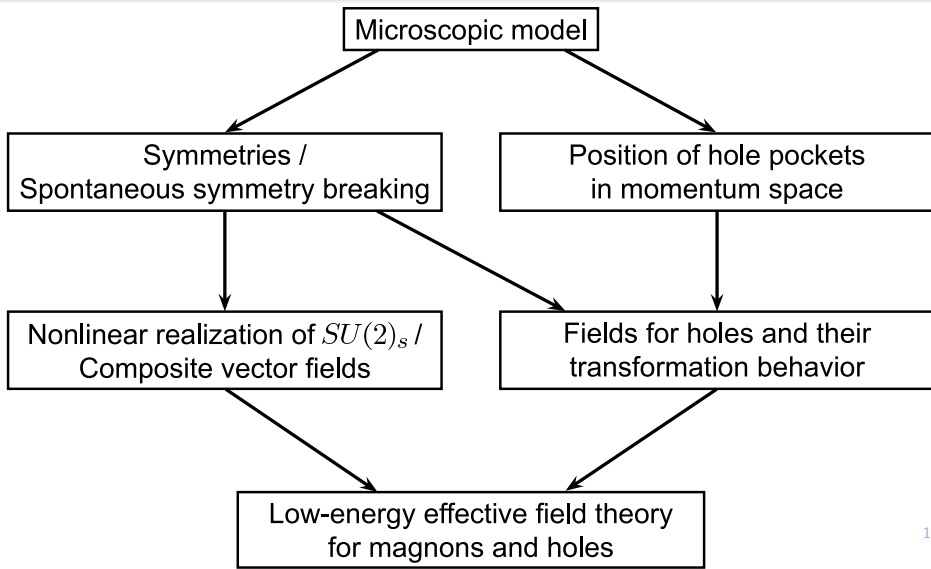
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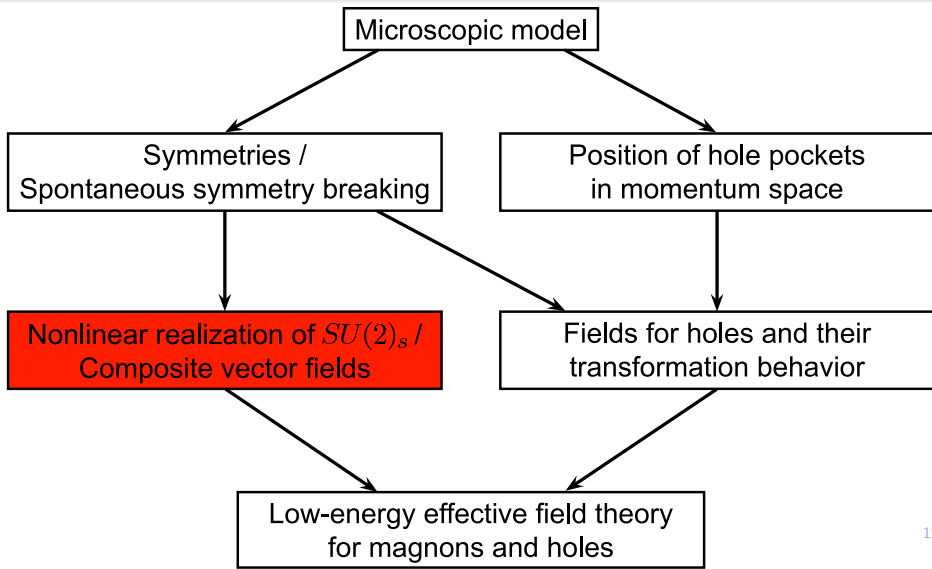
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⇒ Construction of a **systematic low-energy effective field theory** for magnons and holes analogous to baryon chiral perturbation theory

Symmetry-based construction of effective theory



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Nonlinear realization of $SU(2)_S$ symmetry

- $\mathbb{C}P(1)$ representation of magnon field

$$P(x) = \frac{1}{2}(\mathbb{1} + \vec{e}(x) \cdot \vec{\sigma}) \in \mathbb{C}P(1)$$

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$$u(x)P(x)u(x)^\dagger = \begin{pmatrix} 1 & 0 \\ 0 & 0 \end{pmatrix}, \quad u(x) \in SU(2)_s, \quad u_{11}(x) \geq 0$$

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Global $SU(2)_s$ rotation manifests itself as local $U(1)_s$ transformation!

Composite vector fields

- We introduce an anti-Hermitian field

$$v_{\mu}(x) = u(x)\partial_{\mu}u(x)^{\dagger} = \begin{pmatrix} v_{\mu}^3(x) & v_{\mu}^+(x) \\ v_{\mu}^-(x) & -v_{\mu}^3(x) \end{pmatrix}$$

with $\mu \in \{1, 2, t\}$

- Components used to couple magnons to holes

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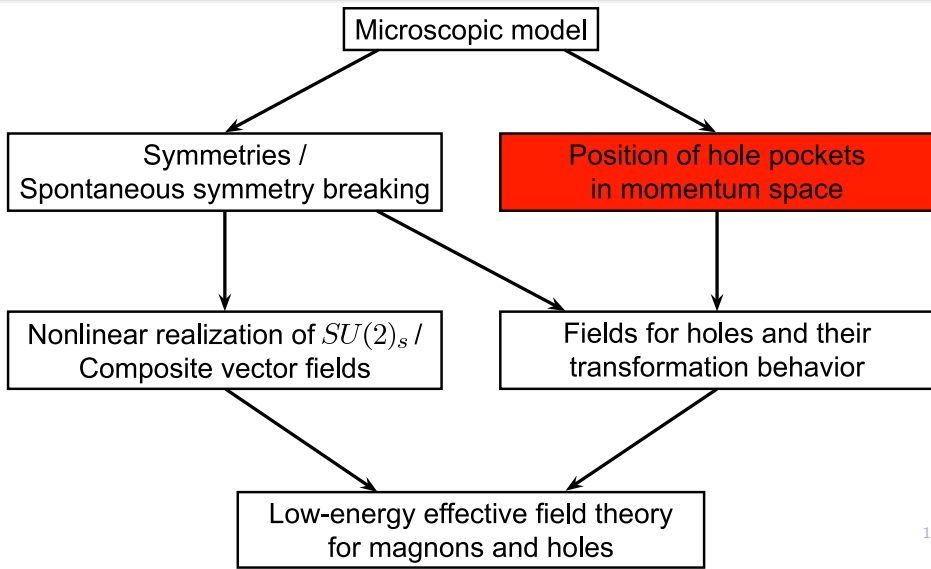
- Components used to couple magnons to holes
- Under global $SU(2)_s$ the components transform as

$$v_\mu^3(x)' = v_\mu^3(x) - \partial_\mu \alpha(x), \quad v_\mu^\pm(x)' = v_\mu^\pm(x) \exp(\pm 2i\alpha(x))$$

$v_\mu^3(x)$: Abelian gauge field

$v_\mu^\pm(x)$: Vector field (“charged” under $U(1)_s$)

Symmetry-based construction of effective theory



Hole pockets \iff Effective fields for holes

Where in momentum space do doped holes reside?

\implies Angle resolved photoemission spectroscopy (ARPES)

\implies Numerical simulations of single hole in AF

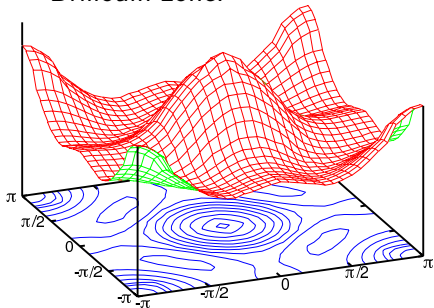
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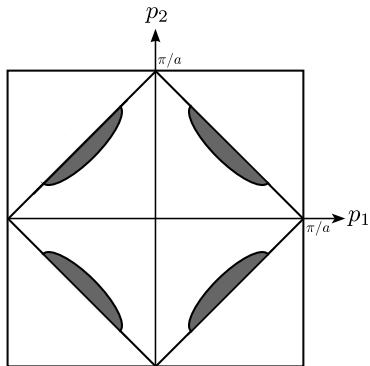
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Single hole (away from half-filling) dispersion relation in the first Brillouin zone:

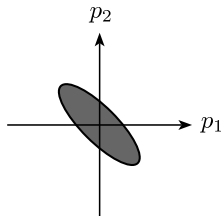


Minima at lattice momenta $\vec{k} = \pm\left(\frac{\pi}{2a}, \pm\frac{\pi}{2a}\right)$

Hole pockets \iff Effective fields for holes



Microscopic theory/ ARPES



Effective theory

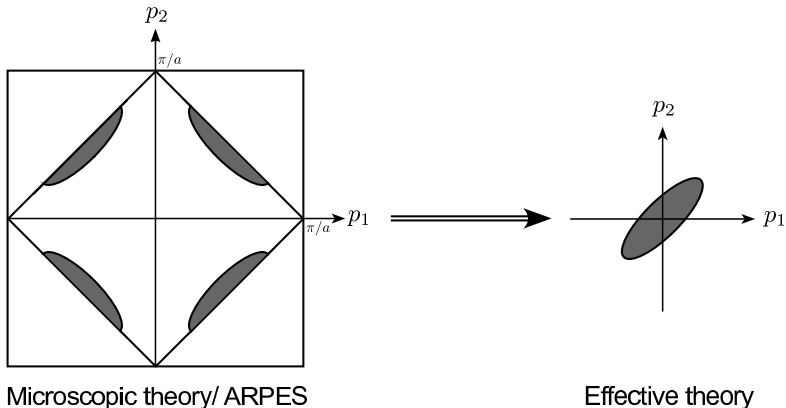
Symmetry considerations: Two **half-pockets** combine to a **full pocket**

In the effective theory:

p_2

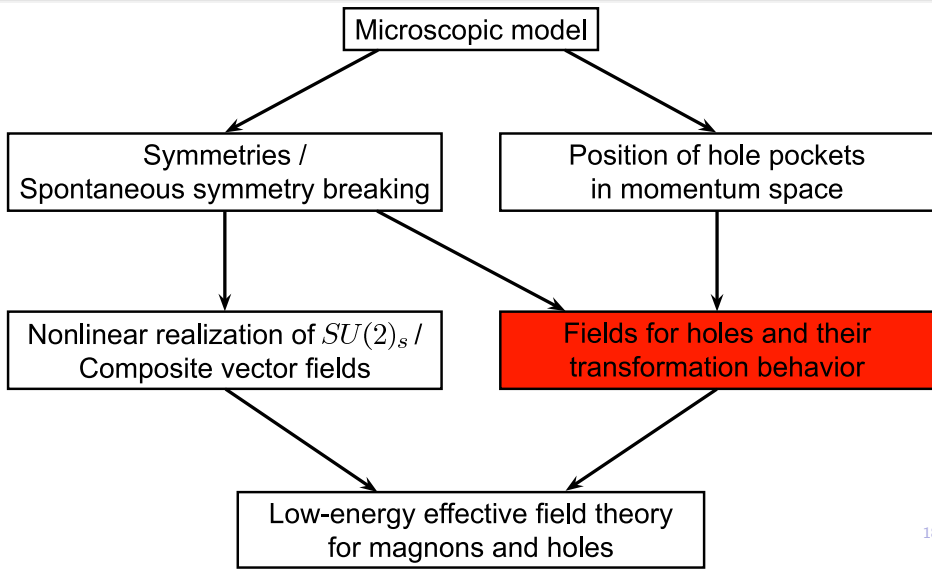
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Transformation behavior of hole fields

- The **symmetry properties** of the underlying system have to be **inherited by the effective theory!**

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$$U(1)_Q : \quad Q\psi_{\pm}^f(x) = \exp(i\omega)\psi_{\pm}^f(x)$$

$$D_i : \quad D_i\psi_{\pm}^f(x) = \mp \exp(ik_i^f a) \exp(\mp i\varphi(x))\psi_{\mp}^f(x)$$

$$O : \quad O\psi_{\pm}^{\alpha}(x) = \mp \psi_{\pm}^{\beta}(Ox) \quad O\psi_{\pm}^{\beta}(x) = \psi_{\pm}^{\alpha}(Ox)$$

$$R : \quad R\psi_{\pm}^{\alpha}(x) = \psi_{\pm}^{\beta}(Rx) \quad R\psi_{\pm}^{\beta}(x) = \psi_{\pm}^{\alpha}(Rx)$$

$$T : \quad T\psi_{\pm}^f(x) = \mp \exp(\mp i\varphi(Tx))\psi_{\pm}^{f\dagger}(Tx)$$

$$T\psi_{\pm}^{f\dagger}(x) = \pm \exp(\pm i\varphi(Tx))\psi_{\pm}^f(Tx)$$

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- In a **systematic derivative expansion**: Construct the most general Lagrangian which respects all symmetries

Effective Lagrangian for magnons and holes

Lagrangian at leading order: [Brügger et al. PRB 74 \(2006\) 224432](#)

$$\begin{aligned}
 \mathcal{L} = & \frac{\rho_s}{2} (\partial_i \vec{e} \cdot \partial_i \vec{e} + \frac{1}{c^2} \partial_t \vec{e} \cdot \partial_t \vec{e}) \\
 & + \sum_{\substack{f=\alpha,\beta \\ s=+,-}} \left[M \psi_s^{f\dagger} \psi_s^f + \psi_s^{f\dagger} D_t \psi_s^f \right. \\
 & + \frac{1}{2M'} D_i \psi_s^{f\dagger} D_i \psi_s^f + \sigma_f \frac{1}{2M''} (D_1 \psi_s^{f\dagger} D_2 \psi_s^f + D_2 \psi_s^{f\dagger} D_1 \psi_s^f) \\
 & + \Lambda (\psi_s^{f\dagger} v_1^s \psi_{-s}^f + \sigma_f \psi_s^{f\dagger} v_2^s \psi_{-s}^f) \\
 & \left. + N_1 \psi_s^{f\dagger} v_i^s v_i^{-s} \psi_s^f + \sigma_f N_2 (\psi_s^{f\dagger} v_1^s v_2^{-s} \psi_s^f + \psi_s^{f\dagger} v_2^s v_1^{-s} \psi_s^f) \right]
 \end{aligned}$$

with

$$D_\mu \psi_\pm^f(x) = [\partial_\mu \pm i v_\mu^3(x)] \psi_\pm^f(x)$$

$$\sigma_\alpha = +1, \quad \sigma_\beta = -1$$

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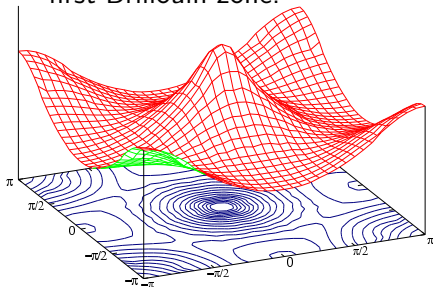
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Electron pockets \iff Effective fields for electrons

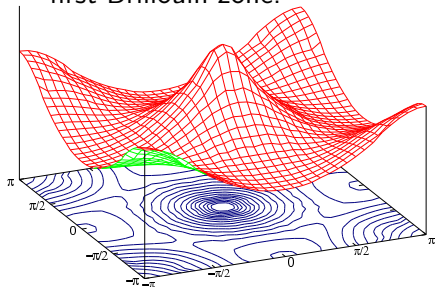
Single electron (away from half-filling) dispersion relation in the first Brillouin zone:



Minima at lattice momenta $\vec{k} = (\frac{\pi}{a}, 0)$ and $\vec{k} = (0, \frac{\pi}{a})$

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Symmetry considerations: **one single** electron pocket!

\implies Fields for electrons: $\psi_{\pm}(x)$ $\psi_{\pm}^{\dagger}(x)$

Effective Lagrangian for magnons and electrons

Lagrangian at leading order:

$$\begin{aligned} \mathcal{L} = & \frac{\rho_s}{2} (\partial_i \vec{e} \cdot \partial_i \vec{e} + \frac{1}{c^2} \partial_t \vec{e} \cdot \partial_t \vec{e}) \\ & + \sum_{s=+,-} \left[M \psi_s^\dagger \psi_s + \psi_s^\dagger D_t \psi_s + \frac{1}{2M'} D_i \psi_s^\dagger D_i \psi_s + N \psi_s^\dagger v_i^s v_i^{-s} \psi_s \right. \\ & \left. + iK (D_1 \psi_s^\dagger v_1^s \psi_{-s} - \psi_s^\dagger v_1^s D_1 \psi_{-s} - D_2 \psi_s^\dagger v_2^s \psi_{-s} + \psi_s^\dagger v_2^s D_2 \psi_{-s}) \right] \end{aligned}$$

with

$$D_\mu \psi_\pm(x) = [\partial_\mu \pm i v_\mu^3(x)] \psi_\pm(x)$$

Brügger, Hofmann, Kämpfer, Moser, Pepe, and Wiese, PRB 75
 (2007) 214405

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$$+ \sum_{s=+,-} \left[M \psi_s^\dagger \psi_s + \psi_s^\dagger D_t \psi_s + \frac{1}{2M'} D_i \psi_s^\dagger D_i \psi_s + N \psi_s^\dagger v_i^s v_i^{-s} \psi_s \right.$$

$$\left. + iK (D_1 \psi_s^\dagger v_1^s \psi_{-s} - \psi_s^\dagger v_1^s D_1 \psi_{-s} - D_2 \psi_s^\dagger v_2^s \psi_{-s} + \psi_s^\dagger v_2^s D_2 \psi_{-s}) \right]$$

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Effective Lagrangian for magnons and electrons

Lagrangian at leading order:

$$\begin{aligned} \mathcal{L} = & \frac{\rho_s}{2} (\partial_i \vec{e} \cdot \partial_i \vec{e} + \frac{1}{c^2} \partial_t \vec{e} \cdot \partial_t \vec{e}) \\ & + \sum_{s=+,-} \left[M \psi_s^\dagger \psi_s + \psi_s^\dagger D_t \psi_s + \frac{1}{2M'} D_i \psi_s^\dagger D_i \psi_s + N \psi_s^\dagger v_i^s v_i^{-s} \psi_s \right. \\ & \left. + iK (D_1 \psi_s^\dagger v_1^s \psi_{-s} - \psi_s^\dagger v_1^s D_1 \psi_{-s} - D_2 \psi_s^\dagger v_2^s \psi_{-s} + \psi_s^\dagger v_2^s D_2 \psi_{-s}) \right] \end{aligned}$$

with

$$D_\mu \psi_\pm(x) = [\partial_\mu \pm i v_\mu^3(x)] \psi_\pm(x)$$

Additional derivatives! \implies One-magnon exchange between electrons is weaker than between holes

Brügger, Hofmann, Kämpfer, Moser, Pepe, and Wiese, PRB 75
 (2007) 214405

Comparison of Holes and Electrons

Holes:

- Carry flavor
- Elliptical pockets in BZ
- Interact with background at $O(p)$

Electrons:

- No flavor
- Circular pockets in BZ
- Interact with background at $O(p^2)$

Outline

- 1 Motivation
 - Condensed matter analog of baryon chiral perturbation theory
 - Effective field theory for magnons
- 2 Hole-Doping
 - Construction of effective field theory
- 3 Electron-Doping
 - Construction of effective field theory
- 4 **Spiral Phases**
 - Uniform Background Field
 - Dispersion Relations of Indoped Fermions
 - Spiral Phases for Holes
 - Homogeneous Phase for Electrons
- 5 Conclusions

Assumptions

To describe the antiferromagnet with finite doping, we assume

- Fermions are indoped homogeneously
- The magnetic background does not vary in time: $v_t = 0$
- Fermion contact interactions are small

- The homogeneous doping of fermions requires a homogeneous magnetic background.

$\Rightarrow v_i = \text{const.}$ up to a $U(1)_s$ “gauge” transformation:

$$v_i^3(x)' = v_i^3(x) - \partial_i \alpha(x) = \sin^2 \frac{\theta(x)}{2} \partial_i \varphi(x) - \partial_i \alpha(x) = c_i^3,$$

$$v_i^\pm(x)' = v_i^\pm(x) \exp(\pm 2i\alpha(x))$$

$$= \frac{1}{2} [\sin \theta(x) \partial_i \varphi(x) \pm i \partial_i \theta(x)] \exp(\mp i(\varphi(x) - 2\alpha(x)))$$

$$= c_i^\pm$$

Theorem

The staggered magnetization $\vec{e}(x)$ configuration formed for uniform background fields c_i, c_i^3 is either homogeneous or a spiral

Brügger, Hofmann, Kämpfer, Pepe, and Wiese PRB 75
 (2007) 014421

Hamiltonian Formulation I

Hole Hamiltonian:

$$H^f = \begin{pmatrix} M + \frac{(p_i - c_i^3)^2}{2M'} + \sigma_f \frac{(p_1 - c_1^3)(p_2 - c_2^3)}{M''} & \Lambda(c_1 + \sigma_f c_2) \\ \Lambda(c_1 + \sigma_f c_2) & M + \frac{(p_i + c_i^3)^2}{2M'} + \sigma_f \frac{(p_1 + c_1^3)(p_2 + c_2^3)}{M''} \end{pmatrix}$$

Electron Hamiltonian:

$$H = \begin{pmatrix} M + \frac{(p_i - c_i^3)^2}{2M'} + Nc_i c_i & 2K(-p_1 c_1 + p_2 c_2) \\ 2K(-p_1 c_1 + p_2 c_2) & M + \frac{(p_i + c_i^3)^2}{2M'} + Nc_i c_i \end{pmatrix}$$

Acting on the 2d spin space $\begin{pmatrix} \psi_+(x) \\ \psi_-(x) \end{pmatrix}$.

Hamiltonian Formulation II

Energy of the indoped fermions through diagonalization of the single-particle Hamiltonians

Hole Energy

$$E_{\pm}^f(\vec{p}) = M + \frac{p_i^2}{2M'} + \sigma_f \frac{p_1 p_2}{M''} \pm \Lambda |c_1 + \sigma_f c_2|$$

Electron Energy

$$E_{\pm}(\vec{p}) = M + \frac{p_i^2}{2M'} + N c_i c_i \pm 2K |p_1 c_1 - p_2 c_2|$$

- \pm now refers to upper and lower energy states.
- Minimizing the energies of the fermions leads to $c_i^3 = 0$

Identification of Parameters

$c_i^3 = 0$ has important consequences. It can be shown, that

- $\theta(x) = \frac{\pi}{2}$
 \Rightarrow The spiral plane lies in the plane CuO plane
- $\varphi(x) = 2c_i x_i$
- $k = 2\sqrt{c_1^2 + c_2^2} = \text{const.}$
 k is the spiral pitch

Background field contribution to the total energy density:

$$\epsilon = 2\rho_s v_i^+ v_i^- = 2\rho_s (c_1^2 + c_2^2)$$

Density of Indoped Holes - Total Energy Density

- Density per flavor:

$$n_{\pm}^f = \frac{1}{(2\pi)^2} \int_{P_{\pm}^f} d^2p = \frac{1}{2\pi} M_{\text{eff}} T_{\pm}^f, \quad M_{\text{eff}} = \frac{M' M''}{\sqrt{M''^2 - M'^2}}$$

- Kinetic energy density:

$$t_{\pm}^f = \frac{1}{(2\pi)^2} \int_{P_{\pm}^f} d^2p \left(\frac{p_i^2}{2M'} + \sigma_f \frac{p_1 p_2}{M''} \right) = \frac{1}{4\pi} M_{\text{eff}} T_{\pm}^f{}^2$$

- Total energy density:

$$\epsilon_h = \sum_{\substack{f=\alpha,\beta \\ s=+,-}} \left[(M + s\Lambda |c_1 + \sigma_f c_2|) n_s^f + t_s^f \right]$$

Minimize $\epsilon_h - \lambda n$: Minimizing the fermion energy density with fixed hole density. λ is a Lagrange multiplier.

Phases of Hole-Doped Antiferromagnets

4 Pocket Phase:

Fact

Bounded from below for $2\pi\rho_s > M_{\text{eff}}\Lambda^2$

$$c_i = 0 \quad \epsilon_4 = \epsilon_0 + Mn + \frac{\pi n^2}{4M_{\text{eff}}}$$

3 Pocket Phase:

Fact

Bounded from below for $2\pi\rho_s > M_{\text{eff}}\Lambda^2$

$$|c_i| = \frac{\pi}{2} \frac{\Lambda n}{6\pi\rho_s - M_{\text{eff}}\Lambda^2} \quad \epsilon_3 = \epsilon_0 + Mn + \frac{\pi}{3M_{\text{eff}}} \left(1 - \frac{1}{2} \frac{M_{\text{eff}}\Lambda^2}{6\pi\rho_s - M_{\text{eff}}\Lambda^2} \right) n^2$$

Phases of Hole-Doped Antiferromagnets

2 Pocket Phase:

Fact

Bounded from below for $2\pi\rho_s > \frac{1}{2}M_{\text{eff}}\Lambda^2$

$$|c_{1,2}| = \frac{\Lambda}{4\rho_s}n, \quad |c_{2,1}| = 0, \quad \epsilon_2 = \epsilon_0 + Mn + \left(\frac{\pi}{2M_{\text{eff}}} - \frac{\Lambda^2}{8\rho_s} \right) n^2.$$

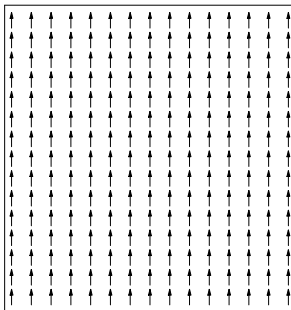
1 Pocket Phase:

Fact

Always bounded from below but unstable against forming inhomogeneities for $2\pi\rho_s < \frac{1}{2}M_{\text{eff}}\Lambda^2$

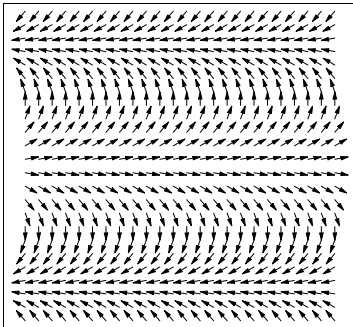
$$|c_1| = |c_2| = \frac{\Lambda}{4\rho_s}n \quad \epsilon_1 = \epsilon_0 + Mn + \left(\frac{\pi}{M_{\text{eff}}} - \frac{\Lambda^2}{4\rho_s} \right) n^2$$

Phases of Hole-Doped Antiferromagnets

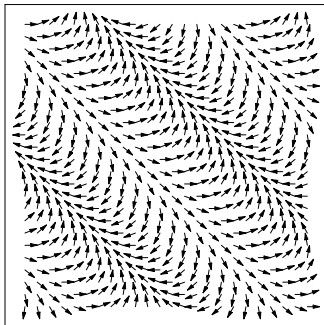


Homogeneous Phase
4 Hole Pockets

Phases of Hole-Doped Antiferromagnets

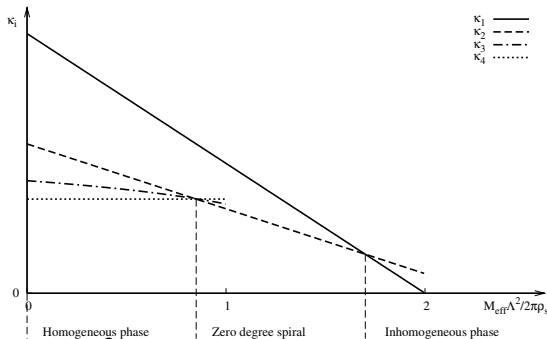


0 Degree Spiral
2 Hole Pockets



45 Degrees Spiral
3 (or 1) Hole Pockets

Stability of Phases for Hole Doping



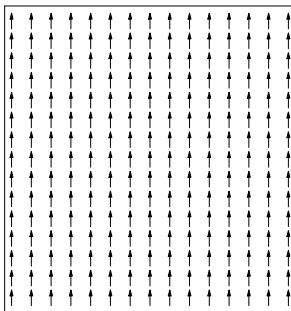
$$\kappa_1 = \left(\frac{\pi}{M_{\text{eff}}} - \frac{\Lambda^2}{\rho_s} \right)$$

$$\kappa_2 = \left(\frac{\pi}{2M_{\text{eff}}} - \frac{\Lambda^2}{8\rho_s} \right)$$

$$\kappa_3 = \frac{\pi}{3M_{\text{eff}}} \left(1 - \frac{1}{2} \frac{M_{\text{eff}}\Lambda^2}{6\pi\rho_s - M_{\text{eff}}\Lambda^2} \right)$$

$$\kappa_4 = \frac{\pi}{4M_{\text{eff}}}$$

Homogeneous Phase for Electron-Doped Antiferromagnets



Homogeneous Phase
No Spiral Phases

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Conclusions

- High-T superconductors: Condensed matter analog of baryon chiral perturbation theory
- We have constructed a systematic low-energy effective field theory for lightly doped antiferromagnets
- Using the effective theory we have investigated spiral phases in hole- and electron-doped cuprates
- While spiral phases do exist for hole-doping, they are absent in electron-doped cuprates
- We also calculated the one-magnon-exchange potential and investigated the possibility of hole-hole and electron-electron bound states

Outlook

- Analysis of materials with other lattice geometries:
Honeycomb and Triangular lattices
- Incorporation of Phonons as low-energy degrees of freedom
- Systematic treatment of loop graphs
- Towards the mysterious Mechanism of high-T
superconductivity

Nonlinear Realization of $SU(2)_S$ on the fermions

$$u(x)' = h(x)u(x)g^\dagger \quad C'_x = gC_x$$

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$$\Psi^X(x)' = h(x)u(x)C_x = h(x)\Psi^X(x)$$

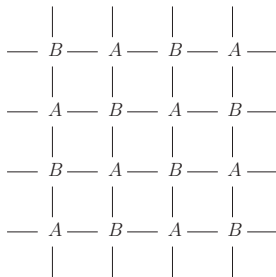
$$\Psi^X(x) = \begin{pmatrix} \psi_+^X(x) & \psi_-^{X\dagger}(x) \\ \psi_-^X(x) & -\psi_+^{X\dagger}(x) \end{pmatrix}, \quad x \in \text{even sublattice}$$

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The global spin rotation symmetry is also realized locally on the fermions.

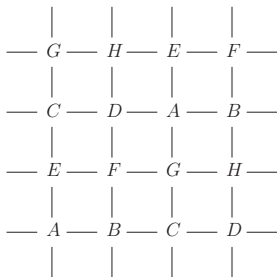
Sublattice Structure 1

$$k = (k_1, k_2) \in \left\{ (0, 0); \left(\frac{\pi}{a}, \frac{\pi}{a} \right) \right\}$$



Sublattice Structure 2

$$k = (k_1, k_2) \in \left\{ (0, 0), \left(\frac{\pi i}{a}, \frac{\pi}{a} \right), \left(\frac{\pi}{a}, 0 \right), \left(0, \frac{\pi}{a} \right), \left(\pm \frac{\pi}{2a}, \pm \frac{\pi}{2a} \right) \right\}$$



From Charge Carriers to Grassmann numbers

From Charge Carriers to Grassmann numbers

Hubbard model in manifestly $SU(2)_{\vec{Q}}$ invariant form (at half filling)

$$\mathcal{H} = -\frac{t}{2} \sum_{x,i} \text{Tr}[C_x^\dagger C_{x+i} + C_{x+i}^\dagger C_x] + \frac{U}{12} \sum_x \text{Tr}[C_x^\dagger C_x C_x^\dagger C_x].$$

$$C_x = \begin{pmatrix} c_{x\uparrow} & (-1)^x c_{x\downarrow}^\dagger \\ c_{x\downarrow} & -(-1)^x c_{x\uparrow}^\dagger \end{pmatrix}$$

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$$\Psi^X(x) = u(x) C_x \quad X := \text{sublatticeindex}$$

Solution for $\vec{e}(x)$ for constant background fields

$$\vec{e}(x) = (\sin\theta(x)\cos\varphi(x), \sin\theta(x)\sin\varphi(x), \cos\theta(x))$$

$$\cos\theta(x) = \frac{1}{\sqrt{1 + \left(\frac{c_i}{c_i^3}\right)^2}} \left[\cos\eta + \frac{c_i}{c_i^3} \sin\eta \cos\left(2\sqrt{1 + \left(\frac{c_i}{c_i^3}\right)^2} c_i^3 x_i\right) \right].$$

$$\varphi(x) = \text{atan} \left(\frac{\frac{c_i}{c_i^3} \sin\left(2\sqrt{1 + \left(\frac{c_i}{c_i^3}\right)^2} c_i^3 x_i\right)}{\sin\eta - \frac{c_i}{c_i^3} \cos\eta \cos\left(2\sqrt{1 + \left(\frac{c_i}{c_i^3}\right)^2} c_i^3 x_i\right)} \right).$$

Antiferromagnetism: Hubbard model

Four possible states at each lattice site: $|0\rangle$, $|\uparrow\rangle$, $|\downarrow\rangle$, $|\uparrow\downarrow\rangle$

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Assumptions:

- Half-filling: **in average one fermion per lattice site**
- Strong coupling limit: $U \gg t$

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How does the ground state order?

- By hopping system can lower its energy
- Hopping only possible for antiparallel spins

\implies **Antiferromagnetic** spin alignment is favored!

Relating microscopic operators to effective fields I

With the matrix-valued operator

$$C_x = \begin{pmatrix} c_{x\uparrow} & (-1)^x c_{x\downarrow}^\dagger \\ c_{x\downarrow} & -(-1)^x c_{x\uparrow}^\dagger \end{pmatrix}$$

the Hubbard Hamiltonian can be written

$$H = -\frac{t}{2} \sum_{x,i} \text{Tr}[C_x^\dagger C_{x+\hat{i}} + C_{x+\hat{i}}^\dagger C_x] + \frac{U}{12} \sum_x \text{Tr}[C_x^\dagger C_x C_x^\dagger C_x] \\ - \frac{\mu}{2} \sum_x \text{Tr}[C_x^\dagger C_x \sigma_3]$$

Relating microscopic operators to effective fields II

- Defining new lattice operators with the help of the diagonalizing matrix $u(x)$:

$$\psi_x^{A,B,\dots,H} = u(x)C_x, \quad x \in A, B, \dots, H$$

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- Replace lattice operators by effective Grassmann fields

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- Replace lattice operators by effective Grassmann fields

$$\psi_x^{A,B,\dots,H} \longrightarrow \Psi^{A,B,\dots,H}(x)$$

- Postulate: Transformation properties are inherited!

Accidental Galilean boost invariance

$$\begin{aligned}
 G : \quad G P(x) &= P(Gx), \quad Gx = (\vec{x} - \vec{v} t, t), \\
 G \psi_{\pm}^f(x) &= \exp(\vec{p}^f \cdot \vec{x} - \omega^f t) \psi_{\pm}^f(Gx), \\
 G \psi_{\pm}^{f\dagger}(x) &= \psi_{\pm}^{f\dagger}(Gx) \exp(-\vec{p}^f \cdot \vec{x} + \omega^f t),
 \end{aligned}$$

with $\vec{p}^f = (p_1^f, p_2^f)$ and ω^f given by

$$p_1^f = \frac{M'}{1 - (M'/M'')^2} \left[v_1 - \sigma_f \frac{M'}{M''} v_2 \right],$$

$$p_2^f = \frac{M'}{1 - (M'/M'')^2} \left[v_2 - \sigma_f \frac{M'}{M''} v_1 \right],$$

$$\omega^f = \frac{p_i^{f2}}{2M'} + \sigma_f \frac{p_1^f p_2^f}{M''} = \frac{M'}{1 - (M'/M'')^2} \left[\frac{1}{2}(v_1^2 + v_2^2) - \sigma_f \frac{M'}{M''} v_1 v_2 \right]$$

Transformation behavior of electron fields

$$SU(2)_s : \quad \psi_{\pm}(x)' = \exp(\pm i\alpha(x))\psi_{\pm}(x)$$

$$U(1)_Q : \quad Q\psi_{\pm}(x) = \exp(i\omega)\psi_{\pm}(x)$$

$$D_i : \quad D_i\psi_{\pm}(x) = \mp \exp(ik_i a) \exp(\mp i\varphi(x))\psi_{\mp}(x)$$

$$O : \quad O\psi_{\pm}(x) = \pm\psi_{\pm}(Ox)$$

$$R : \quad R\psi_{\pm}(x) = \psi_{\pm}(Rx)$$

$$T : \quad T\psi_{\pm}(x) = \exp(\mp i\varphi(Tx))\psi_{\pm}^{\dagger}(Tx)$$

$$T\psi_{\pm}^{\dagger}(x) = -\exp(\pm i\varphi(Tx))\psi_{\pm}(Tx)$$