

Title: Electronic Liquid Crystalline Phases of Highly Correlated Electronic Systems

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Abstract: In one extreme, where the interactions are sufficiently weak compared to the interactions, electrons form a “Fermi liquid” the state that accounts for the properties of simple metals. In the other extreme, where the interactions are dominant, the electrons form various “Mott” insulating or “Wigner crystalline” phases, often characterized by broken spatial and/or magnetic symmetries. Corresponding charge and/or magnetically ordered insulating phases are common in nature. Between these two extremes lie highly correlated electronic fluids, and correspondingly a host of interesting and perplexing materials, including such diverse systems as the cuprate and iron-based high temperature superconductors, the failed metamagnet $\text{Sr}_3\text{Ru}_2\text{O}_7$, and a variety of quantum Hall fluids. Some insight into electron fluids in this rich intermediate coupling regime can be obtained from viewing them as partially melted electron solids, rather than as strongly interacting gases. Here, analogies with the liquid crystalline phases of complex classical fluids provide useful guidance for a new approach to this key problem in condensed matter physics.

Electronic liquid crystals

Waterloo - 2013

Strongly Interacting Electrons

$$r_s \sim V/K \sim h^{-2} n^{-1/3}$$

$$V \sim e^2 n^{1/3} \quad K \sim h^2 n^{2/3} / 2m$$

Strongly Interacting Electrons

$$r_s \sim V/K \sim \hbar^{-2} n^{-1/3}$$

$$\text{Quantum "temperature"} \sim 1/r_s$$

Strongly Interacting Electrons

$$r_s \sim V/K \sim \hbar^{-2} n^{-1/3}$$

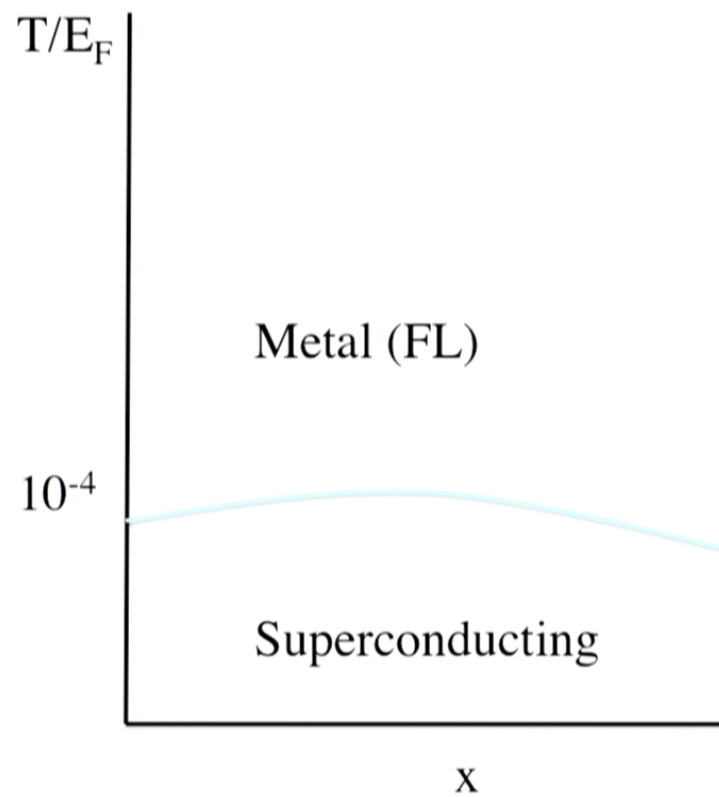
Quantum “temperature” $\sim 1/r_s$

$r_s =$ “small” Fermi Gas (simple metal)

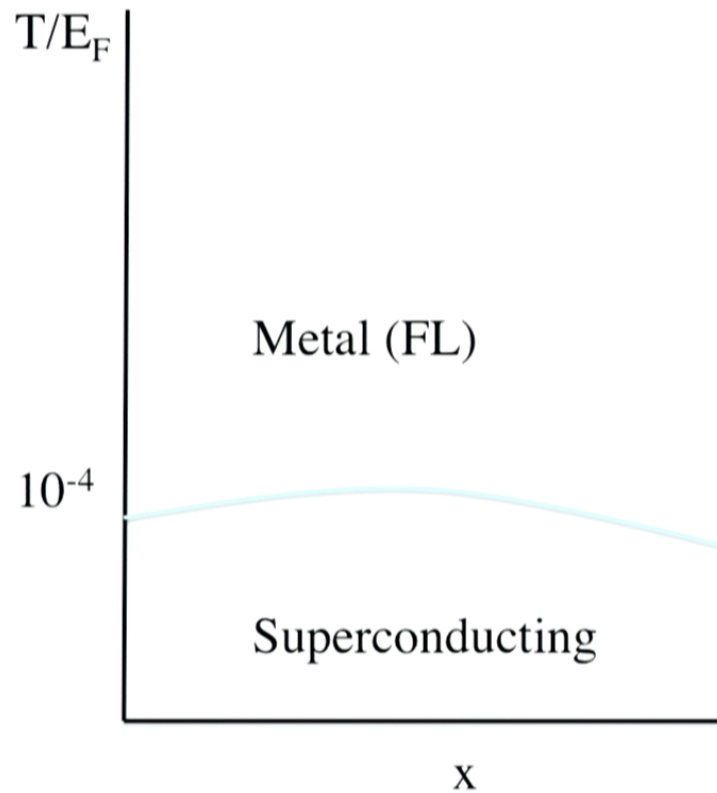
$r_s =$ “very large” Wigner-Mott Insulator

$r_s =$ “pretty darn large” Highly Correlated Fluid

Phase diagram of simple metals – $A_{1-x}B_x$ - with $r_s \sim 1$



Phase diagram of simple metals – $A_{1-x}B_x$ - with $r_s \sim 1$



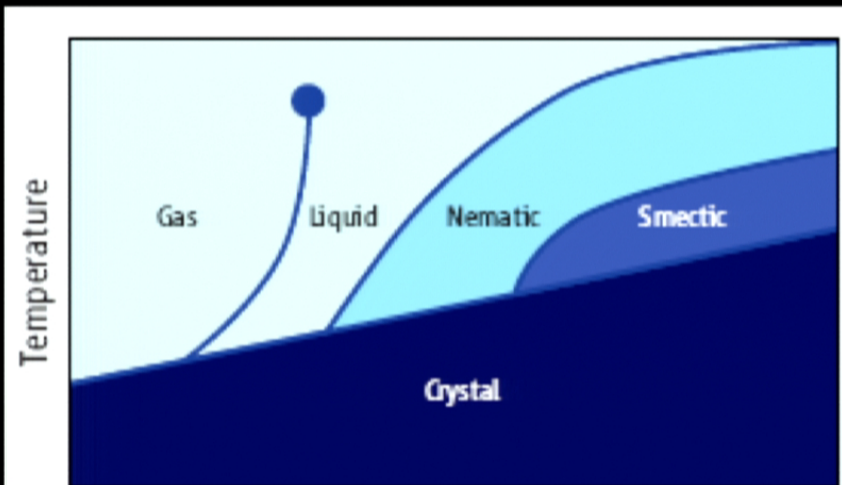
$r_s \sim 1$ is “natural”

The unnatural stability
of the Fermi liquid

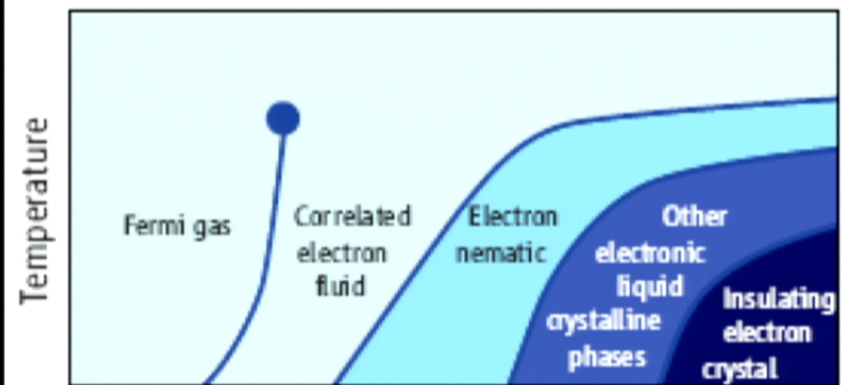
Anderson’s theorem

Traditional approach: View highly correlated fluid as a strongly interacting gas, figure out approximate ways of resumming perturbation series, and invoke the magic of adiabatic continuity (*i.e.* Fermi liquid theory).

Correlated
Fluids
Viewed as
Partially
Melted
Crystals



Another physical parameter



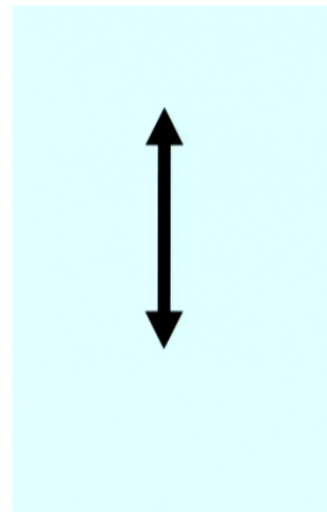
Strength of interactions

Classical Liquid Crystals: the broken symmetries

Molecular
liquid



Isotropic



Nematic

Traceless symmetric
tensor

(Headless vector)



Smectic

Unidirectional
“density-wave”

Electronic liquid crystalline phases are “unnatural” if we start from an interacting Fermi gas.

However, electron liquid crystalline phases seem reasonable when thought of as partially quantum melted electron crystalline phases.

Spin-Stripe Crystal in a Doped Antiferromagnet



Breaks τ , T_x , T_y , $R_{\pi/2}$, mirrors, (spin-rotation)
 Incommensurate or high order commensurate
 Bond-centered or site centered, diagonal or vertical

Some “practical” considerations:

1) **Crystals are imperfect** ... quenched disorder

In case of a broken symmetry, this can lead to the presence of “random fields:”

$$\Delta H = \int d\vec{r} \mathbf{h}(\vec{r}) \cdot \phi(\vec{r})$$

In 2d, and under many conditions in 3d, this spoils most macroscopic manifestations of broken symmetry, and “rounds” the phase transition.

Some “practical” considerations:

2) Often spatial symmetries are only approximate

In case of a broken symmetry, this leads to a “small” symmetry breaking field:

$$\Delta H = \int d\vec{r} \mathbf{h} \cdot \phi(\vec{r})$$

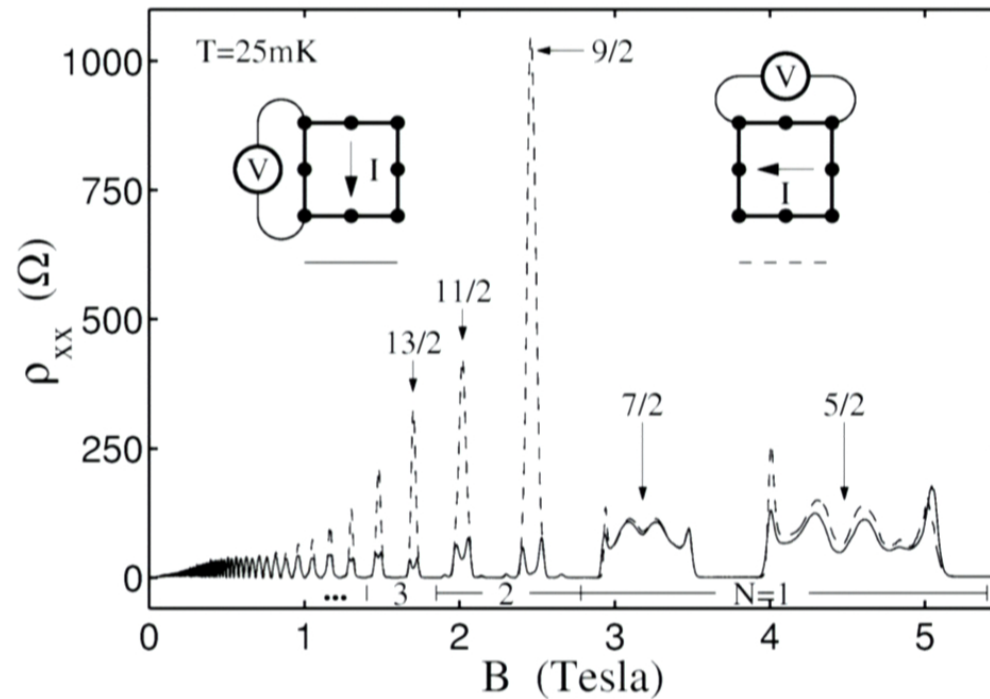
This spoils the sharp distinction between the broken symmetry state and the symmetric phase, and so “rounds” the phase transition

However, field cooling circumvents vexing non-equilibrium effects of domain formation.

Anisotropic State in Higher Landau Levels

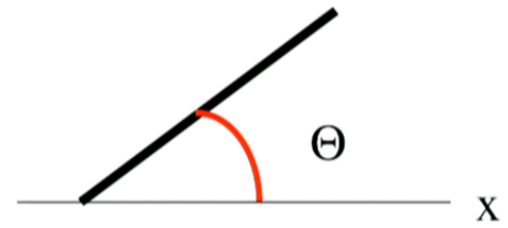
Eisenstein and co. PRL 1999 etc.

$$\nu = n_{el} \phi_0$$



Ultra high mobility 2DEG in GaAs-GaAlAs heterostructures

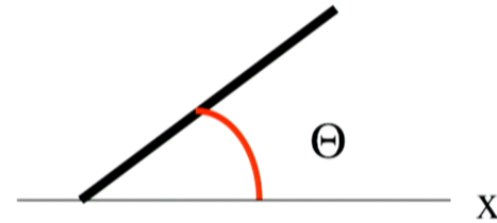
Symmetry considerations:



Symmetry considerations:

$$\text{Nematic } \Theta = \pi - \Theta$$

$$m = \langle \cos(2\Theta) \rangle$$

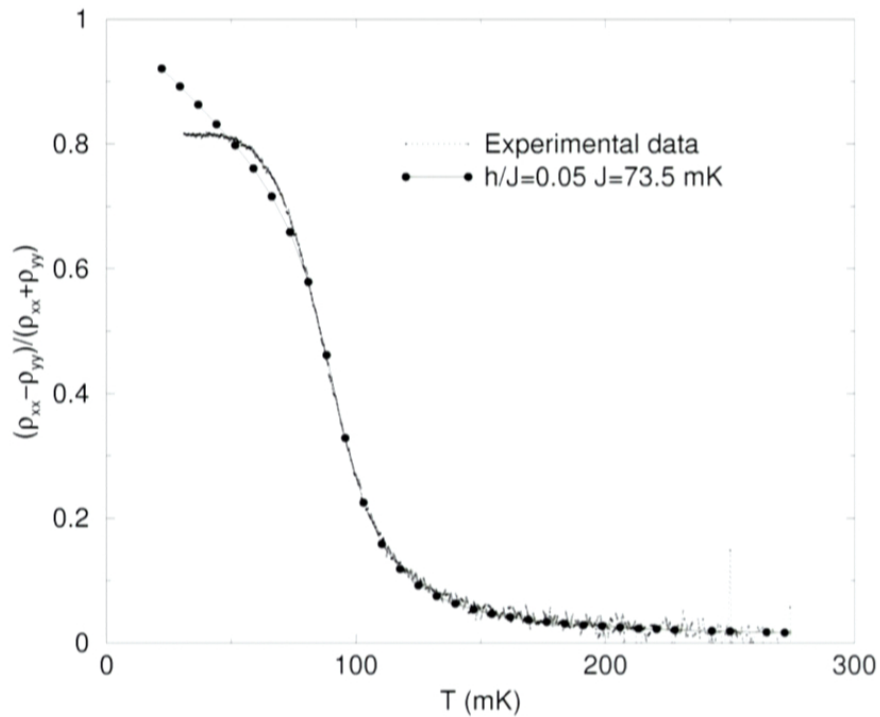


Within effective mass approximation -
full rotational symmetry.

$$\text{Crystal field effects in bulk } V_{\text{CF}} = V_4 \cos(4\Theta)$$

$$V_4 \sim \text{Ry}^* (a_{\text{B}}/a_{\text{B}}^*)^2 \sim 10^{-4} \text{Ry}^*$$

Comparison with classical XY model of Nematic transition

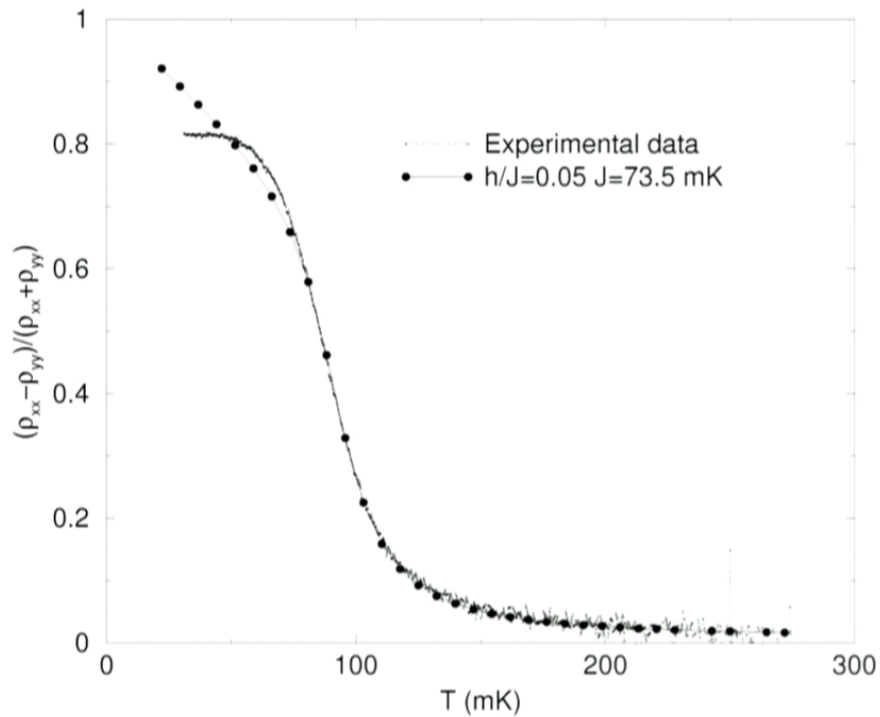


$$H = -J \cos(2\theta_i - 2\theta_j) - h \cos(2\theta_i)$$

$$\frac{\rho_{xx} - \rho_{yy}}{\rho_{xx} + \rho_{yy}} \sim \langle \cos(2\theta) \rangle$$

Fradkin and SAK (1999)

Comparison with classical XY model of Nematic transition



$$H = -J \cos(2\theta_i - 2\theta_j) - h \cos(2\theta_i)$$

$$\frac{\rho_{xx} - \rho_{yy}}{\rho_{xx} + \rho_{yy}} \sim \langle \cos(2\theta) \rangle$$

Fradkin and SAK (1999)

Quantum Hall Nematic:

Finite T phase transition to a conducting state
with a large resistivity anisotropy

Transition fits well to a KT transition in a small
symmetry breaking field

No depinning field or similar signatures of a CDW state
in non-linear I - V 's

Symmetries (available for breaking) of the square lattice electronic system

Point group symmetry including

Discrete rotations $R_{\pi/2}$, R_{π} ,

Reflection planes,...

Translations, T_x , T_y , ...

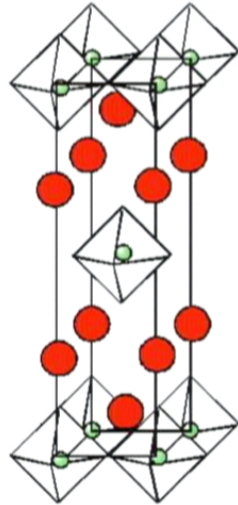
Time reversal, τ

Large scale gauge symmetry (number conservation)

An electron liquid is a conducting (or superconducting) state of the electrons in a solid.

Introduction: Materials

The Ruddlesden-Popper Series: $Sr_{1+n}Ru_nO_{1+3n}$

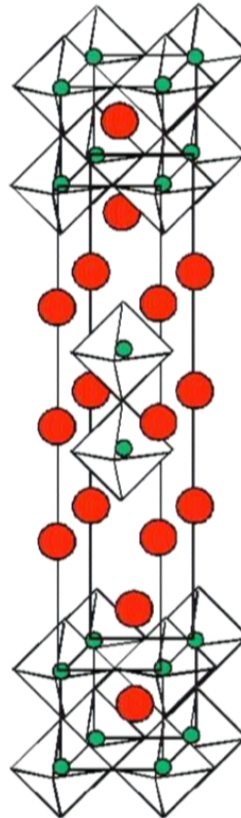


n=1

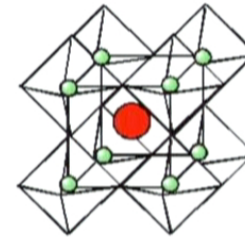
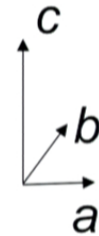
quasi-2d

Paramagnetic Fermi liquid

Triplet superconductor



n=2



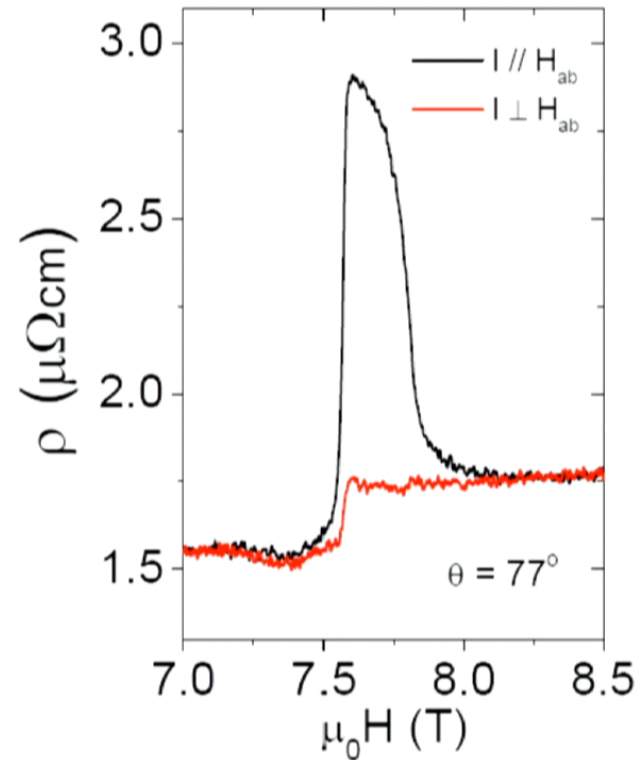
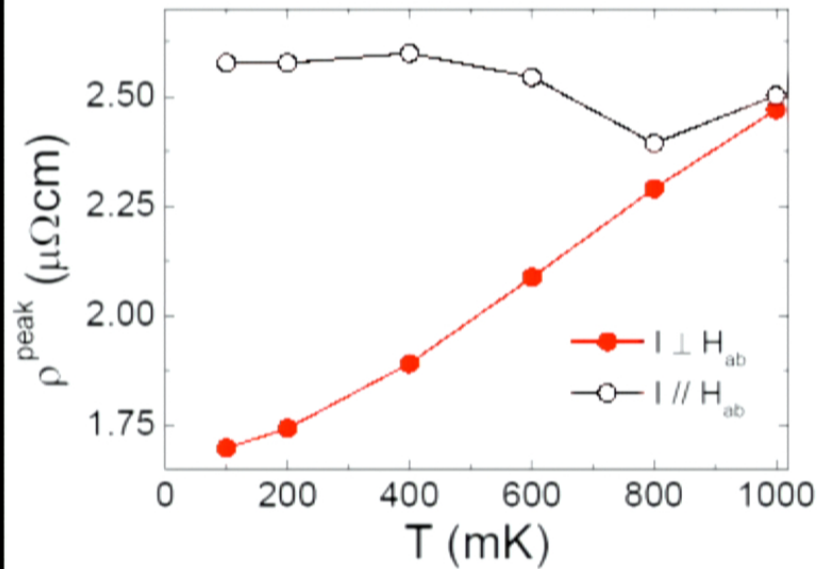
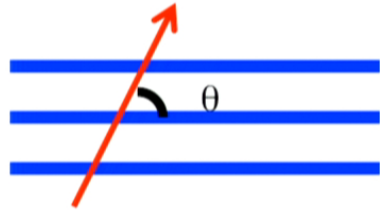
n=∞

3d

Ferromagnetic,

$T_c=160K$

Anisotropic resistivity: temperature dependence



No orthorhombicity has been detected in X-ray scattering

O(N) quantum rotor model: $\vec{S}_i \cdot \vec{S}_i = N$

$$H_{2D} = J_1 \sum_{\langle i,j \rangle_1} \vec{S}_i \cdot \vec{S}_j + J_2 \sum_{\langle i,j \rangle_2} \vec{S}_i \cdot \vec{S}_j - \frac{K}{N} \sum_{\langle i,j \rangle_1} [\vec{S}_i \cdot \vec{S}_j]^2$$

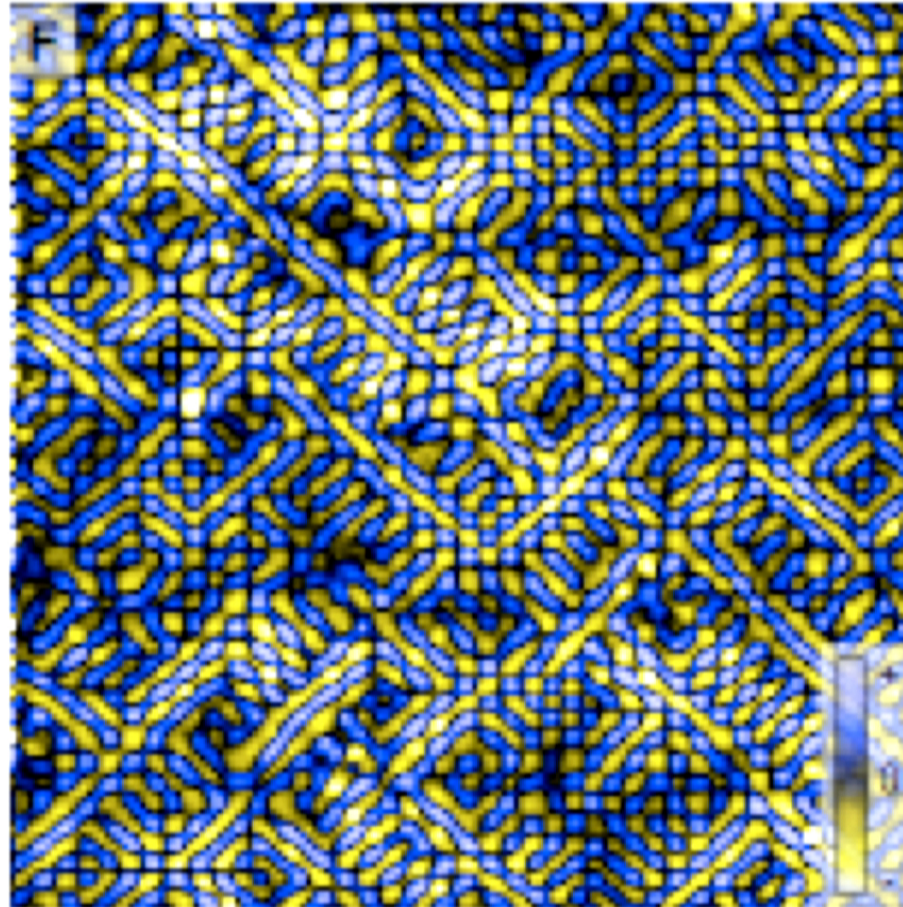
“Weak” interlayer coupling: $H_{\perp} = -J_{\perp} \sum_{\langle i,j \rangle_{\perp}} \vec{S}_i \cdot \vec{S}_j$

Experimental “proof” of the existence of electron nematic (and other electronic liquid crystalline phases) has been adduced primarily in certain, very clean, very intensely studied materials.

Disorder is a disaster, especially in quasi 2D:

Inevitably leads to a glassy (domain) state.

STM
image
of a
“Nematic
Glass”
phase on
the
surface of
BSCCO
with
 $T_c \sim 90\text{K}$



From Kohsaka *et al*, Nature (2007)
STM on BSCCO or NaCCOC

The End

Electronic liquid crystalline phases, with new and unexpected properties exist in a variety of materials

Both nematic and various forms of electronic smectic order appear prominently in the cuprate phase diagram