

Frankfurt, June 24, 2014

Exercises for Computational Methods in Solid State Theory
SS 2014

Exercise Set 6

(Due date: Monday, July 7, 2014)

Exercise 8 (Hartree-Fock mean field calculations) (10 points)

The physical properties of κ -(BEDT-TTF)₂X charge transfer salts can be described by a Hubbard model on the anisotropic triangular lattice which is given as

$$(1) \quad H = \sum_{\langle ij \rangle, \sigma} t_{ij} c_{i\sigma}^\dagger c_{j\sigma} + U \sum_i n_{i\uparrow} n_{i\downarrow} - \mu \sum_{i\sigma} c_{i\sigma}^\dagger c_{i\sigma},$$

where $c_{i\sigma}^\dagger$ ($c_{i\sigma}$) are creation (annihilation) operators for electrons of spin σ , $n_{i\sigma} = c_{i\sigma}^\dagger c_{i\sigma}$ is the density of σ spin electrons and t_{ij} is the hopping amplitude. We choose to study here the case where $t_{ij} = t$ in two of the directions of the triangular lattice, and $t_{ij} = t'$ in the third direction (see Figure 1).

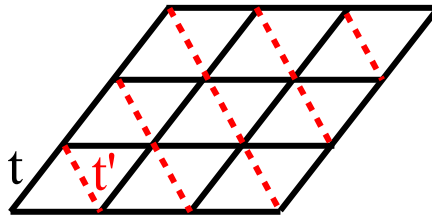


Figure 1: Anisotropic triangular lattice.

To study this model in the Hartree-Fock approximation we can consider two particular mean fields:

$$(2) \quad H^{\text{MF1}} = H_{\uparrow} + H_{\downarrow} + C,$$

with:

$$\begin{aligned} H_{\uparrow} &= -t \sum_{\langle i,j \rangle} c_{i,\uparrow}^{\dagger} c_{j,\uparrow} + U \sum_i n_{i,\uparrow} \langle n_{i,\downarrow} \rangle, \\ H_{\downarrow} &= -t \sum_{\langle i,j \rangle} c_{i,\downarrow}^{\dagger} c_{j,\downarrow} + U \sum_i \langle n_{i,\uparrow} \rangle n_{i,\downarrow}, \\ C &= -U \sum_i \langle n_{i,\uparrow} \rangle \langle n_{i,\downarrow} \rangle. \end{aligned}$$

and

$$(3) \quad H^{\text{MF2}} = C^{\dagger} \begin{pmatrix} H^{\uparrow\uparrow} & H^{\uparrow\downarrow} \\ H^{\downarrow\uparrow} & H^{\downarrow\downarrow} \end{pmatrix}_{2N \times 2N} C,$$

with:

$$\begin{aligned} C^{\dagger} &= (c_{1,\uparrow}^{\dagger}, \dots, c_{N,\uparrow}^{\dagger}, c_{1,\downarrow}^{\dagger}, \dots, c_{N,\downarrow}^{\dagger}) \\ H_{i,j}^{\uparrow\uparrow} &= (U \langle n_{i,\downarrow} \rangle - \mu) \delta_{i,j} + T_{i,j} \\ H_{i,j}^{\uparrow\downarrow} &= -U \langle S_i^- \rangle \delta_{i,j} \\ H_{i,j}^{\downarrow\uparrow} &= -U \langle S_i^+ \rangle \delta_{i,j} \\ H_{i,j}^{\downarrow\downarrow} &= (U \langle n_{i,\uparrow} \rangle - \mu) \delta_{i,j} + T_{i,j} \end{aligned}$$

$$\begin{aligned} H^{\text{MF2}} &= -t \sum_{\langle i,j \rangle, \sigma} c_{i,\sigma}^{\dagger} c_{j,\sigma} + U \sum_i [\langle n_{i,\downarrow} \rangle n_{i,\uparrow} + \langle n_{i,\uparrow} \rangle n_{i,\downarrow} - \langle n_{i,\uparrow} \rangle \langle n_{i,\downarrow} \rangle] \\ &\quad -U \sum_i \left[\langle S_i^+ \rangle c_{i,\downarrow}^{\dagger} c_{i,\uparrow} + \langle S_i^- \rangle c_{i,\uparrow}^{\dagger} c_{i,\downarrow} - \langle S_i^+ \rangle \langle S_i^- \rangle \right] \end{aligned}$$

and in this exercise we will use the particular case where all the spin are in the xz -plane: $\langle S_i^+ \rangle = \langle S_i^- \rangle = \langle S_i^x \rangle$

- Using MF1 compute: the ground state energy per site, the gap and the local magnetization $\langle S^z \rangle$ at $U/t = 1$ and $U/t = 16$ at $t'/t = 0, 0.5$ for a system of size 10×10 .
- Using MF2 compute: the ground state energy, the local magnetizations $\langle S^z \rangle$ and $\langle S^x \rangle$ at $U/t = 16$ at $t'/t = 0.5, 0.8, 1$ for a system of size 10×10 and 12×12 .
- How could you define/describe the phase corresponding to each of the cases of question a) and b)?
- Discuss the difference between results for the systems of size 10×10 and 12×12 .

The computation should be realized in real space and at half-filling for a total magnetization equal to zero.