

Phy220 Group theory final

①

Problem 1: Cubic anisotropy

Symmetry analysis can help to build up field theory for the system you want to study. Now let me aside to write down a free energy for a magnetic system, which is a functional of magnetic density $\vec{m}(\vec{r})$ denoted as $F[\vec{m}(\vec{r})]$. Let us

expand $F[\vec{m}(\vec{r})]$ in terms of polynomials, and only consider the homogenous distribution, i.e. $\vec{m}(\vec{r}) = \vec{m}$ which is a const over space. Then $F(m) = a_i m_i + a_{ij} m_i m_j + a_{ijk} m_i m_j m_k$

$$+ a_{ijkl} m_i m_j m_l m_k + \dots, \quad (i, j, k, l = x, y, z)$$

and we truncate at the quartic level.

① For an isotropic space, figure out the structures of the coefficients $a_i, a_{ij}, a_{ijk}, a_{ijkl}$.

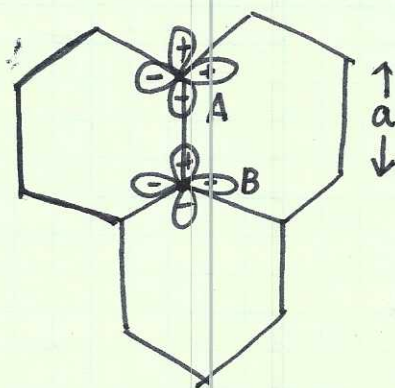
Hint, $F(m)$ needs to be invariant under all the symmetries of an isotropic space, and also possibly other symmetries due to physical requirement such as time-reversal symmetry.

(3)

② If the system does not have all the symmetric of a free space, but has the cubic symmetry O . Try to do ~~problem~~ part ① again.

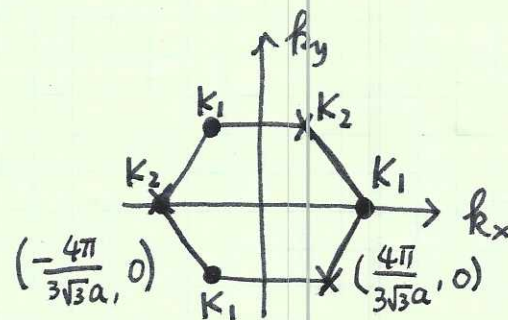
Problem 2: p-orbital band structure in the honeycomb lattice.

① One each site, there are a pair of degenerate P_x and P_y -orbitals. Each unit cell has a pair of sites A and B.



The Brillouine zone is plotted below.

~~The honeycomb lattice has the~~



① What's the group of wavevector K_1 ? You only need to figure out the ~~rotation~~ part of point group.

② There are four bases for the group of wavevector K_1 .

$$\psi_{A, P_{x,y}}(\vec{r}) = \frac{1}{\sqrt{N}} \sum_{\vec{r}_A^0} e^{i\vec{K}_1 \cdot \vec{r}} \phi_{P_{x,y}}(\vec{r} - \vec{r}_A^0)$$

$$\psi_{B, P_{x,y}}(\vec{r}) = \frac{1}{\sqrt{N}} \sum_{\vec{r}_B^0} e^{-i\vec{K}_1 \cdot \vec{r}} \phi_{P_{x,y}}(\vec{r} - \vec{r}_B^0)$$

③

Basing on this set of basis, decompose them into irreducible representations of the group of wavevector $\vec{G}_K$. Can you understand them in terms of angular momentum numbers? Can you compare with the case of graphene?

③ open part: for extra credit. You may further consider to add spin-orbit coupling, and try to figure out more physics.

Problem 3: Consider a ^{3D} diamond lattice. Let us use the convention that from each A site there are four bonds towards its four B-neighbours along (111) , $(\bar{1}\bar{1}1)$, $(\bar{1}1\bar{1})$ and $(1\bar{1}\bar{1})$. Figure out the elements of its space group.

Is it symmetric or non-symmetric?