

Condensed Matter Physics I
II intermediate test a.y. 2025/26 - 18 December 2025
(Time: 1:30 h)

Exercise 1: Bloch functions

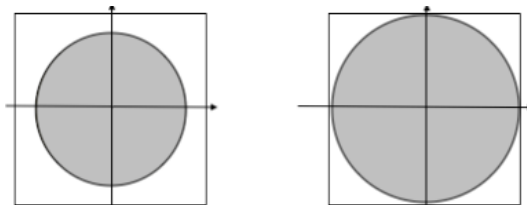
1. Does a Bloch function have the same periodicity of the Bravais lattice or not? Justify your answer. **No, since, upon translation of \mathbf{R} , it is multiplied by a phase $\exp(i\mathbf{k} \cdot \mathbf{R})$**
2. Does the spacial probability density for Bloch electrons have the periodicity of the corresponding Bravais lattice or not? Justify your answer. **Yes, since the multiplying phase has modulus=1**
3. What is $u_{n,\mathbf{k}}(\mathbf{r})$ in case of free electrons? **A constant, the choice depending on the normalization**

Exercise 2: Density of electronic states

A two-dimensional material has a "relativistic" dispersion relation $E(\mathbf{k}) = \alpha|\mathbf{k}|$. An example of such a system is graphene.

1. Calculate the velocity in the semiclassical model. **$\mathbf{v} = \frac{\alpha}{\hbar} \hat{\mathbf{k}}$**
2. Calculate the density of electronic states. **$g(E) = \frac{E}{\pi\alpha^2}$**

Exercise 3: Nearly free electrons



Consider a 2D square Bravais lattice of parameter a . The two figures represents the first Brillouin zone with Fermi circles ("empty" lattice model) corresponding to two different electron densities.

1. Focus on the left plot: estimate the (integer) number of electrons per unit cell. **$Z=1$**
2. Then consider the plot on the right and give the exact value of the number of electrons (not necessarily integer) per unit cell in this case. **$Z=\pi/2$**
3. If such number is between two integers, Z_A and Z_B , this means that it could represent the average electron number per cell of an alloy $A_{1-x}B_x$. Calculate x . **$x=0.57$ or $x=0.43$, according whether one considers $Z_A \leq Z_B$ or viceversa**
4. Consider a magnetic field perpendicular to the lattice. Which are the crystalline momentum velocities $\dot{\mathbf{k}}$ and the periods T of the orbit in the two cases? **$\dot{\mathbf{k}} = \omega_c(-k_F|_y, k_F|_x)$, $|\dot{\mathbf{k}}| = \omega_c k_F$ with $k_F = \sqrt{2\pi}/a$ for $Z = 1$ and π/a for $Z = 2$; period = $2\pi/\omega_c$ in any case**

5. Now consider to have a small perturbing potential, tight-binding like but small:

$$V(x, y) = 2V_1(\cos(2\pi x/a) + \cos(2\pi y/a)) + 4V_2 \cos(2\pi x/a)\cos(2\pi y/a)$$

Use the weak potential model to find the energy of the lowest electronic state at $\mathbf{k} = (\pi/a, 0)$. Writing sin and cos in terms of complex exponentials, it is evident that the Fourier components of the potential include only the following \mathbf{K} reciprocal lattice vectors: $\pm\mathbf{b}_1 = \pm 2\pi/a(1, 0)$, $\pm\mathbf{b}_2 = \pm 2\pi/a(0, 1)$, $\pm(\mathbf{b}_1 \pm \mathbf{b}_2)$. At $\mathbf{k} = (\pi/a, 0)$ only the component for \mathbf{b}_1 is involved, therefore $E_-(\mathbf{k} = (\pi/a, 0)) = E^0(\mathbf{k} = (\pi/a, 0)) - V_1 = \frac{\hbar^2}{2m} \left(\frac{\pi}{a}\right)^2 - V_1$