

Exercise 8.1 Bohr–van Leeuwen theorem

Prove the Bohr–van Leeuwen theorem, which states that there is no magnetism in classical physics.

Hint: $\mathcal{H}(\mathbf{p}_1, \dots, \mathbf{p}_N; \mathbf{q}_1, \dots, \mathbf{q}_N)$ is the Hamiltonian of the N -particle system with vanishing external magnetic field. In comparison, the Hamiltonian with applied magnetic field \mathbf{B} is then given by $\mathcal{H}(\mathbf{p}_1 - e/c\mathbf{A}_1, \dots, \mathbf{p}_N - e/c\mathbf{A}_N; \mathbf{q}_1, \dots, \mathbf{q}_N)$, where $\mathbf{B} = \nabla \times \mathbf{A}$ and $\mathbf{A}_i \equiv \mathbf{A}(\mathbf{q}_i)$. The (thermal average of the) magnetization can be calculated using

$$M = \left\langle -\frac{\partial \mathcal{H}}{\partial B} \right\rangle = \frac{1}{\beta} \frac{\partial \log Z}{\partial B}, \quad (1)$$

with the partition function Z of the system in the magnetic field.

Exercise 8.2 Landau Diamagnetism

Calculate the orbital part of the magnetization (e.g. ignore the Zeeman-term) of the free electron gas in 3D in the limits of low temperature and small external field ($T \rightarrow 0$, $B \rightarrow 0$). In addition, show that the magnetic susceptibility at $T = 0$ and $B = 0$ is given by

$$\chi = -\frac{1}{3} \frac{m^2}{m^{*2}} \chi_P, \quad (2)$$

where χ_P is the Pauli susceptibility.¹

Hint: Calculate the free energy,

$$F = N\mu - k_B T \sum_i \ln [1 + e^{-(\epsilon_i - \mu)/k_B T}], \quad (3)$$

at $T = 0$ to second order in B using the Euler-Maclaurin formula,

$$\sum_0^{n_0} f(n) \approx \int_{-1/2}^{n_0+1/2} f(n) dn - \frac{1}{24} [f'(n_0 + 1/2) - f'(-1/2)]. \quad (4)$$

Exercise 8.3 Landau Levels in Graphene

Graphene is defined as a single two-dimensional layer of graphite, the C -atoms are arranged on a two-dimensional honeycomb lattice (cf. exercise 5). The latter is not a Bravais-lattice, but a triangular lattice with a diatomic basis. Consequently, the reciprocal lattice, which is a honeycomb lattice as well, has two inequivalent points called K - and K' -points (see Fig. 1). The two atoms per unit cell create a valence and a conduction band which cross linearly in one point (called the Dirac point) at the K - and K' -points and form the so-called Dirac cones (see Fig. 1). In undoped graphene, the Fermi energy is exactly at the Dirac point.

¹The Pauli susceptibility $\chi_P = \mu_B^2 \rho(\epsilon_F)$ (at $T = 0$) is a consequence of the Zeeman energy-term in the Hamiltonian for particles with non-zero spin; $\mu_B = \frac{e\hbar}{2mc}$; ρ is the density of states (including the spin degeneracy).

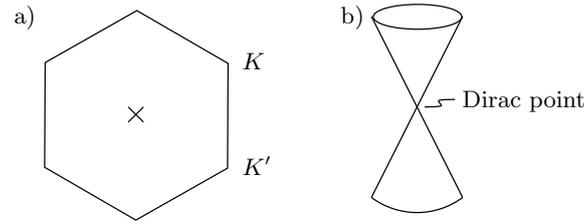


Figure 1: a) First Brillouin zone of graphene with K - and K' -points. b) Band structure of graphene at the K - and K' -points: the Dirac cones.

To a good approximation, the spectrum in graphene is linear at the Fermi energy and described by the Hamiltonian²

$$\hat{\mathcal{H}}_{\text{eff}} = v_F \begin{pmatrix} \hat{p}_x \sigma_x + \hat{p}_y \sigma_y & 0 \\ 0 & \hat{p}_x \sigma_x - \hat{p}_y \sigma_y \end{pmatrix}, \quad (5)$$

where the Pauli matrices σ act on a pseudo-spin³ and the two parts refer to the inequivalent points K and K' .

- a) Using the Peierls-substitution $\mathbf{p} \rightarrow \mathbf{p} - (e/c)\mathbf{A}$, find the Landau levels in graphene for a magnetic field perpendicular to the plane (ignore the Zeeman-term).

Hint: As the 2 parts of the effective Hamiltonian Eq. (5) are independent on each other, it is enough to consider only the first part $\mathcal{H}_K = v_F(p_x \sigma_x + p_y \sigma_y)$; the second part has the same spectrum. Take the “square” of the Schrödinger equation. Note that not only the σ_μ have non-trivial (anti-)commutation relations but also \mathbf{p} and \mathbf{r} do not commute.

- b) Determine the degeneracy of the Landau levels.
- c)* Will the magnetization of graphene oscillate when changing the magnetic field? What is the dependence of the ground state energy and the magnetization on a small magnetic field?

Office hour:

Monday, April 23th, 2012 - 10:00 to 12:00 am

HIT G 33.1

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²One possible way to come to the Hamiltonian given in Eq. (5) is to recall the Eq. (7) from Solution Sheet 5 with the formula for \mathcal{H} within the tight-binding model on a hexagonal lattice. If you would expand it near the Fermi point K , then substitute there $\hbar\mathbf{k}$ by $\hat{\mathbf{p}}$, you would (after a rotation of axes x , y) get the first part of the Hamiltonian given in Eq. (5). Similarly one could obtain the second part by expanding near K' .

³Recall exercise 5: $\begin{pmatrix} 1 \\ 0 \end{pmatrix}$ does correspond to a particle on sublattice A and $\begin{pmatrix} 0 \\ 1 \end{pmatrix}$ to a particle on the sublattice B.