Goethe-Universität Frankfurt Fachbereich Physik



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Frankfurt, 16.01.2020

## Einführung in die Theoretische Festkörperphysik Winter term 2019/2020

# Exercise 11

(Due date: 27.01.2020)

## Problem 1 (Specific heat of a semiconductor) (3 points)

Consider a semiconductor model where the density of states for valence and conducting band are constant, i.e.

$$\rho(E) = \begin{cases} \frac{1}{E_0} &, \ 0 < E < E_0 \text{ and } E_0 + \Delta < E < 2E_0 + \Delta \\ 0 &, \ \text{else} \end{cases}$$

Moreover, we assume one electron per unit cell in the system. How does the specific heat depend on the temperature in the limit of low temperatures?

#### Problem 2 (Extrinsic Semiconductor) (4 points)

Consider a semiconductor with gap  $\Delta$ , doped with donor atoms with concentration  $n_D$ . The doping introduces energy levels within the band gap, with distance  $\epsilon_D$  from the conduction band edge. Let the chemical potential be located close to the conduction band edge,  $\mu \approx \Delta$  (with respect to the upper valence band edge which defines E = 0 here).

*Hint*: We consider temperatures  $k_B T \ll \Delta$ , which is fulfilled at room temperature for semiconductors with a large gap  $\Delta$ . Electrons and holes are treated as free particles with same effective mass  $m_e = m_L$ . Furthermore, we approximate the dispersive bands as constant for simplicity.

- a) How many electrons are provided by the donor atoms on average (assuming single ionization)?
- b) From the neutrality condition  $n_e = n_L$ , derive the relation

(1) 
$$n_D = 2A \left( A e^{\frac{\epsilon_D}{k_B T}} + 1 \right) \left( \frac{m_e k_B T}{2\pi \hbar^2} \right)^{3/2}$$

*Hint*: Show that the valence band almost doesn't provide any conduction electrons.

### Problem 3 (Slater determinant) (3 points)

We consider the fermionic creation operator  $c_k^{\dagger} = \sum_{r_i} e^{ikr_i} c_{r_i}^{\dagger}$  that creates states  $|k\rangle$  with wave function  $\psi_k(r) = \langle r|k\rangle = \langle 0|c_r c_k^{\dagger}|0\rangle$ . Rewrite the wave function

$$\psi_{k_1,k_2,k_3}(r_1,r_2,r_3) = \langle 0|c_{r_1}c_{r_2}c_{r_3}c_{k_1}^{\dagger}c_{k_2}^{\dagger}c_{k_3}^{\dagger}|0\rangle$$

of a system of 3 fermions in terms of a Slater determinant.