Examination in Condensed Matter Physics I, FK7042, 7.5 hp

Thursday, March 19, 2015, 09.00-14.00.

Allowed help:

- periodic table and fundamental constants (distributed)

- formula sheet (distributed)

- pocket calculator, BETA / mathematics handbook or similar

Instructions:

All solutions should be easy to read and have enough details to be followed. The use of nontrivial formulas from the formula sheet should be explained. *Summarize each problem* before its solution, so that the solution becomes self-explained. State any assumptions or interpretation of a problem formulation and define any introduced variables.

Good luck! / A.R.

1. Consider two different ionic crystals, sodium chloride (NaCl) and caesium chloride (CsCl). Their lattice parameters are $a_{\text{NaCl}} = 5.65$ Å and $a_{\text{CsCl}} = 4.12$ Å, respectively. In the structures, the Na⁺ ions arrange in an *fcc* lattice, while the Cs⁺ ions arrange in a simple cubic (*sc*) lattice. The Cs⁺ ion is about 60% larger than the Na⁺ ion.

a) Use this information to make a crude estimate of the ionic radii for Na⁺, Cs⁺, and Cl⁻. (2p)

b) Find how much the density of NaCl would change if NaCl had the CsCl structure. (2p)

2. a) The relaxation time τ for electrons in the Drude model is defined to give a probability dt/τ for an electron to collide during a short time dt. Use this definition to find an expression for the probability of the electron not colliding during a time t. Interpret the result. (2p)

b) How does this relaxation time affect the thermal conductivity of a metal? (0.5p)

c) The electronic specific heat is linear in temperature, $c_{v,el} = \gamma T$. Use the Lorenz number $L = \pi^2 k_B^2 / 3e^2$ and Wiedemann-Franz law to express γ in terms of the Fermi energy ε_F . Describe used relations. (1.5p)

3. Transversal, optical lattice vibrations in germanium have a frequency $f \approx 9 \cdot 10^{12}$ Hz, which can be assumed independent of wavelength and propagation direction. Germanium has diamond structure and lattice parameter a = 5.66 Å.

a) Calculate the average number of transversal, optical phonons in a 1 cm^3 Ge-crystal at room temperature. Hint: the number of allowed wave vectors in the 1st Brillouin zone equals the number of primitive cells. (2p)

b) Calculate the thermal energy associated with these lattice vibrations for the given crystal. (1p)

c) Estimate the contribution from these lattice vibration to the heat capacity of the crystal at room temperature. (1p)

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4. The following three expressions can be found in the formula collection:

$$g_c(\varepsilon) = \frac{1}{2\pi^2} \left(\frac{2m_e}{\hbar^2}\right)^{3/2} (\varepsilon - \varepsilon_c)^{1/2}$$
$$n_c(T) = \int_{\varepsilon_c}^{\infty} \frac{1}{e^{(\varepsilon - \mu)/k_B T} + 1} g_c(\varepsilon) d\varepsilon$$
$$N_c(T) \approx \frac{1}{4} \left(\frac{2m_e k_B T}{\pi \hbar^2}\right)^{3/2}$$

Explain what they describe and and under what conditions/assumptions they apply. Also explain variables etc. in the expressions. (4p)

5. a) In an experiment, x-rays with a wave length $\lambda = 3.1$ Å are diffracted on a crystal with monoatomic simple cubic (*sc*) structure. The lattice parameter is a = 3.50 Å. Find all the possible diffraction angles of the experiment. (2.5p)

b) Suppose that the structure in a) would be modified by adding a light atom with a relatively small atomic form factor to the center of the *sc* cell. This would cause the diffraction intensity of all the diffraction peaks to change slightly. However, the relative change would be different for one of the peaks. Which one, and how would it change? (1.5p)

6. a) Pauli paramagnetism and free spin paramagnetism are two types of paramagnetism of quite different origin. Compare the two by explaining their origin, what main physical property they depend on, and their resulting temperature dependence of magnetic susceptibility. (2p)

b) State two formulations of the Bloch theorem and explain with words. (2p)