PHY 752 Solid State Physics 11-11:50 AM MWF Olin 103

Plan for Lecture 33:

Chapter 15 in GGGPP -- continued:

Electron Gas in Magnetic Fields

- 1. Landau diamagnetism and de Haasvan Alphen effect
- 2. Hall effect

Lecture notes prepared with materials from GGGPP textbook 11/16/2015 PHY 752 Fail 2015 - Lecture 33 1

22 Wed, 10/21/2015	Chap. 10	Scattering of particles by crystals	#20
23 Fri, 10/23/2015	Chap. 11	Optical and transport properties of metals	#21
24 Mon. 10/26/2015	Chap. 11	Optical and transport properties of metals	#22
25 Wed, 10/28/2015	Chap, 11	Transport in metals	#23
26 Fri, 10/30/2015	Chap 12	Optical properties of semiconductors and insulators	
27 Mon, 11/02/2015	Chap. 7 & 12	Excitons	#24
28 Wed, 11/04/2015	Chap 9	Lattice vibrations	#25
29 Fri, 11/06/2015	Chap. 9	Lattice vibrations	#26
30 Mon. 11/09/2015	Chap 13	Defects in semiconductors	#27
31 Wed, 11/11/2015	Chap. 14	Transport in semiconductors	#28
32 Fri, 11/13/2018	Chap. 15	Electron gas in Magnetic fields	#29
33 Mon, 11/16/2015	Chap. 15	Electron gas in Magnetic fields	Prepare presentation
34 Wed, 11/18/2015	Chap. 17	Magnetic ordering in crystals	Prepare presentation
35 Fri, 11/20/2015	Chap. 18	Superconductivity	Prepare presentation
36 Mon, 11/23/2015	Chap 18	Superconductivity	Prepare presentation
Wed, 11/25/2015		Thanksgiving Holiday	
Fri, 11/27/2015		Thanksgiving Holiday	
37 Mon, 11/30/2015	Chap 18	Superconductivity	Prepare presentation
Wed, 12/02/2015		Student presentations i	
Fri, 12/04/2015		Student presentations II	
Mon. 12/07/2015		Begin Take-home final	

















Some steps in the evaluation of the free energy of the electron gas in a magnetic field; summation of the contributions of the Landau levels (Thanks to appendix of Chapter 15 of GGGPP.)

Poisson sum formula

$$f\left(n+\frac{1}{2}\right) = \sum_{s=-\infty}^{+\infty} (-1)^s \int_n^{n+1} f(x) e^{2\pi i x x} dx.$$

$$\sum_{n=0}^{+\infty} f\left(n+\frac{1}{2}\right) = \int_0^{\infty} f(x) dx + 2\sum_{s=1}^{\infty} (-1)^s \int_0^{\infty} f(x) \cos(2\pi s x) dx$$
For a well-behaved function $f(x)$



For convenience, assume that we have shifted the energy scale so that the range of energies of interest are 0<E< μ where μ denotes the chemical potential. In terms of the density of states D(E), we can define:

$$P_1(E) = \int_0^E D(E') dE'$$
 with $P_1(0) = 0$

where $P_1(\mu) = N$, (the number of electrons)

Similarly, we can define

$$P_2(E) = \int_0^E P_1(E') dE' = \int_0^E dE' \int_0^E D(E'') dE''$$
 with $P_2(0) = 0$.

The ground state energy of the system of N electrons is given by

$$E_0 = \int_0^r ED(E) \, dE$$

 $E_0 = \int_0^\mu E \, dP_1(E) = \mu P_1(\mu) - \int_0^\mu P_1(E) \, dE = N \mu - P_2(\mu);$ 11/16/2015 PHY 752 Fall 2015 – Lecture 33

Evaluating the free energy of our system

$$F(T) = N\mu - k_B T \sum_{n} \ln \left[1 + e^{(\mu - E_n)/k_B T} \right],$$
In terms of the density of states:

$$F(T) = N\mu - k_B T \int_0^\infty \ln \left[1 + e^{(\mu - E)/k_B T} \right] D(E) dE.$$
which can be integrated by parts:

$$F(T) = N\mu - \int_0^\infty P_2(E) \left(-\frac{\partial f}{\partial E} \right) dE.$$

$$\approx N\mu - P_2(\mu, B)$$





Using Poisson sum formula with

$$P_{2}(E, B) = \frac{2}{3}\hbar\omega_{c}A\sum_{n=0}^{\infty} f\left(n + \frac{1}{2}\right) \text{ with } f(x) = (E - x\hbar\omega_{c})^{3/2}\Theta(E - x\hbar\omega_{c}).$$

$$\sum_{n=0}^{+\infty} f\left(n + \frac{1}{2}\right) = \int_{0}^{\infty} f(x) \, dx + \frac{1}{24}f'(0)$$

$$-\sum_{s=1}^{\infty} \frac{(-1)^{s}}{2\pi^{2}s^{2}} \int_{0}^{\infty} f''(x) \cos(2\pi sx) \, dx.$$

$$\int_{0}^{\infty} f(x) \, dx = \int_{0}^{E/\hbar\omega_{c}} (E - x\hbar\omega_{c})^{3/2} \, dx = \frac{2}{5} \frac{1}{\hbar\omega_{c}} E^{5/2}.$$

$$\frac{1}{24}f'(0) = \frac{1}{24}\frac{3}{2}(-\hbar\omega_{c})E^{1/2} = -\frac{1}{16}\hbar\omega_{c}E^{1/2}.$$

$$\int_{0}^{\infty} f''(x) \cos(2\pi sx) \, dx = \frac{3}{4}\hbar^{2}\omega_{c}^{2} \int_{0}^{E/\hbar\omega_{c}} \frac{1}{\sqrt{E - x\hbar\omega_{c}}} \cos(2\pi sx) \, dx$$

$$= \frac{3}{4}(\hbar\omega_{c})^{3/2} \frac{1}{\sqrt{2s}} \cos\left(\frac{2\pi s}{\hbar\omega_{c}}E - \frac{\pi}{4}\right)$$
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J. Phys. Chem. Solids Pergamon Press 1968. Vol. 29, pp. 1485–1502. Printed in Great Britain.

THE DE HAAS-VAN ALPHEN EFFECT AND THE FERMI SURFACE OF TUNGSTEN*

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(Received 8 April 1968)

(Received Appril 1990) Abstract – The de Haus-wan Alphen effect in tanguates has been studied in detail using large impulsive magnetic fields, and the frequencies have been determined to high accuracy with the aid of digital data recording, digital analysis and dynamic calibration of the entire upparture. These improvements in experimental technique have resulted in the detection of several new frequency branches which have not been previously reported. A quantitative description of all beets of the Ferrei surface is given in terms of simple analytical functions, and this geometrical model is in excellent agreement with the observed angular variations of the de Haas-van Alphen frequencies. In most respects, the model surface is found to be quite similar to that deduced by Loucks from first-principles relativistic band calculations, and the predictions of the empirical model are compared with the results of other experiments relating to the Fermi surface.

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