Condensed Matter Physics

PHY 5111

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The Free Fermi Gas and Single Electron Model

Condensed Matter Physics – Michael P. Marder

Chapter 6

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2

Much of condensed matter physics lies within a Hamiltonian that one easily can write down in a single line. It is

 $\hat{\mathcal{H}} = \sum_{l} \frac{P_{l}^{2}}{2M_{l}} + \frac{1}{2} \sum_{l \neq l'} \frac{q_{l}q_{l'}}{|\hat{R}_{l} - \hat{R}_{l'}|}.$

The single-electron model

$$
\sum_{l=1}^N (\frac{-\hbar^2 \nabla_l^2}{2m} + U(\vec{r}_l)) \Psi(\vec{r}_1 \dots \vec{r}_N) = \mathcal{E} \Psi(\vec{r}_1 \dots \vec{r}_N)
$$

$$
\left| \left(\frac{-\hbar^2 \nabla^2}{2m} + U(\vec{r}) \right) \psi_I(\vec{r}) = \mathcal{E}_I \psi_I(\vec{r}) \right|
$$

The free Fermi gas

$$
\frac{-\hbar^2}{2m}\sum_{l=1}^N \nabla_l^2 \Psi(\vec{r}_1 \dots \vec{r}_N) = \mathcal{E}\Psi(\vec{r}_1 \dots \vec{r}_N)
$$

To simplify further we imposes periodic boundary conditions

$$
\Psi(x_1 + L, y_1, z_1 \ldots, z_N) = \Psi(x_1, y_1, z_1 \ldots z_N)
$$

$$
\Psi(x_1, y_1 + L, z_1 \ldots, z_N) = \Psi(x_1, y_1, z_1 \ldots z_N)
$$

One Free Fermion

 $|l_x, l_y,$ and l_z are integers ranging from $-\infty$ to ∞

$$
\varepsilon_{\vec{k}}^0 = \frac{\hbar^2 k^2}{2m}
$$

The ground state of electrons obeying free Fermi gas assumption is constructed from products of the one-electron wave functions. The Pauli exclusion principle forbids any given state from being occupied more than once, and therefore any given state indexed by \vec{k} is able to host no more than two electrons, one for each value of spin.

Many Free Fermions

 $|l_x, l_y,$ and l_z are integers ranging from $-\infty$ to ∞

Density of States

Volume occupied by 2 states $=$

Total number of states $=$

States per unit volume =

Density of States

$$
D_{\vec{k}} = 2 \frac{1}{(2\pi)^3}
$$

For each wave vector Pauli's exclusion principle allows two electrons, one with spin up and the other with spin down.

$$
\int [d\vec{k}] \equiv \frac{2}{\overline{v}} \sum_{\vec{k}} = \int d\vec{k} D_{\vec{k}} = \frac{2}{(2\pi)^3} \int d\vec{k}
$$

Energy Density of States

$$
D(\mathcal{E}) = \int [d\vec{k}] \delta(\mathcal{E} - \mathcal{E}_{\vec{k}})
$$

The units of densities of states are able to change without much warning. Often they are expressed in units of 1/[eV atom], which means they are related to the function defined by above equation by a factor of density *n*.

Energy Density of States

$$
D(\mathcal{E}) = \int [d\vec{k}] \ \delta(\mathcal{E} - \mathcal{E}_{\vec{k}}^0)
$$

= $4\pi \frac{2}{(2\pi)^3} \int_0^\infty dk \ k^2 \delta(\mathcal{E} - \mathcal{E}_{\vec{k}}^0)$
= $\frac{1}{\pi^2} \int_0^\infty \frac{d\mathcal{E}^0}{|d\mathcal{E}^0/dk|} \frac{2m\mathcal{E}^0}{\hbar^2} \delta(\mathcal{E} - \mathcal{E}^0)$
= $\frac{m}{\hbar^3 \pi^2} \sqrt{2m\mathcal{E}}$
= 6.812 \cdot 10²¹ $\sqrt{\mathcal{E}/eV} eV^{-1}$ cm⁻³.

Fermi gas

For the free

Electron density

The number of electrons that can fit into a sphere of radius k_F is

$$
N = \sum_{\vec{k}\sigma} f_{\vec{k}}
$$

= $\sqrt{\int [d\vec{k}] f_{\vec{k}}},$
= $\frac{\sqrt{2}}{4\pi^3} \frac{4\pi}{3} k_F^3 = \frac{\sqrt{2}}{3\pi^2}$

Because $f_{\vec{k}}$ is 1 only if the state is occupied.

$$
n = N/\mathcal{V}
$$

Electron density

$$
4\pi \int_0^{k_F} g(k)k^2 dk = \frac{N}{V_g} = n
$$

$$
k_F = (3\pi^2 n)^{1/3}, \qquad E_F = \frac{\hbar^2}{2m} (3\pi^2 n)^{2/3}
$$

$$
k_F = (3\pi^2 n)^{1/3} = 3.09 [n \cdot \text{\AA}^3]^{1/3} \text{\AA}^{-1}
$$

$$
\mathcal{E}_F = \frac{\hbar^2 k_F^2}{2m} = 36.46 [n \cdot \text{\AA}^3]^{2/3} \text{eV}
$$

For the free Fermi gas

$$
v_F = \hbar k_F / m = 3.58 [n \cdot \text{\AA}^3]^{1/3} \cdot 10^8 \text{ cm s}^{-1}
$$

$$
D(\mathcal{E}_F) = \frac{3}{2} \frac{n}{\mathcal{E}_F} = 4.11 \cdot 10^{-2} [n \cdot \text{\AA}^3] \text{ eV}^{-1} \text{ \AA}^{-3}
$$

Grand Partition Function

$$
Z_{\rm gr} = \sum_{\rm states} e^{\beta(\mu N - \mathcal{E})}
$$

$$
= \sum_{n_1=0}^1 \sum_{n_2=0}^1 \sum_{n_3=0}^1 \ldots e^{\beta \sum_l n_l(\mu - \mathcal{E}_l)}
$$

$$
= \prod_{l} \left\{ \sum_{n_l=0}^{1} e^{\beta n_l [\mu - \mathcal{E}_l]} \right\}
$$

=
$$
\prod_{l} \left[1 + e^{\beta[\mu - \mathcal{E}_l]} \right].
$$

$$
\sum_{n_1=0}^N \sum_{n_2=0}^N \ldots \sum_{n_M=0}^N \prod_{l=1}^M A_{n_l} = \prod_{l=1}^M \left\{ \sum_{n_l=0}^N A_{n_l} \right\}
$$

 $\Pi \equiv -k_B T \ln Z_{gr}$ $= -k_BT \sum_{l} \ln \left[1 + e^{\beta[\mu - \mathcal{E}_l]} \right].$ $= -k_B T \mathcal{V} \int d\mathcal{E} D(\mathcal{E}) \ln \left[1 + e^{\beta[\mu - \mathcal{E}]} \right].$

Fermi Distribution Function

$$
N = -\frac{\partial \Pi}{\partial \mu}
$$

= $\sqrt[n]{\int d\mathcal{E}' D(\mathcal{E}') \frac{e^{\beta \mu - \beta \mathcal{E}'}}{1 + e^{\beta \mu - \beta \mathcal{E}'}}}$

$$
\Rightarrow n = \frac{N}{\sqrt[n]{\mu}} = \int d\mathcal{E}' D(\mathcal{E}') f(\mathcal{E}'),
$$

$$
f(\mathcal{E}) = \frac{1}{e^{\beta(\mathcal{E} - \mu)} + 1}
$$

18

Fermi Distribution Function

Fermi Distribution Function

Fermi Surface

The derivative of the Fermi function

Specific Heat of Noninteracting Electrons at Low T

$$
c_{\mathcal{V}} = \frac{1}{\mathcal{V}} \frac{\partial \mathcal{E}}{\partial T} \big|_{\mathcal{NV}}
$$

$$
c_V = \frac{\pi^2}{3} k_B^2 T D(\mathcal{E}_F)
$$

