

## Degenerate electron gas

↓  
quantum statistics [Fermi-Dirac]  
must be applied

$$k_B T \ll E_F = \frac{P_F^2}{2m} = (3\pi^2)^{2/3} \frac{\hbar^2}{m} n^{2/3}$$

Chemical potential  $\mu = E_F - \text{corrections}$   
 $O\left[\left(\frac{k_B T}{E_F}\right)^2\right]$

Occupation numbers

$$n = \frac{1}{e^{\frac{\epsilon - \mu}{k_B T}} + 1} \approx 1 - e^{-\frac{\mu - \epsilon}{k_B T}} \approx 1 \quad \begin{matrix} \text{large} \\ \forall \epsilon \end{matrix}$$

$$+ \frac{\mu - \epsilon}{k_B T} \gg 1$$

Plasma: electrons of density  $n$   
 (neutral)  $+ \text{ions } \oplus Ze \quad n_i = \frac{1}{Z} n$

e-ion interaction  $\frac{Ze^2}{a} \ll E_F$  kinetic energy per one electron

$$a \sim n_i^{-1/3} = (Z/n)^{1/3}$$

mean distance between ions

$$ze \cdot \left(\frac{n}{Z}\right)^{1/3} \ll \frac{\hbar^2}{m} n^{2/3}$$

$$Z^{2/3} \ll \frac{\hbar^2}{me^2} n^{1/3}$$

Bohr radius

$$n \gg Z^2 \cdot \left(\frac{\hbar^2}{me^2}\right)^{-3} = Z^2 a_0^{-3}$$

$$n \cdot a_0^3 \gg Z^2$$

The higher density of degenerate electron gas, the more ideal it is:

$$K \gg U_{e\text{-ion}}, U_{e\text{-e}}$$

Thermal properties of degenerate electron gas



$$k_B T \ll E_F$$

ground state

$$T=0$$

$$+ \text{ corrections } \sim \left(\frac{k_B T}{E_F}\right)$$

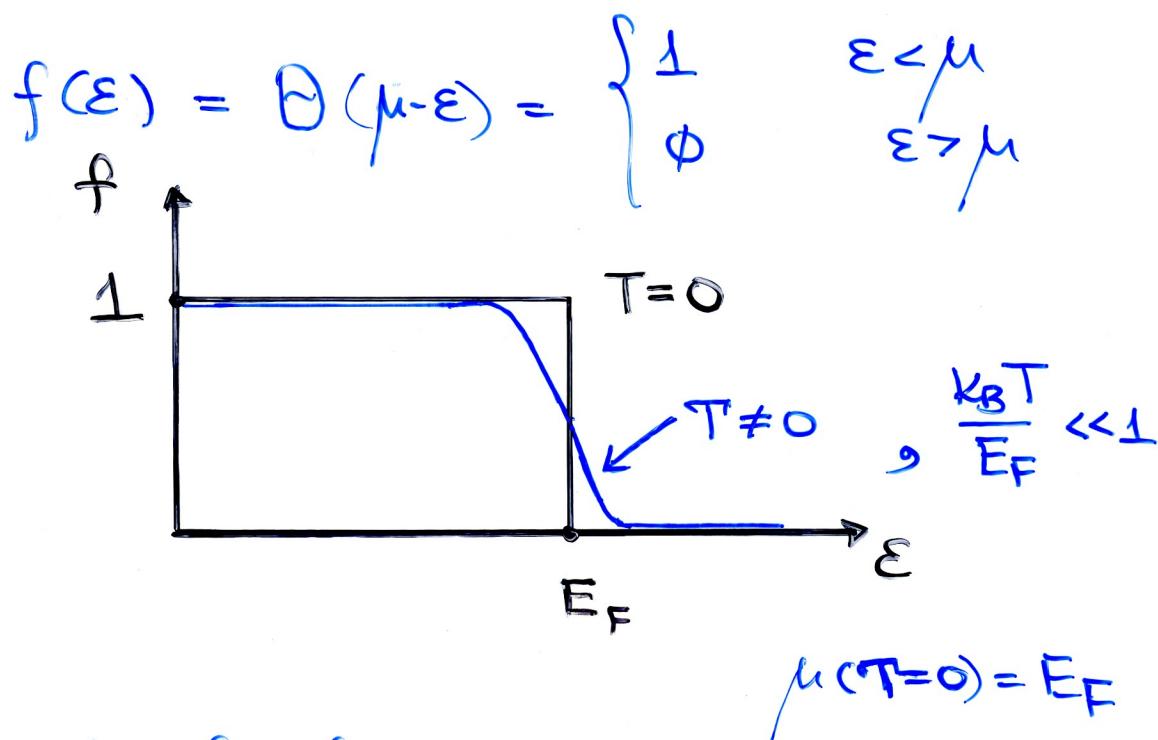


some power

the only source  
of  $C_V$  ...

$$T \rightarrow 0$$

$$f(\epsilon) = \frac{1}{\exp[(\epsilon - \mu)/k_B T] + 1}$$



Condition for finding  $\mu$ :

$$n = \int_{-\infty}^{\infty} d\epsilon f(\epsilon) \cdot g(\epsilon)$$

↑ Density of states

$$n = 2 \cdot \int \frac{d^3 k}{(2\pi)^3} f(\epsilon_k)$$

↑ electron density

$$d^3 k = 4\pi k^2 dk = 4\pi \cdot \left(\frac{2m\epsilon}{\hbar}\right) \cdot \frac{1}{2} \left(\frac{2m}{\hbar}\right)^{1/2} \epsilon^{1/2} dk$$

$$\frac{\hbar^2 k^2}{2m} = \epsilon \quad k = \left(\frac{2m\epsilon}{\hbar}\right)^{1/2} \quad dk = \frac{1}{2} \left(\frac{2m}{\hbar}\right)^{-1/2} \epsilon^{-1/2} d\epsilon$$

$$g(\epsilon) = \frac{(2m)^{3/2}}{2\pi^2 \hbar^3} \sqrt{\epsilon}$$

DOS per unit volume  
 $v=1$  for both  $T \downarrow$   
 directions of spin

DOS

$$g(\epsilon) = \begin{cases} \frac{3}{2} \cdot \frac{n}{E_F} \left( \frac{\epsilon}{E_F} \right)^{1/2}, & \epsilon > 0 \\ 0 & \epsilon < 0 \end{cases}$$

$$E_F = \frac{\hbar}{2m} \cdot (3\pi^2 n)^{2/3} \uparrow_{\text{used}}$$

We can calculate now various thermodynamic quantities:

Internal energy density

$$U = \int_{-\infty}^{\infty} d\epsilon f(\epsilon) g(\epsilon) \cdot \epsilon \quad \text{etc.}$$

For metals  $T_{\text{room}} \ll E_F \sim (2-25) \times \underbrace{12 \times 10^3 \text{ K}}_{1 \text{ eV}}$

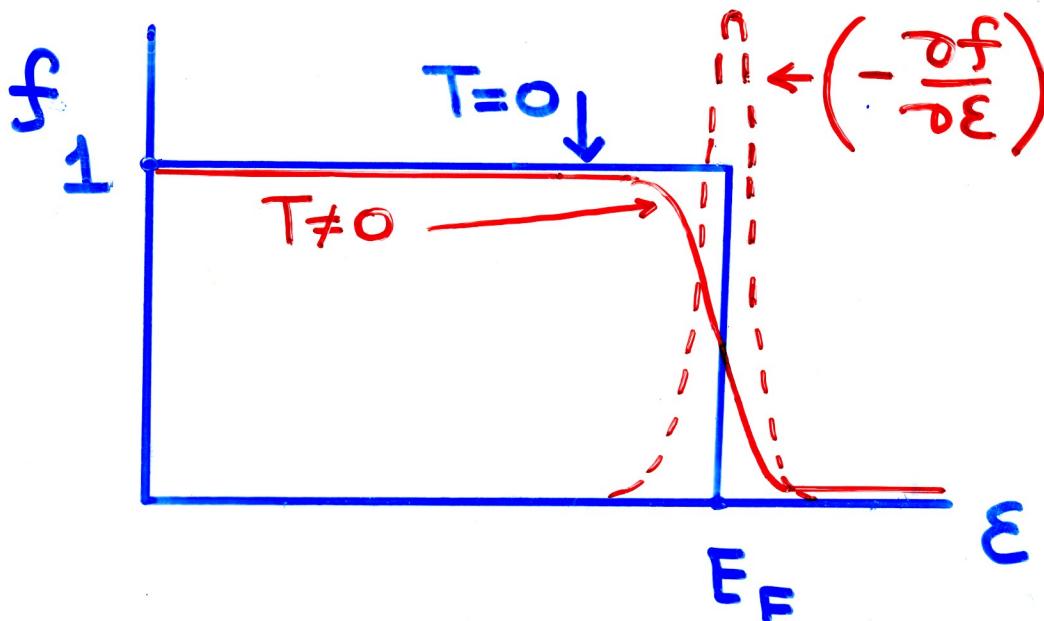
Low-temperature expansion Sommerfeld

$$\int_{-\infty}^{\mu} d\epsilon \cdot f(\epsilon) \cdot H(\epsilon) = \int_{-\infty}^{\mu} d\epsilon \cdot H(\epsilon) +$$

$$+ \sum_{n=1}^{\infty} (k_B T)^{2n} \cdot a_n \left. \frac{d^{2n-1} H(\epsilon)}{d \epsilon^{2n-1}} \right|_{\epsilon=\mu}$$

dimensionless constants of order unity

$$a_1 = \frac{\pi^2}{6} \quad a_2 = \frac{7\pi^4}{360}$$



$$f(\varepsilon) = \frac{1}{\exp\left(\frac{\varepsilon-\mu}{k_B T}\right) + 1}$$

$$x = \frac{\varepsilon - \mu}{k_B T}$$

$$-\frac{\partial f}{\partial \varepsilon} = \frac{1}{k_B T} \cdot \frac{e^x}{[e^x + 1]^2} = \frac{1}{k_B T} \cdot \frac{1}{4 \cosh^2 \frac{x}{2}}$$

even

$$\int_{-\infty}^{\infty} d\varepsilon \left( -\frac{\partial f}{\partial \varepsilon} \right) K(\varepsilon) = K(\mu) + \text{corrections}$$

↑       $k_B T \ll E_F$

behaves like  $\delta(\varepsilon - \mu) + \text{corrections}$

$$T=0: f(\varepsilon) = \Theta(\mu - \varepsilon) \quad -\frac{\partial f}{\partial \varepsilon} = \delta(\varepsilon - \mu)$$

$$\int_{-\infty}^{\mu} d\epsilon f(\epsilon) H(\epsilon) = \int_{-\infty}^{\mu} d\epsilon H(\epsilon) + \frac{\pi^2}{6} (k_B T)^2 H'(\mu) + \dots$$

If  $H(\epsilon)$  is a smooth function with variations on the energy scale  $\sim \mu$

$$H' \sim \frac{H}{\mu}$$

$$\Rightarrow \text{expansion in } \frac{k_B T}{E_F} \ll 1$$

Examples: T-corrections to

- ① internal energy
- ② chemical potential

$$\begin{aligned} \textcircled{1} \quad u &= \int_{-\infty}^{\mu} d\epsilon f(\epsilon) \left[ g(\epsilon) \epsilon \right] \xrightarrow{\substack{\text{--- ---} \\ \text{--- ---}}} H(\epsilon) & H'(\epsilon) = g + g' \cdot \epsilon \\ \textcircled{2} \quad n &= \int_{-\infty}^{\mu} d\epsilon f(\epsilon) \left[ g(\epsilon) \right] \xrightarrow{\substack{\text{--- ---} \\ \text{--- ---}}} H(\epsilon) & H'(\epsilon) = g'(\epsilon) \end{aligned}$$

$$g(\epsilon) = \frac{(2m)^{3/2}}{2\pi^2 \hbar^3} \sqrt{\epsilon} \quad \text{DOS}$$

$$\int_0^{\mu} d\epsilon H(\epsilon) \simeq \int_0^{E_F} d\epsilon H(\epsilon) + H(E_F) \cdot (\mu - E_F)$$

Chemical potential

$$\mu = E_F \left[ 1 - \frac{\pi^2}{12} \cdot \left( \frac{k_B T}{E_F} \right)^2 \right] \checkmark$$

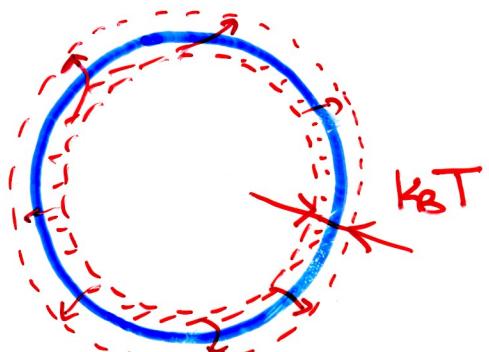
$$U = U_0 + \frac{\pi^2}{6} (k_B T)^2 \cdot g(E_F)$$

thermal energy density

$$g(E_F) = \frac{3}{2} \cdot \frac{n}{E_F}$$

DOS at the Fermi energy

correction  $\delta U \sim k_B T \cdot \frac{k_B T}{E_F}$



Physical meaning:

$$\Delta N \sim k_B T \cdot g(E_F)$$

# of excited electrons

$$\delta U \sim k_B T \cdot \Delta N \sim (k_B T)^2 \cdot g(E_F)$$

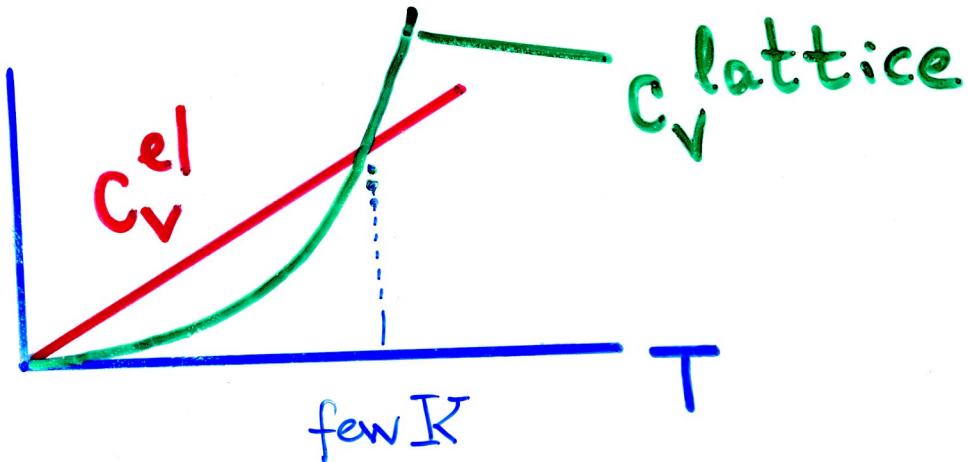
$$C_V = \left( \frac{\partial U}{\partial T} \right)_V = \frac{\pi^2}{3} k_B^2 T \cdot g(E_F)$$

$$C_V = \frac{\pi^2}{2} \cdot \frac{k_B T}{E_F} \cdot n k_B \sim \frac{k_B T}{E_F} \cdot C_V^{\text{class}}$$

$$C_V^{\text{el}} \sim T$$

$$C_V = C_V^{\text{el}} + C_V^{\text{lattice}}, \quad C_V^{\text{lattice}} \sim T^3$$

$$C_V = \gamma \cdot T + A \cdot T^3$$



Specific heat effective mass  $m^*$

$$C_V^{\text{el}} \sim k_B^2 T \cdot g(E_F) \sim n/E_F \sim \left(\frac{\hbar^2}{m r_s^2}\right)^{-1} \xrightarrow{\text{electron mass}}$$

$$C_V^{\text{el}} = \frac{m^*}{m} \cdot \gamma T$$

$\uparrow$  for free electrons of density  $n$

$m^* \sim 1$	Alkali, Noble metals	$m^* \sim 10^{-2} - 10^{-1}$	Bi, SB
		$m^* \sim 10^{-2}$	22
		$m^* \sim 10^{-1}$	..
		$m^* \sim 10$	Mn, Fe

## Conduction in metals

Boltzmann distribution  $\rightarrow$  Fermi-Dirac distr.

$$\textcircled{1} \quad f(\vec{v}) = \frac{(m/h)^3}{4\pi^3} \cdot \frac{1}{\exp([\frac{mv^2}{2} - \mu]/k_B T) + 1}$$

\textcircled{2}. (Quasi)classical description of electron motion

$$\Delta x \cdot \Delta p \sim h$$

$$\Delta x \ll l, \lambda$$

$$\Delta p \gg \frac{\hbar}{e}, \frac{\hbar}{\lambda} \quad \ll \hbar k_F = p_F$$

$k_F \sim r_s^{-1}$

mean free path

wave length,  
charact. length

$l, \lambda \gg r_s$

Mean free path

$$l = v_F \cdot \tau \sim 10^8 \frac{\text{cm}}{\text{sec}} \times 10^{-14} \text{ sec}$$

$$l \sim 10^{-6} \text{ cm} = 100 \text{ \AA}$$

Large compared to lattice spacing!

## Thermal conductivity

$$K = \frac{1}{3} V^2 C_V$$

$$\downarrow V_F^2 \quad \downarrow \frac{k_B T \cdot C_V^{\text{class}}}{E_F} = \frac{\frac{1}{3} \frac{m V_{\text{class}}^2}{2}}{\frac{m V_F^2}{2}} \cdot C_V^{\text{class}}$$

quantum contributions  
compensate

## Electrical conductivity

$$\sigma = \frac{n e^2 \tau}{m}$$

does not change  
if  $\tau$  does not  
depend on energy  $\Leftrightarrow$   
on velocity distribution

## Lorenz number

Sommer.  $\frac{K}{\sigma T} = \left[ \frac{\pi^2}{3} \left( \frac{k_B}{e} \right)^2 \right] = 2.4 \times 10^{-8} \frac{\text{watt} \cdot \text{ohm}}{\text{K}^2}$

excellent agreement  
with exp.

Drude  $\frac{K}{\sigma T} = \frac{3}{2} \cdot \left( \frac{k_B}{e} \right)^2 = 1.1 \times 10^{-8}$

## Thermopower Q

$$\vec{E} = Q \vec{\nabla} T$$

$$Q = - \frac{C_V}{3ne}$$

$$Q = - \frac{\pi^2}{6} \frac{k_B}{e} \cdot \frac{k_B T}{E_F}$$

Somm.  $C_V = \frac{\pi^2}{2} \left( \frac{k_B T}{E_F} \right) \cdot n k_B$

Drude  $C_V^{\text{class}} = \frac{3}{2} n k_B$  :  $\left[ \frac{k_B T}{E_F} \right]^{-1} \approx 10^2$  overestimation stat.

Solids: Is classical description of electron kinetics worthless?

No! : Non-degenerate electron gas  
classical

low densities of electrons

$$\frac{\hbar^2}{m r_s^2} \approx \frac{\hbar^2}{m} \cdot 12^{2/3} \ll k_B T$$

$$r_s \gg \left( \frac{\hbar^2}{m k_B T} \right)^{1/2}$$

Semiconductors