

Condensed Matter Physics I

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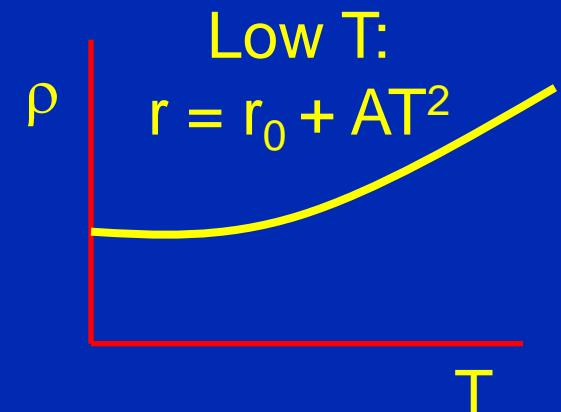
Website: <http://www.loosdrecht.net/>

Metals

What is a metal ?

Electrical conductivity:

$$\rho_{300\text{ K}} \sim 1.7 \text{ (Cu)} - 153 \mu\Omega\cdot\text{cm} \text{ (Pu)}$$



Thermal conductivity:

$$\text{Cu: } K_{300\text{ K}} \sim 3.9 \text{ W/Kcm} \quad \text{Pu: } K_{300\text{ K}} \sim 0.049 \text{ W/Kcm}$$

$$\text{Wiedemann-Franz: } K/\sigma = \alpha T$$

$$\text{Quartz: } K \sim 0.13 \text{ W/Kcm} \quad \text{NaCl: } K \sim 0.27 \text{ W/Kcm}$$

Reflectivity:

Highly reflecting upto plasma-frequency

$$\omega < \omega_p \quad \omega_p^2 = 4 \pi n e^2 / m$$

FREE ELECTRON MODEL

FEM, overview

- Free electron model (Drude, Sommerfeld theory)
- Statistics and density of states
- Heat capacity
- Electrical conductivity (Ohm's law)
- Influence of a magnetic field (Hall effect)
- Thermal conductivity and Wiedemann-Franz law

Electrons in metals

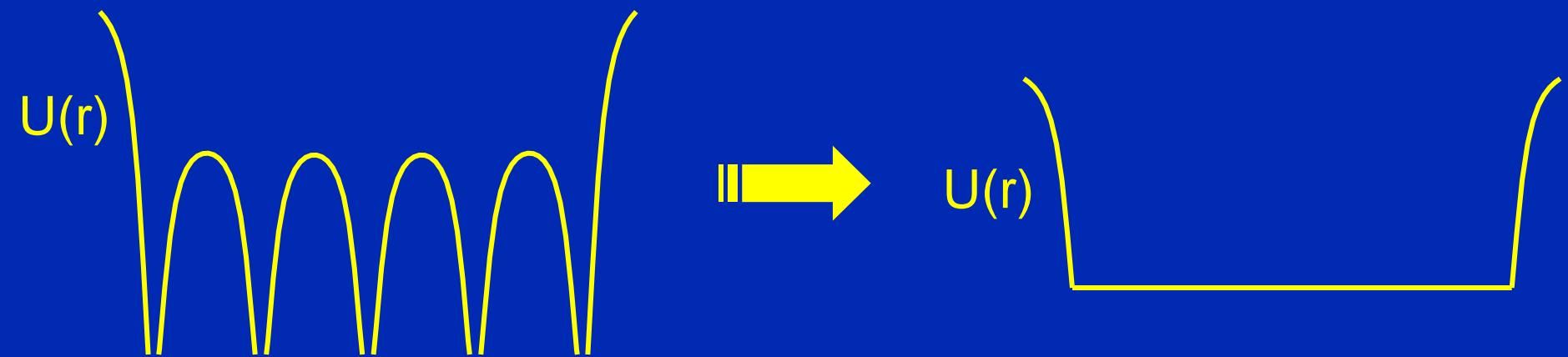
P. Drude: 1900 kinetic gas theory of electrons, classical
Maxwell-Boltzmann distribution
independent electrons
free electrons
scattering from ion cores (relaxation time approx.)

A. Sommerfeld: 1928
Fermi-Dirac statistics

F. Bloch's theorem: 1928
Bloch electrons

L.D. Landau: 1957
Interacting electrons (Fermi liquid theory)

Free electron approximation



Neglect periodic potential & scattering (Pauli)

Reasonable for “simple metals” (Alkali Li,Na,K,Cs,Rb)

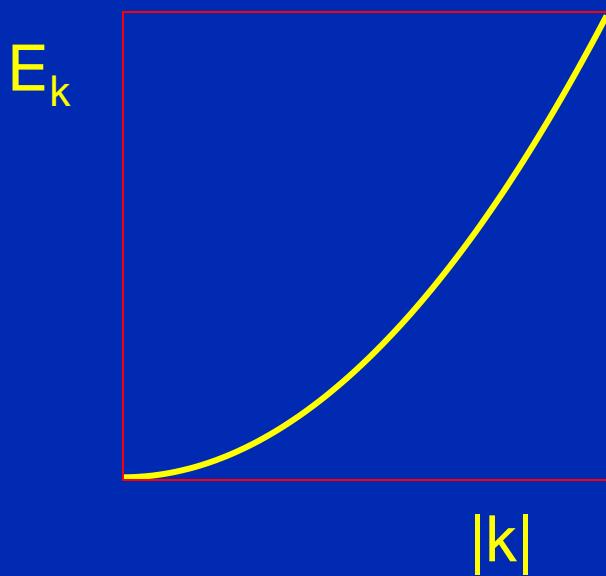
Eigenstates & energies

$$\left(\frac{-\hbar^2}{2m} \nabla^2 + \chi \right) \Psi = i\hbar \frac{d\Psi}{dt}$$

$$\Psi_{\vec{k}}(r, t) = \Psi_0 e^{i(\omega t - \vec{k} \cdot \vec{r})}$$

$$E_{\vec{k}} = \frac{\hbar^2}{2m} |\vec{k}|^2$$

$$\vec{k} = 2\pi(n_x/L_x, n_y/L_y, n_z/L_z)$$

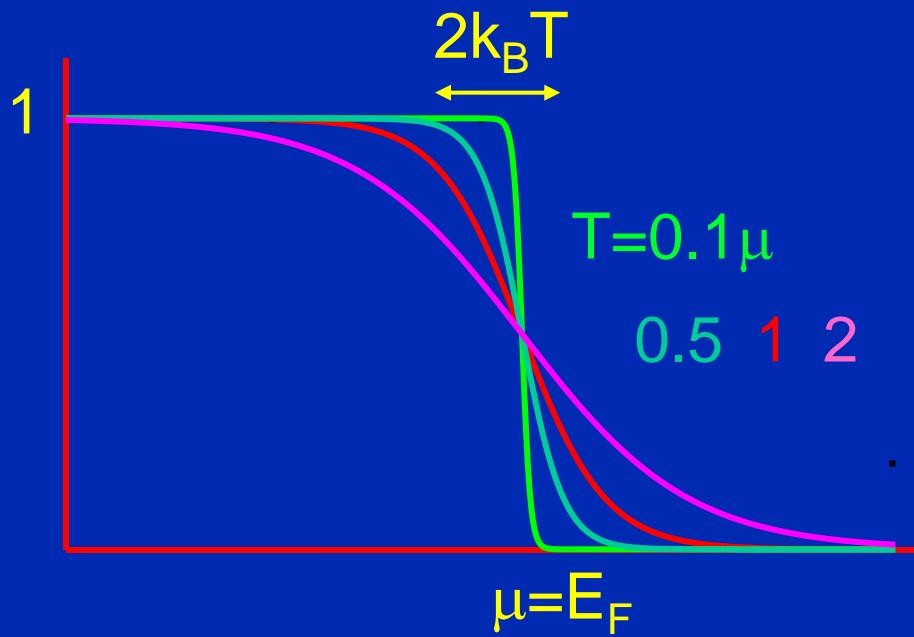


$$D_d(k) \propto L^d \cdot k^{d-1}$$

Statistics & DOS

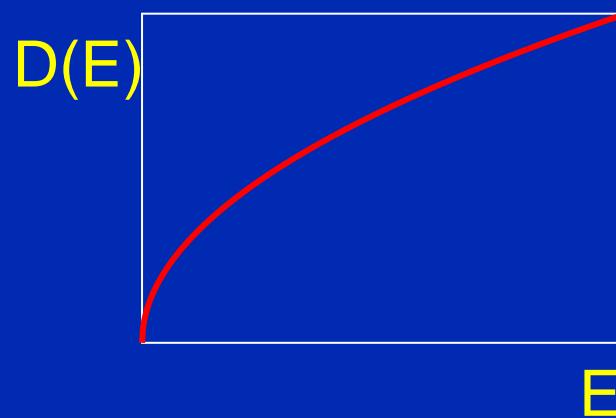
Fermi-Dirac statistics:

$$f_{FD}(E) = \frac{1}{e^{\frac{E-\mu}{kT}} + 1}$$

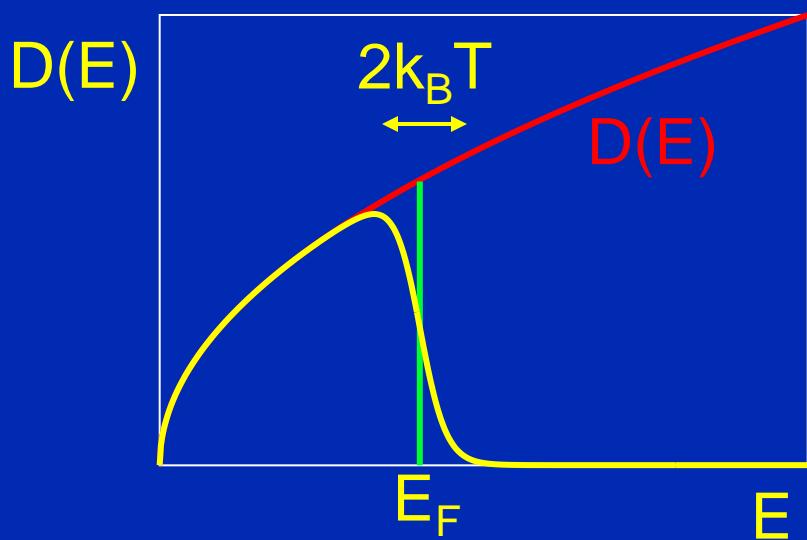
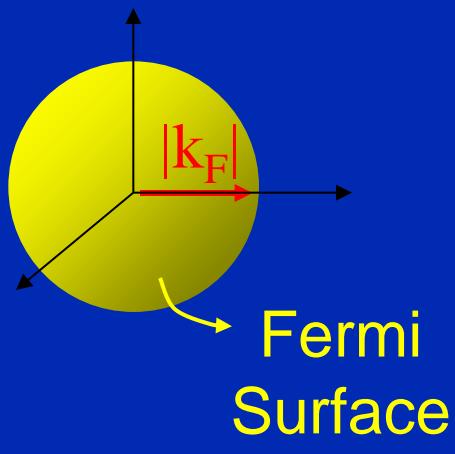


Density of states:

$$D(E) = 2 \cdot D(k) \cdot \frac{dk}{dE} = \frac{V_m}{\pi^2 \hbar^3} \sqrt{2mE}$$



Occupation of states



Free electron gas parameters

$$N = \int_0^{\infty} D(E) \cdot f_{FD}(E) dE$$

$$E_F = \frac{\hbar^2 k_F^2}{2m}$$

$$E_F = \frac{1}{2} m v_F^2$$

$$T_F = E_F / k_B$$

$$D(E_F) = \frac{V}{\pi^2 \hbar^3} \sqrt{2m E_F}$$

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

$$k_F = (3\pi^2 n)^{1/3}$$

$$v_F = \frac{\hbar k_F}{m}$$

$$T_F = \frac{\hbar^2}{2mk_B} (3\pi^2 n)^{2/3}$$

$$D(E_F) = \frac{V}{\pi \hbar^2} \cdot \left(\frac{3}{\pi} n \right)^{1/3} m$$

Sodium

Na $1s^2 2s^2 2p^6 3s = [\text{Ne}]3s$

$r_{\text{ion}} = 0.98 \text{ \AA}$ $d_{\text{nn}} = 1.83 \text{ \AA}$

$$n = 2.65 \cdot 10^{22} \text{ cm}^{-3}$$

$$E_F = 3.23 \text{ eV}$$

$$k_F = 2.65 \cdot 10^{22} \text{ cm}^{-1}$$

$$v_F = 1.07 \cdot 10^8 \text{ cm/s}$$



Non relativistic

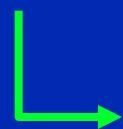
$$T_F = 3.75 \cdot 10^4 \text{ K}$$



Degenerate quantum gas
Chemical potential
hardly depends on T

So far

- Free electrons, i.e. no periodic potential
- Independent electrons, i.e. no $e^- - e^-$ interactions
- Relaxation time approximation (scattering time τ)
- Classical statistics (Drude)
- Fermions (Sommerfeld)



- Density of states (1D, 2D, 3D)
- Fermi energy (E_F , k_F v_F , T_F)

t.b.d.

- Sommerfeld (=Drude+quantum)
 - Compressibility
 - Heat capacity
 - Conductivity, Hall effect, 1D conduction
 - Thermal conductivity, Wiedemann-Franz law
- FAILURES of the free electron models
- Including the periodic potential the e^- live in.

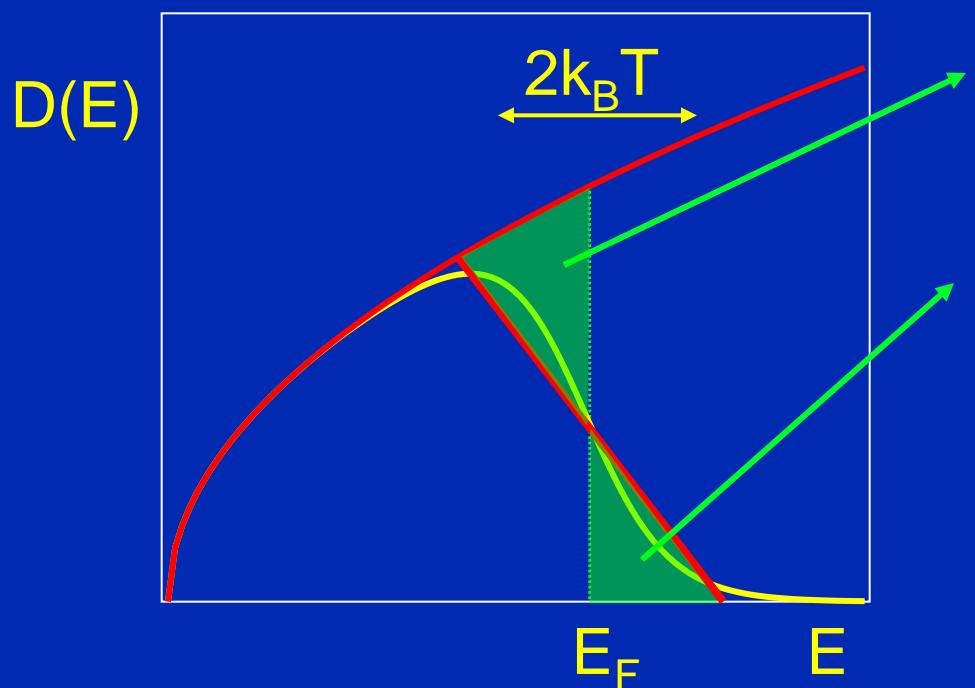
Compressibility

$$B = 1/K = -V \frac{dP}{dV} = V \frac{d^2U}{dV^2}$$

	$B_{f.e.}$ (Gpa)	B_{obs} (Gpa)
Li	24	12
Na	9	6.5
K	3	3
Rb	2	2
Cs	1.5	1.5
Cu	64	134
Ag	35	100
Al	228	76

Heat capacity: Quick&Dirty

$$C_{\text{el}} = \frac{dU_{\text{el}}}{dT} = \frac{d}{dT} \int_0^{\infty} E \cdot D(E) \cdot f_{\text{FD}}(E, T) dE$$



$$\begin{aligned} & -\frac{1}{2} k_B T \times \frac{1}{2} D(E_F) \cdot \left(E_F - \frac{k_B T}{2} \right) \\ & + \frac{1}{2} k_B T \times \frac{1}{2} D(E_F) \cdot \left(E_F + \frac{k_B T}{2} \right) \end{aligned}$$

$$\Delta U = \frac{k_B^2 T^2}{2} D(E_F)$$

$$\begin{aligned} C_{\text{el}} &= k_B^2 D(E_F) \cdot T \\ \Rightarrow C_{\text{el}} &\propto T \end{aligned}$$

Heat capacity

Electronic contribution

$$C_{\text{el}} = \frac{1}{3} \pi^2 \cdot D(E_F) \cdot k_B^2 T = \frac{1}{2} \pi^2 N k_B \frac{T}{T_F} \ll \frac{3}{2} N k_B$$

Electrons + lattice (low T): $C_v = \gamma \cdot T + A \cdot T^3$

$$\gamma = \frac{1}{3} \pi^2 \cdot D(E_F) \cdot k_B^2 \propto m$$

$$\frac{\gamma_{\text{exp}}}{\gamma_{f.e.}} \neq 1 \quad \Rightarrow \quad m^{*}_{th} \equiv \frac{\gamma_{\text{exp}}}{\gamma_{f.e.}} m_0$$



Periodic potential (band mass)
e-p interaction (polarons)
e-e interaction

Heat capacity: Na & K

PHYSICAL REVIEW

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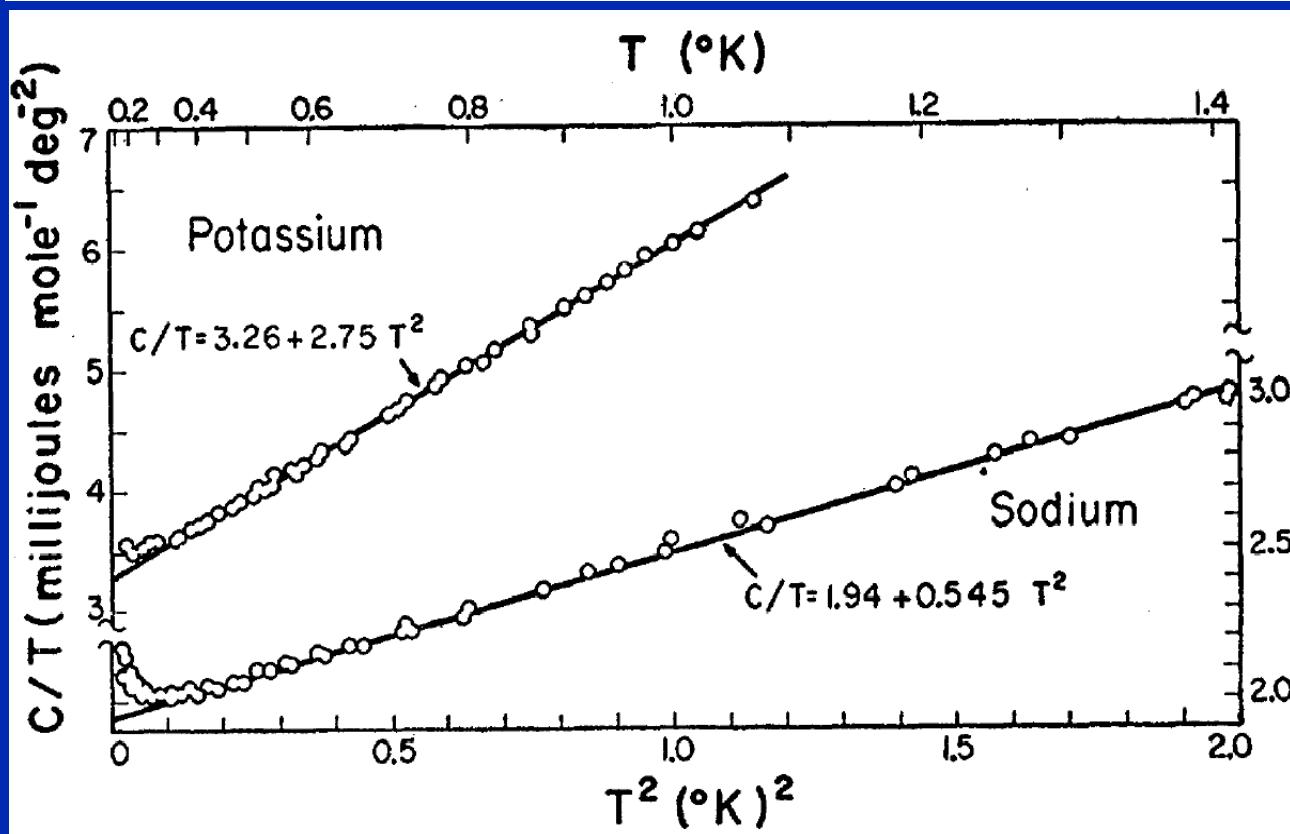
MAY 15, 1960

Heat Capacity of Sodium and Potassium at Temperatures below 1°K

WILLIAM H. LIEN AND NORMAN E. PHILLIPS

Department of Chemistry and Lawrence Radiation Laboratory, University of California, Berkeley, California

(Received December 17, 1959)



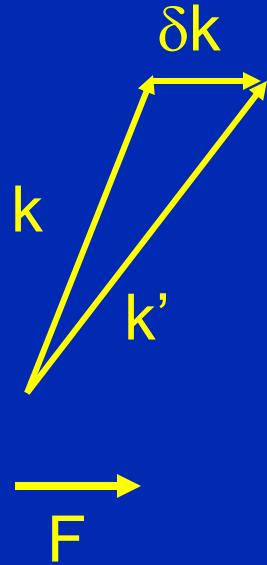
$$C = \gamma T + A T^3$$

$$m_{\text{th}} = 1.25 m_0$$

γ & thermal effective mass

Element	Free e ⁻ γ	Expt. γ	m_{th}^*/m_0
	10^{-4} cal/mol K ²		
Li	1.8	4.2	2.3
Na	2.6	3.5	1.3
K	4.0	4.7	1.2
Cu	1.2	1.6	1.3
Be	1.2	0.5	0.42
Fe	1.5	12	8
Mn	1.5	40	27
Bi	4.3	0.2	0.047

Electrical conductivity



$$2^{\text{nd}} \text{ Newton: } m \frac{dv}{dt} = F \quad \text{or} \quad \delta k = \frac{1}{\hbar} \int F dt$$

Relaxation time τ : $\delta k \rightarrow 0$

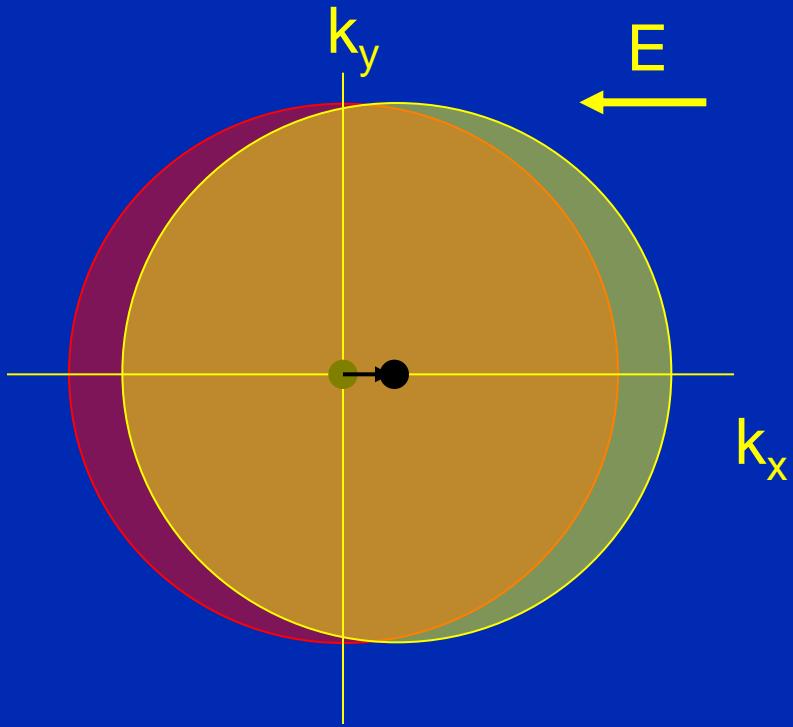
e-p,e-e scattering (appendix J, Ch. 10)
Impurity scattering

$$\delta k = \frac{1}{\hbar} F \tau$$

Equation of motion:

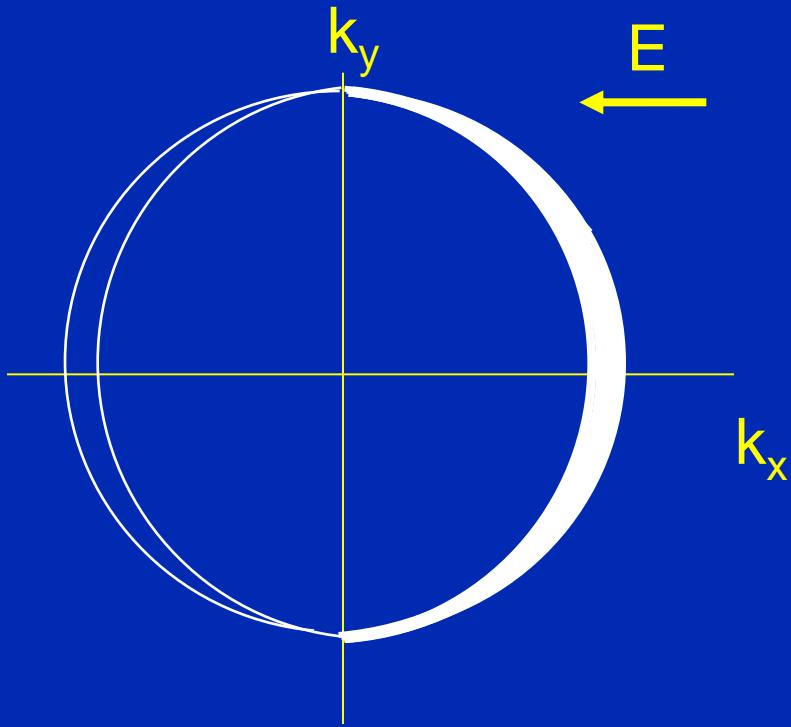
$$\hbar \left(\frac{d}{dt} + \frac{1}{\tau} \right) k = F$$

Electrical conductivity



$$\delta k = \frac{1}{\hbar} F \tau$$

Electrical conductivity



$$\delta k = \frac{1}{\hbar} F \tau$$

Ohm's law

$$\delta k = \frac{1}{\hbar} F \tau$$

$$\delta k = \frac{q E \tau}{\hbar};$$

$$v_{\text{drift}} = \delta v = \frac{q E \tau}{m}$$

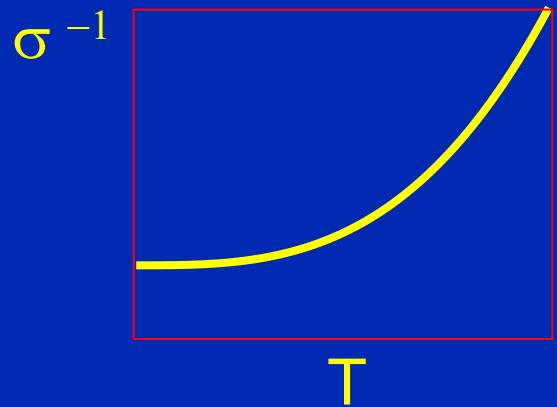
$$F = qE$$

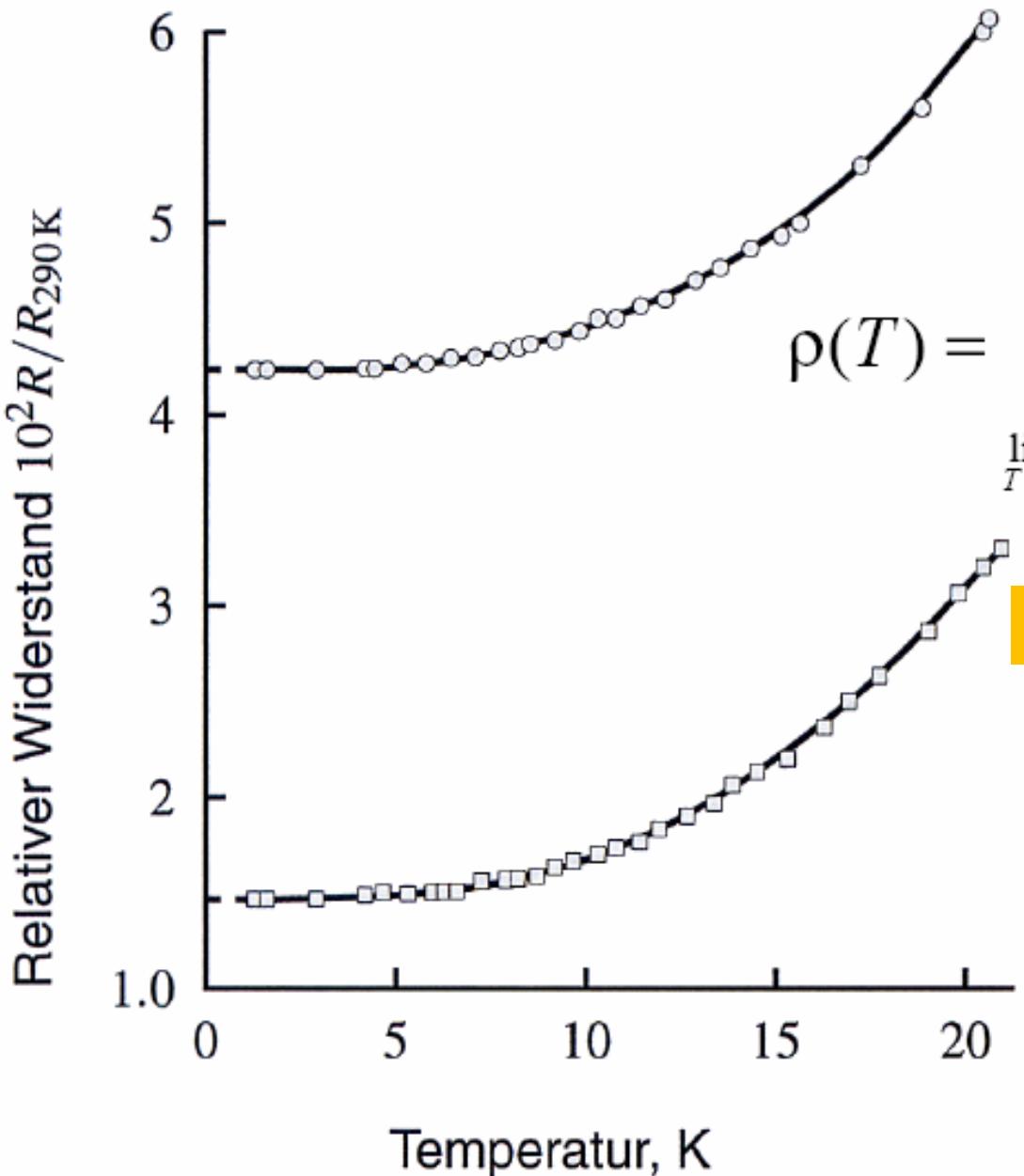
Current density $j = nq\delta v = \frac{n e^2 \tau E}{m}$

Ohm's law $\sigma = \frac{j}{E} = \frac{n e^2 \tau}{m}$

$$\left. \begin{array}{l} \sigma \sim 10^7 \Omega^{-1} m^{-1} \\ n \sim 10^{28} m^{-3} \end{array} \right\} \quad \tau \sim 10 \text{ fs} \quad \Rightarrow \quad \ell = v_F \tau \sim 10 \text{ nm}$$

Very pure metals, low T: $\ell > 1 \text{ cm} !!$





Resistance of potassium at 20 K for two different samples.

$$\rho(T) = \underbrace{\rho_{\text{Ph}}(T)}_{\lim_{T \rightarrow 0} \rho_{\text{Ph}}(T) = 0} + \underbrace{\rho_i(T)}_{\lim_{T \rightarrow 0} \rho_i(T) = \rho_i(0)}$$

Residual resistance ration RRR

$$\frac{\rho(T = 293 \text{ K})}{\rho(T = 0 \text{ K})} = 1.1 \dots 1000$$

An impurity causes a residual resistance of:

$1 \cdot 10^{-6} \Omega \text{cm}$ pro Atomprozent der Verunreinigung.

Temperature dependence of the electric resistance

(1) Very low temperatures

$$\rho_{\text{Phonon}} \propto T^5 / \Theta_D^5$$

Umklapp-processes are not possible
for $T < 2K$ in the case of K

(2) Low temperatures

Umlkapp-processes become possible

$\langle n \rangle \propto e^{-\Theta_U/T}$ characteristic Umklapp-temperature

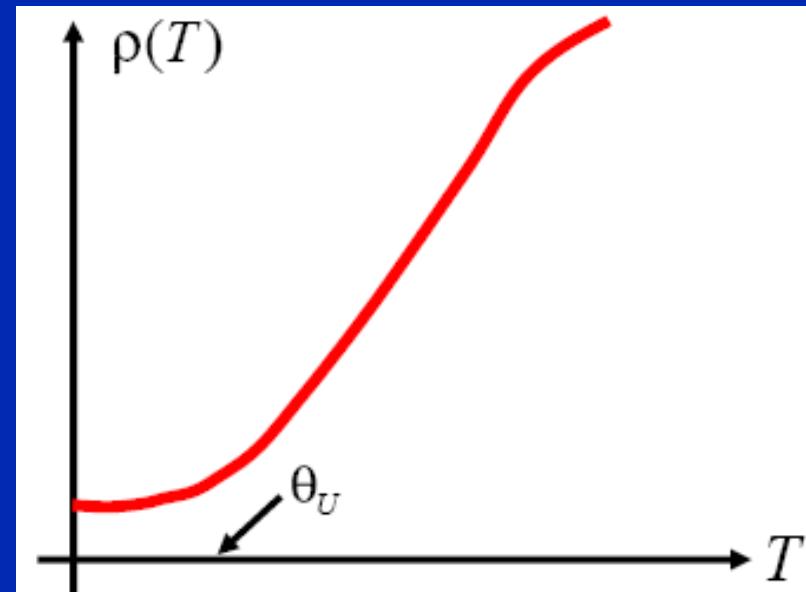
Potassium: $\Theta_U=23\text{K}$; $\Theta_D=91\text{K}$

(3) High temperatures

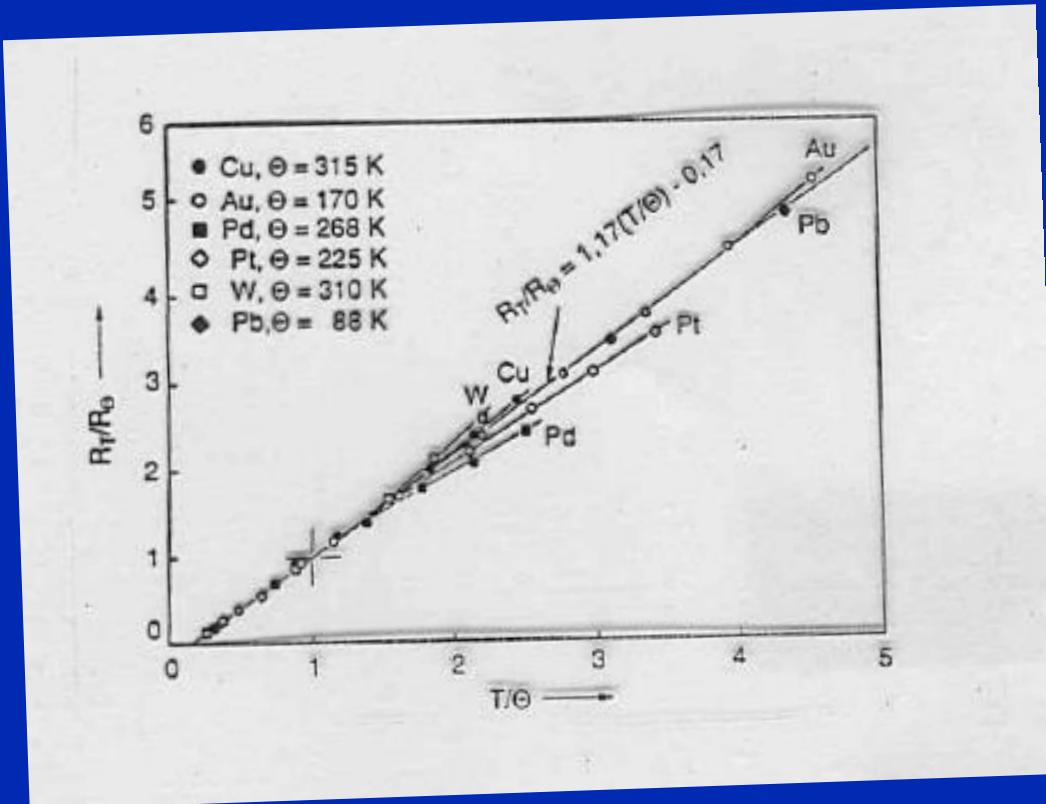
$$\langle n \rangle \propto T \rightarrow \rho_{\text{Phonon}} \propto T$$

(4) Very high temperatures

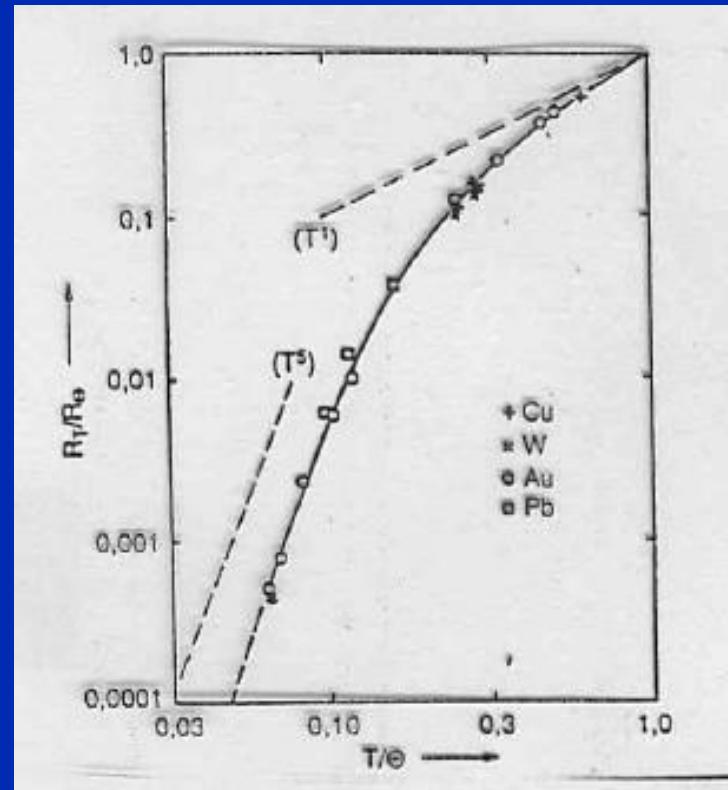
saturation as the phonon scattering cannot lead to mean free path-lengths shorter than the lattice constant (Ioffe rule)



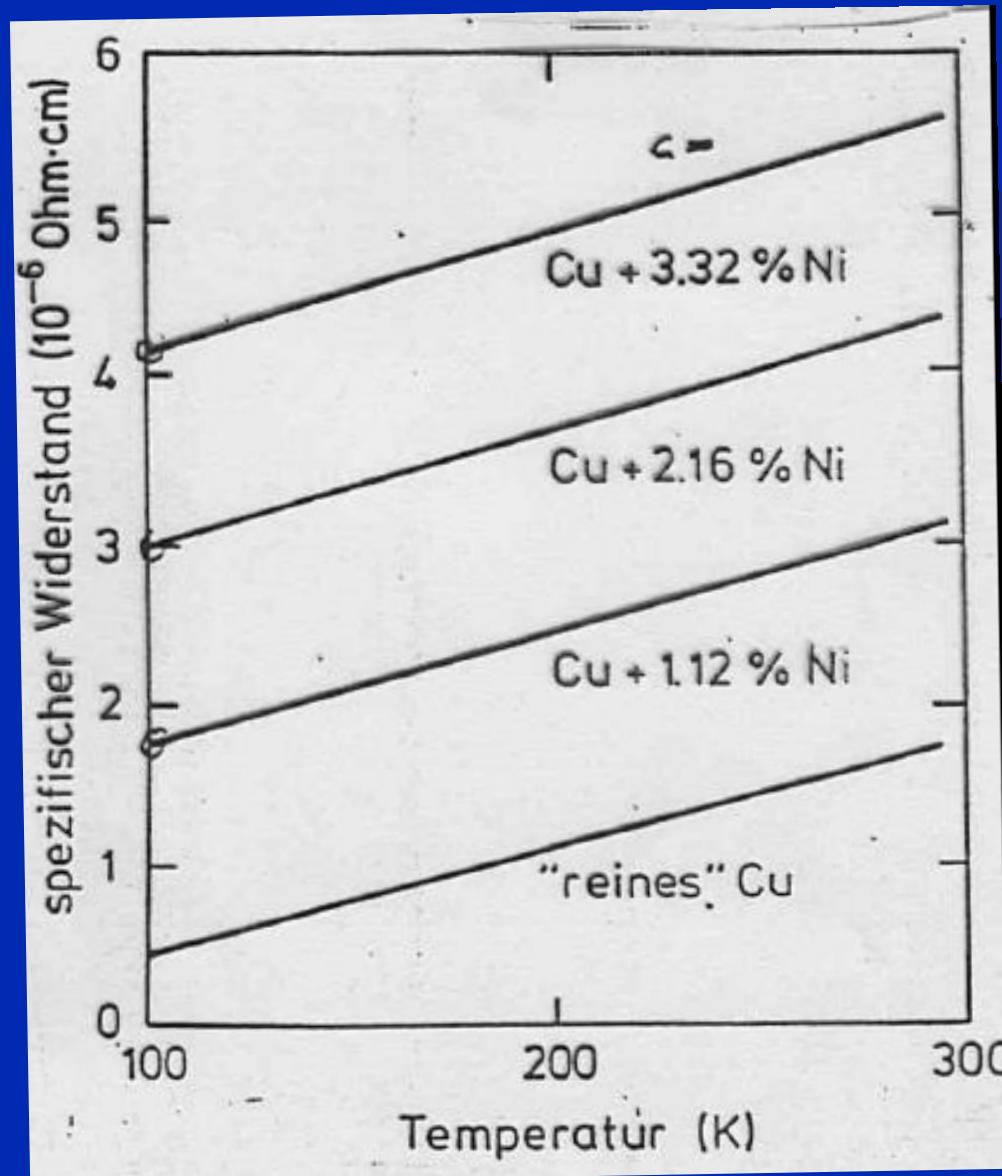
The resistance roughly scales with the Debye temperature !



Transition from T^5 to T behavior !



Matthiessen rule: alloys \rightarrow vertical displacement of $\rho(T)$ curves



Quantized Conductance of Point Contacts in a Two-Dimensional Electron Gas

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(Received 31 December 1987)

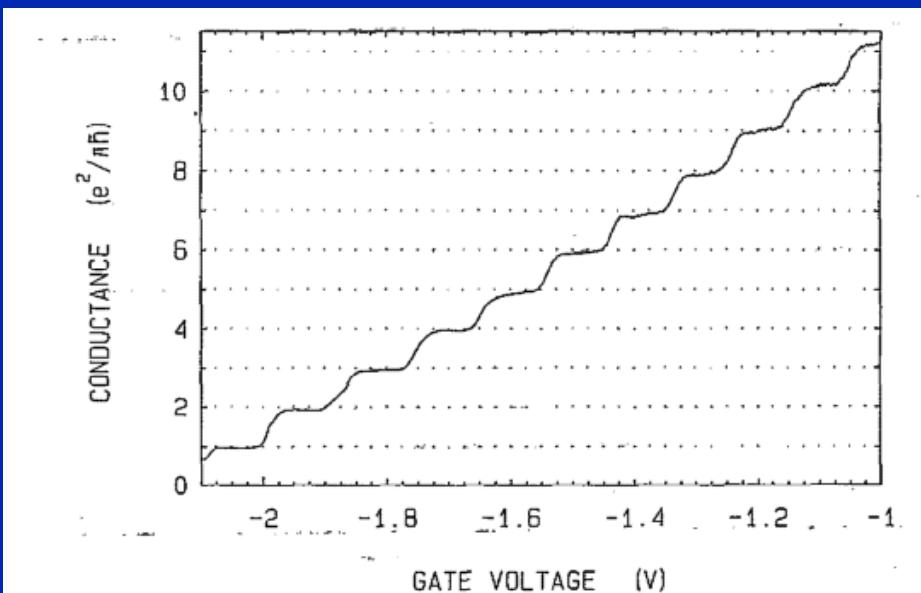


FIG. 2. Point-contact conductance as a function of gate voltage, obtained from the data of Fig. 1 after subtraction of the lead resistance. The conductance shows plateaus at multiples of $e^2/\pi\hbar$.