



Drude (1900) → Sommerfeld (1927) → Bloch (1940)

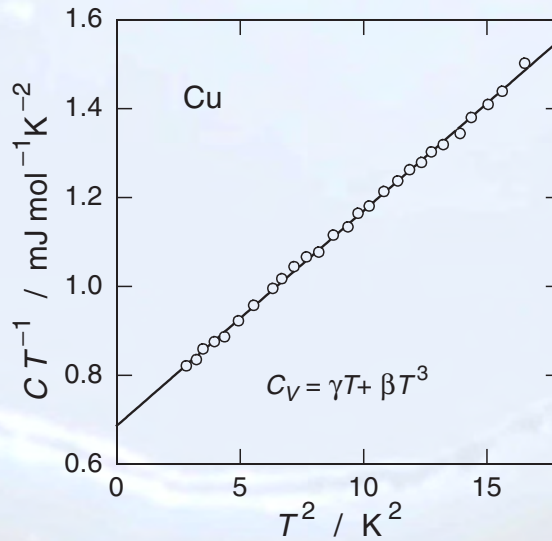
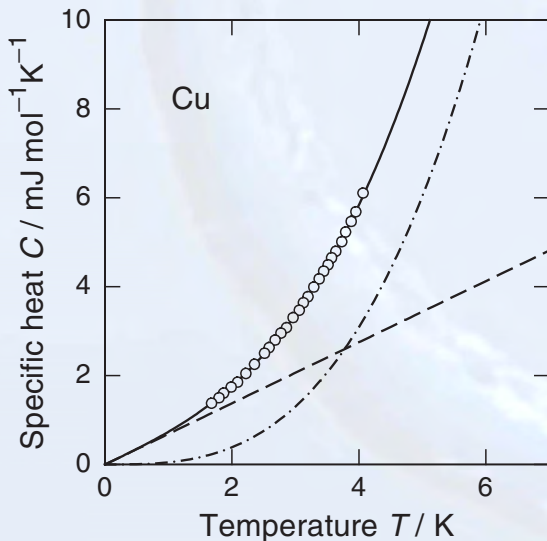
7.1 Specific heat

a) simple metals → free electrons gas

free electrons gas: $c_V = \gamma T + \beta T^3$

/ phonons

electrons: $\gamma = \frac{\pi^2 n k_B^2}{2E_F} \propto n^{1/3} m_{th}^*$



- ▶ electrons dominate below ~ 4 K
- ▶ very good qualitative agreement



7.1 Specific Heat



Element	γ_{exp}	γ_{theo}	m_{th}^*/m	Element	γ_{exp}	γ_{theo}	m_{th}^*/m
Ag	0.64	0.64	1.00	Cu	0.69	0.50	1.37
Al	1.35	0.91	1.48	Ga	0.60	1.02	0.59
Au	0.69	0.64	1.08	In	1.66	1.26	1.31
Ba	2.70	1.95	1.38	K	2.08	1.75	1.19
Be	0.17	0.49	0.35	Li	1.65	0.75	2.19
Ca	2.73	1.52	1.80	Mg	1.26	1.00	1.26
Cd	0.69	0.95	0.73	Na	1.38	1.3	1.22
Cs	3.97	2.73	1.46	Pb	2.99	1.50	1.99

- good qualitative agreement for simple metals
- $\gamma_{\text{exp}}/\gamma_{\text{theo}} = m_{\text{th}}^*/m$ for quantitative agreement

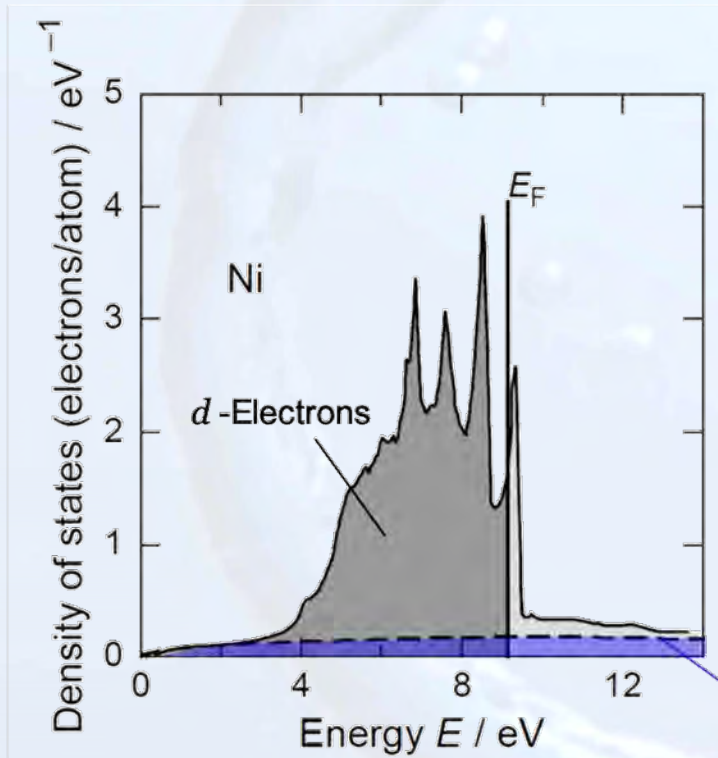


but: **transition series metals**

example nickel: $m_{th}^* \approx 15 m$ \longrightarrow reason is **d-electrons contribute**, which are not (completely) free

\longrightarrow involved in **covalent bond**, highly oriented

\longrightarrow **no spherical** Fermi surface



► **d-electrons** with large density of state **dominate** at E_F

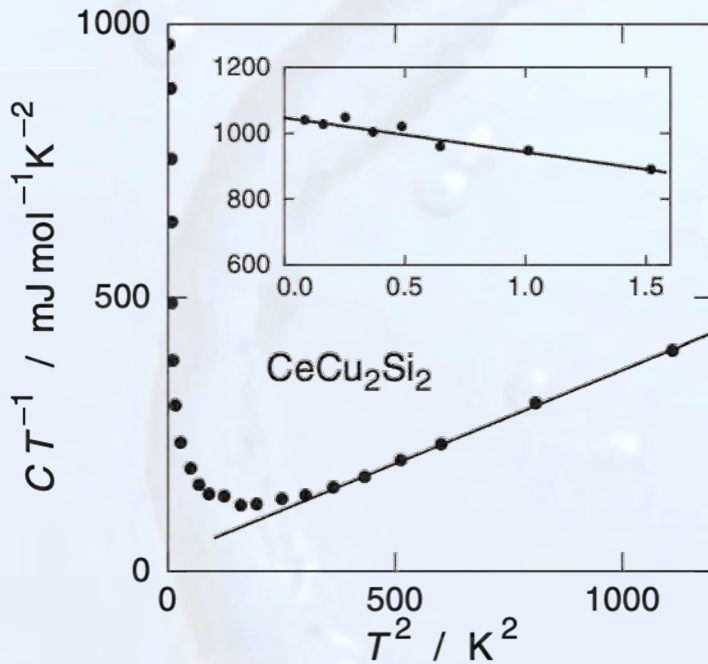
► **d-electrons** are localized

\longleftarrow s-Electrons



b) metal with **heavy** electrons

examples: CeCu2Si2 cer electronic configuration $[\text{Xe}] 5d^1 4f^1 6s^2$



→ γ not constant below 15 K

$\gamma = C_{el}/T$

- $\gamma \approx 30 \text{ mJ (mol K}^2\text{)}^{-1}$ extrapolated from high T
- $\gamma \approx 1050 \text{ mJ (mol K}^2\text{)}^{-1}$ at low temperatures $T \rightarrow 0$

→ **4f** electrons are localized at **high T** and form a **conduction band** at **low T**

→ effective mass: $m^* \approx 100 m_e$

- ▶ $T > 15 \text{ K}$, $D(E)$ and m^* are constant
- ▶ $T < 15 \text{ K}$, C/T **increase** strongly with **decreasing** temperature



Heavy fermion systems

- ▶ interesting class of solids with **strongly correlated electrons**
- ▶ **effective masses m^*** up to **2000 m_e** observed
- ▶ origin: interaction with localized spins

$$\chi = \mu_0 \mu_B^2 D(E_F) \propto n^{1/3} m^*$$

Wilson ratio:

$$R = \frac{\chi}{\gamma} \frac{\pi^2 k_B^2}{\mu_0 \mu_{\text{eff}}^2}$$

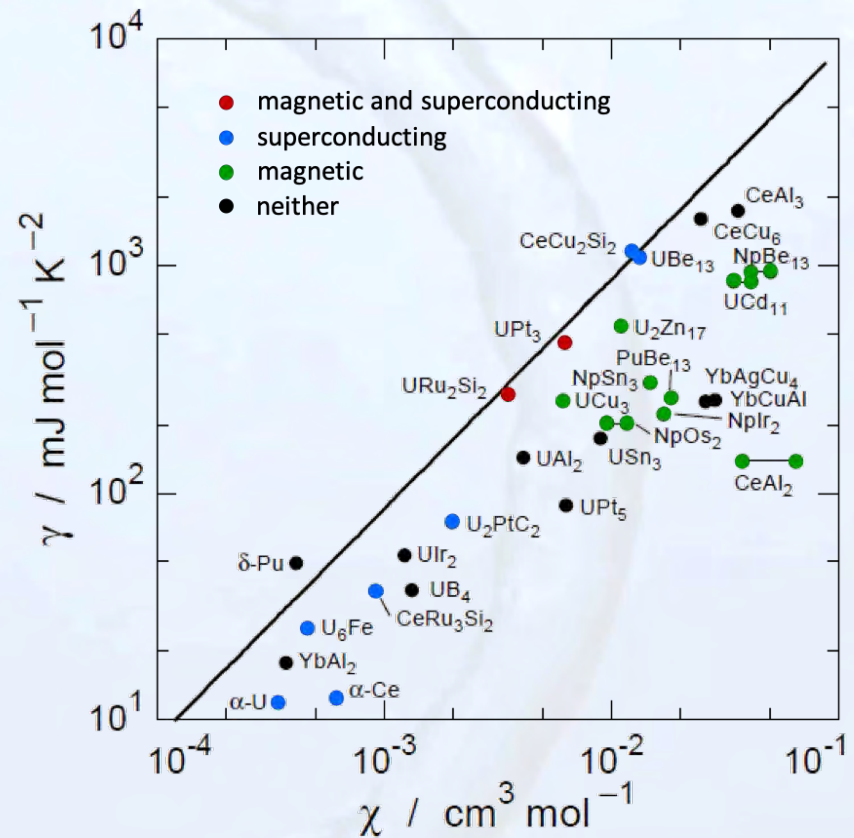
$$\gamma \propto n^{1/3} m_{\text{th}}^*$$

important: **Fermi liquid theory**

$$C = \frac{m^*}{m} C_{\text{FG}} = \left(1 + \frac{1}{3} F_1\right) C_{\text{FG}}$$

analogy to ^3He reaches even further

→ some heavy fermion systems show **unconventional superconductivity** ($S \neq 0$) : UPt_3 , URu_2Si_2 ...





metals, no superconductors, no semiconductors

Boltzmann equation \longleftrightarrow kinetic gas theory

- ▶ starting point: equilibrium distribution without external fields $f_0(\mathbf{k})$
Fermi-Dirac distribution
- ▶ with field: stationary non-equilibrium value of $f(\mathbf{k}, \mathbf{r}, t)$
- ▶ expand $f_0(\mathbf{k}) - f(\mathbf{k}, \mathbf{r}, t)$ in linear order + relaxation ansatz for collisions

→ linearized Boltzmann equation $f(\mathbf{k}) \approx f_0(\mathbf{k}) + \frac{e\tau(\mathbf{k})}{\hbar} \boldsymbol{\varepsilon} \cdot \frac{\partial f_0(\mathbf{k})}{\partial \mathbf{k}}$
scattering time electric field

→ $j_x = -e \int D(k) v_x(k) f(k) dk = -\frac{e}{\pi^2} \int k^2 v_x(k) f(k) dk$

→ $\sigma = \frac{1}{3} e^2 D(E_F) v_F^2 \tau(E_F) \longrightarrow \sigma = \frac{ne^2}{m} \tau(E_F)$



scattering time determined by:

- ▶ defect scattering
- ▶ phonon scattering
- ▶ magnon scattering (in ferromagnets)
- ▶ electron-electron scattering (can be neglected in most cases)

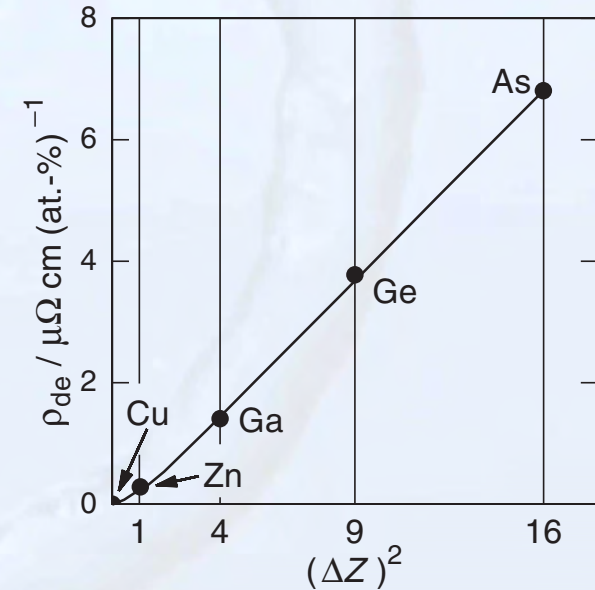
a) defect scattering

- ↗ local charge density variations
- ↘ local strain fields (less important)

Local charge variations

- ▶ **Rutherford scattering** on ionic cores of impurity atoms
- ▶ scattering **cross section** : $\sigma_{\text{cross}} \propto (\Delta Z)^2$
- ▶ resistivity $\rho_D \propto (\Delta Z)^2$

- ▶ residual resistance of copper with 1 at% impurities with **different valence** electrons **configurations**
- ▶ agrees well with: $\rho_D \propto (\Delta Z)^2$





Concentration dependence

example: binary mixture A_xB_{1-x} \longrightarrow **mixing increases the resistance**

average potential $U_0 = xU_A + (1 - x)U_B$

deviations at atoms A and B: $(U_0 - U_A) = (1 - x)(U_B - U_A)$

$(U_0 - U_B) = x(U_A - U_B)$

scattering probability:

$$w_A = (1 - x)^2 \left| \int \psi^*(\mathbf{k}) (U_B - U_A) \psi(\mathbf{k}') d^3k' \right|^2$$

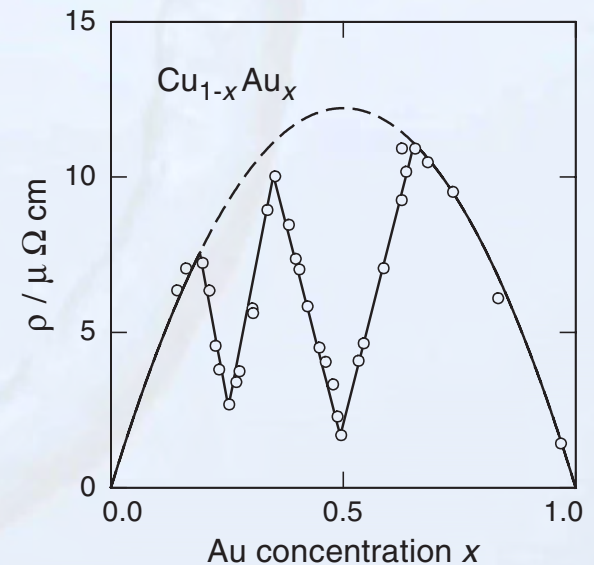
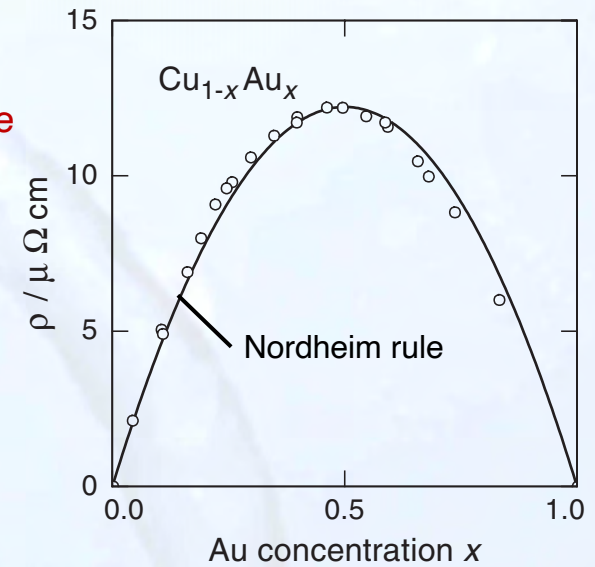
$$w_B = x^2 \left| \int \psi^*(\mathbf{k}) (U_A - U_B) \psi(\mathbf{k}') d^3k' \right|^2 = \frac{x^2}{(1 - x)^2} w_A$$

resistivity: $\rho_{de} \propto xw_A + (1 - x)w_B$

Nordheim rule: $\rho_{de} \propto x(1 - x)$

► $Cu_{1-x}Au_x$ **data agree** with Nordheim rule

► **tempering** of sample \longrightarrow formation of **ordered compounds**

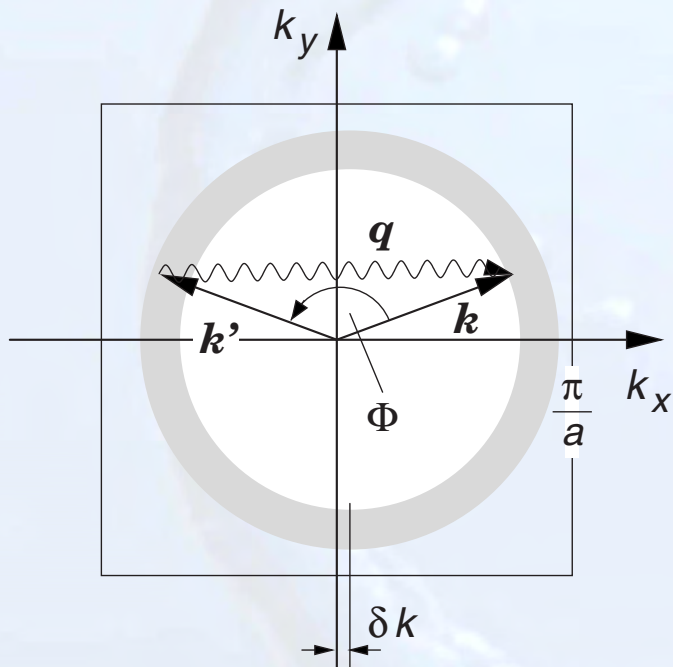




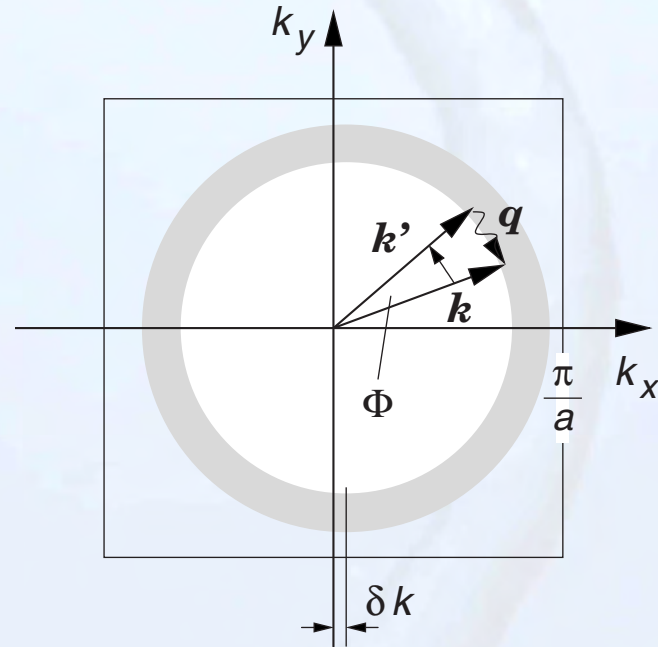
Electron-Phonon scattering

only electrons at the **Fermi surface** can participate $|\mathbf{k}| \approx |\mathbf{k}'|$.

high temperatures



low temperatures





Electron-Phonon scattering

a) high temperatures ($T > \theta_D$) $l \propto \tau v_F \longrightarrow l^{-1} \propto n_{ph} \propto T \longrightarrow \rho \propto T$

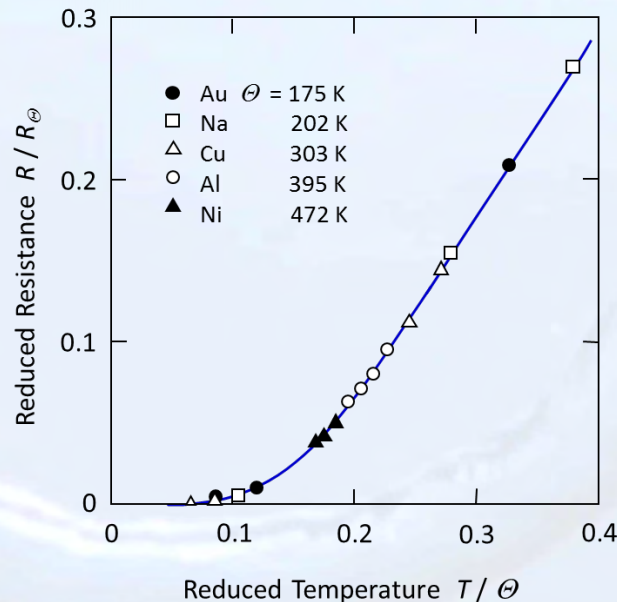
b) intermediate temperatures ($T < \theta_D$)

- ▶ cross-section depends on temperature $e_0^2 \propto \omega/\theta \propto T/\theta$
 - ▶ number of scattering centers (phonons) reduces $n_{ph} \propto T^2/\theta^2$
 - ▶ effectiveness of scattering process goes down $(1 - \cos \phi) \propto T^2/\theta^2$
- $\left. \begin{array}{l} \phantom{\text{▶}} \\ \phantom{\text{▶}} \\ \phantom{\text{▶}} \end{array} \right\} \tau^{-1} \propto n\Sigma \propto T^3/\theta^3$

→ $\rho_{ph} \propto \tau_{eff}^{-1} \propto \left(\frac{T}{\theta}\right)^5$

Bloch-Grüneisen law

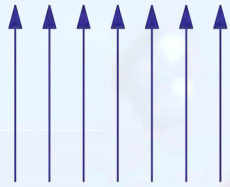
- ▶ reduced plot → material independent
- ▶ defect scattering subtracted
- ▶ good agreement with Bloch-Grüneisen



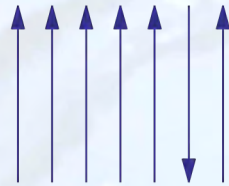
Small angles ϕ
 $\rightarrow (1 - \cos \phi) \propto \phi^2$
 with $q \approx \left(\frac{T}{\theta}\right) q_{max}$
 and $k_F = \frac{\pi}{a}$
 $\phi \approx \frac{q}{k_F} \approx \frac{T}{\theta} \frac{\pi}{a} \frac{1}{k_F} = \frac{T}{\theta}$
 $\rightarrow \phi^2 \propto \left(\frac{T}{\theta}\right)^2$



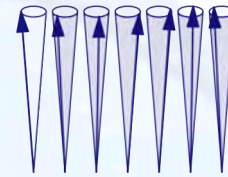
Electron magnon scattering
spin waves in ferromagnets



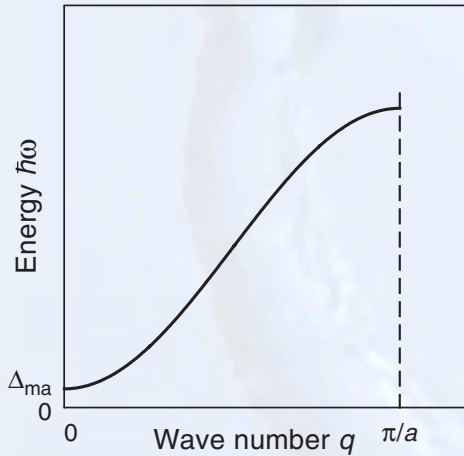
ground state



single spin excitation
(cost too much energy $E \approx k_B T_C$)



collective excitations
spin waves $\hat{=}$ magnons



dispersion curve: (Terbium)

$$\hbar\omega = 2JS [1 - \cos(qa)] + \Delta_{ma}$$

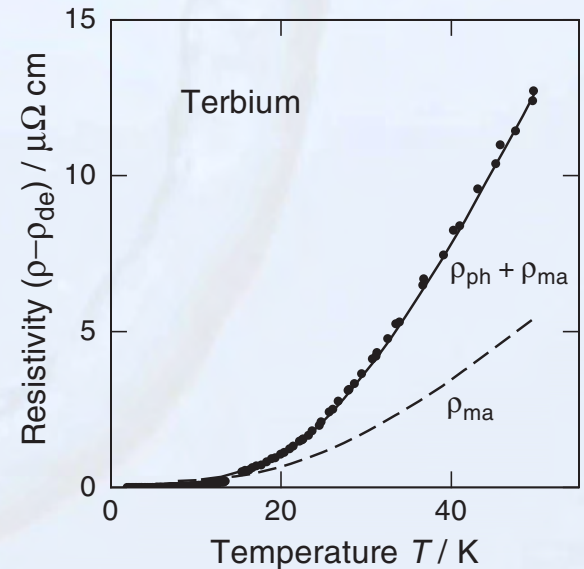
anisotropic exchange
interaction of spins

$$\Delta_{ma}/k_B \approx 20 \text{ K}$$

resistivity: (Terbium)

$$\tau^{-1} \propto n_{ma} \propto e^{-\Delta_{ma}/k_B T}$$

$$\longrightarrow \rho_{ma} \propto e^{-\Delta_{ma}/k_B T}$$





Thermal conductivity of a free Fermi gas

$$\Lambda_{el} = \frac{1}{3} c_V^{el} v \ell = \frac{1}{3} \frac{\pi^2 n k_B^2 T}{m v_F^2} v_F \ell$$

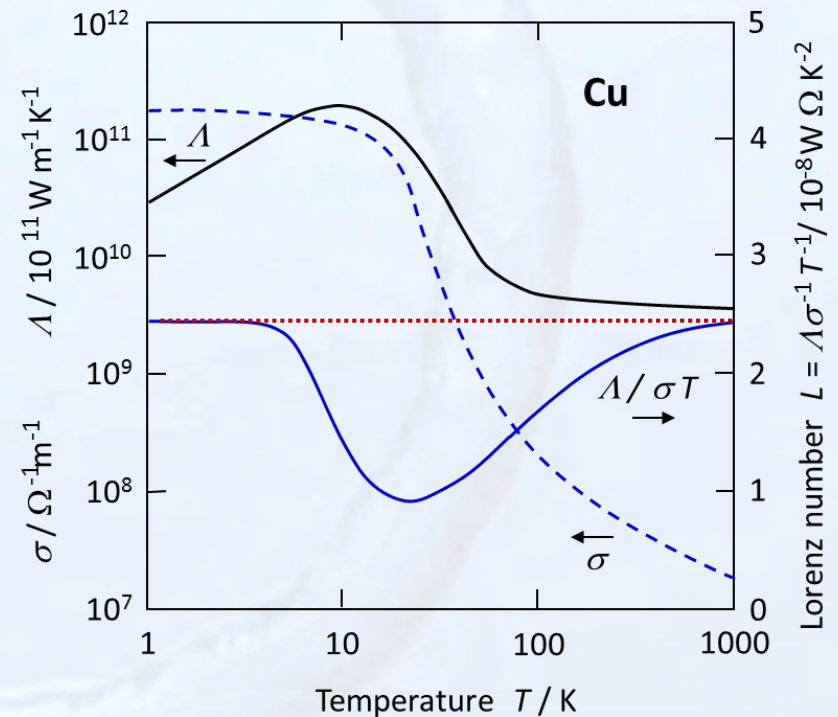
if **scattering processes** are **identical** for electrical and thermal transport

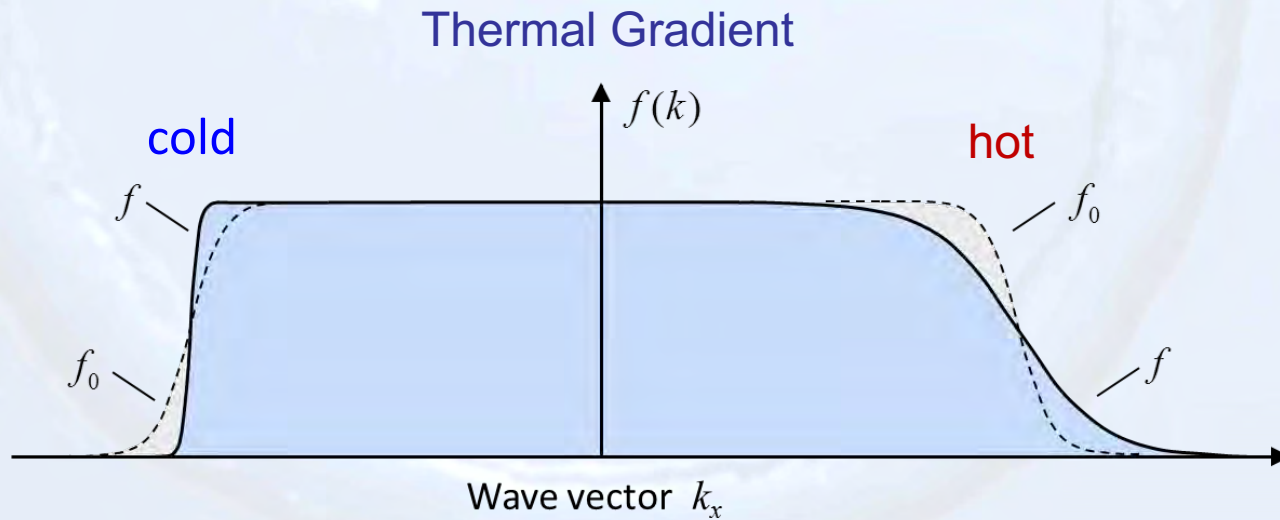
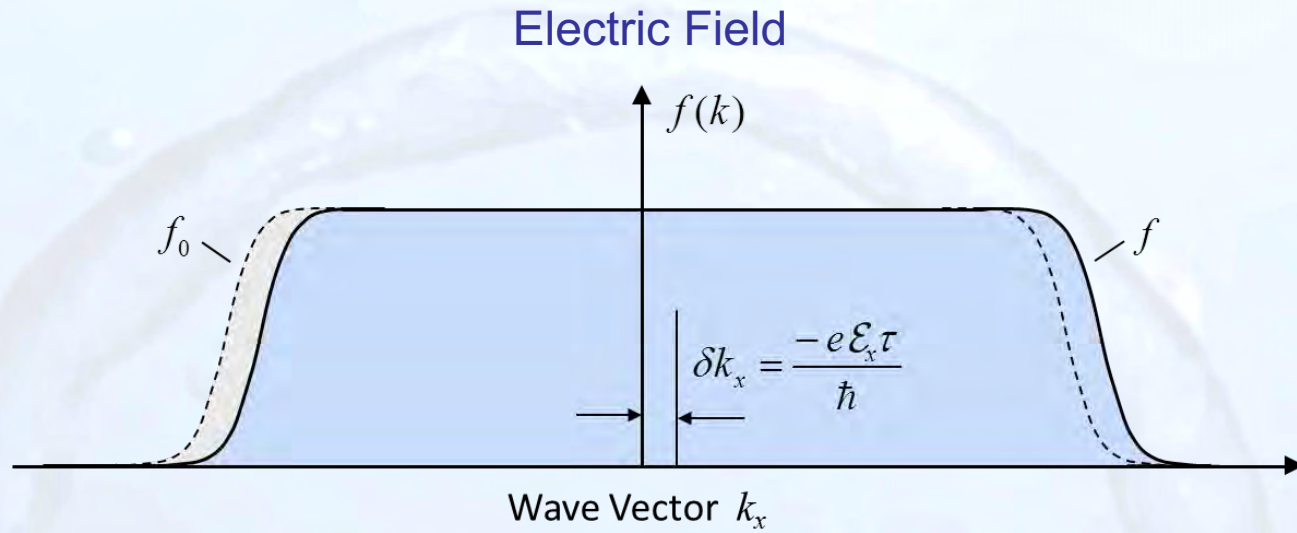
➔ Widemann-Franz law

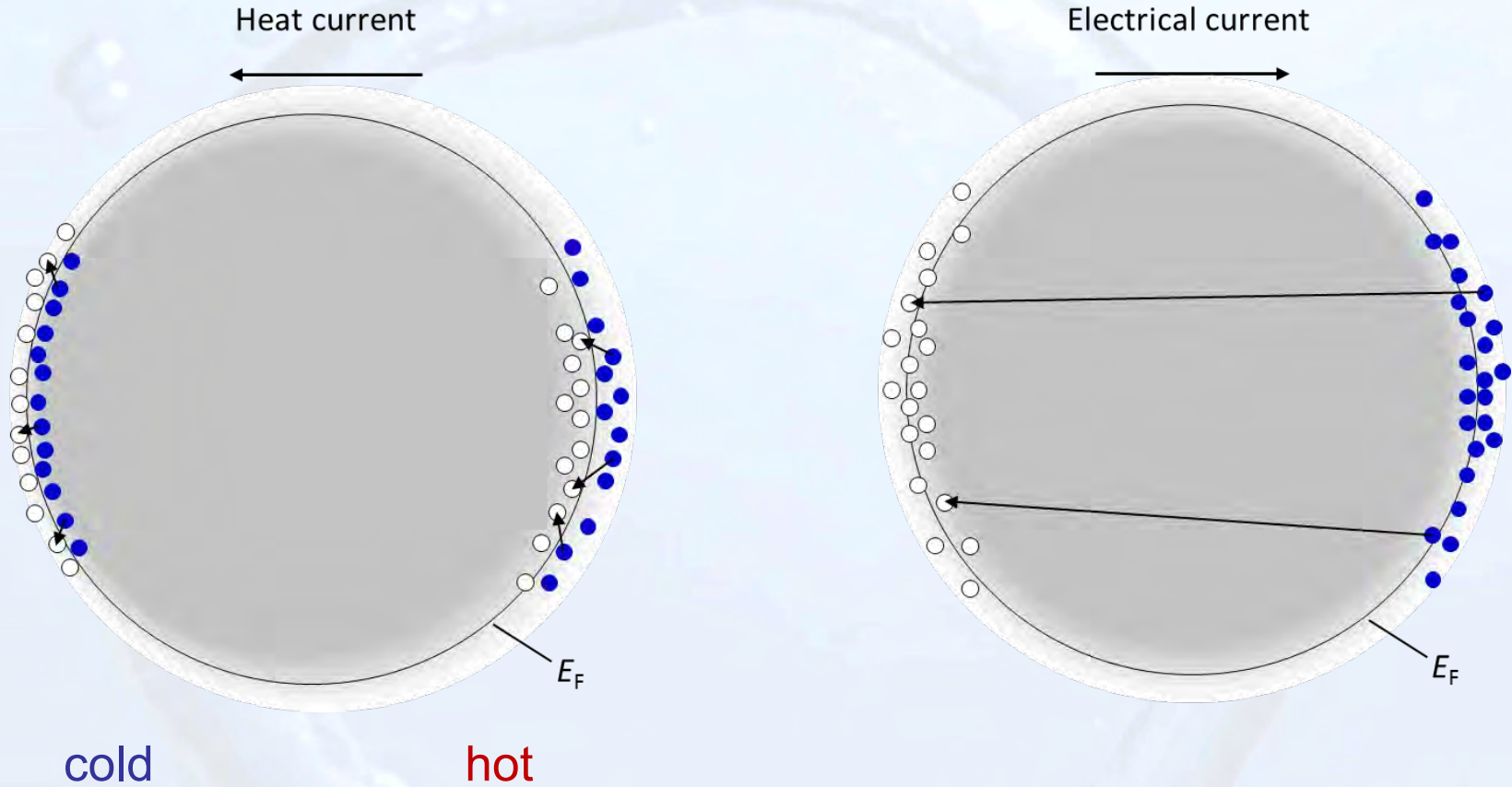
$$\frac{\Lambda_{el}}{\sigma} = \frac{\pi^2}{3} \left(\frac{k_B}{e} \right)^2 T = \mathcal{L} T$$

$$\mathcal{L} = 2.45 \times 10^{-8} \text{ V}^2 \text{ K}^{-2}$$

- ▶ Lorenz number depends on temperature
- ▶ works well at **very low** and **very high temperatures**





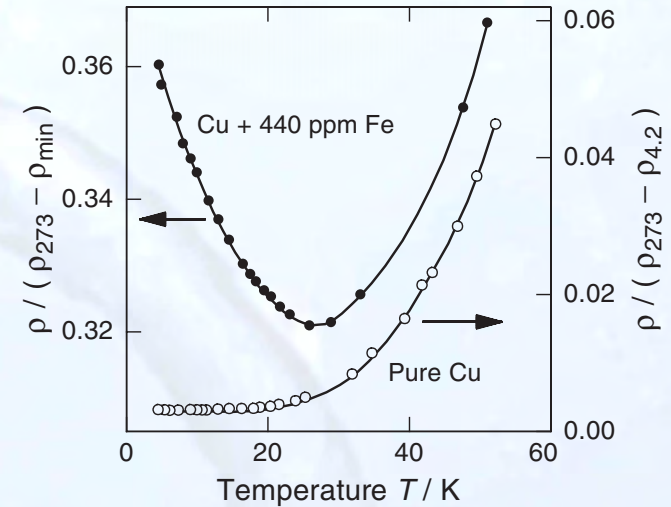




1930 Meissner and Voigt observe a **resistance minimum** for Au and Cu with **magnetic impurities**

- ▶ example: Cu + 440 ppm Fe
- ▶ resistance minimum at 27 K

1964 explanation by **Kondo** by spin dependent scattering of electrons on magnetic impurities



a) Influence of conduction electrons on localized magnetic moments

example: **d-levels** of **transition metals** in simple metals

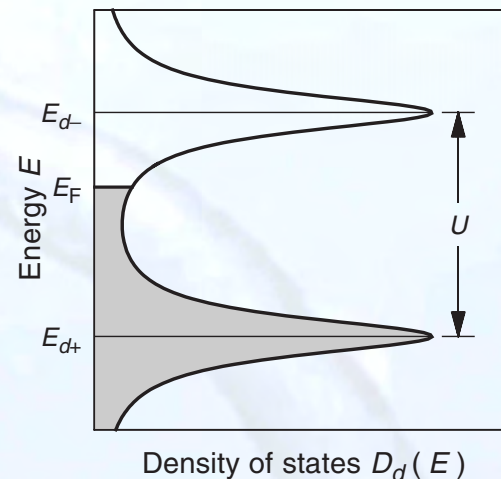
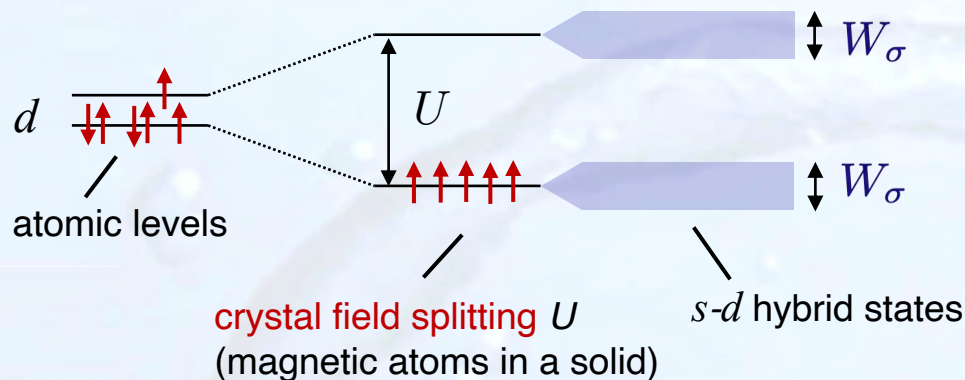
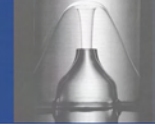
- ▶ **d-d** interaction \longrightarrow **splitting** and **polarization** of **d-levels**, because of **crystal field**
- ▶ interaction of **d-electrons** with **conduction electrons** (*s*) \longrightarrow hybrid states

width determined by **$s \leftrightarrow d$ transition rate**

matrix element

\longrightarrow golden rule:
$$\frac{W_{\sigma}}{\hbar} = \frac{\pi}{\hbar} \nu^2 D_s(E_{d\sigma})$$

Density of states of *s*-electrons at $E_{d\sigma}$



AgMn:

$$W_\sigma = 0.5 \text{ eV}$$

$$U = 4.8 \text{ eV}$$

localized moments remain if interaction
(W_σ) is not too strong $W_\sigma < U$

spin-dependent interaction of s electrons with d -electrons

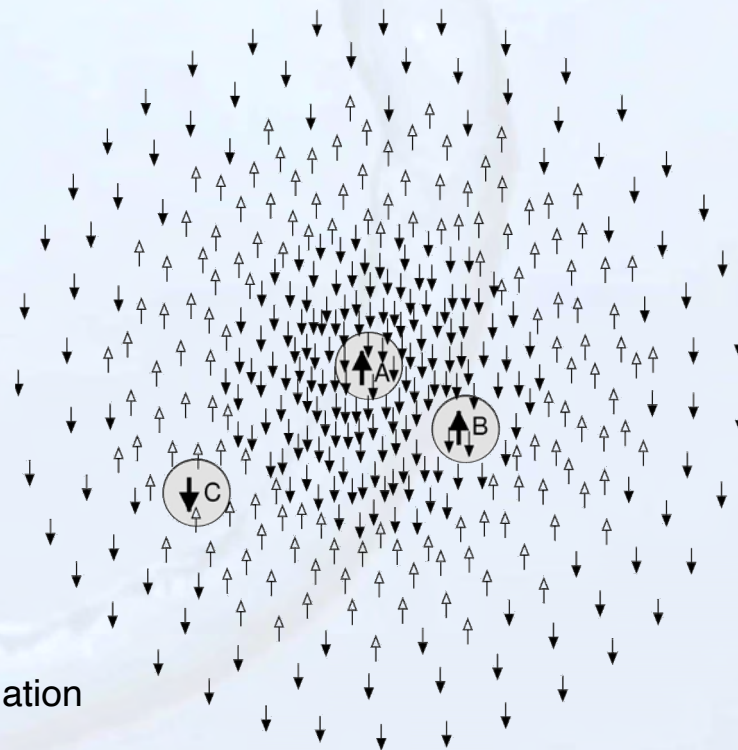
$$J \approx -\mathcal{V}^2/U$$

position of localized spin

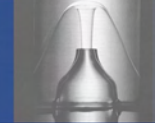
$$\mathcal{H}_{sd} = -J \mathbf{S} \cdot \mathbf{s} \delta(\mathbf{r} - \mathbf{R}) \quad (\star)$$

localized spins

conduction electron



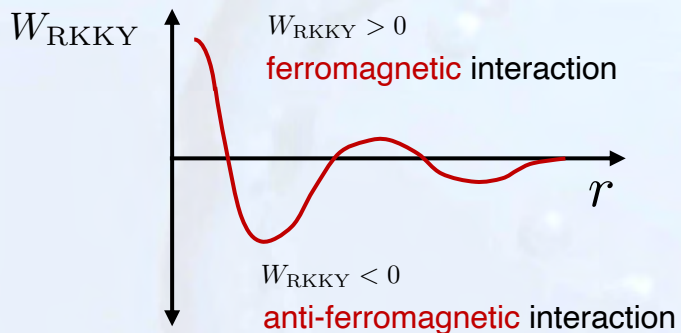
Friedel oscillation



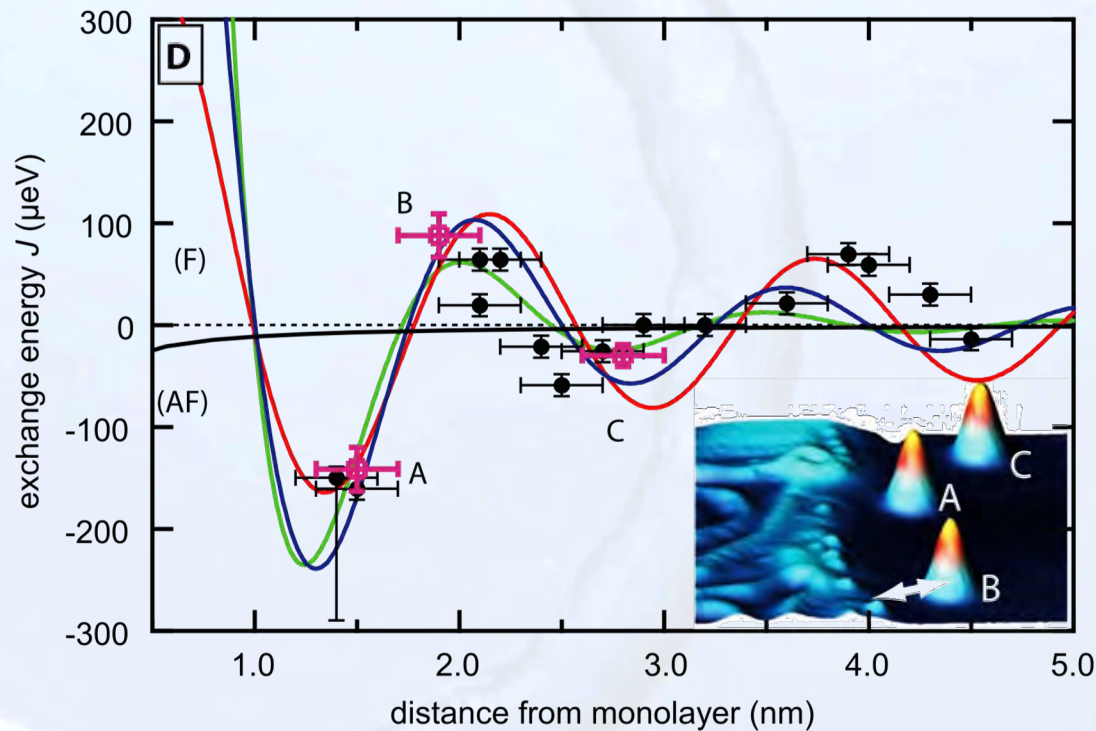
Indirect exchange interaction (*d-d*) — RKKY interaction

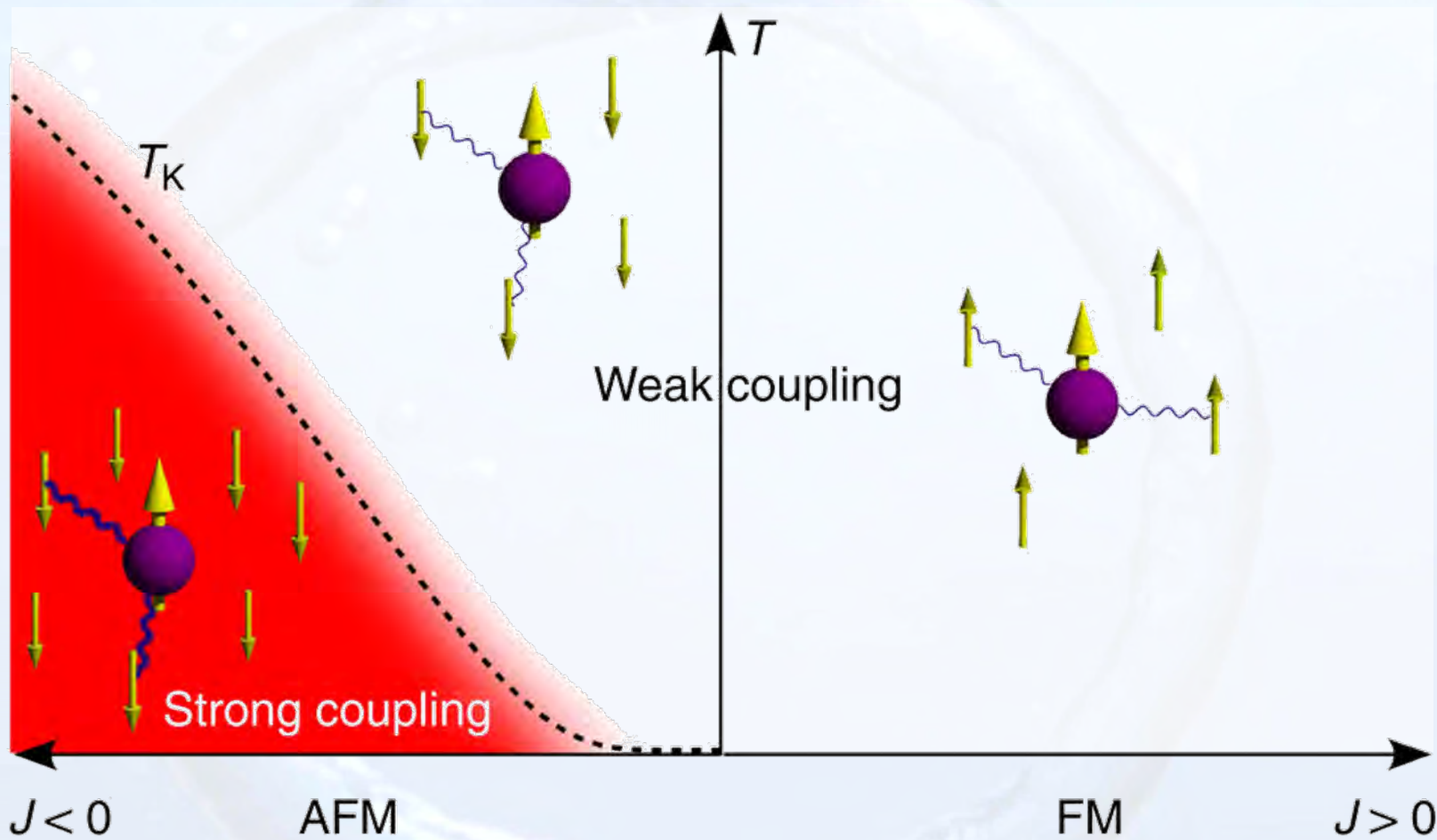
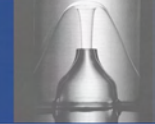
Ruderman
Kittel
Kusuya
Yosida

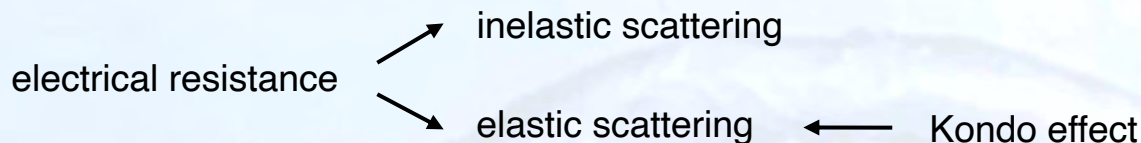
$$W_{\text{RKKY}} \propto J_{\text{RKKY}}^2 D(E_F) \frac{\cos(2k_F r)}{r^3}$$



MFM measurement using magnetic moments on a noble metal surface







consider (★) in “N representation” $\left| \mathcal{H}_{sd} = -J \mathbf{S} \cdot \mathbf{s} \delta(\mathbf{r} - \mathbf{R}) \right.$

$$\mathcal{H}_{sd} = -J \sum_{\mathbf{k}, \mathbf{k}'} S_z (c_{\mathbf{k}'\uparrow}^+ c_{\mathbf{k}\uparrow} - c_{\mathbf{k}'\downarrow}^+ c_{\mathbf{k}\downarrow}) + S_+ c_{\mathbf{k}'\downarrow}^+ c_{\mathbf{k}\uparrow} + S_- c_{\mathbf{k}'\uparrow}^+ c_{\mathbf{k}\downarrow}$$

const. < 0!

$c_{\mathbf{k}}^+$ creation operator
 $c_{\mathbf{k}}$ annihilation operator
 } act on conduction electrons

$S_+ = S_x + iS_y$
 $S_- = S_x - iS_y$
} spin states

\mathbf{k} wave vector of conduction electrons



Jun Kondo

Harmonic Oscillator

$$H = \hbar\omega \left(a^\dagger a + \frac{1}{2} \right) \quad \left| \quad H = \frac{p^2}{2m} + \frac{m\omega^2 Q^2}{2} \right.$$

$$Q = \sqrt{\frac{\hbar}{2m\omega}} (a + a^\dagger)$$

$$P = \frac{1}{i} \sqrt{\frac{m\hbar\omega}{2}} (a - a^\dagger)$$

$$a^\dagger = \sqrt{\frac{m\omega}{2\hbar}} \left(Q - \frac{i}{m\omega} P \right)$$

$$a = \sqrt{\frac{m\omega}{2\hbar}} \left(Q + \frac{i}{m\omega} P \right)$$

$$a^\dagger a = N$$

$$a^\dagger | \psi_n \rangle = \sqrt{n+1} | \psi_{n+1} \rangle$$



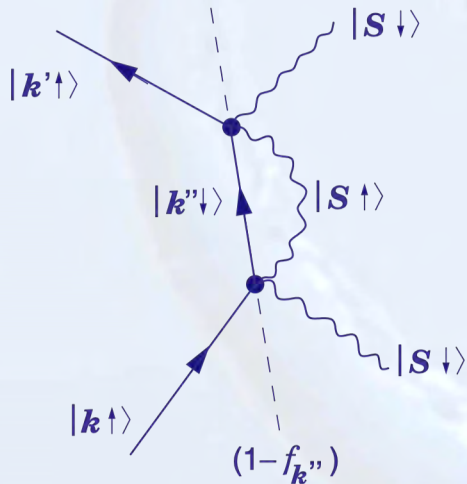
(i) **first-order** perturbation theory $|\mathbf{k} \uparrow\rangle \longrightarrow |\mathbf{k}' \uparrow\rangle$

$$t^{(1)} = \langle \mathbf{k}' \uparrow | \mathcal{H}_{sd} | \mathbf{k} \uparrow \rangle = -J S_z$$

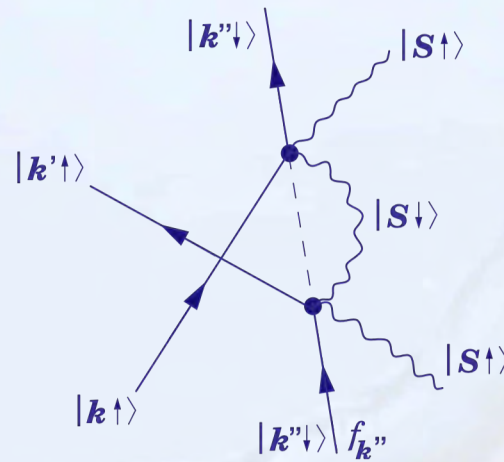
scattering amplitude \searrow temperature independent constant

(ii) **second-order** perturbation theory: two possible process $\begin{cases} \text{direct process (normal order)} \\ \text{exchange process (reverse order)} \end{cases}$

processes **without** spin flip



direct process



exchange process



all first and second order processes

