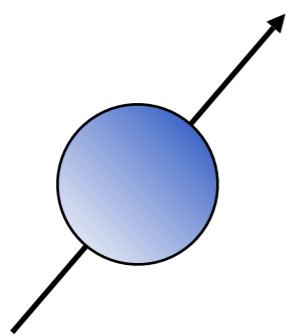
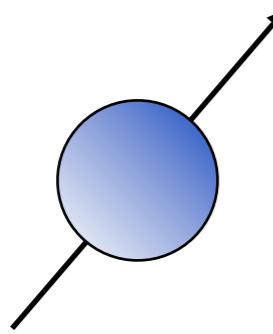


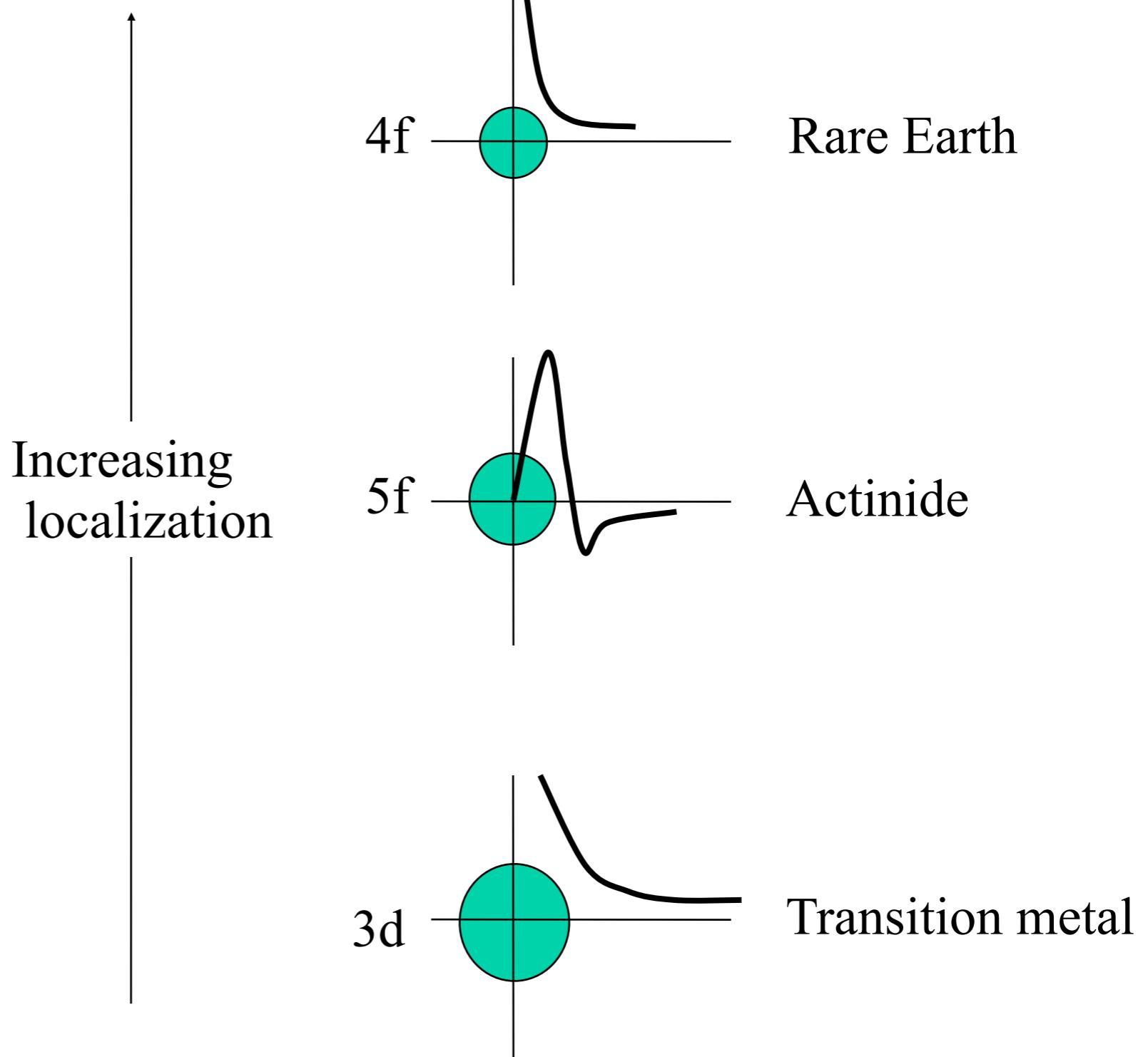
$$\vec{M} = g\mu_B \vec{S}$$



$$\overrightarrow{M} = g\mu_B \overrightarrow{S}$$



Localized Moment

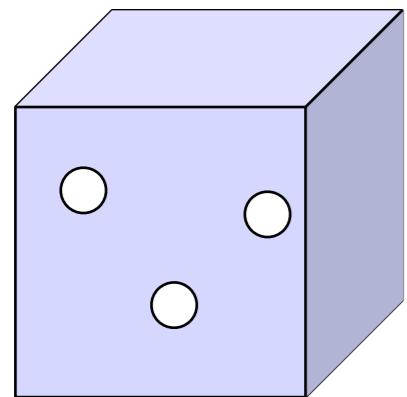
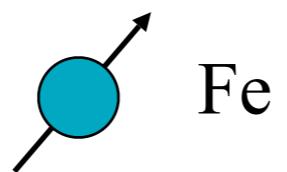


## Outline:

1. Chemistry of Kondo.
2. Key Properties.
3. Tour of Heavy Fermions.
4. Anderson model: the atomic limit.
5. Adiabaticity: the heavy Fermi liquid.
6. Scaling: the Kondo effect.
7. Doniach's Kondo Lattice Concept.

# Moment Formation

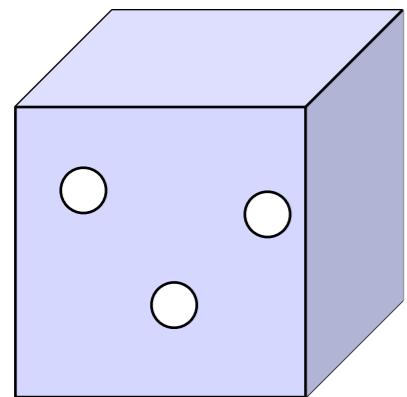
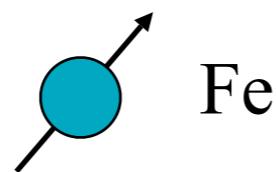
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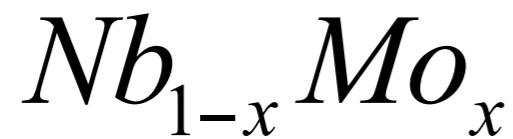
$$x < 0.4$$

# Moment Formation

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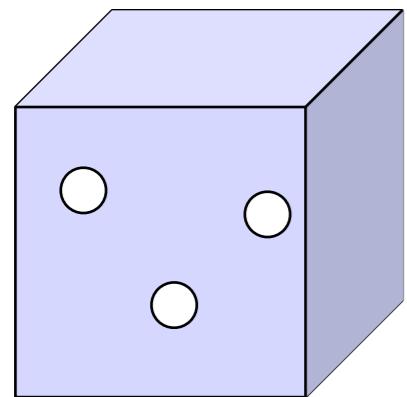
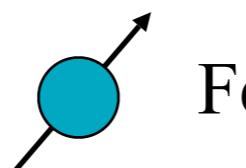
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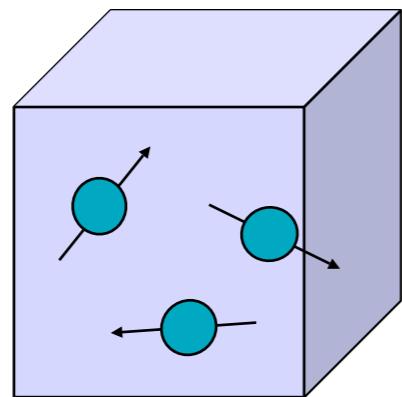
Clogston, Mathias et al, 1962

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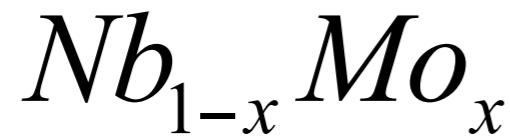
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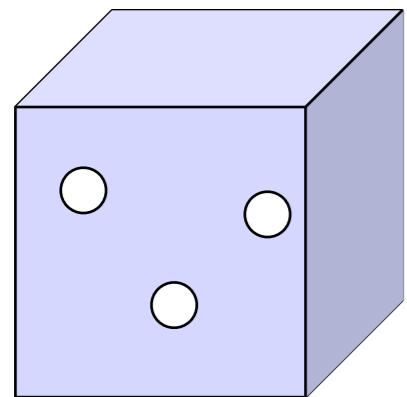
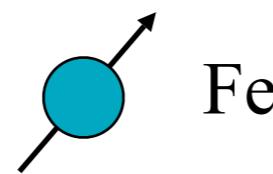
$x > 0.4$



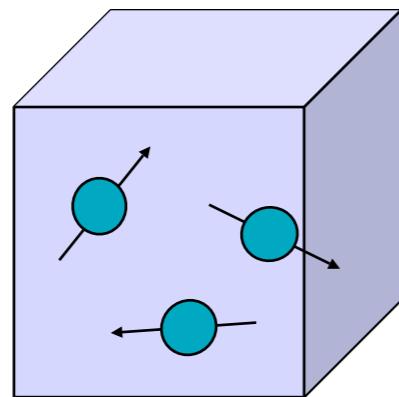
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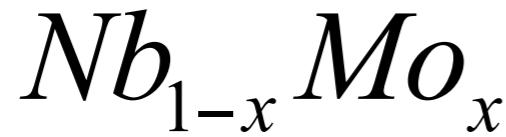
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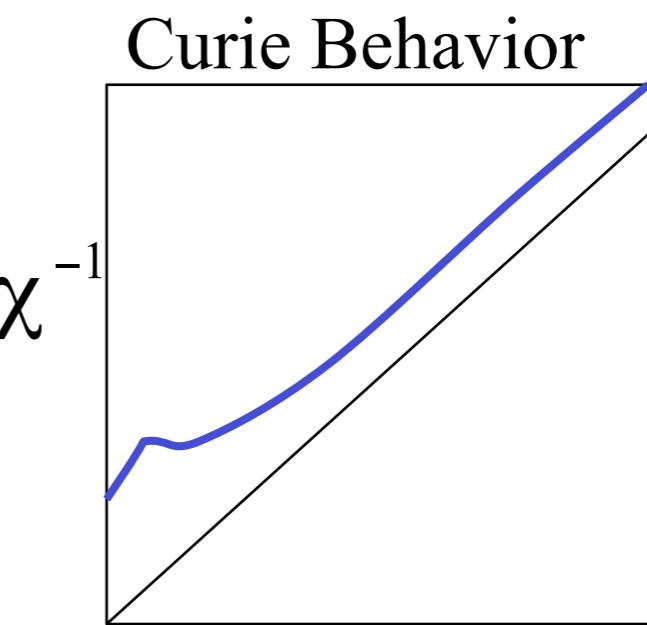
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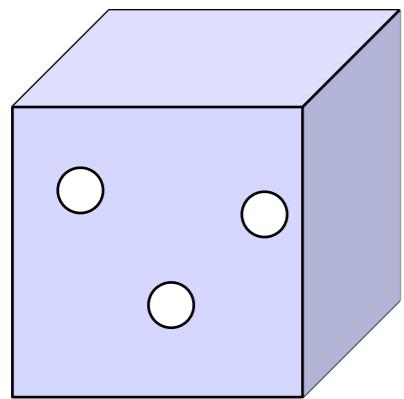
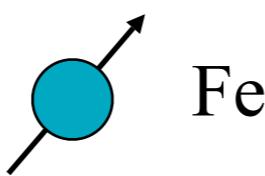


Clogston, Mathias et al, 1962

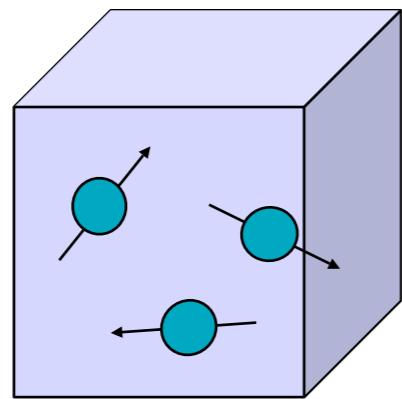


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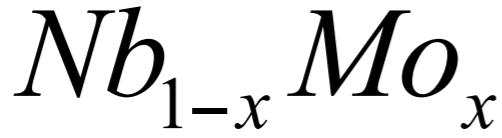
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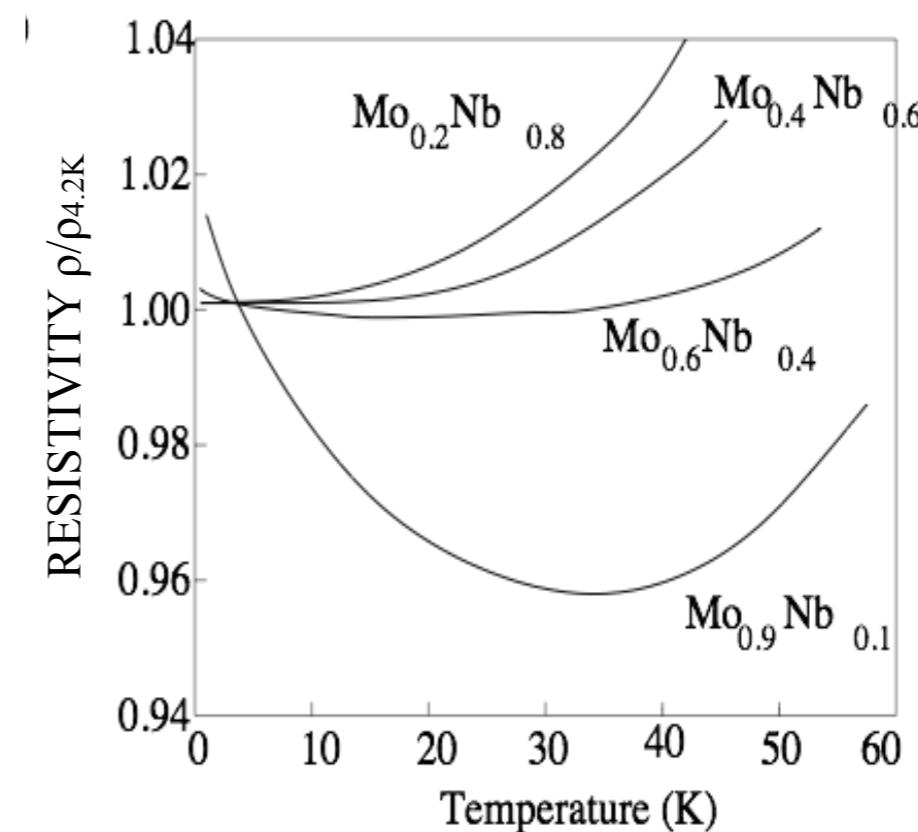
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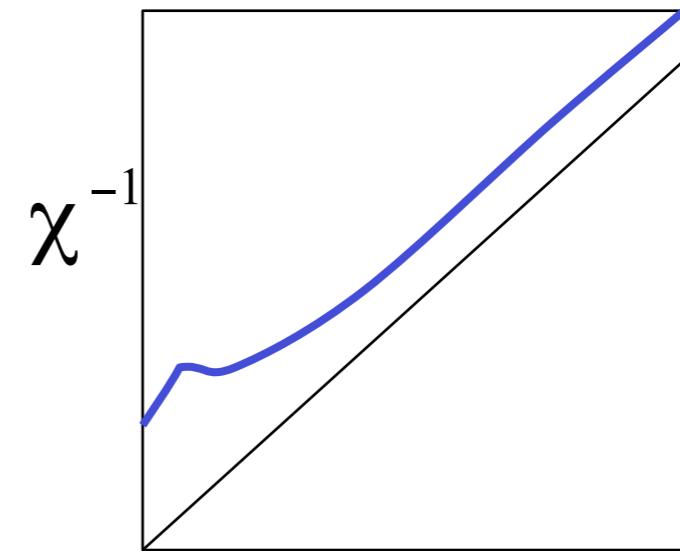
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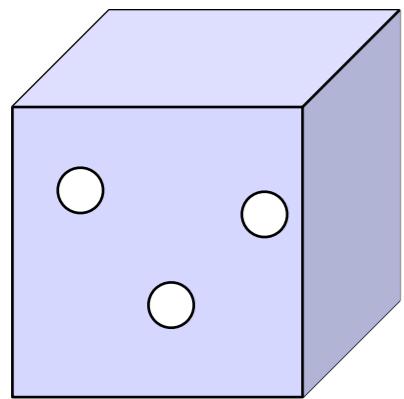
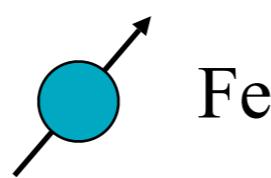


Curie Behavior

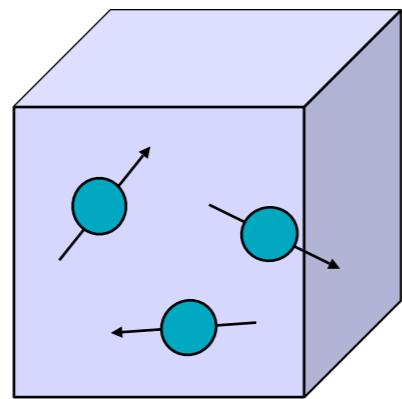


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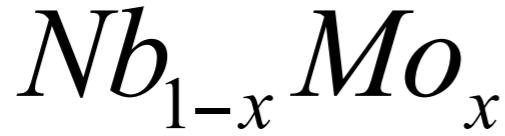
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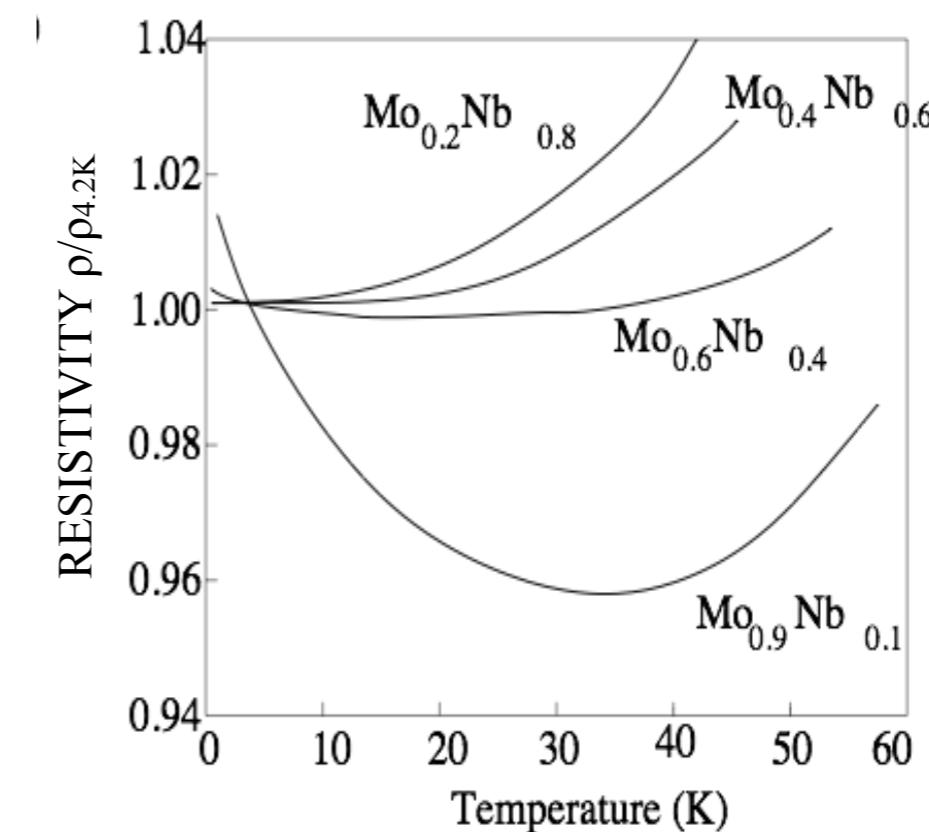
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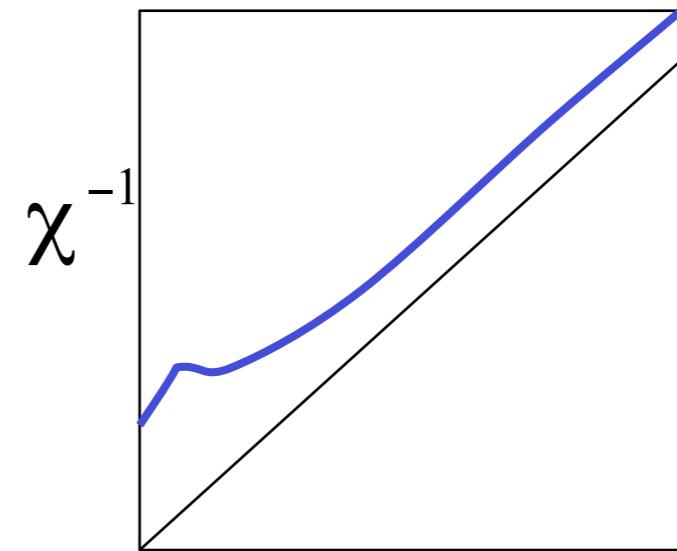
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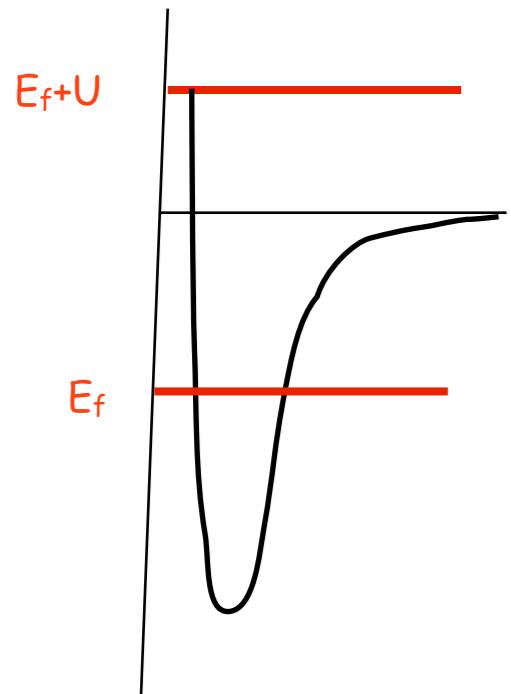


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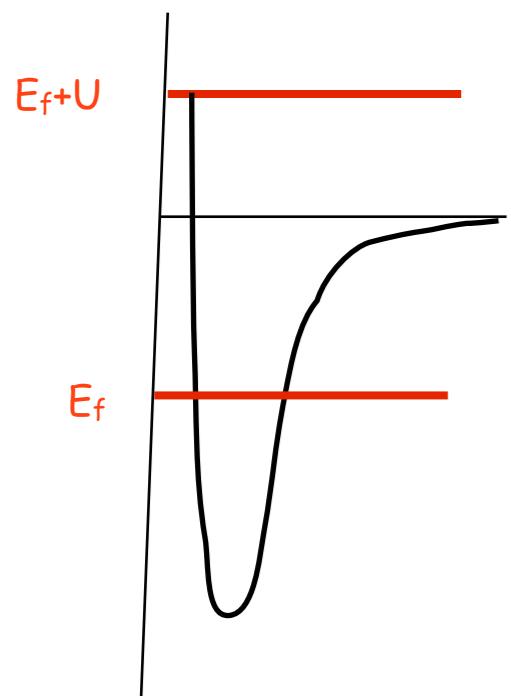
How?

# Anderson Model



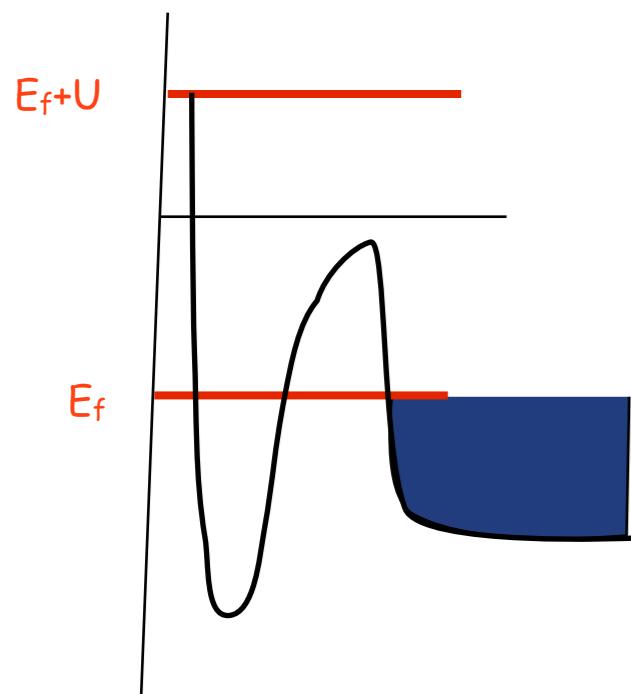
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$$H = \underbrace{E_f n_f + U n_{f\uparrow} n_{f\downarrow}}_{H_{atomic}}$$



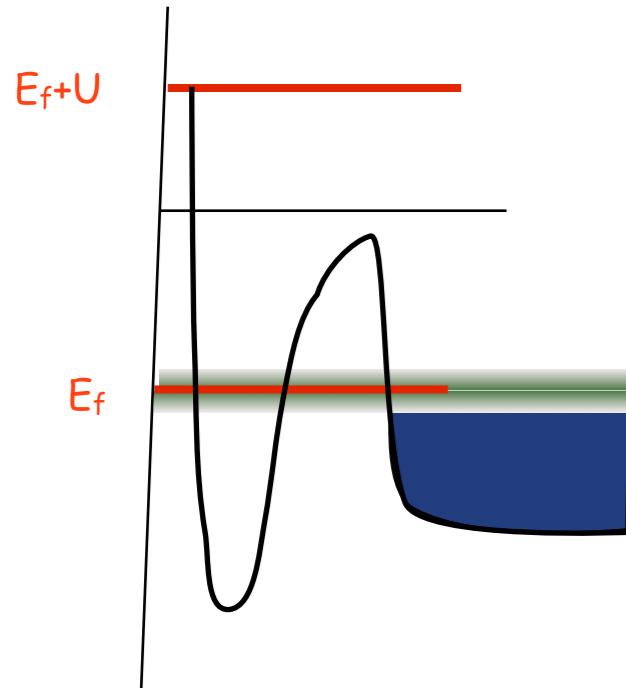
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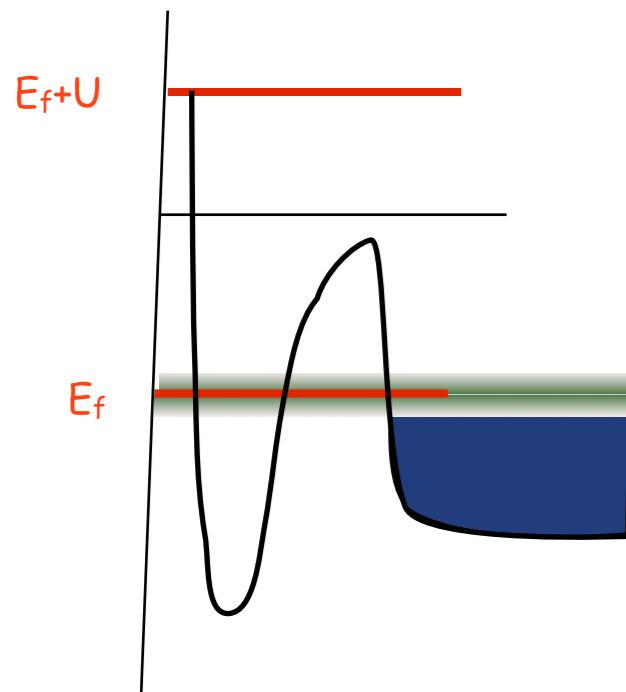
$$H = \underbrace{\sum_{k,\sigma} \epsilon_k n_{k\sigma} + \sum_{k,\sigma} V(k) [c_{k\sigma}^\dagger f_\sigma + f_\sigma^\dagger c_{k\sigma}]}_{H_{resonance}} + \underbrace{E_f n_f + U n_{f\uparrow} n_{f\downarrow}}_{H_{atomic}}$$



# Anderson Model

$$V(k) = \langle k | V_{atomic} | f \rangle = 4\pi i^l \int_0^\infty r^2 dr j_l(kr) V(r) R_f(r)$$

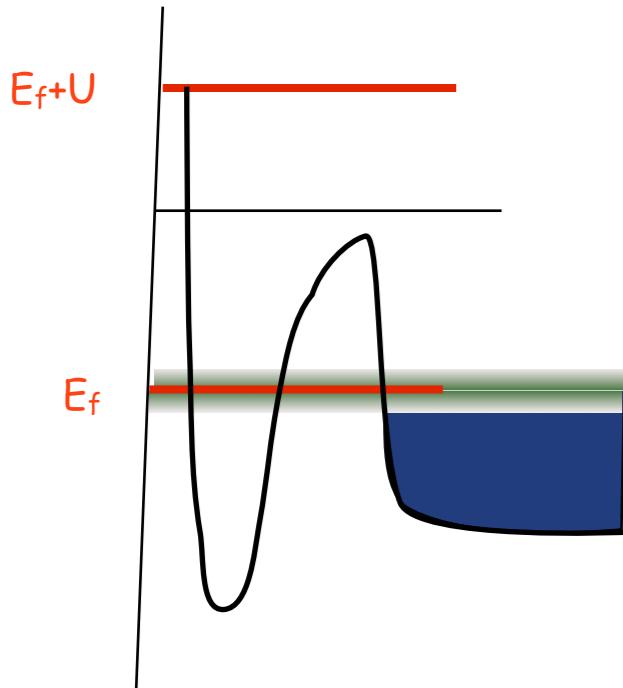
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$$U = \int d^3x d^3x' V(\mathbf{x} - \mathbf{x}') |\psi_f(\mathbf{x})|^2 |\psi_f(\mathbf{x}')|^2$$



- Atomic approach : Start with  $V(k)=0$ , then dial up the hybridization

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Local moment, states at  $E_f$  and  $E_f+U$

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Friedel-Abrikosov-Suhl Resonance.

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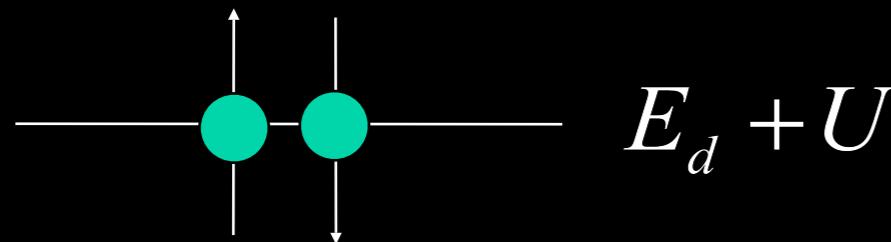
How to reconcile the two approaches?

# Anderson Model of Moment Formation

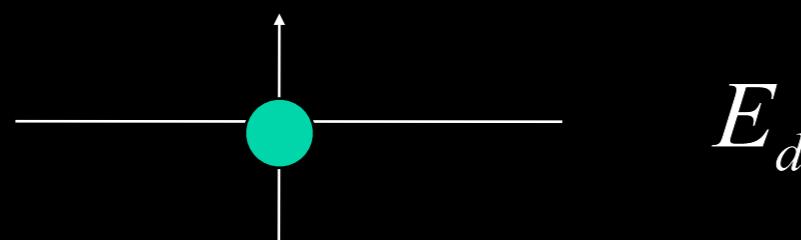
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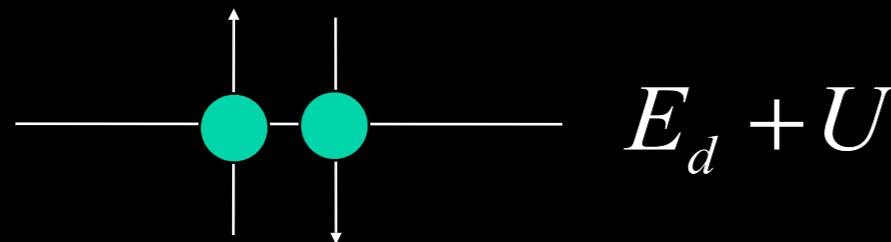
$$E_d + U$$



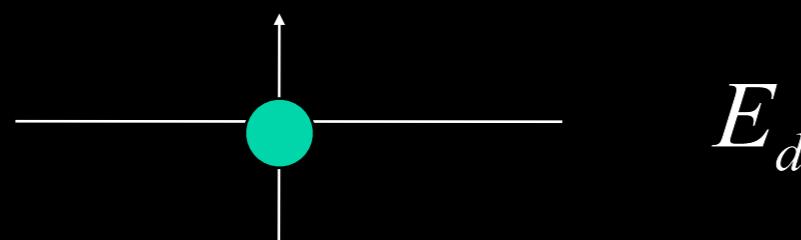
$$E_d$$

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$$E_d + U$$

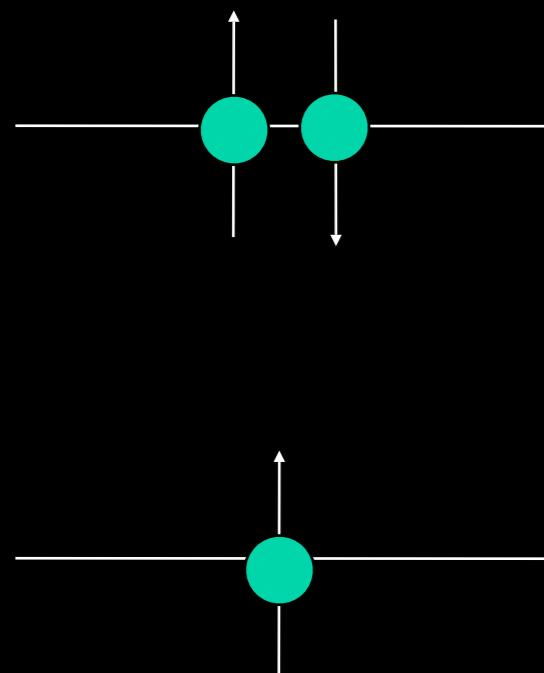


$$E_d$$

$$U/2 > |E_d + U/2|$$

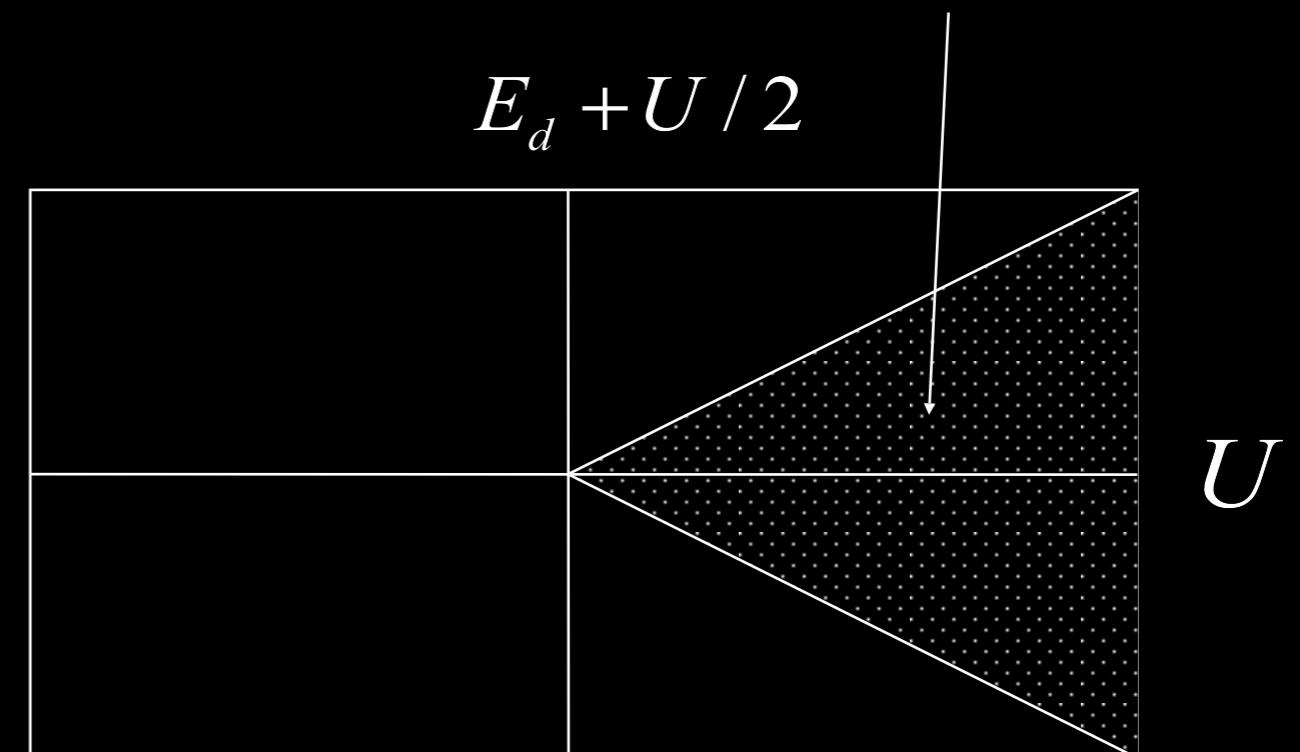
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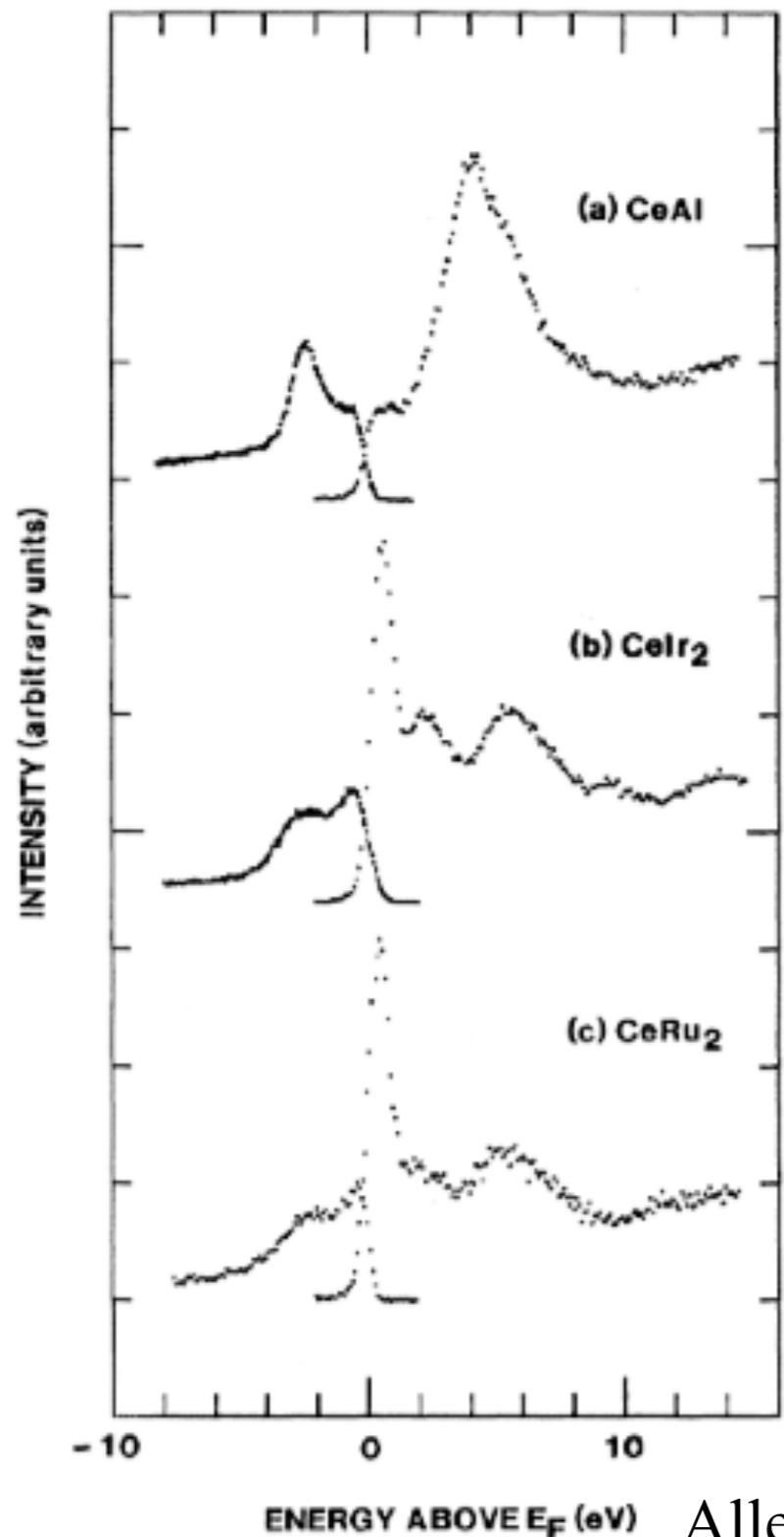


$$E_d + U$$
$$E_d$$

Local Moment  
Forms

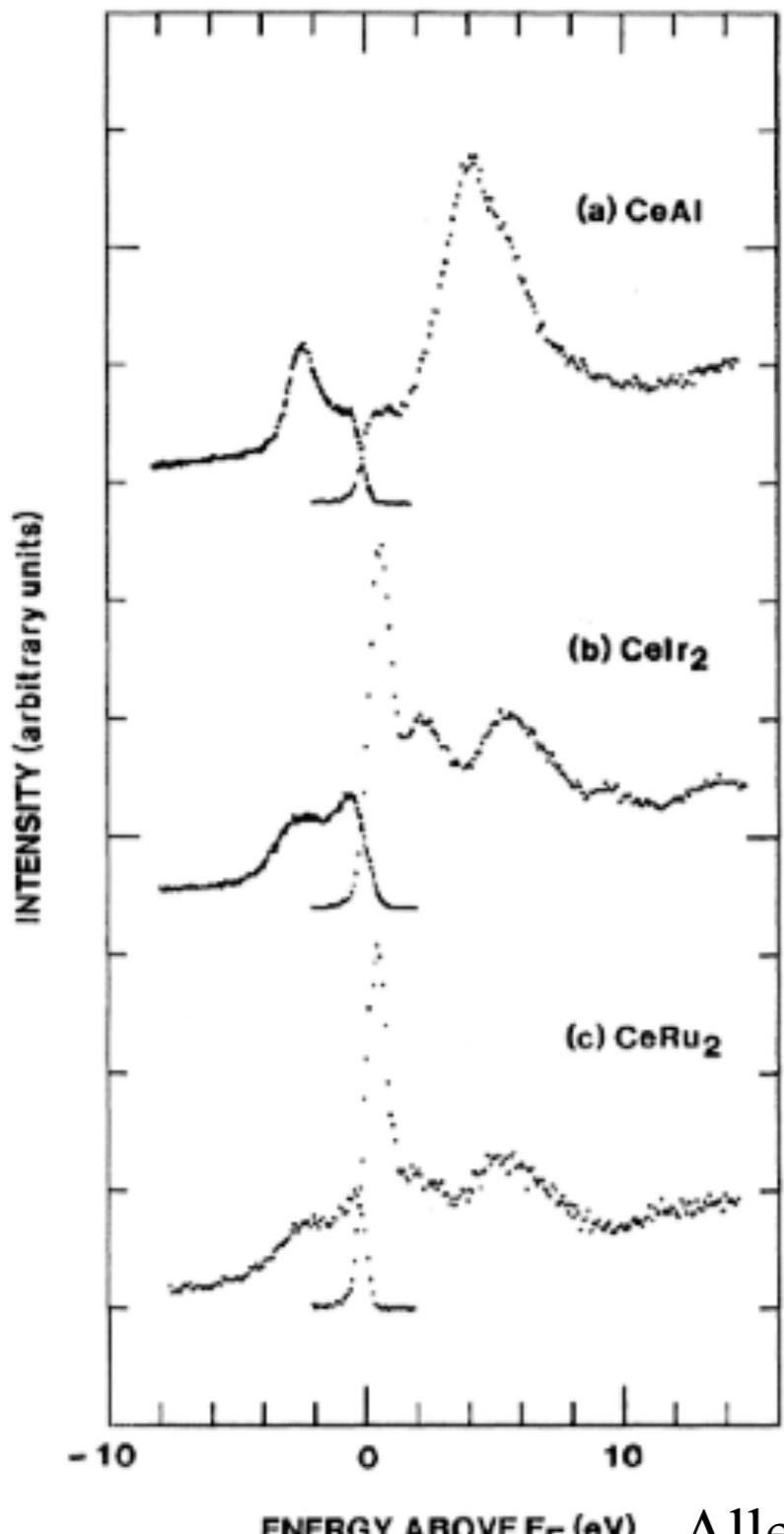


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$$A_f(\omega) = \frac{1}{\pi} \text{Im}G_f(\omega - i\delta)$$

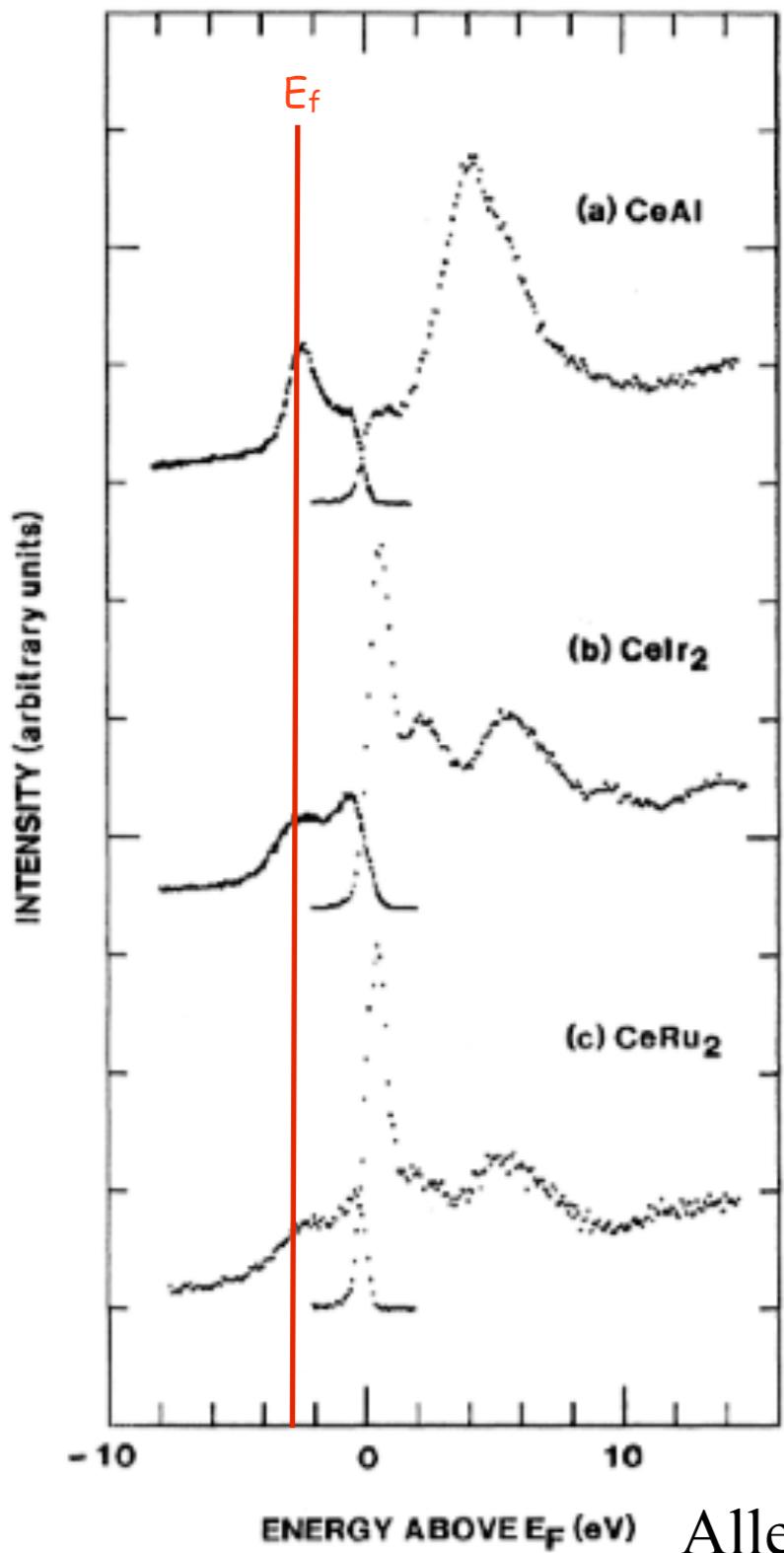
Allen et al (1983)



ENERGY ABOVE  $E_F$  (eV) Allen et al (1983)

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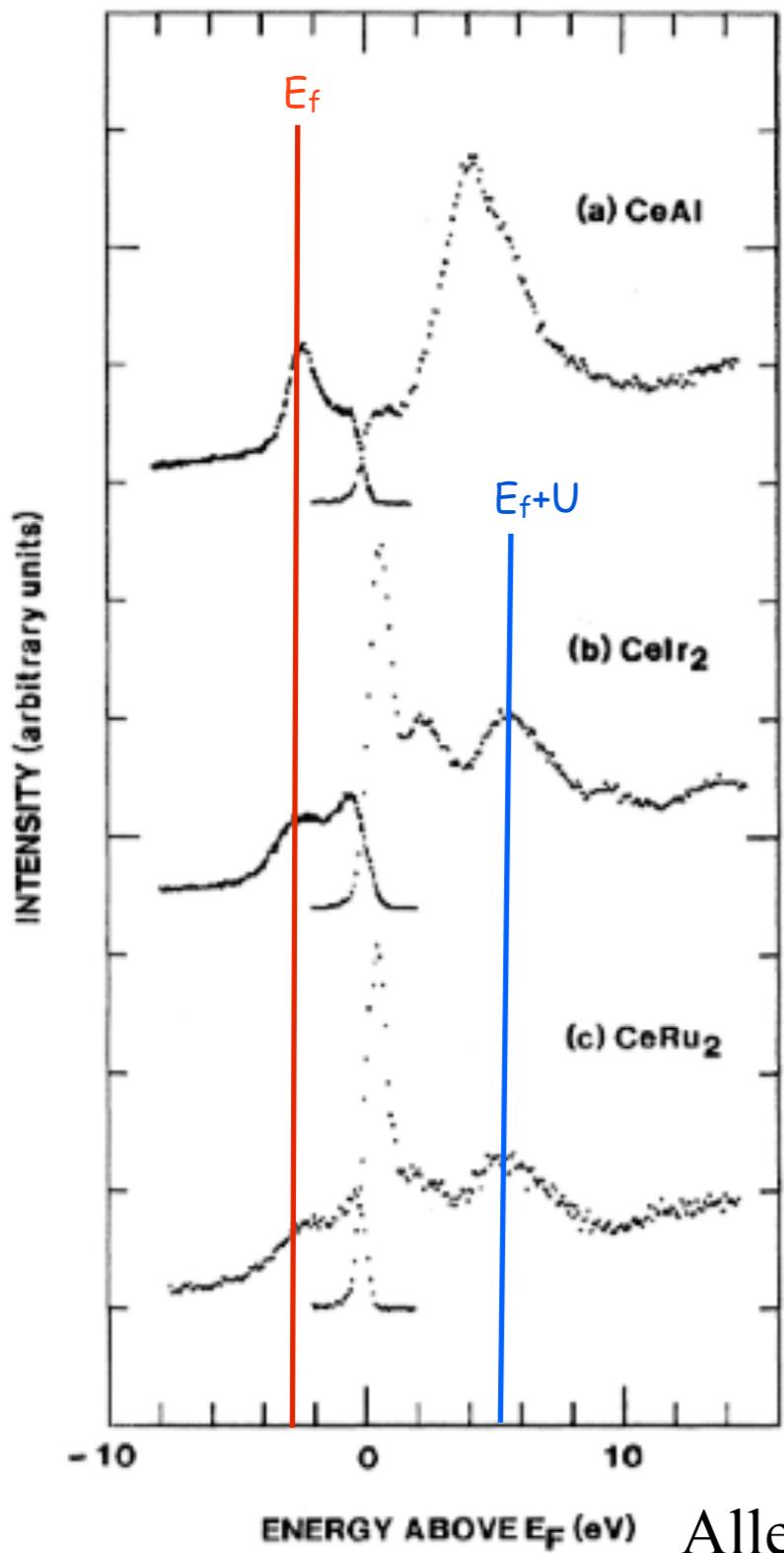
$$G_f(\omega) = -i \int_{-\infty}^{\infty} dt \langle T f_{\sigma}(t) f_{\sigma}^{\dagger}(0) \rangle e^{i\omega t}$$



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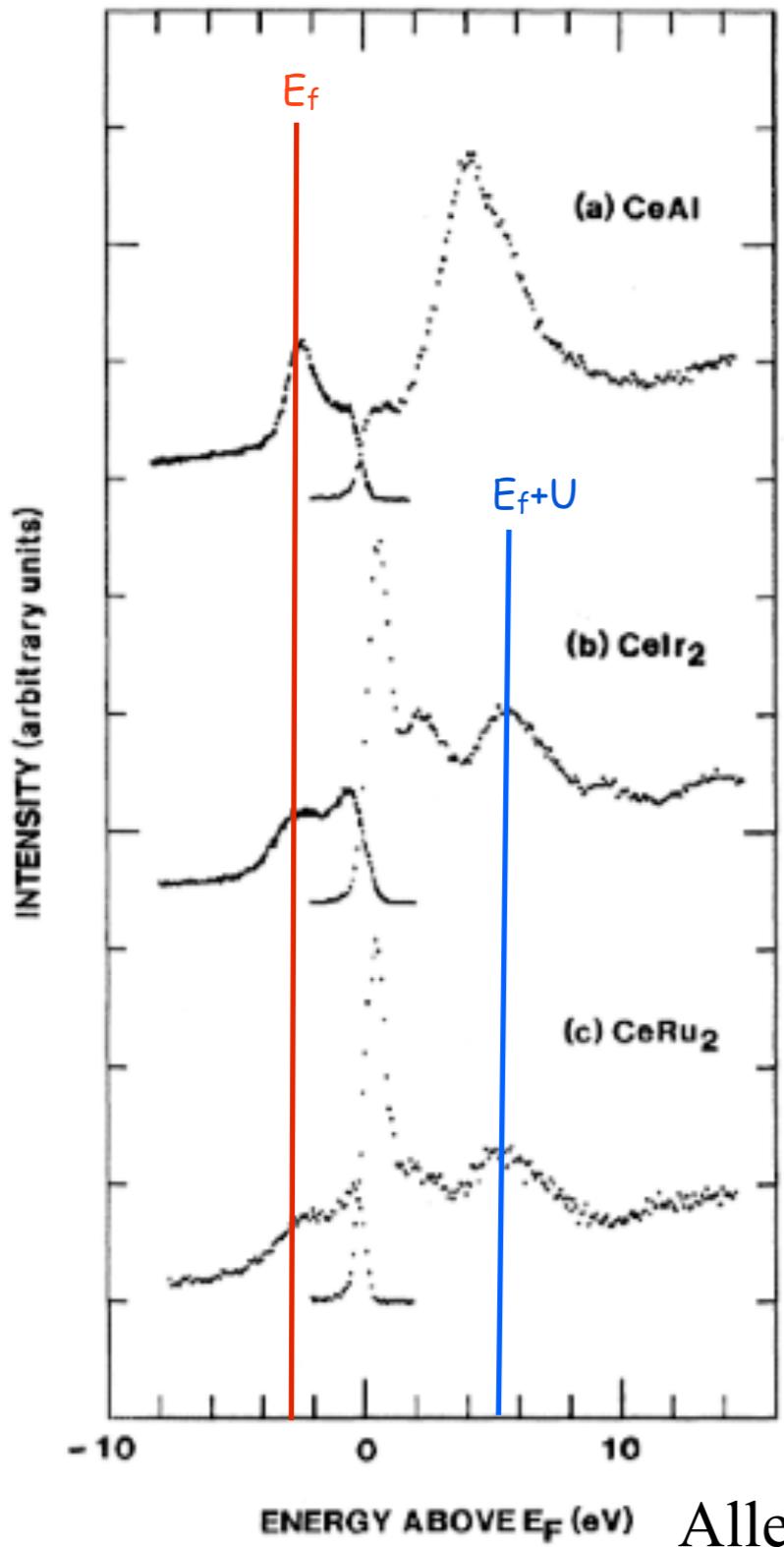
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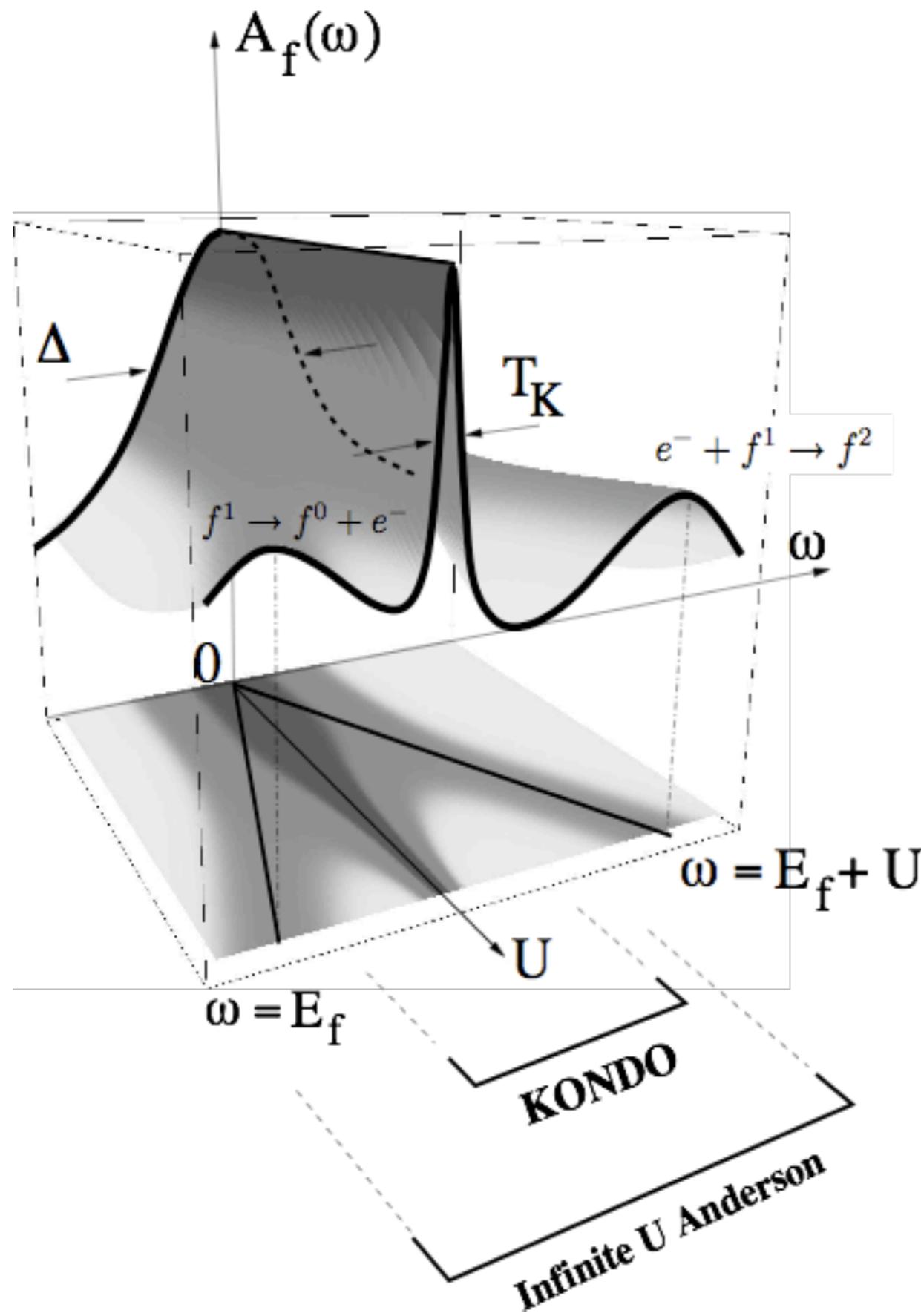
$$G_f(\omega) = -i \int_{-\infty}^{\infty} dt \langle T f_{\sigma}(t) f_{\sigma}^{\dagger}(0) \rangle e^{i\omega t}$$

$$A_f(\omega) = \begin{cases} \overbrace{\sum_{\lambda} |\langle \lambda | f_{\sigma}^{\dagger} | \phi_0 \rangle|^2 \delta(\omega - [E_{\lambda} - E_0])}, & (\omega > 0) \\ \overbrace{\sum_{\lambda} |\langle \lambda | f_{\sigma} | \phi_0 \rangle|^2 \delta(\omega - [E_0 - E_{\lambda}])}, & (\omega < 0) \end{cases}$$

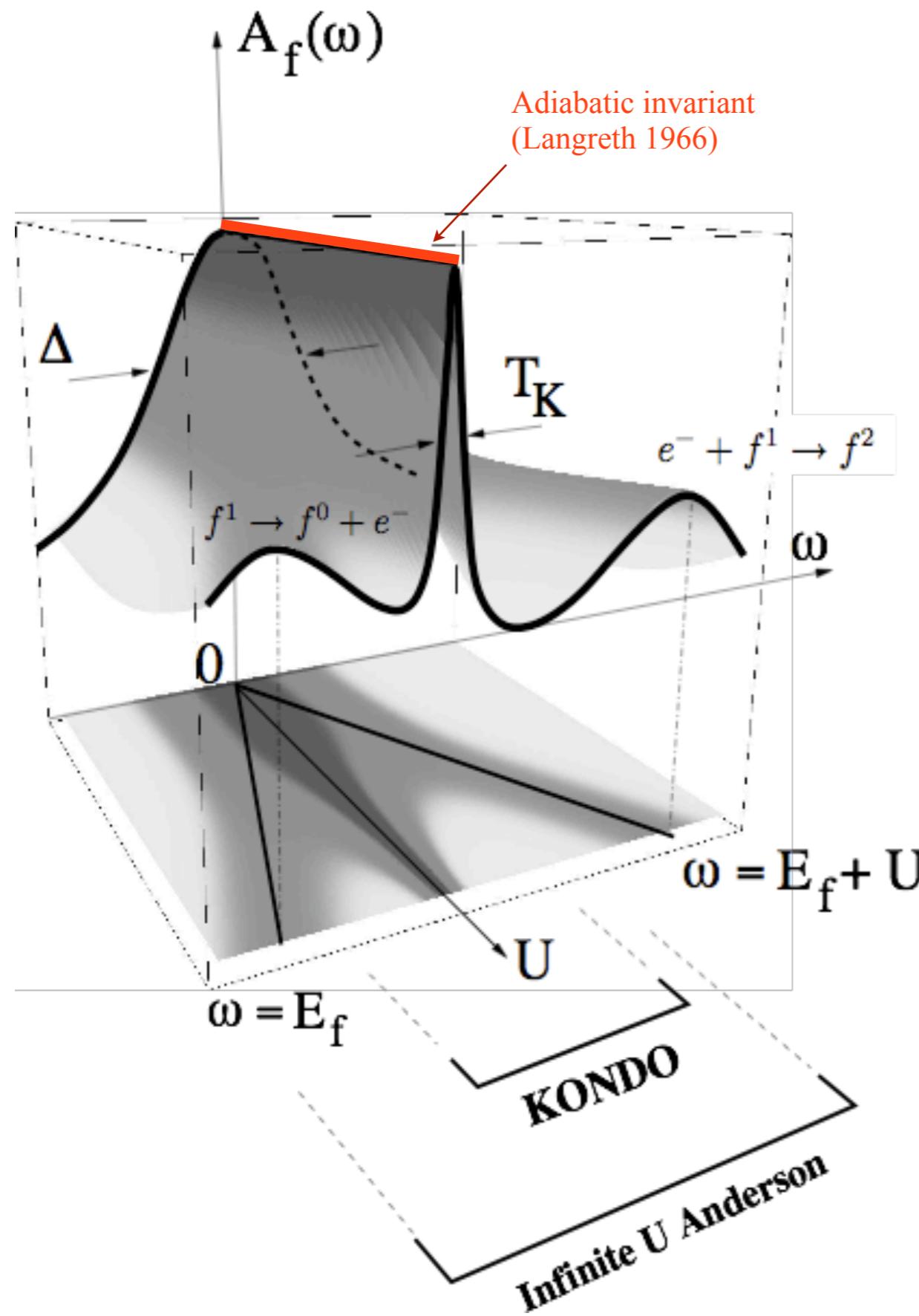
Energy distribution of state formed by adding one f-electron.

Energy distribution of state formed by removing an f-electron

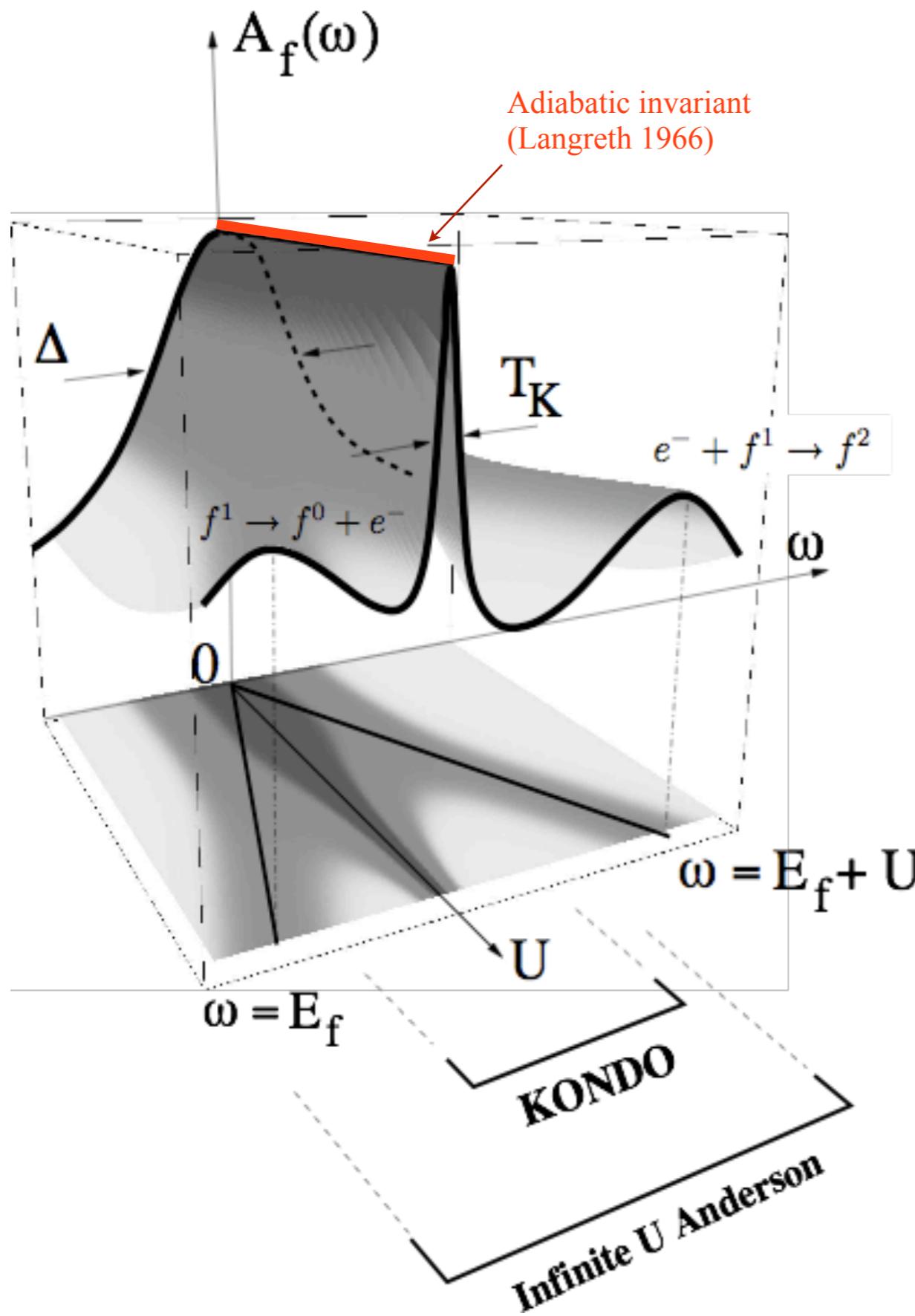
# Adiabatic approach & Kondo Resonance



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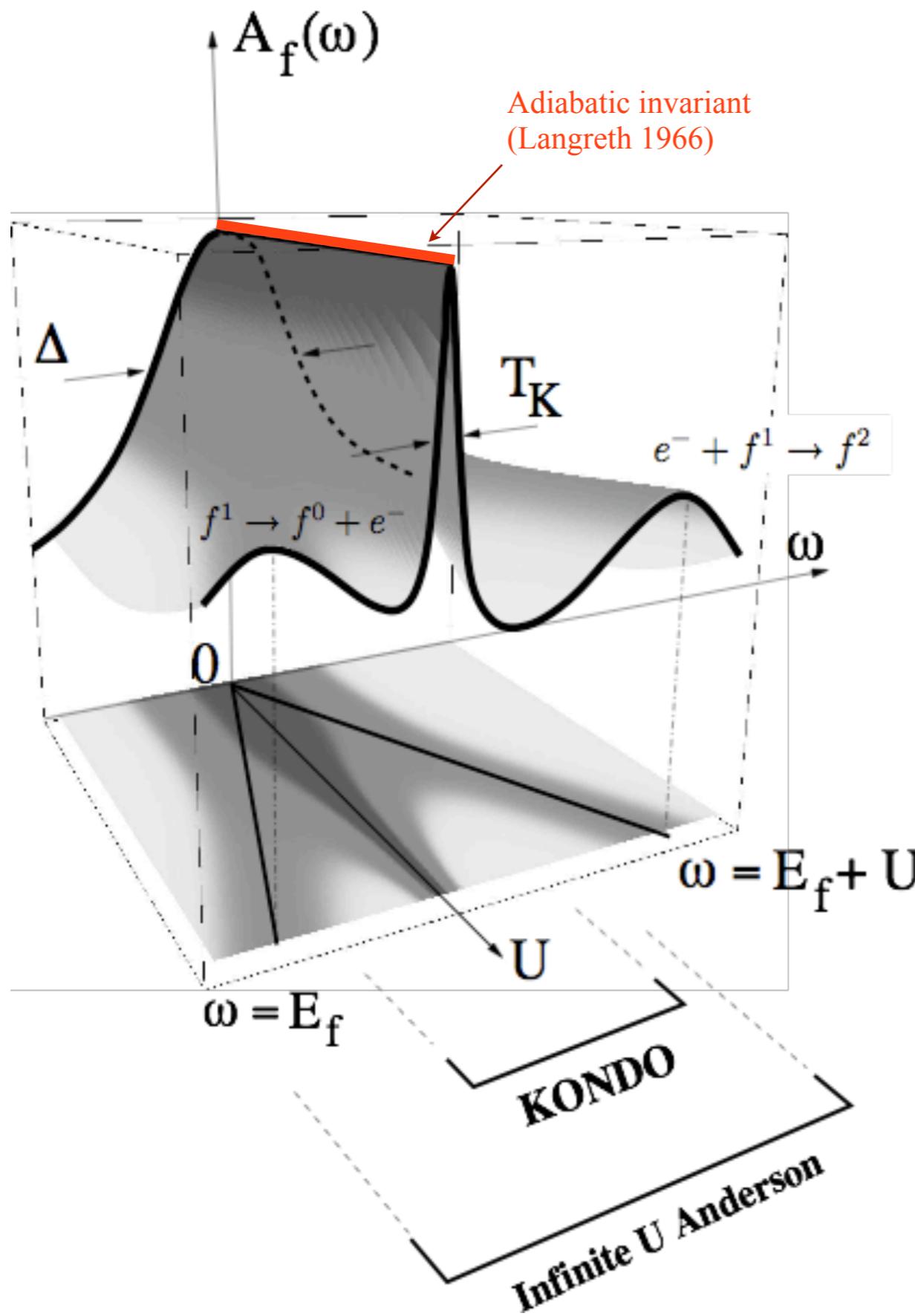


# Adiabatic approach & Kondo Resonance



$$A_f(\omega = 0) = \frac{\sin^2 \delta_f}{\pi \Delta}$$

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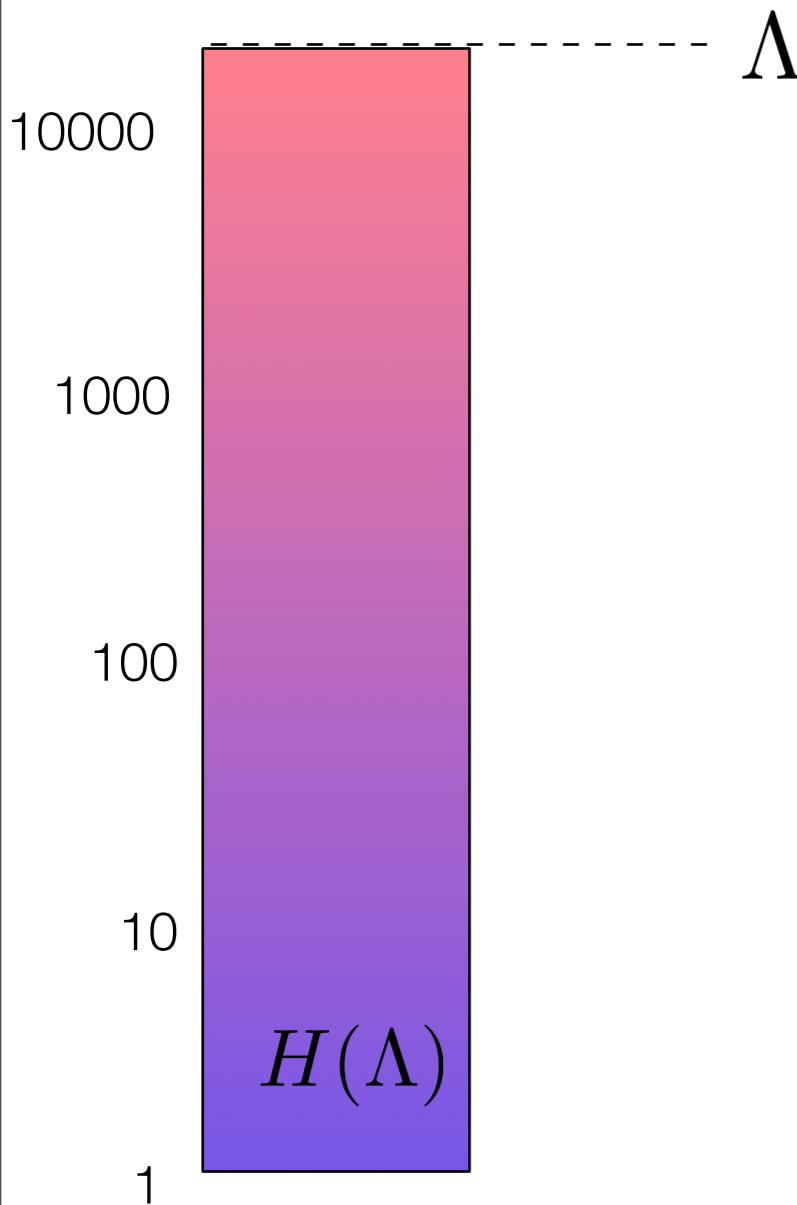


$$A_f(\omega = 0) = \frac{\sin^2 \delta_f}{\pi \Delta}$$

$$\sum_{\sigma} \frac{\delta_{f\sigma}}{\pi} = 2 \frac{\delta}{\pi} = n_f$$

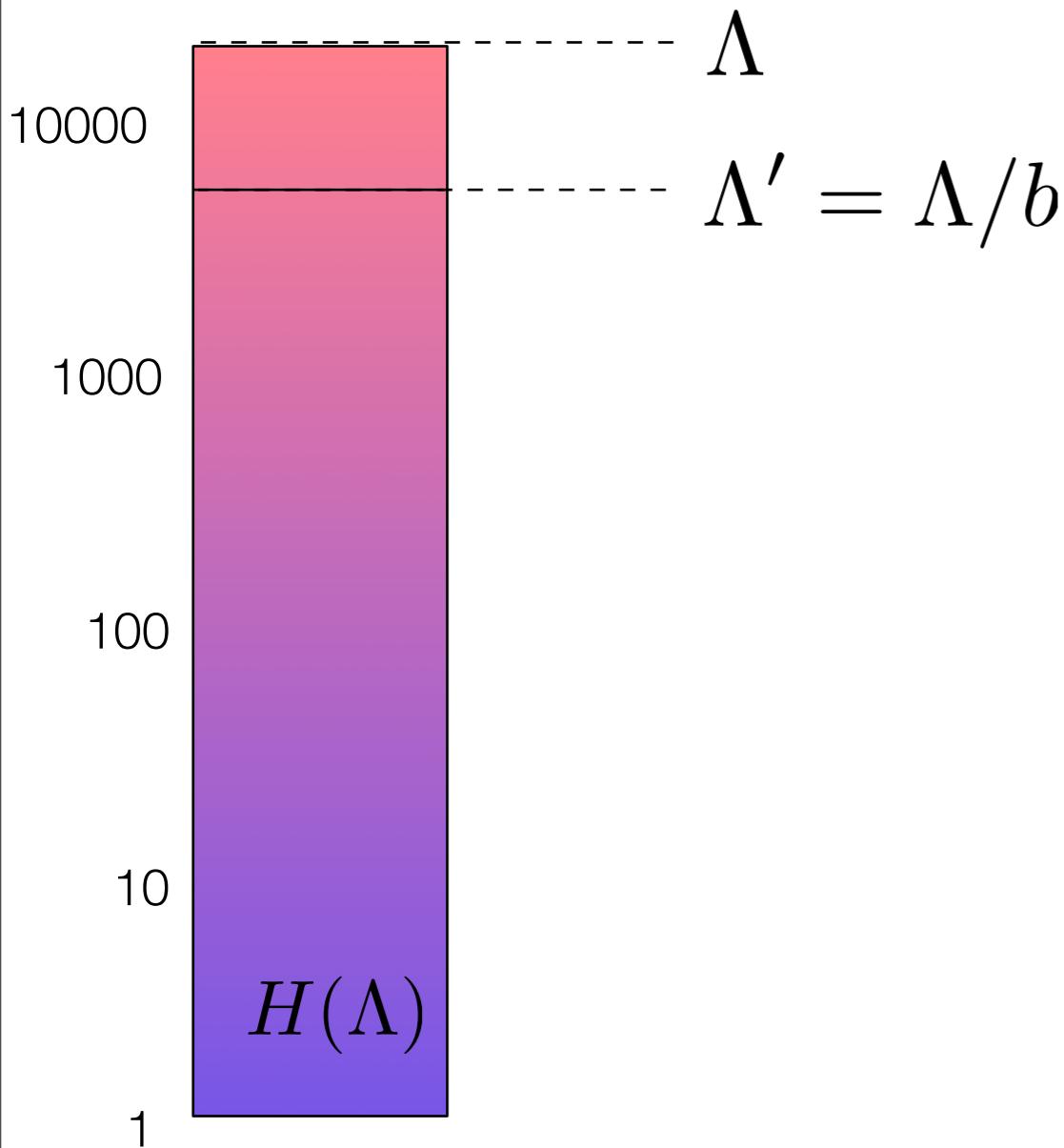
# Hierachies of Energy scales: Renormalization concept.

(Anderson, Wilson, ....)



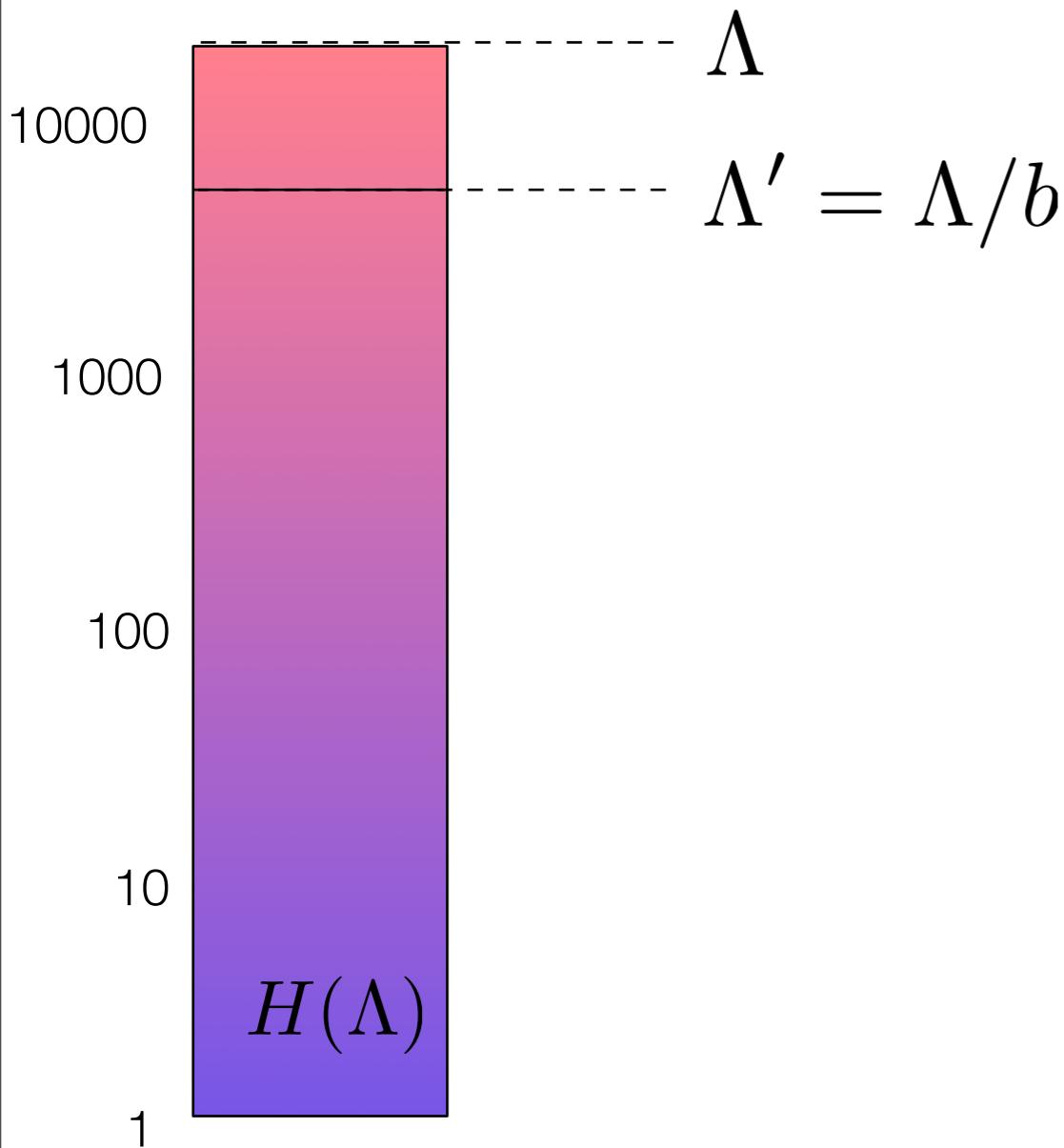
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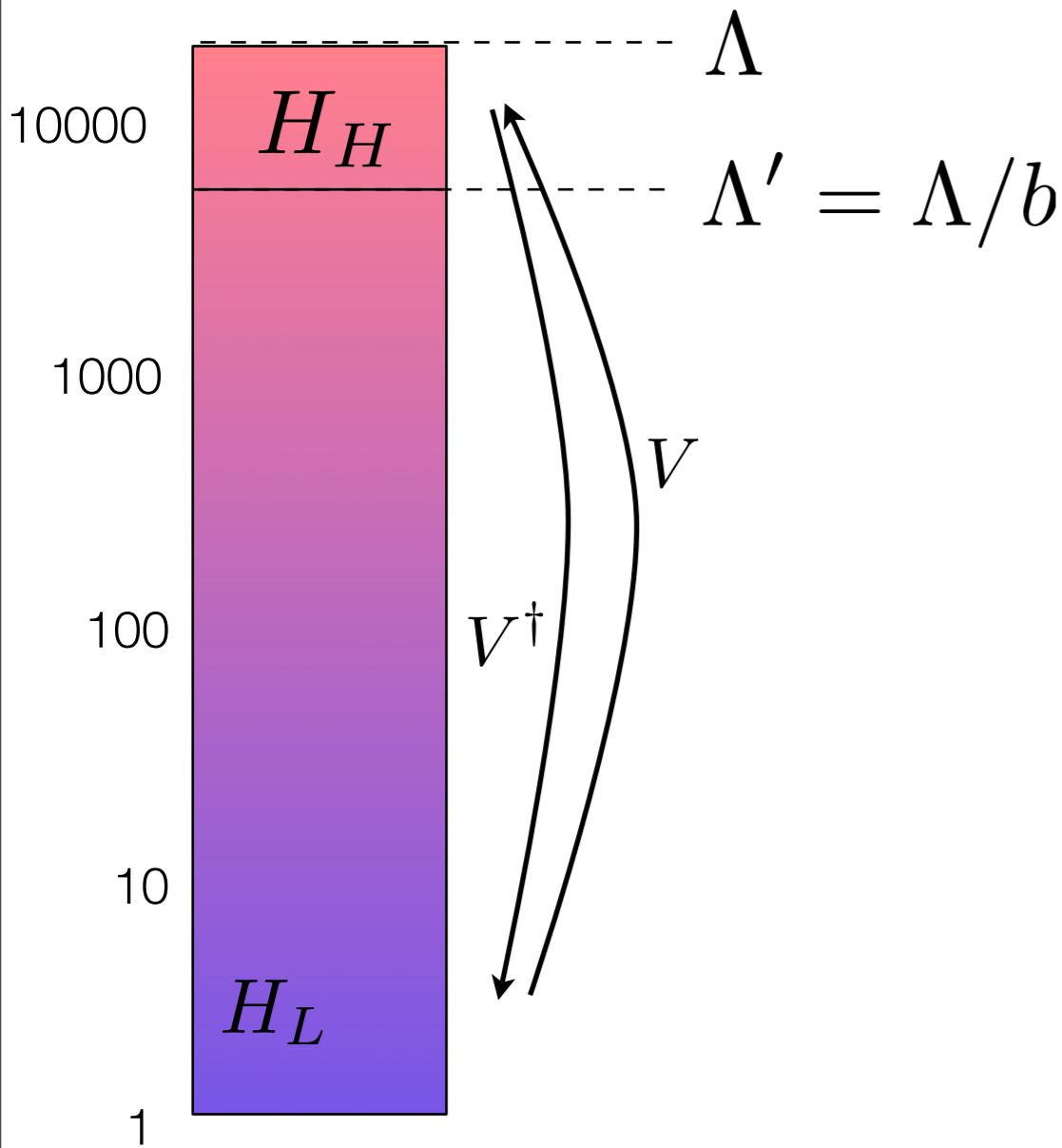
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$$H(\Lambda) = \left[ \frac{H_L}{V} \middle| \frac{V^\dagger}{H_H} \right],$$

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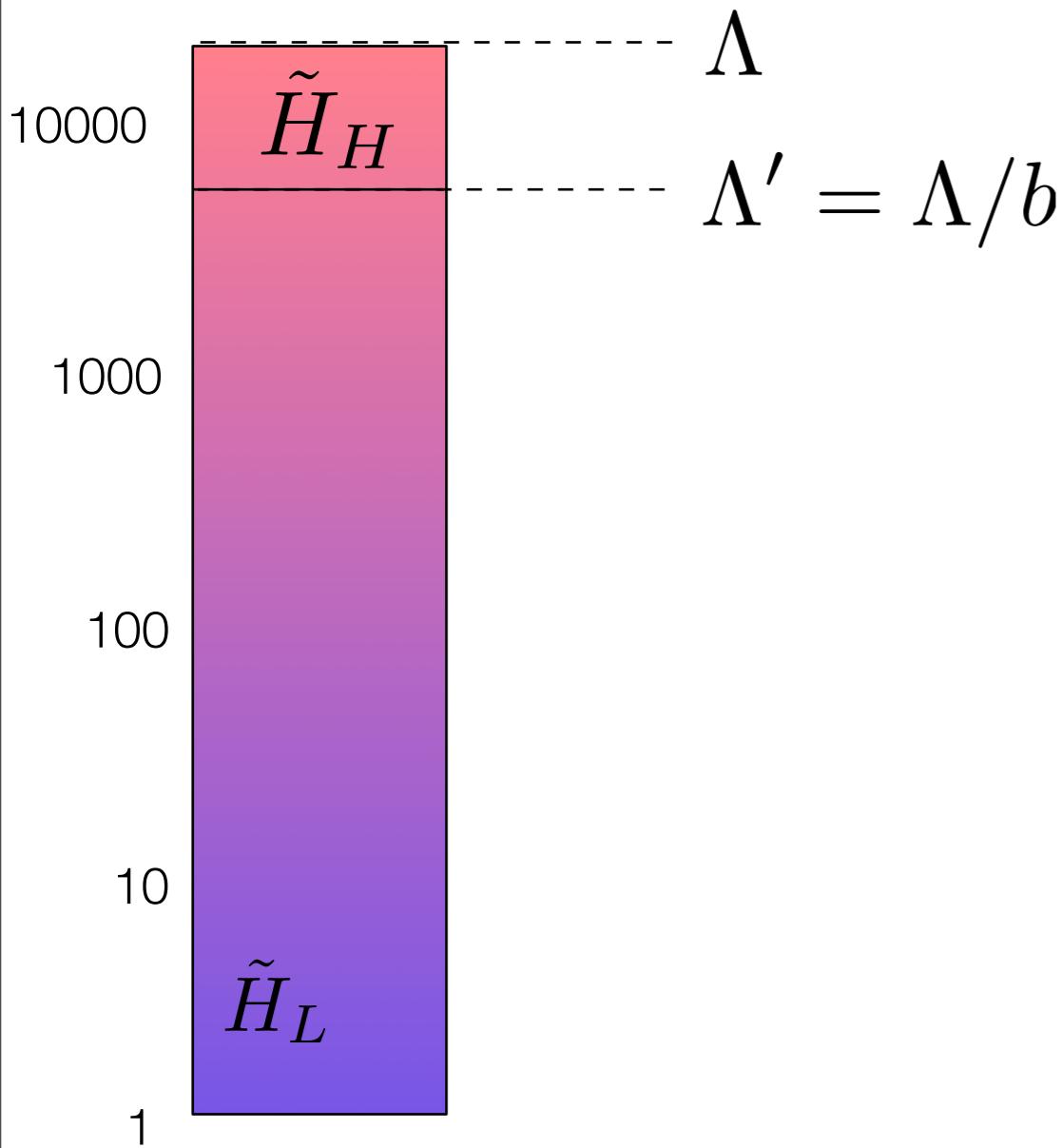
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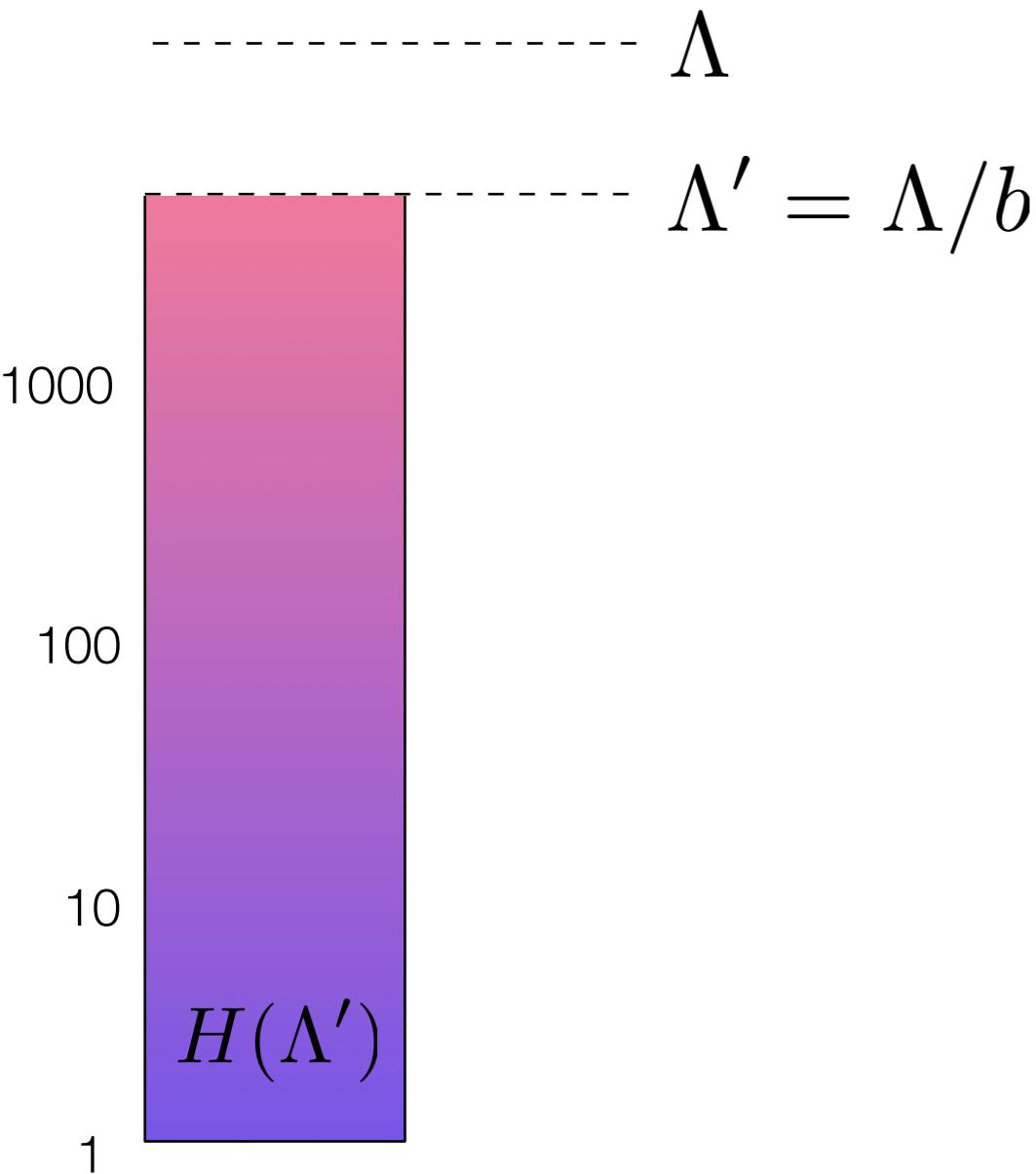


$$H(\Lambda) = \left[ \begin{array}{c|c} H_L & V^\dagger \\ \hline V & H_H \end{array} \right],$$

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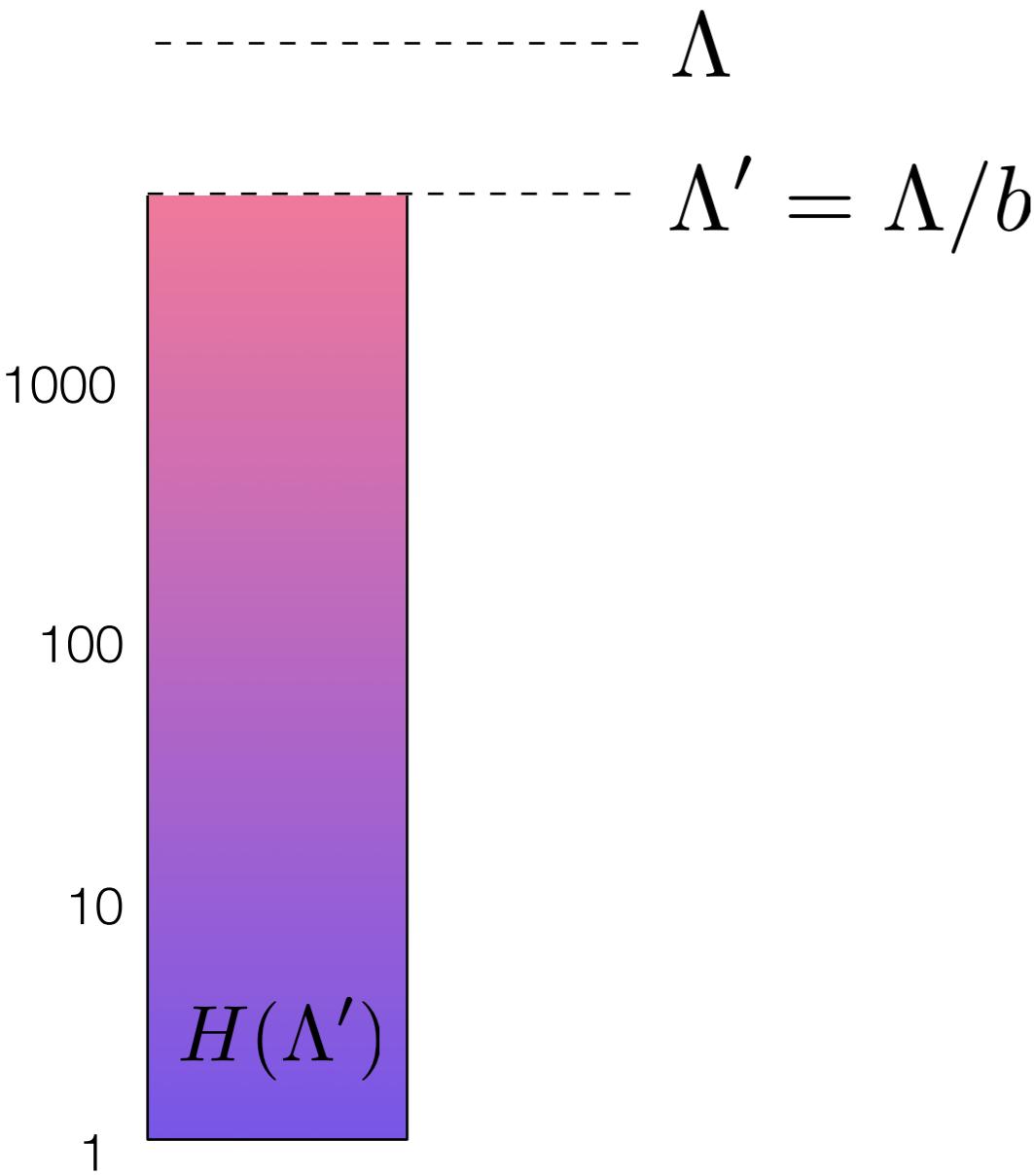
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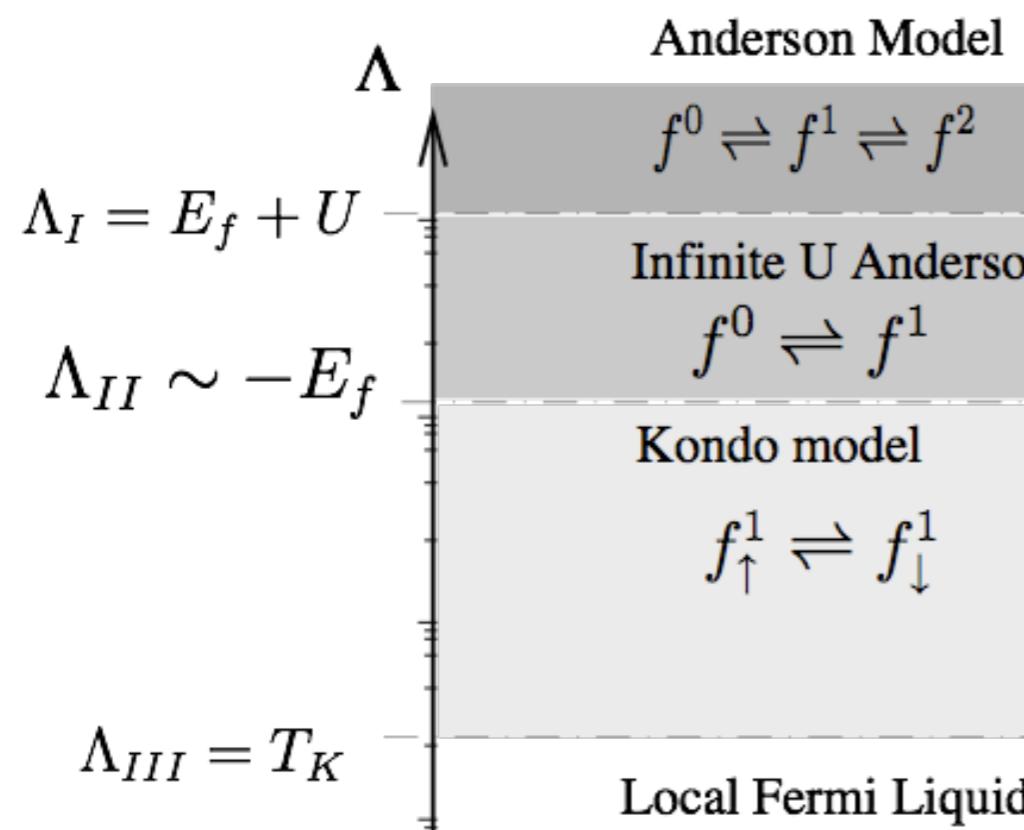
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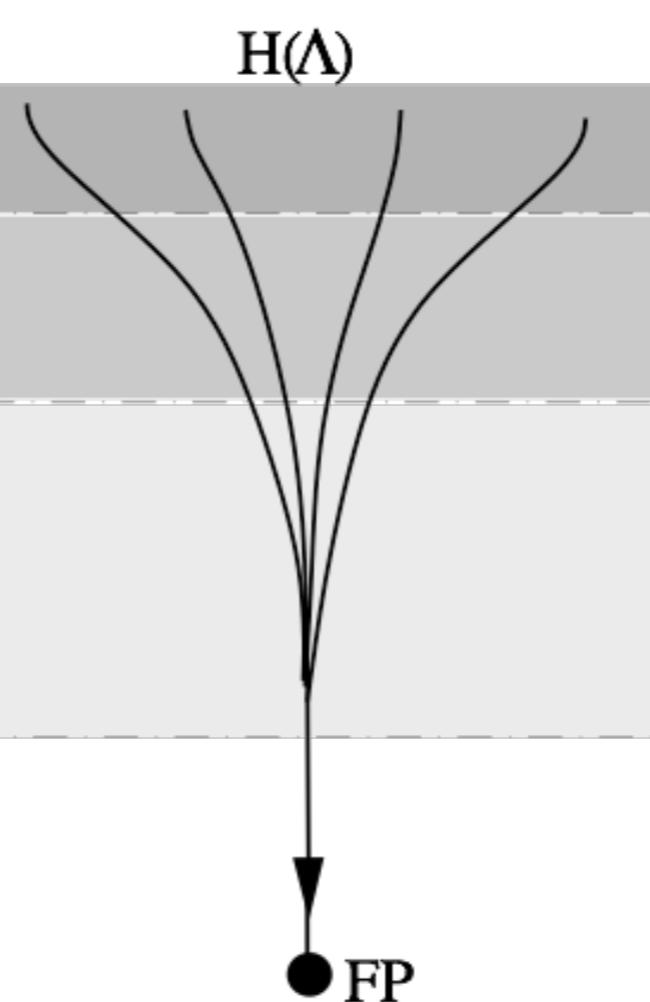
$$\tilde{H}_L = H_L + \delta H = H(\Lambda')$$

$$\delta H = P_L V^\dagger \left( \frac{1}{E - \tilde{H}_H} \right) V P_L \Big|_{E \sim E_L}$$

(a)

**Hamiltonian**

(b)

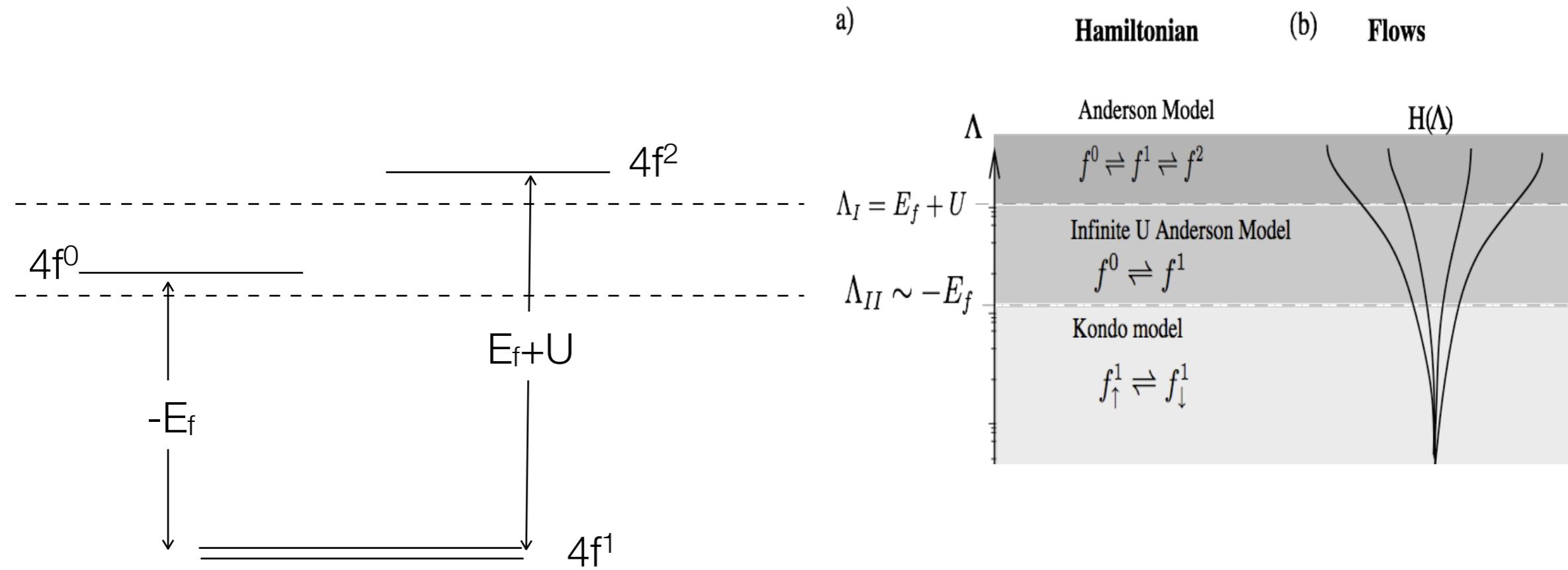
**Flows****Excitations**

- ↑ valence fluctuations
- ↓ moment formation
- ↑ local moments
- ↓ quasiparticles

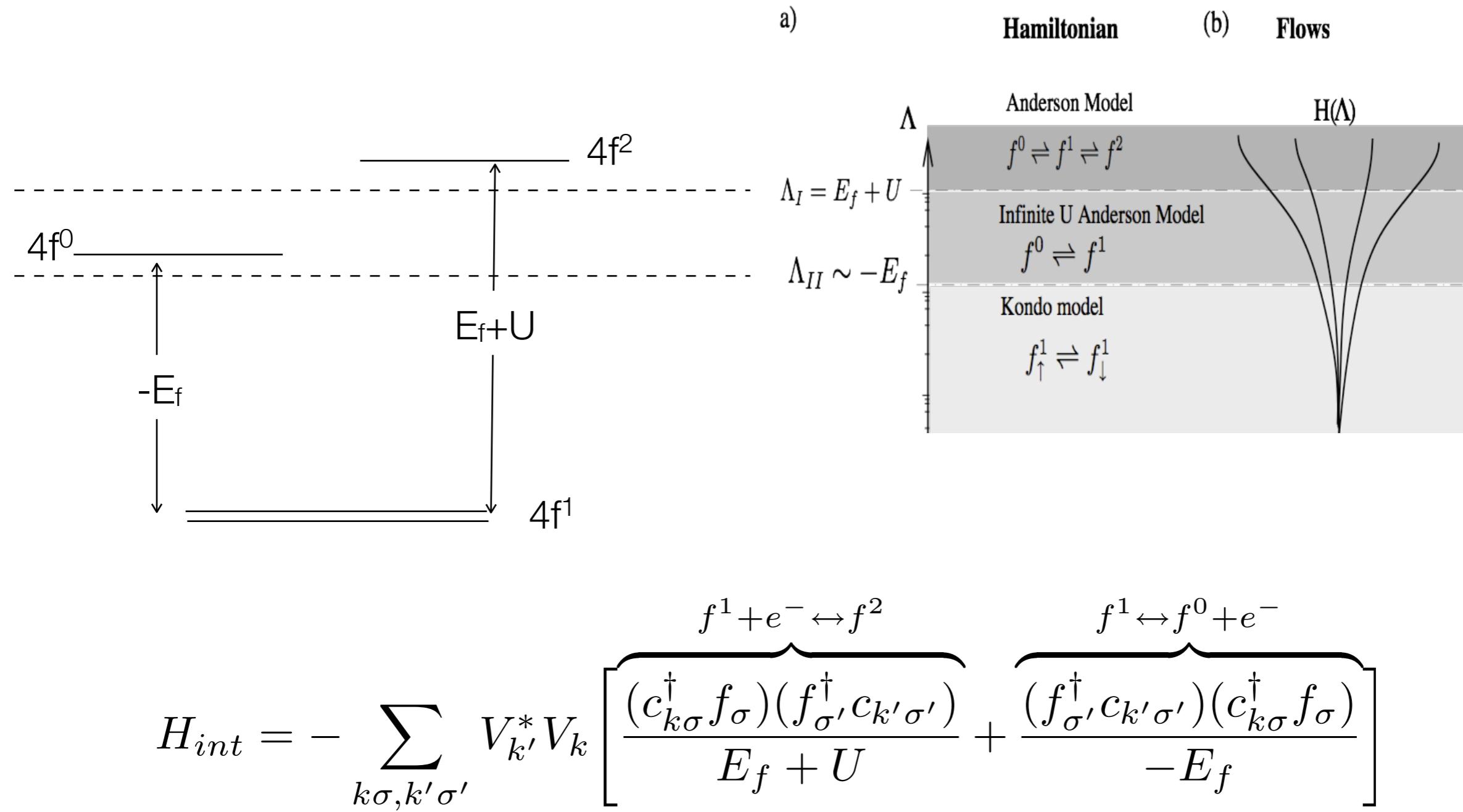
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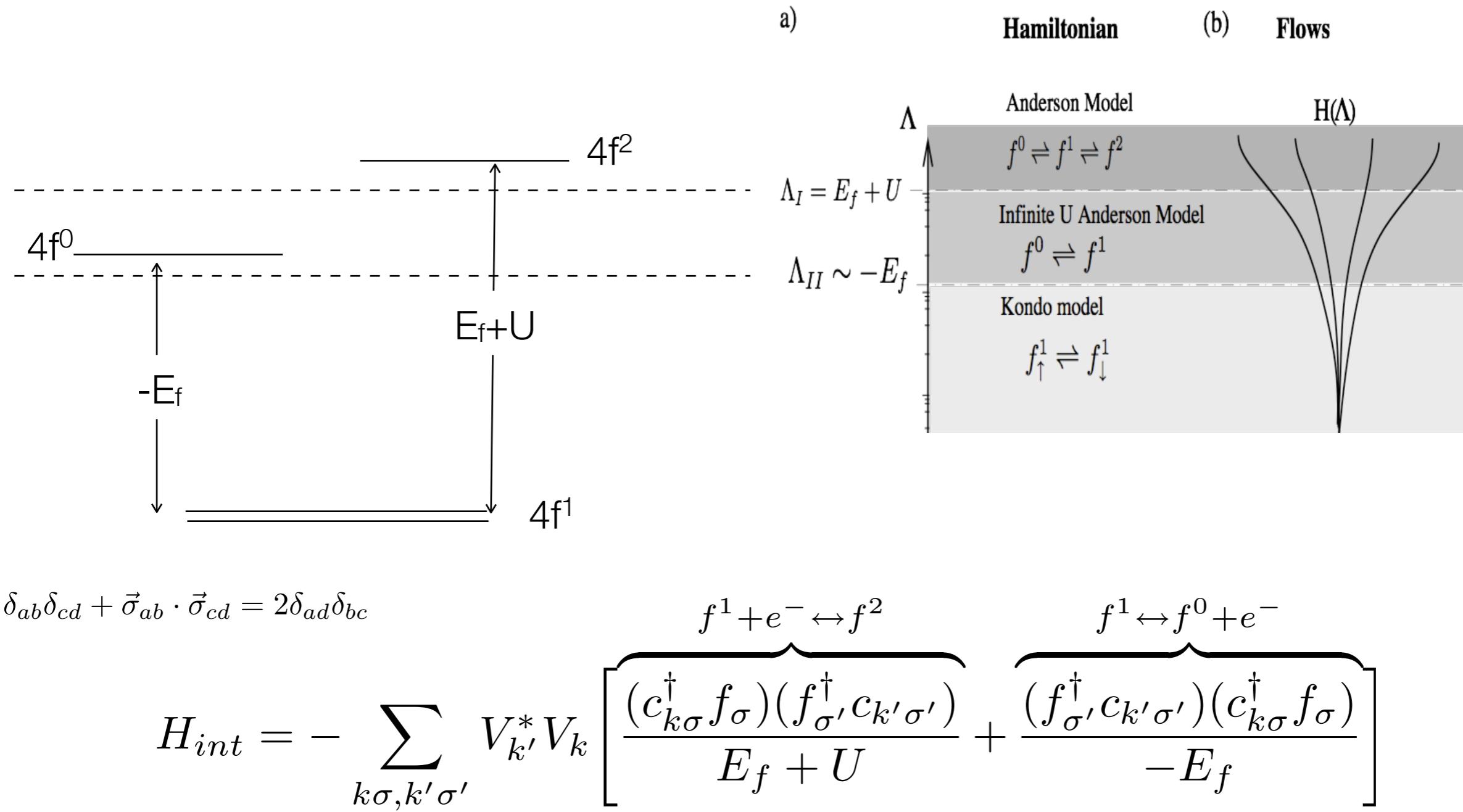
# Schrieffer-Wolff Transformation.



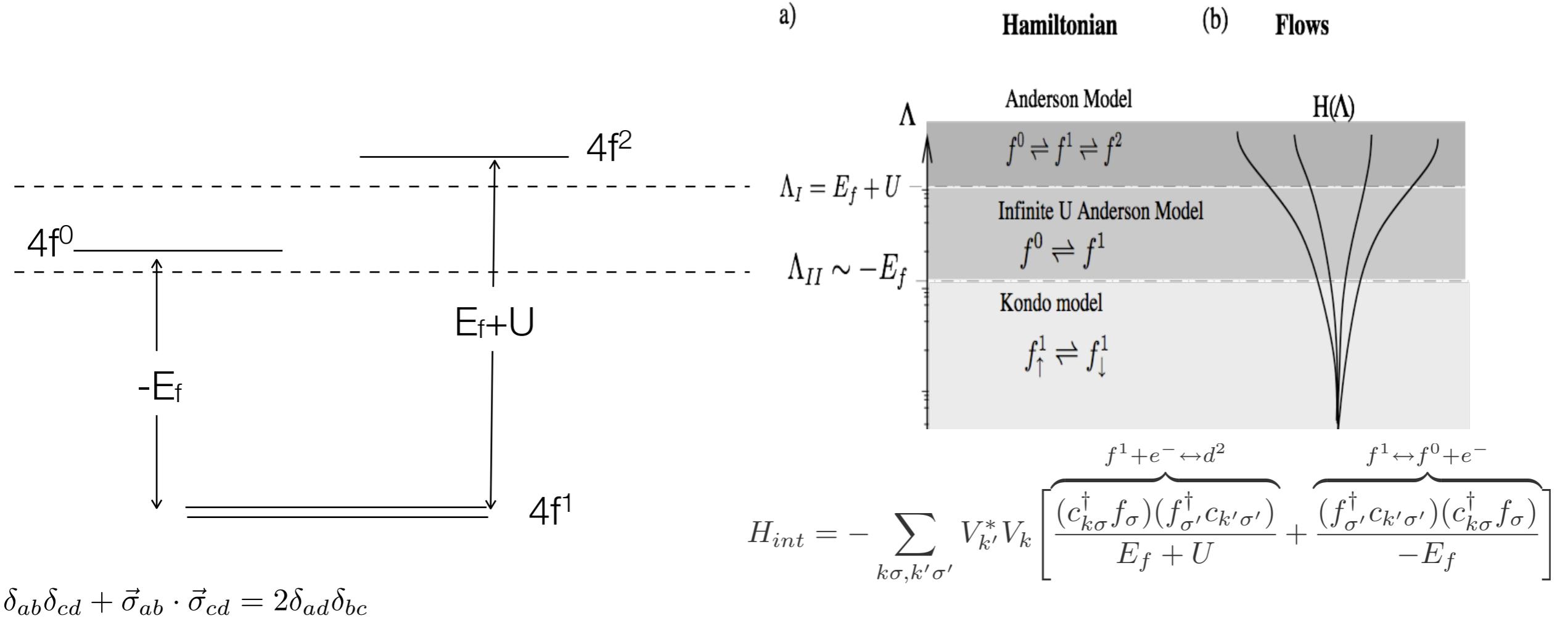
# Schrieffer-Wolff Transformation.



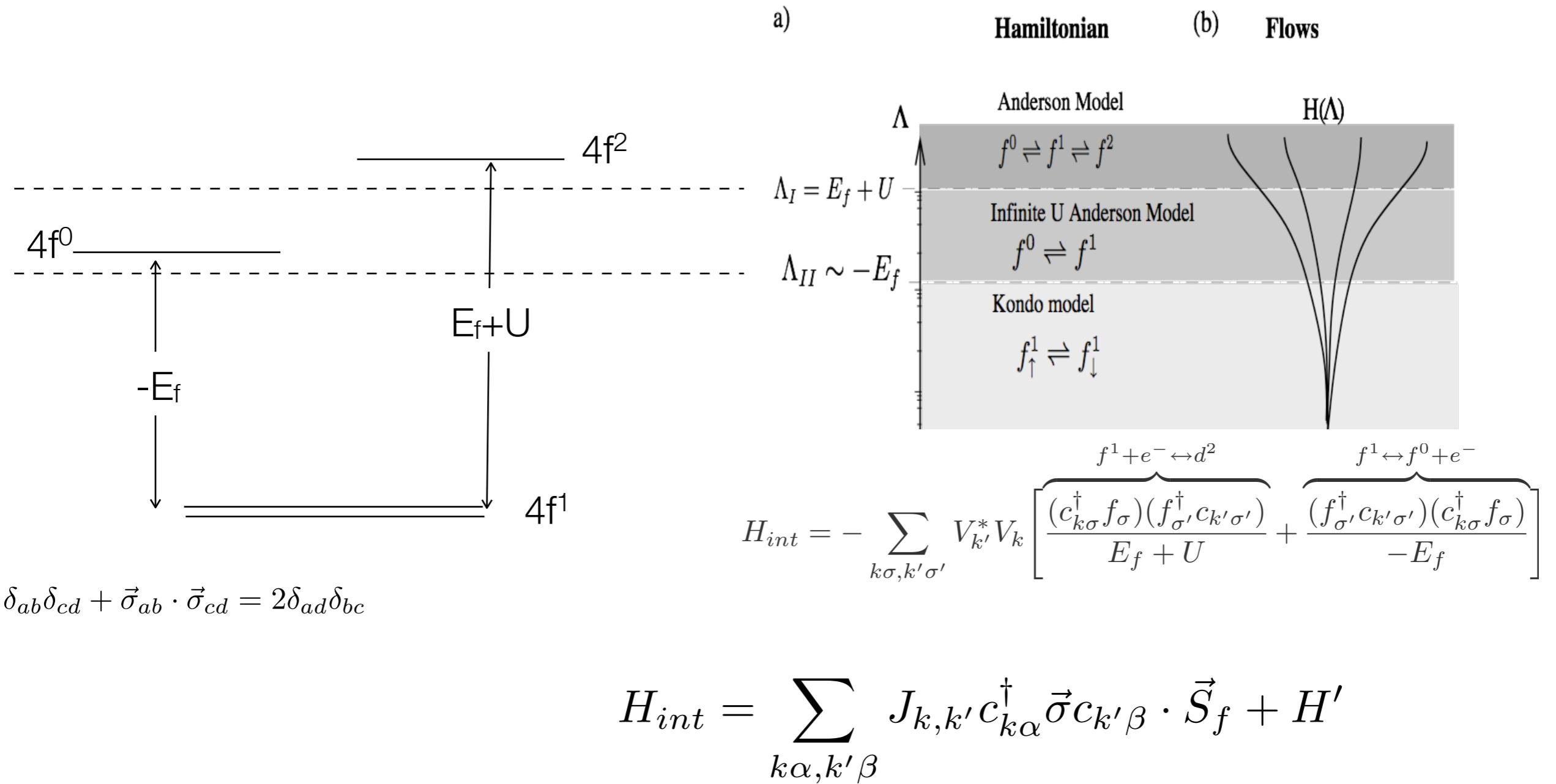
# Schrieffer-Wolff Transformation.



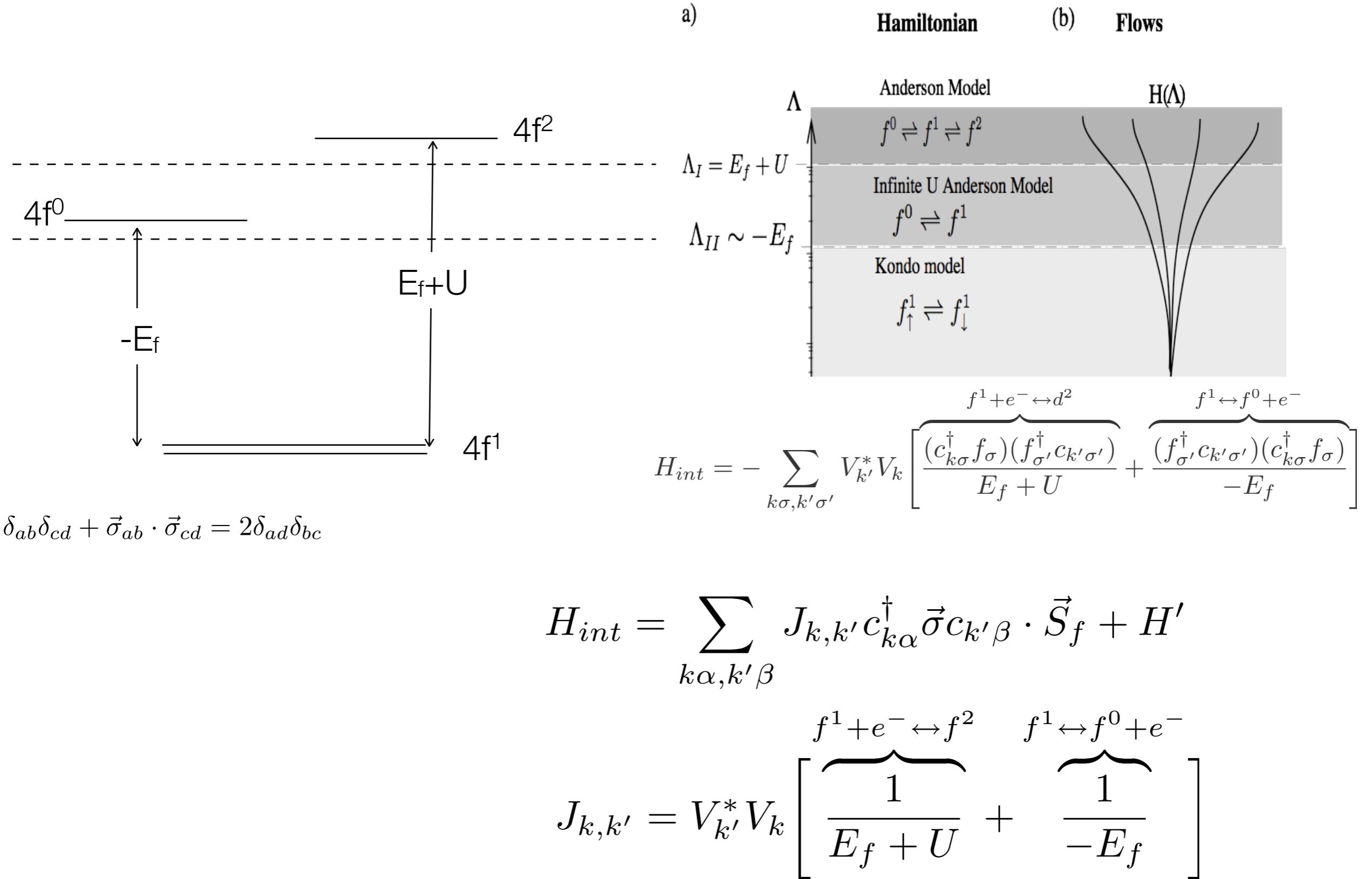
# Schrieffer-Wolff Transformation.

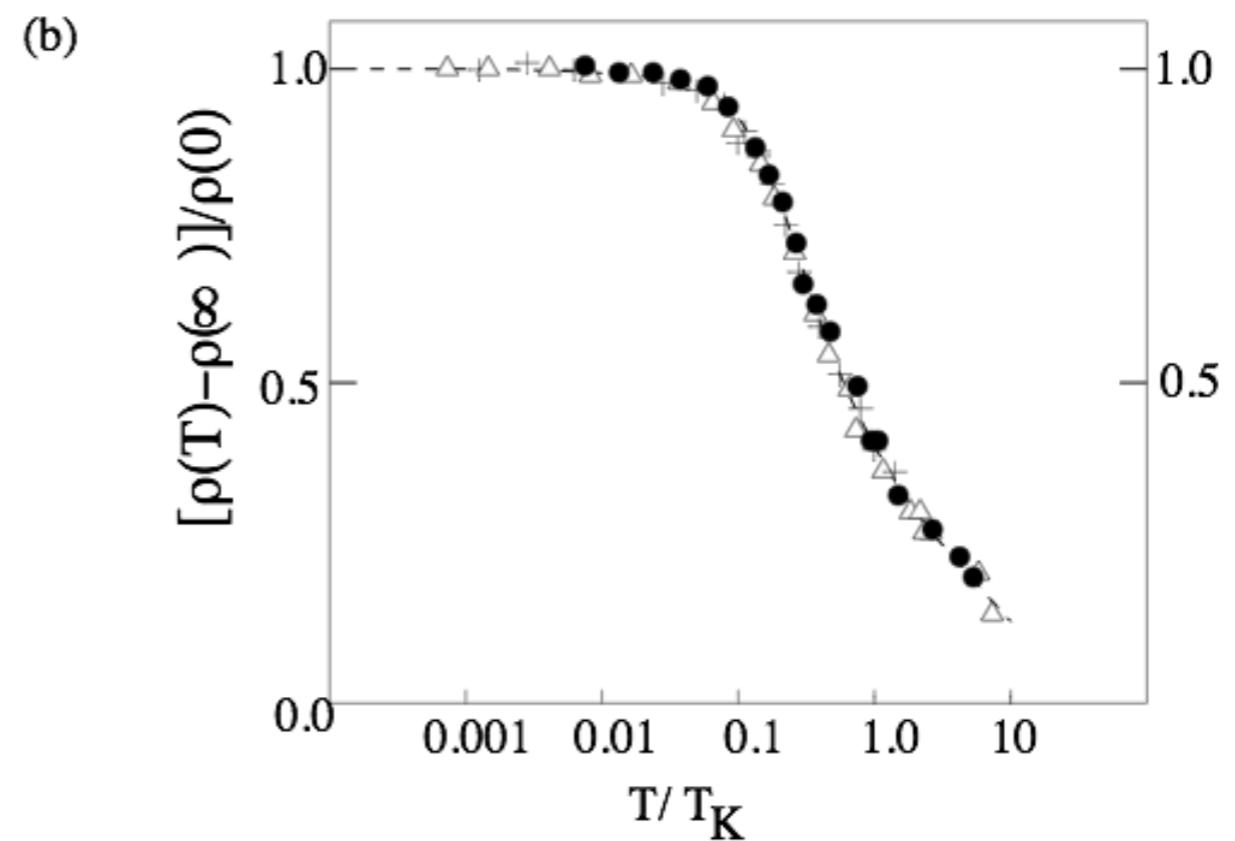
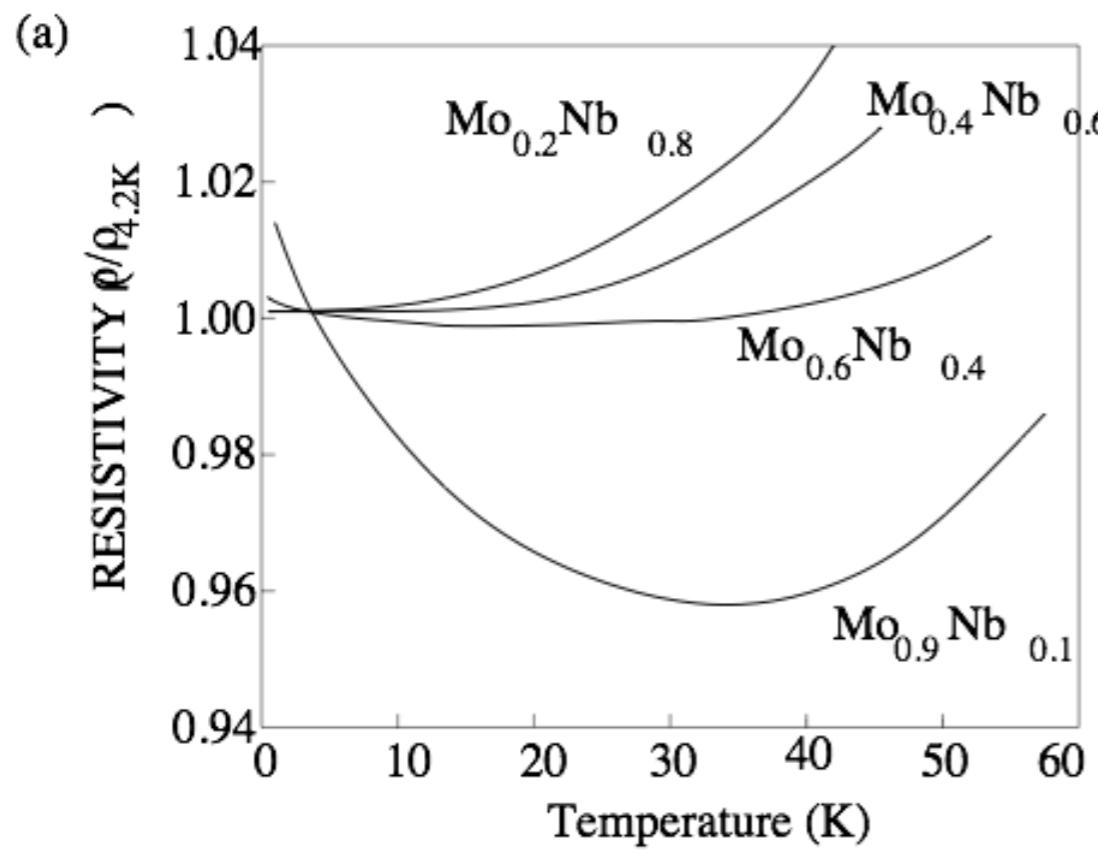


# Schrieffer-Wolff Transformation.



# Schrieffer-Wolff Transformation.





Mathias and Clogston (62)  
Sarachik et al, (1964)

White and Geballe, (1979)