

# From single-particle to many-body localisation in disordered systems.

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**NEC**



The 4th Windsor Summer School on Condensed Matter Theory  
**Quantum Transport and Dynamics in Nanostructures**  
Great Park, Windsor, UK, August 6 - 18, 2007

# *Lecture 3*

## *Many Body Localization*

# Previous Lectures:

1. Anderson Localization as Metal-Insulator Transition
2. Anderson model. Cayley tree problem.
3. Spectral Statistics and Localization.  
Poisson versus Wigner-Dyson.  
Invariant definition of the localization
4. Chaos and Localization.  
Localization in the space of quantum numbers.

Wigner-Dyson random matrix statistics follows from the delocalization.

**Q** ■ *Why the random matrix theory (RMT) works so well for nuclear spectra* ■ **?**

Many-Body excitations are delocalized !

What does it mean ?

Consider a finite system of quantum particles, e.g., fermions. Let the **one-particle spectra** be **chaotic** (Wigner-Dyson).

**Q** ■ What is the statistics of the **many-body** spectra? **?**

a. The particles **do not interact** with each other.

**Poisson:**

individual energies are conserving quantum numbers.

a. The particles **do interact**. **????**

*Part 4.*

*Many-Body  
excitation in finite  
systems.*

# Decay of a quasiparticle with an energy $\varepsilon$ in Landau Fermi liquid

$\varepsilon$  ●

$\varepsilon - \omega$  ●

$\varepsilon_1 + \omega$  ●

$\varepsilon_1$  ●

Fermi Sea

Landau Fermi liquid in a clean bulk system:

$$\gamma(\varepsilon) \propto \frac{\varepsilon^2}{\varepsilon_F}$$

Fermi energy

# Decay of a quasiparticle with an energy $\varepsilon$ in Landau Fermi liquid

$\varepsilon$  ●

Quantum dot – zero-dimensional case ?

$\varepsilon - \omega$  ●

$\varepsilon_1 + \omega$  ●

Fermi golden rule

( U.Sivan, Y.Imry & A.Aronov, 1994 ):

$$\gamma(\varepsilon) \propto \delta_1 \left( \frac{\varepsilon}{E_T} \right)^2$$

Mean  
level  
spacing

Thouless  
energy

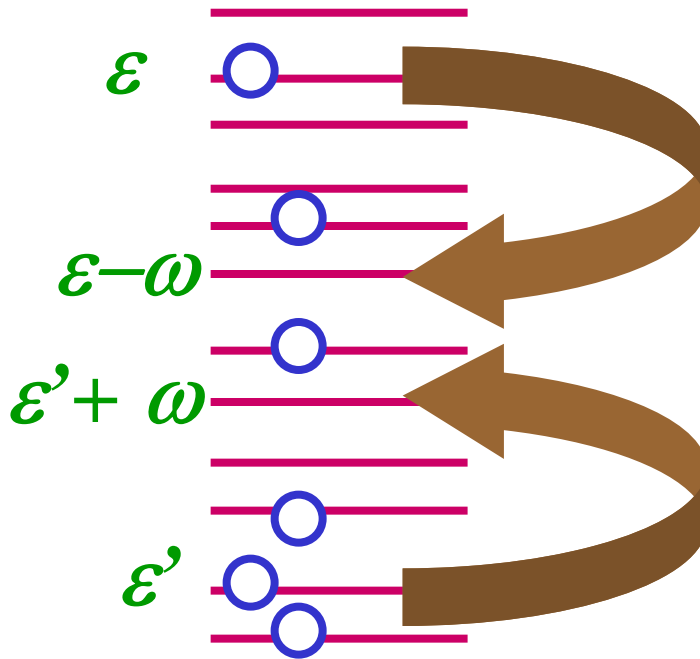
$\varepsilon_1$  ●

Fermi Sea



*Inelastic relaxation rate in 0D case*

$T=0$



**Offdiagonal**  
matrix  
element

$$M(\omega, \varepsilon, \varepsilon') \propto \frac{\delta_1}{g} \ll \delta_1$$

# Decay of a quasiparticle with an energy $\varepsilon$ in Landau Fermi liquid

$\varepsilon$  ●

$\varepsilon - \omega$  ●

$\varepsilon_1 + \omega$  ●

$\varepsilon_1$  ●

Fermi Sea

zero-dimensional case



one-particle spectrum is  
**discrete**



equation

$$\varepsilon_1 + \varepsilon_2 = \varepsilon'_1 + \varepsilon'_2$$

can not be satisfied exactly

???

# Decay of a quasiparticle with an energy $\varepsilon$ in Landau Fermi liquid

$\varepsilon$  ●

$\varepsilon - \omega$  ●

$\varepsilon_1 + \omega$  ●

$\varepsilon_1$  ●

Fermi Sea

zero-dimensional case



one-particle spectrum is  
**discrete**



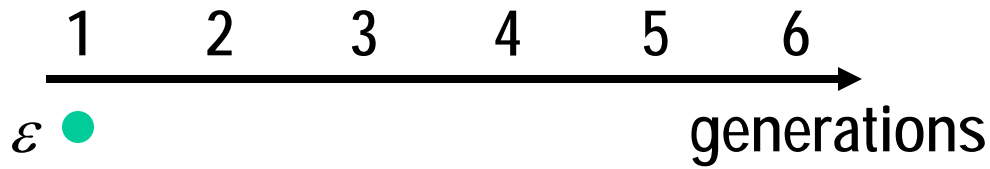
equation

$$\varepsilon_1 + \varepsilon_2 = \varepsilon'_1 + \varepsilon'_2$$

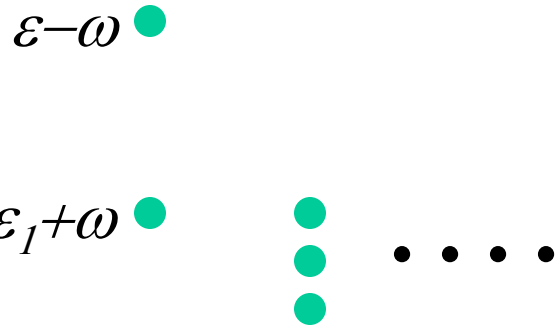
can not be satisfied exactly

**Recall:** in the Anderson model the site-to-site hopping **does not** conserve the energy

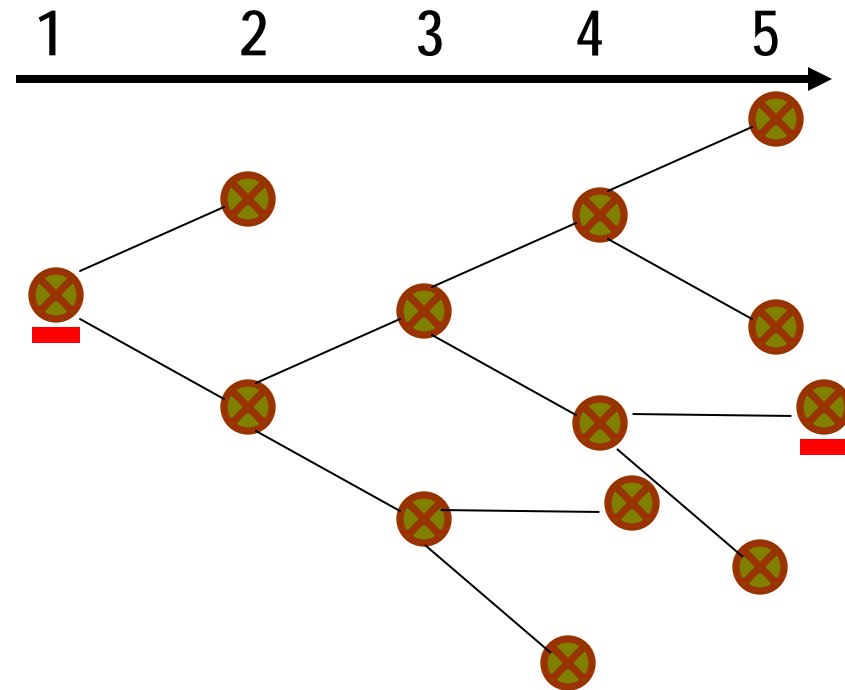
# Chaos in Nuclei – Delocalization?

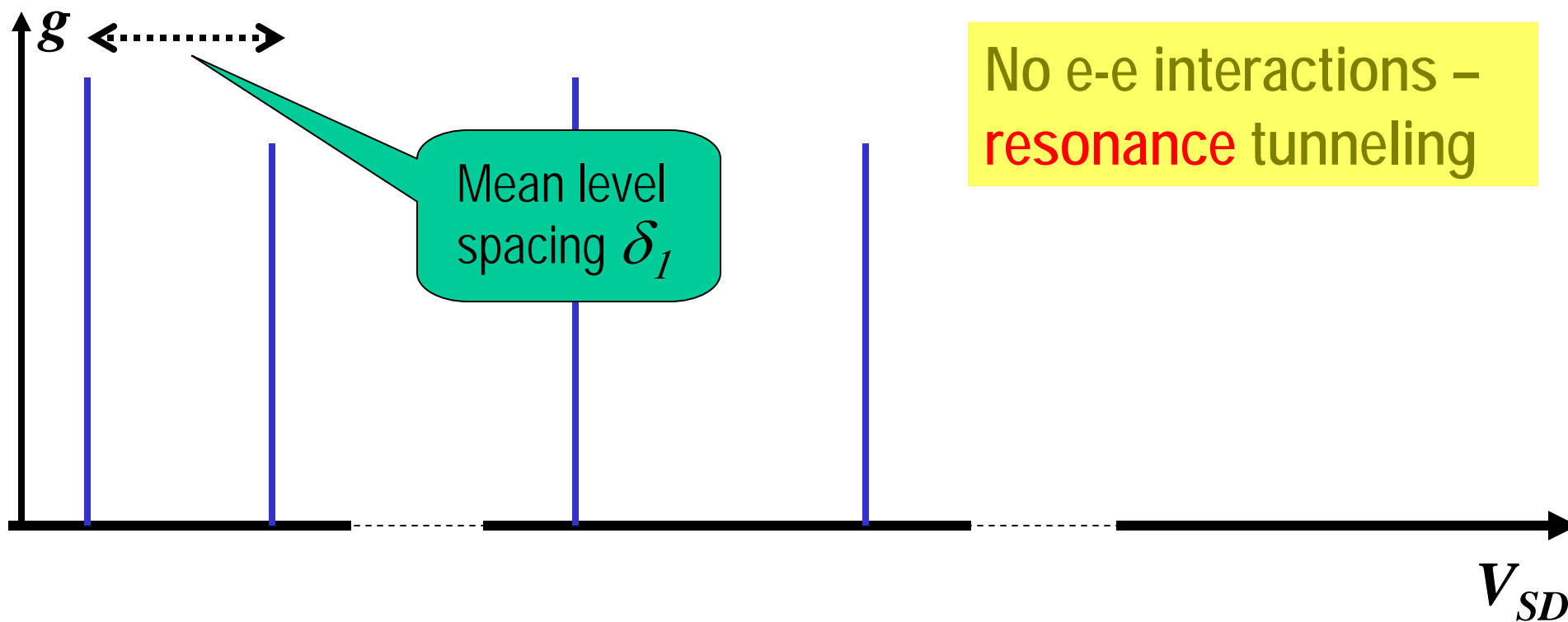
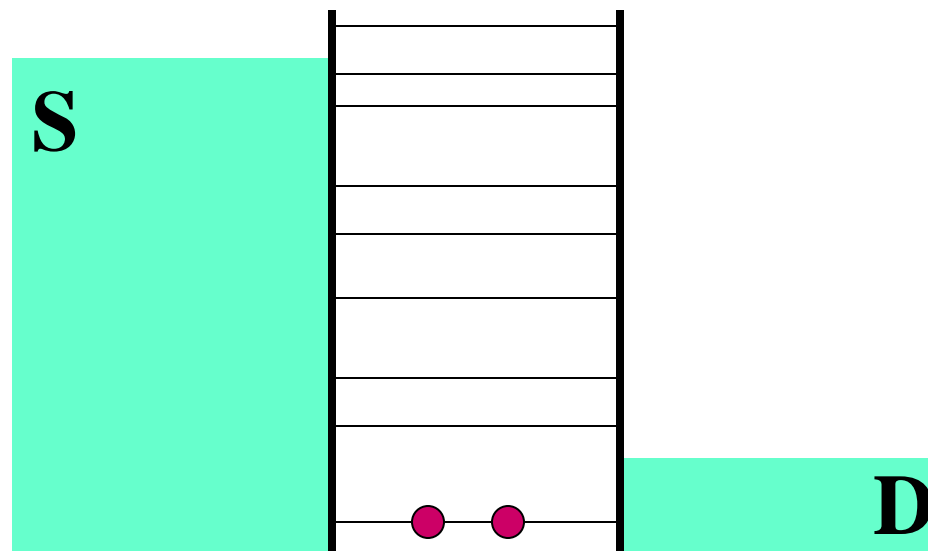
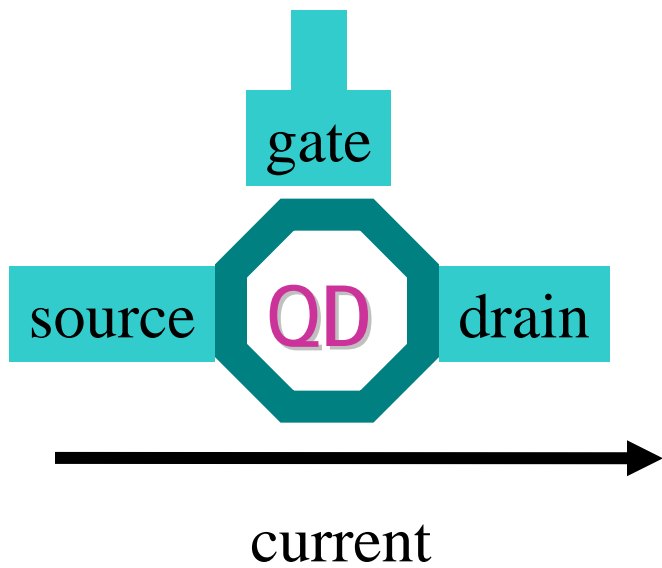


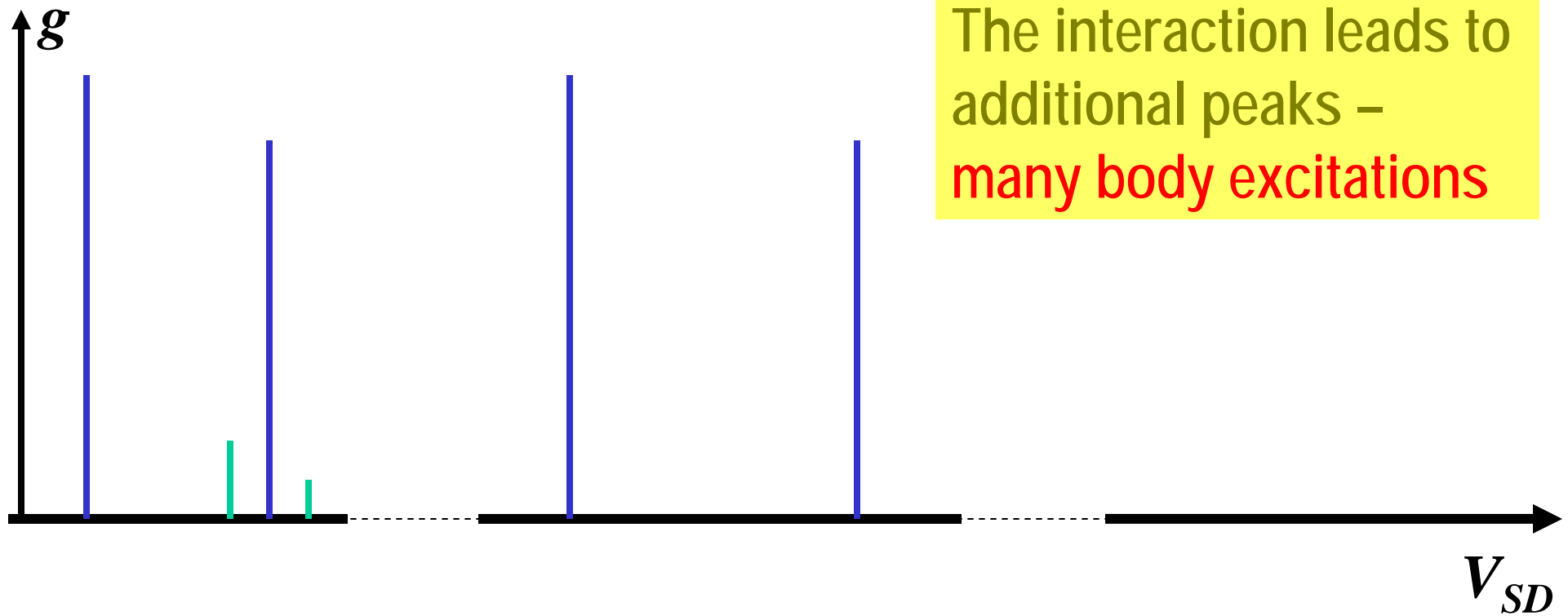
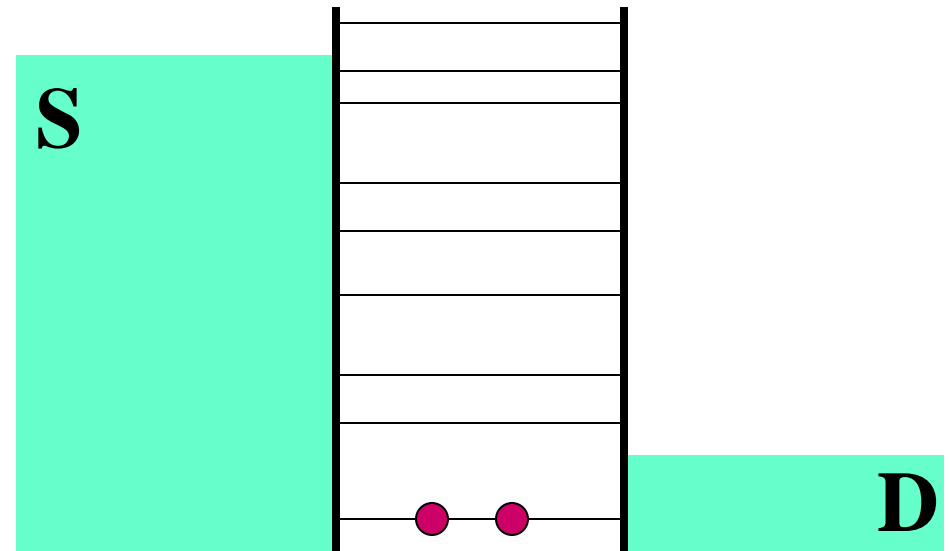
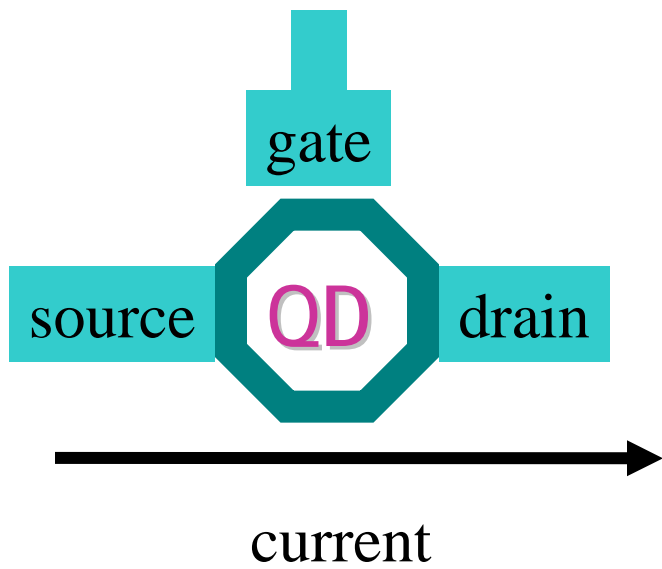
Delocalization  
in Fock space



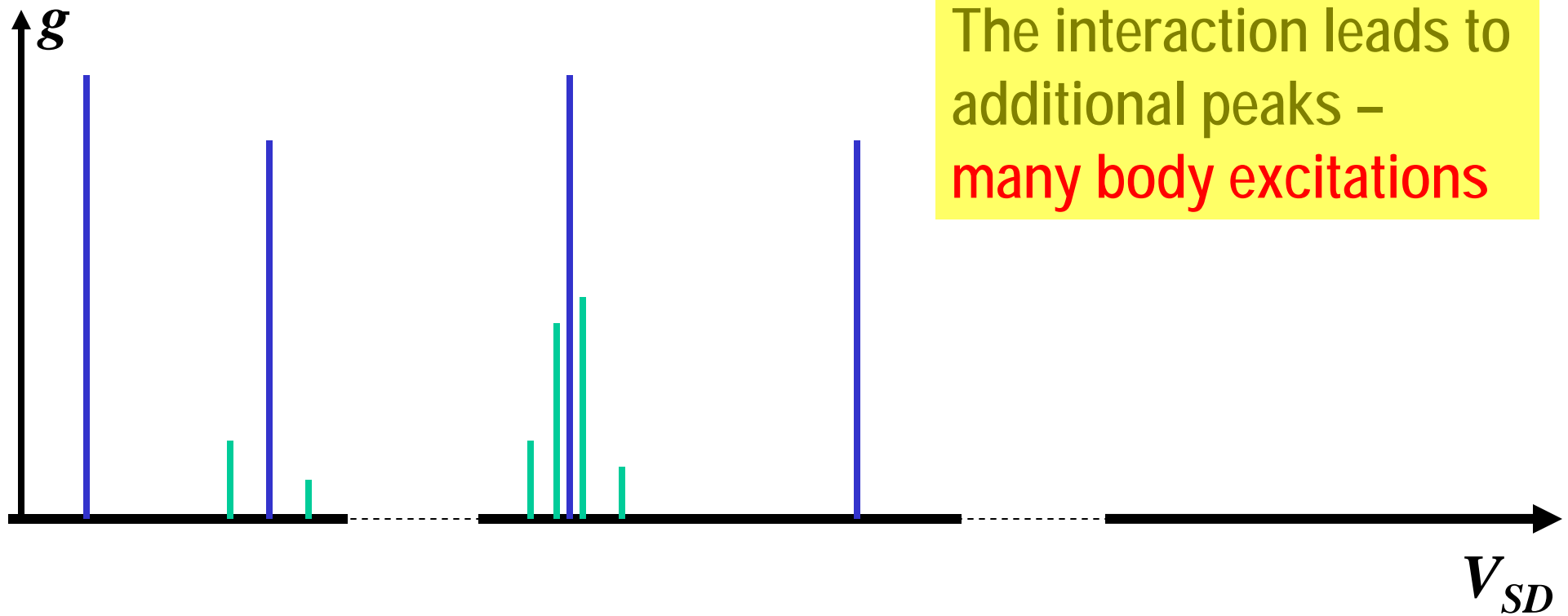
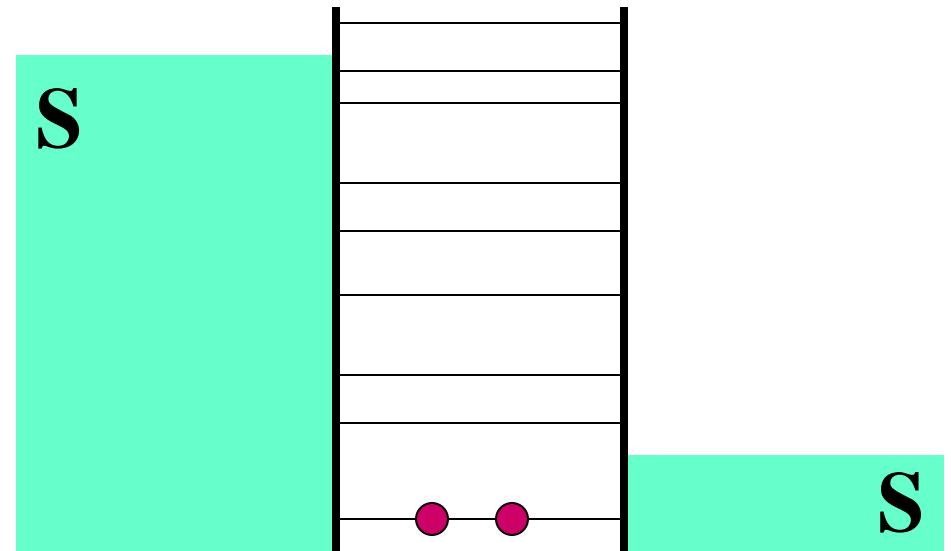
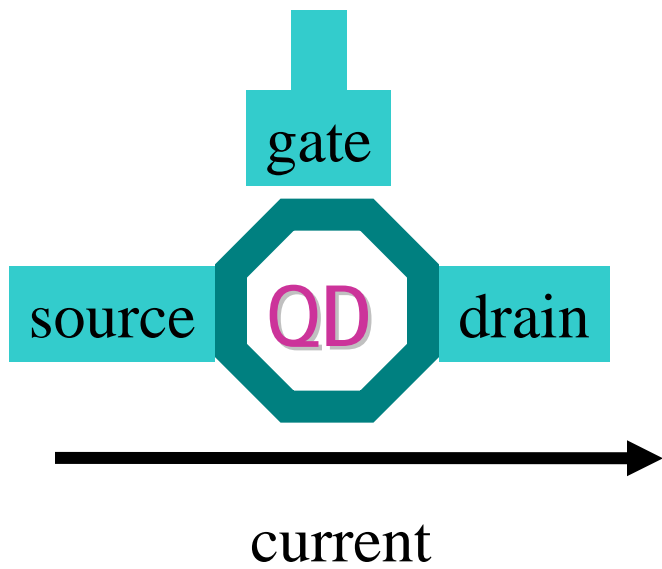
Can be mapped (approximately)  
to the problem of localization  
on Cayley tree



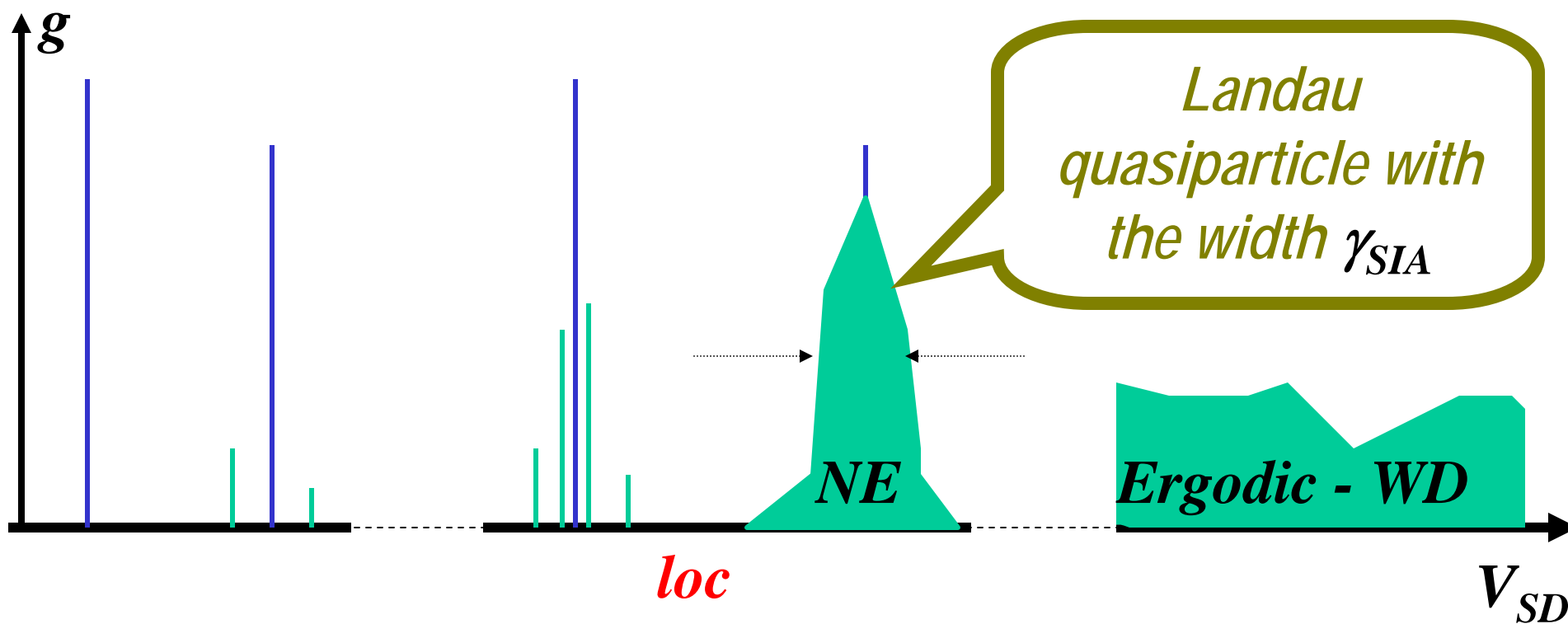
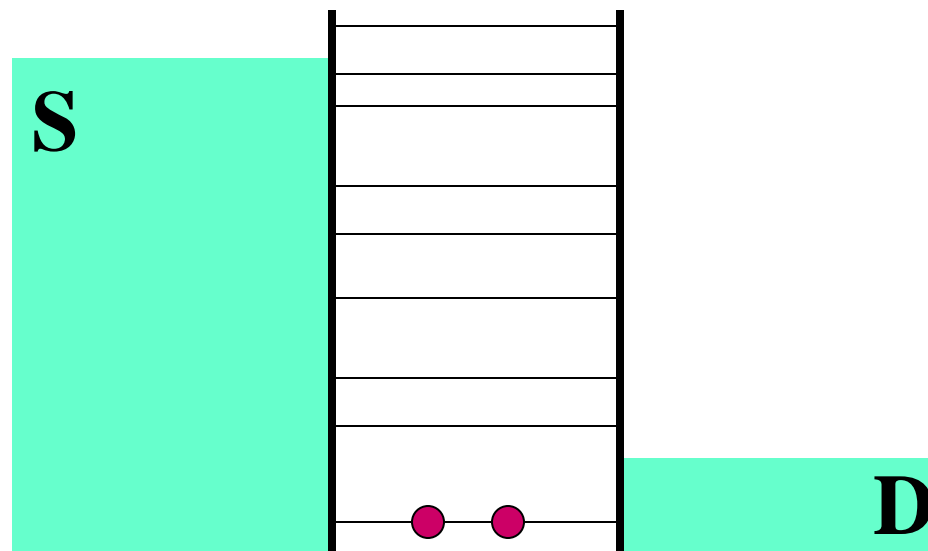
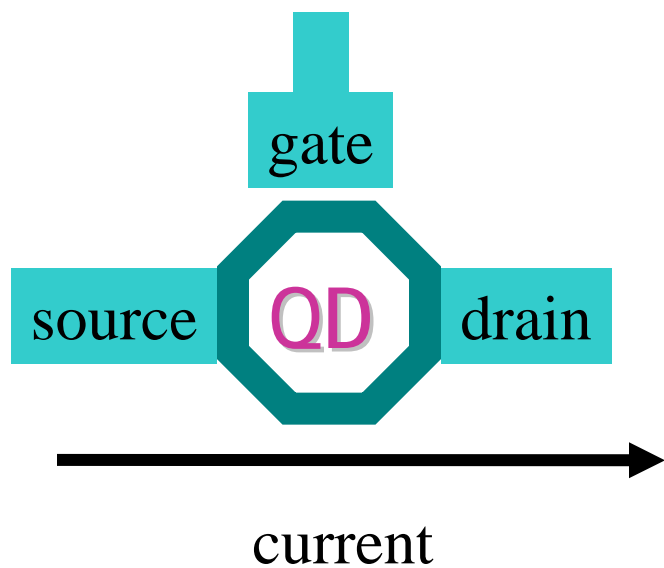




The interaction leads to additional peaks – many body excitations



The interaction leads to additional peaks – many body excitations





## nonergodic states

Such a state occupies infinitely many sites of the Anderson model but still negligible fraction of the total number of sites

$N$

total  
number of  
sites in the  
system

$n$

support of a  
given wave  
function

$N \rightarrow \infty$

$n \rightarrow \text{const}$

localized

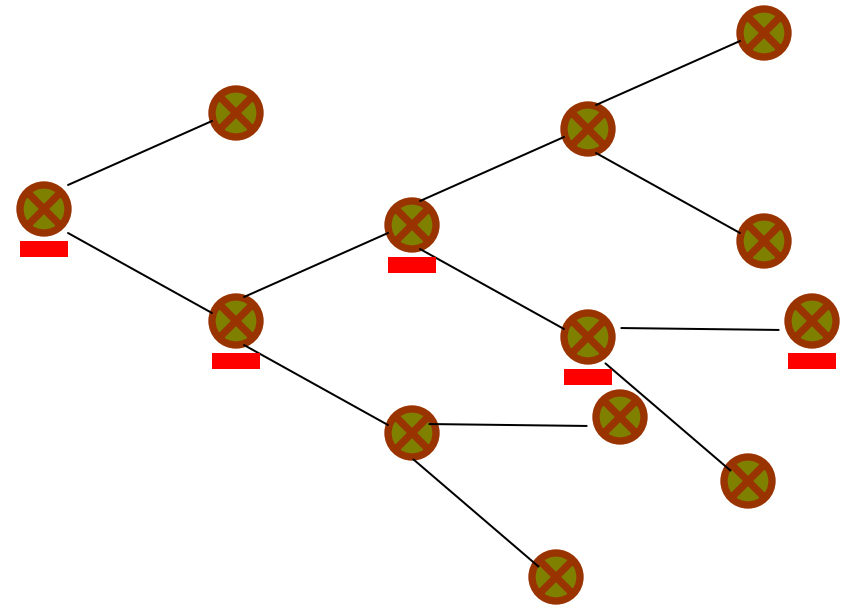
$n \rightarrow \infty$

extended

$$W > I > W/K$$

There is of order one  
resonance at every step

$$n \sim \ln N \quad \text{nonergodic}$$



*Part 5.*

*Many-Body  
Localization*

# Metal-insulator transition in a weakly interacting many-electron system with localized single-particle states

Annals of Physics, v. 321, p. 1126-1205 (2006)

D. M. Basko & I.L. Aleiner

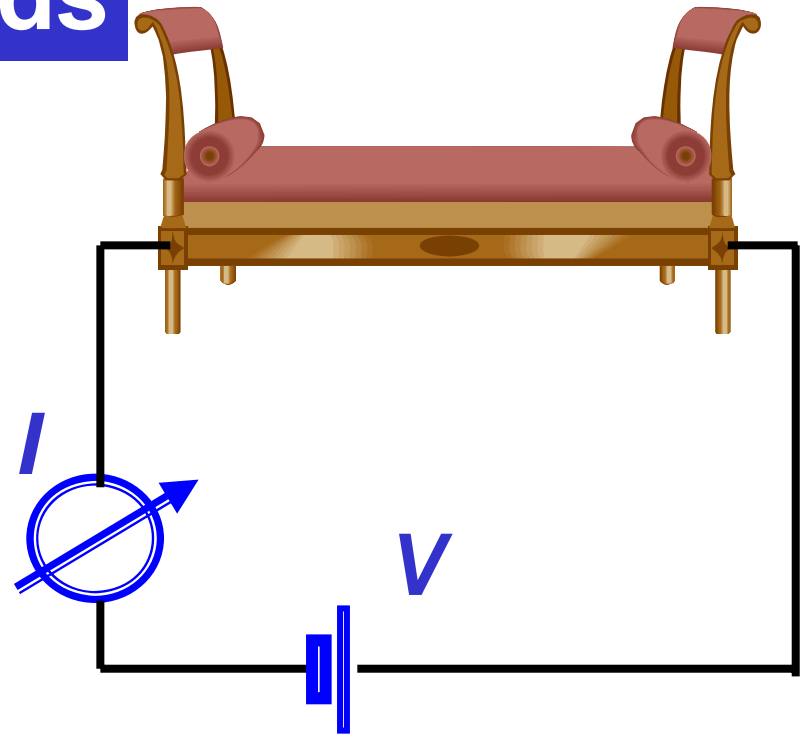
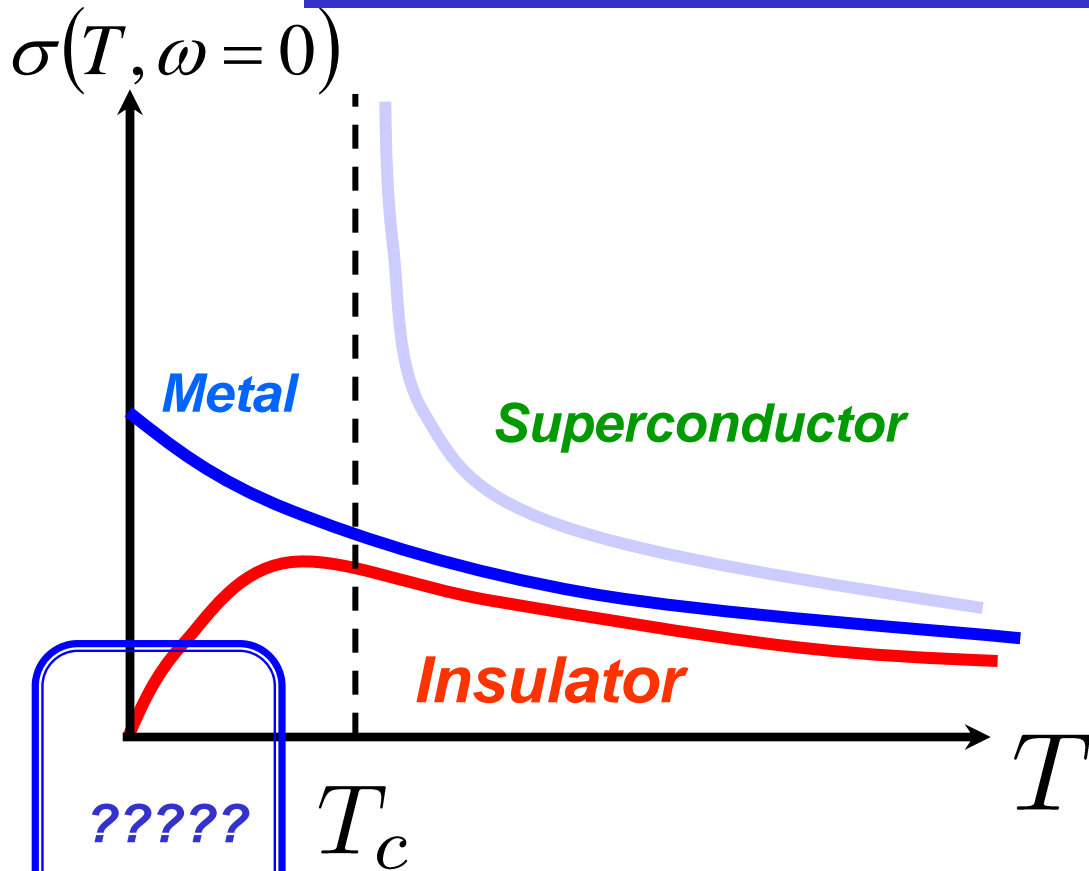
*Columbia University*



B. L. Altshuler

*Columbia University, NEC-Laboratories America*

# Transport in solids

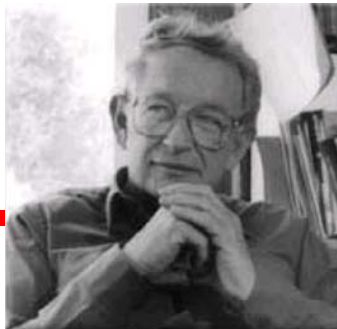


**Conductance:**  $G(\omega, T) = \left. \frac{I}{V} \right|_{V \rightarrow 0}$

**Conductivity:**  $G(\omega, T) = \sigma(\omega, T) \frac{L_x L_y}{L_z}$

# Disorder + interactions

strength  
of the  
disorder



strength  
of the  
interaction

Fermi liquid

Wigner crystal

$r_s$

Can hopping conductivity  
exist **without phonons**



**Question:** can  $e$ - $e$  interaction **alone** sustain hopping conduction in a localized system?

- Given:**
1. All one-electron states are localized
  2. Electrons interact with each other
  3. The system is closed (no phonons)
  4. Temperature is low but finite

**Find:** DC conductivity  $\sigma(T, \omega=0)$

**zero or finite?**



“All states are **localized**”

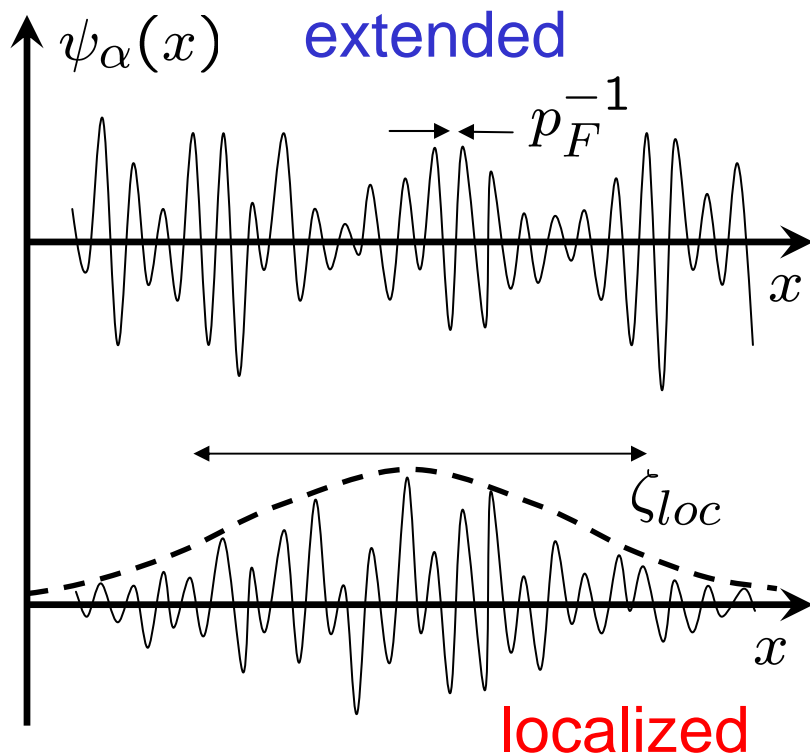
means that

probability to find a state extended over the system size  $L$  is

$$\mathcal{P}_{ext} \propto \exp\left(-\# \frac{L}{\zeta_{loc}}\right)$$

# 1. Localization of single-electron wave-functions:

$$\left[ -\frac{\nabla^2}{2m} + U(\mathbf{r}) - \epsilon_F \right] \psi_\alpha(\mathbf{r}) = \xi_\alpha \psi_\alpha(\mathbf{r})$$



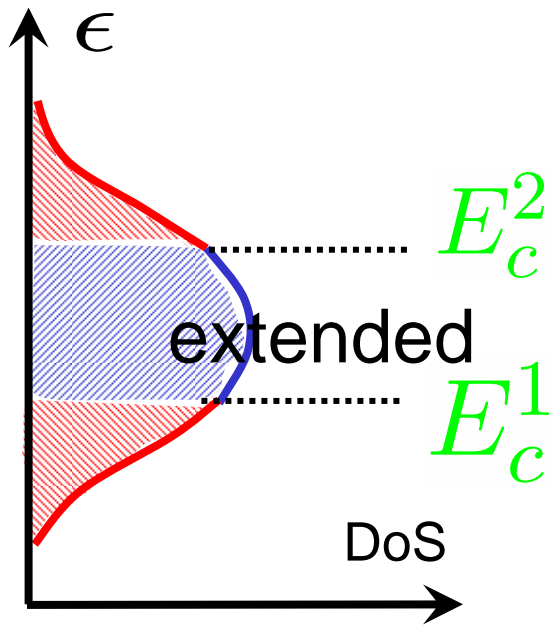
$d=1$ ; All states are localized

$d=2$ ; All states are localized

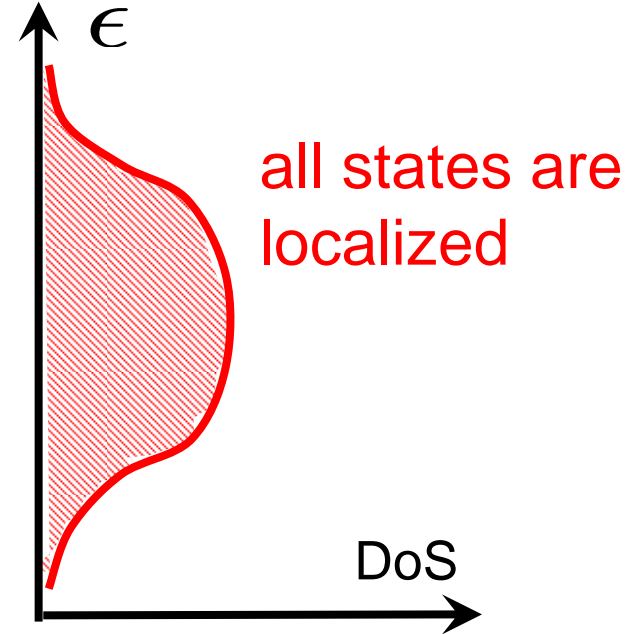
$d>2$ ; Anderson transition

# Anderson Transition

$$I > I_c$$



$$I < I_c$$

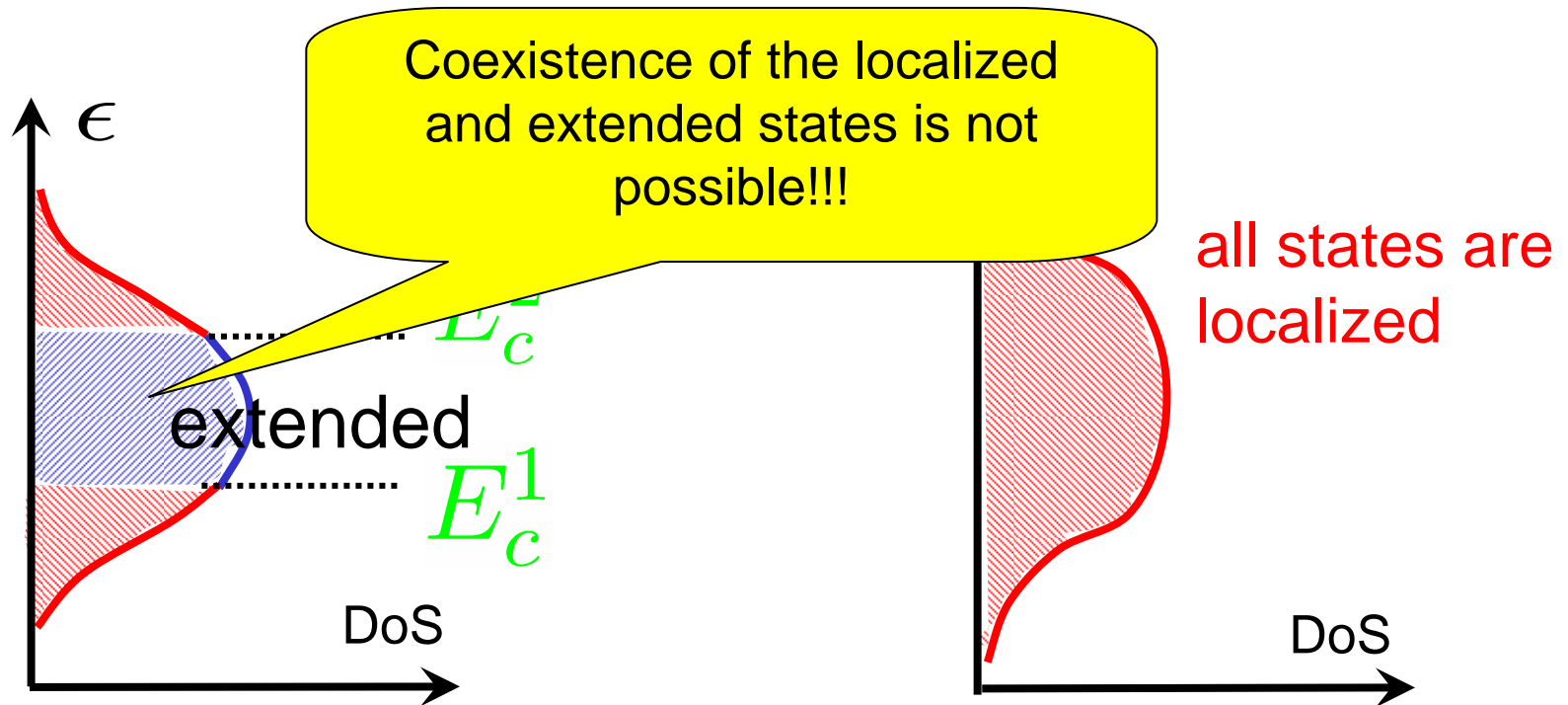


$E_c$  - mobility edges (one particle)

# Anderson Transition

$$I > I_c$$

$$I < I_c$$



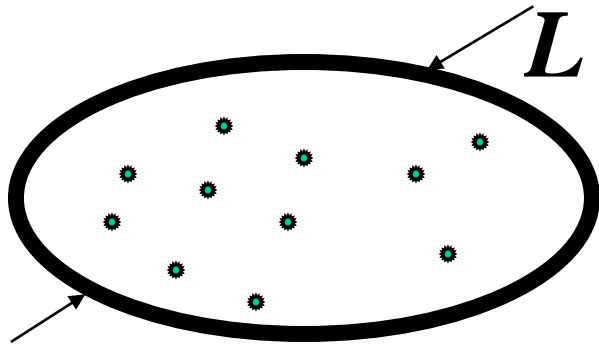
$E_c$  - mobility edges (one particle)

# Quantum particle in a random potential (*Thouless, 1972*)

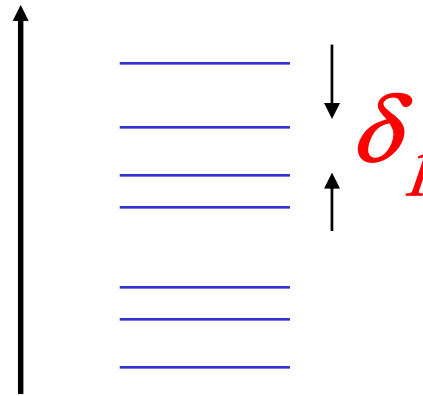
## Energy scales

### 1. Mean level spacing

$$\delta_1 = 1/\nu \times L^d$$



energy



$L$  is the system size;

$d$  is the number of dimensions

### 2. Thouless energy

$$E_T = \hbar D / L^2$$

$D$  is the diffusion const

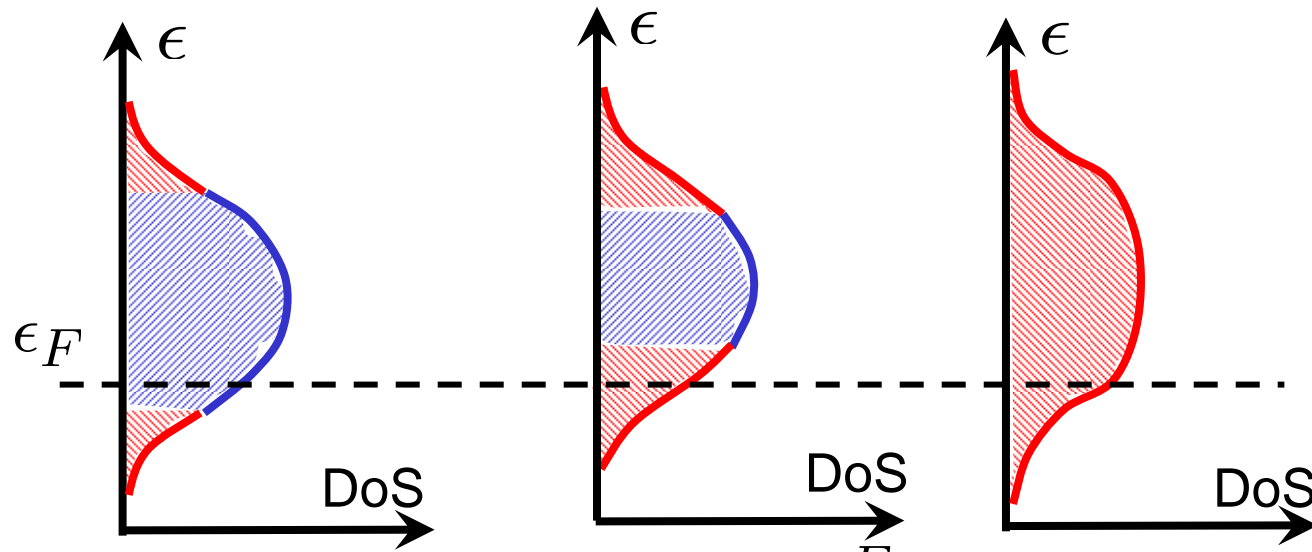
$E_T$  has a meaning of the *inverse diffusion time* of the traveling through the system or the *escape rate* (for open systems)

$$g = E_T / \delta_1$$

dimensionless  
*Thouless*  
conductance

$$g = Gh/e^2$$

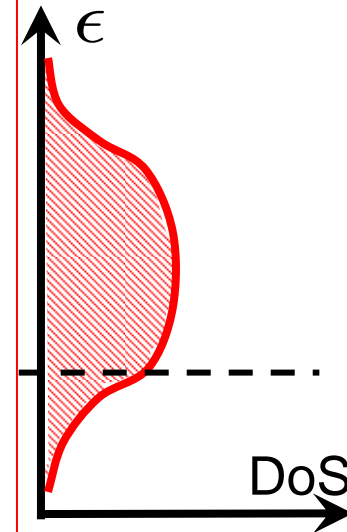
# *Temperature dependence of the conductivity of noninteracting electrons*



$$\sigma(T \rightarrow 0) > \sigma(T) \propto e^{-\frac{E_c - \epsilon_F}{T}} \quad \sigma(T) = 0$$

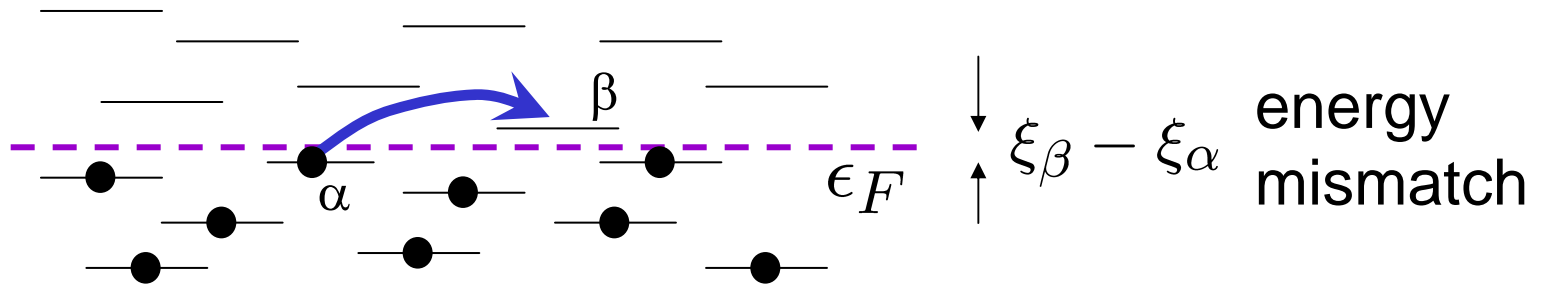
# *Temperature dependence of the conductivity of noninteracting electrons*

Assume that all the states  
are localized



$$\underline{E} \quad \sigma(T) = 0$$

# Inelastic processes - transitions between localized states



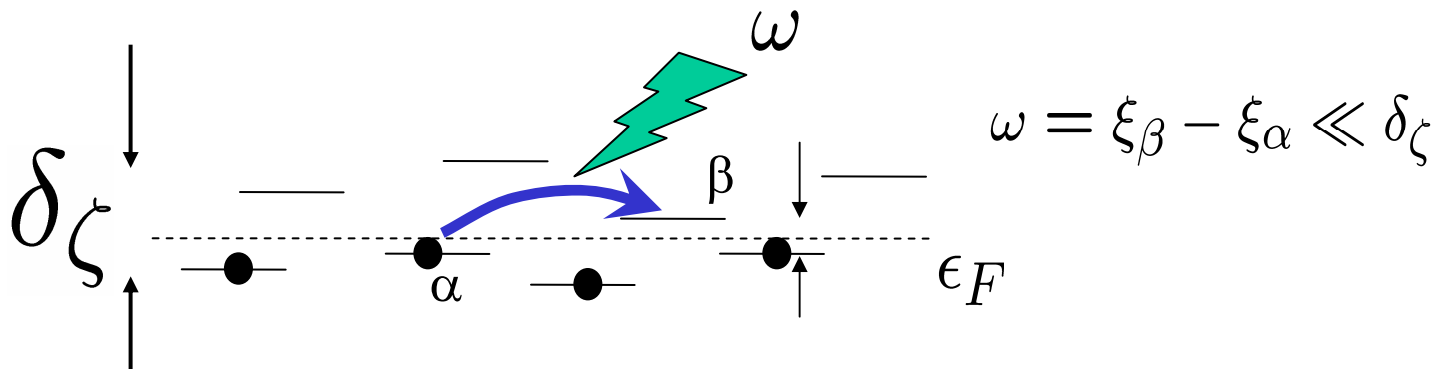
$$\sigma(T) \propto \Gamma_\alpha \text{ (inelastic lifetime)}^{-1}$$

$$T = 0 \Rightarrow \sigma = 0 \text{ (any mechanism)}$$

$$T > 0 \Rightarrow \sigma = ?$$



# Phonon-induced hopping



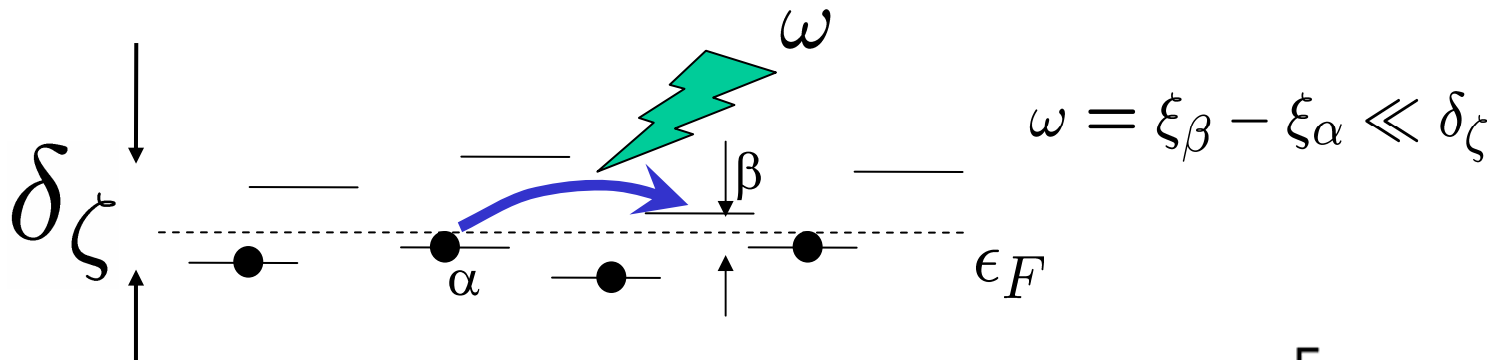
$$\omega = \xi_\beta - \xi_\alpha \ll \delta_\zeta$$

$$\delta_\zeta \equiv \frac{1}{v \zeta_{loc}^d}$$

localization spacing

energy difference can be matched by a phonon

# Phonon-induced hopping



Variable Range Hopping  
N.F. Mott (1968)

$$\sigma(T) \propto T^\gamma \exp \left[ - \left( \frac{\delta_\zeta}{T} \right)^{\frac{1}{d+1}} \right]$$

Mechanism-dependent  
prefactor

Optimized  
phase volume

Any bath with a continuous spectrum of **delocalized excitations** down to  $\omega = 0$  will give the same exponential

**Q** ■ Can e-h pairs replace phonons and lead to **phonon-less** *Variable Range Hopping* ?

**A#1:** Sure

**Easy steps:**

1) Recall phonon-less AC conductivity:

Sir N.F. Mott (1970)

$$\sigma(\omega) \simeq \frac{e^2 \zeta_{loc}^{d-2}}{\hbar} \left( \frac{\hbar\omega}{\delta\zeta} \right)^2 \ln^{d+1} \left| \frac{\delta\zeta}{\hbar\omega} \right|$$

2) Calculate the Nyquist noise.

3) Use the electric noise instead of phonons.

4) Do self-consistency (whatever it means).

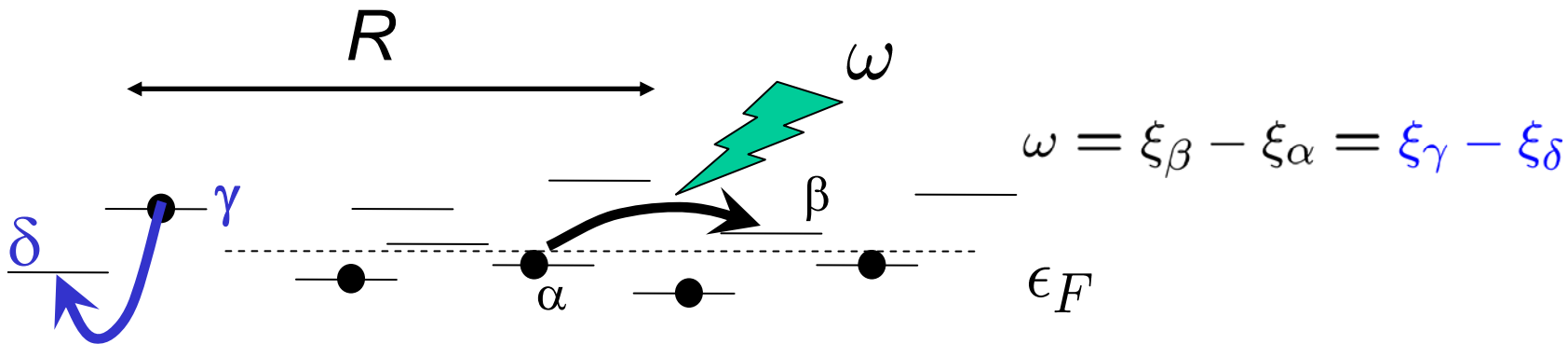
**Q** ■ Can e-h pairs replace phonons and lead to **phonon-less Variable Range Hopping** ?

A#1: Sure

A#2: No way [L. Fleishman, P.W. Anderson (1980)]  
 (for Coulomb interaction in 3D – may be)

$$\sigma(\omega) \simeq \frac{e^2 \zeta_{loc}^{d-2}}{\hbar} \left( \frac{\hbar\omega}{\delta\zeta} \right)^2 \ln^{d+1} \left| \frac{\delta\zeta}{\hbar\omega} \right|$$

is contributed by rare resonances

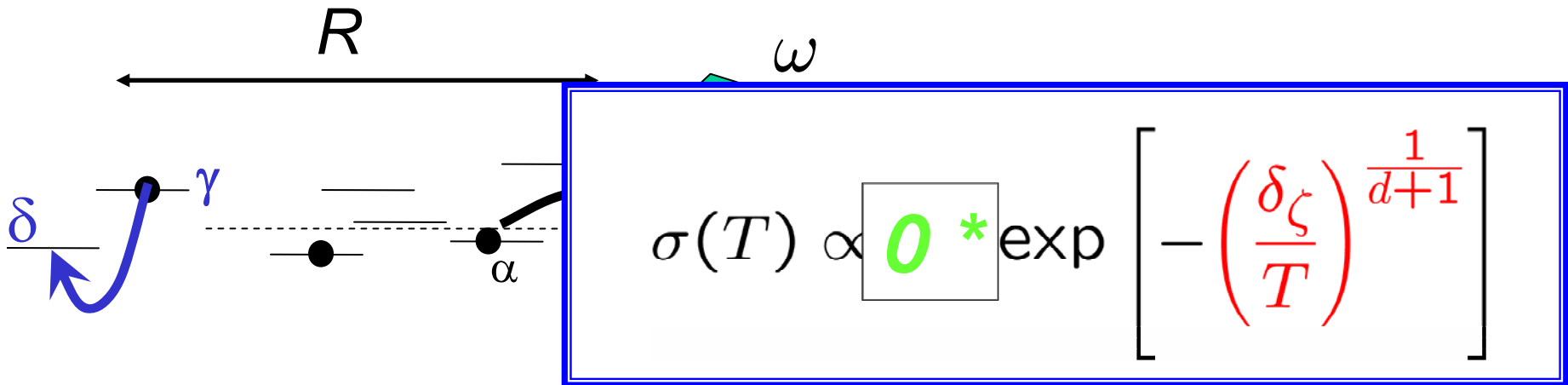


**Q** ■ Can e-h pairs replace phonons and lead to **phonon-less Variable Range Hopping** ?

A#1: Sure

A#2: No way [L. Fleishman, P.W. Anderson (1980)]  
 (for Coulomb interaction in 2D, may be)

$R \rightarrow \infty$  Thus, the matrix element vanishes !!!



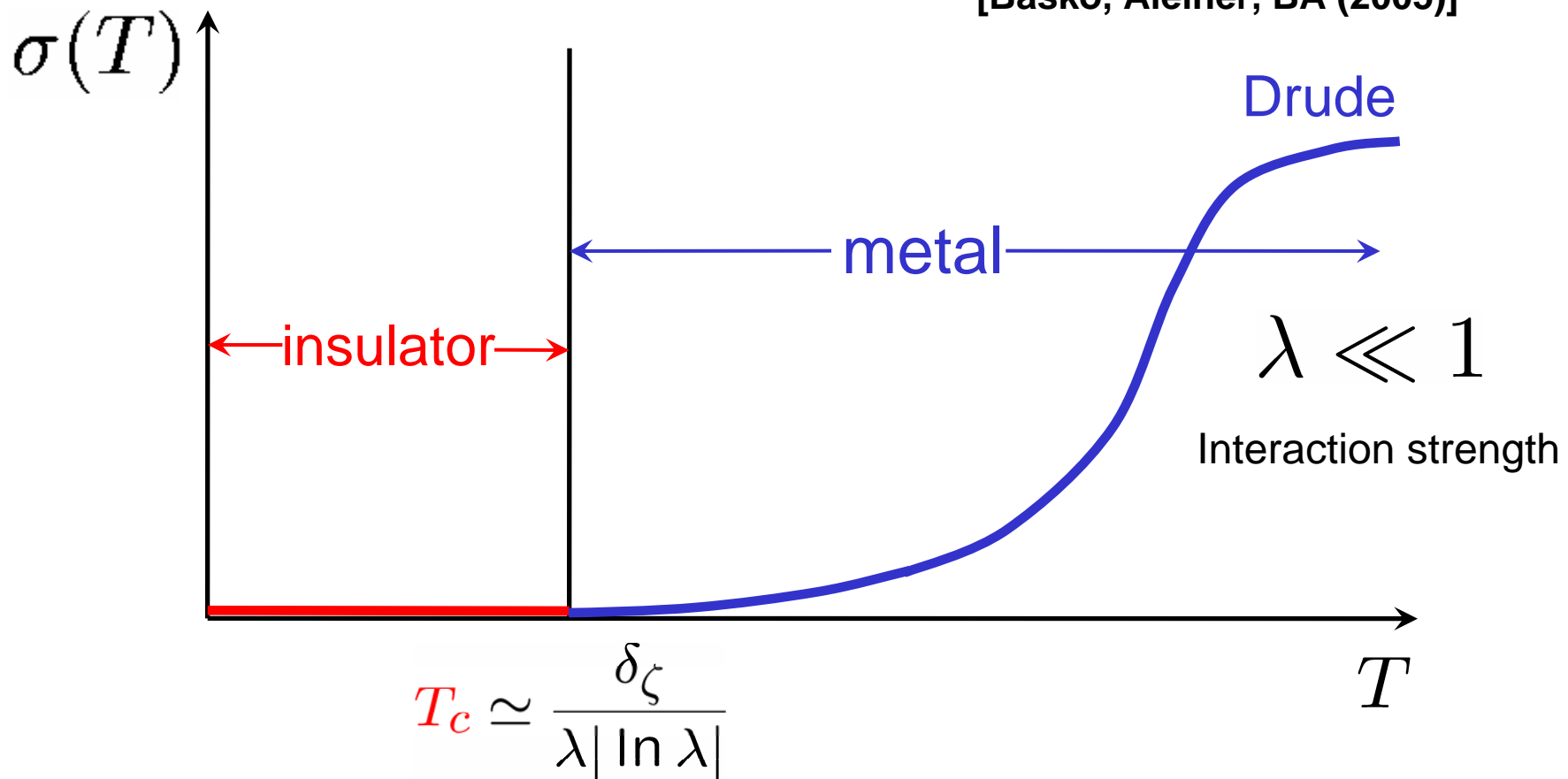
Q: Can we replace phonons with e-h pairs and obtain phonon-less VRH?

A#1: Sure [a person from the street (2005)]:

A#2: No way [L. Fleishman, P.W. Anderson (1980)]

A#3: Finite  $T$  Metal-Insulator Transition

[Basko, Aleiner, BA (2005)]



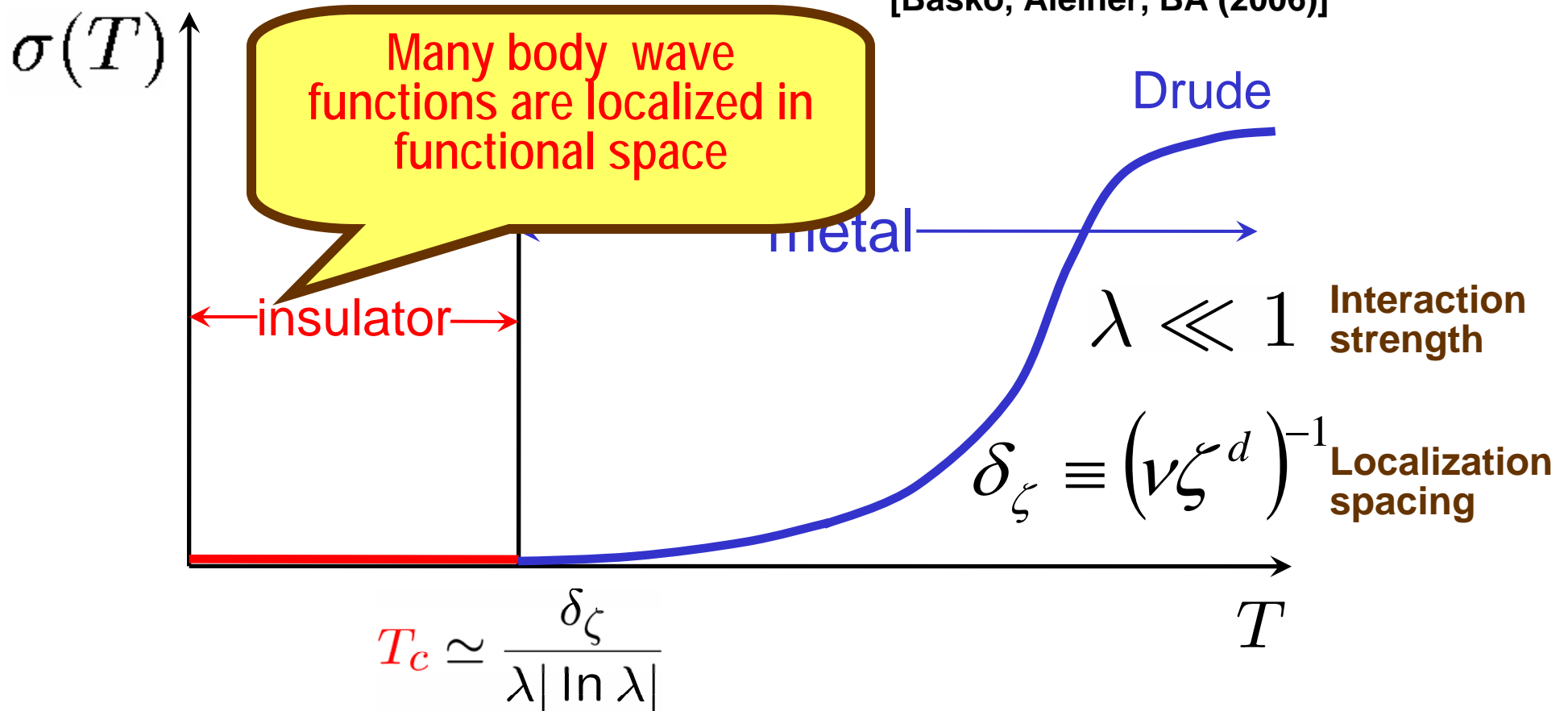
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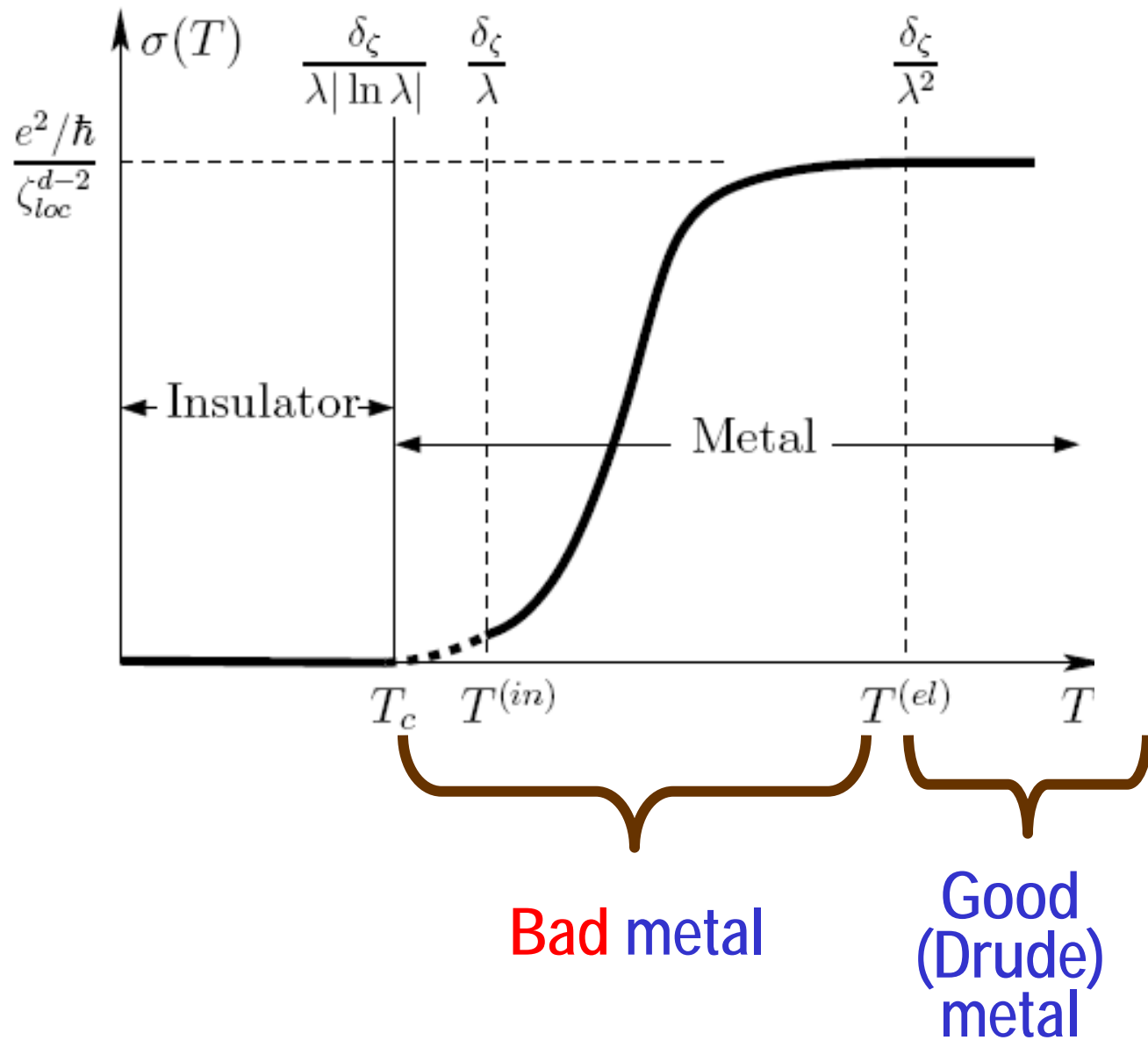
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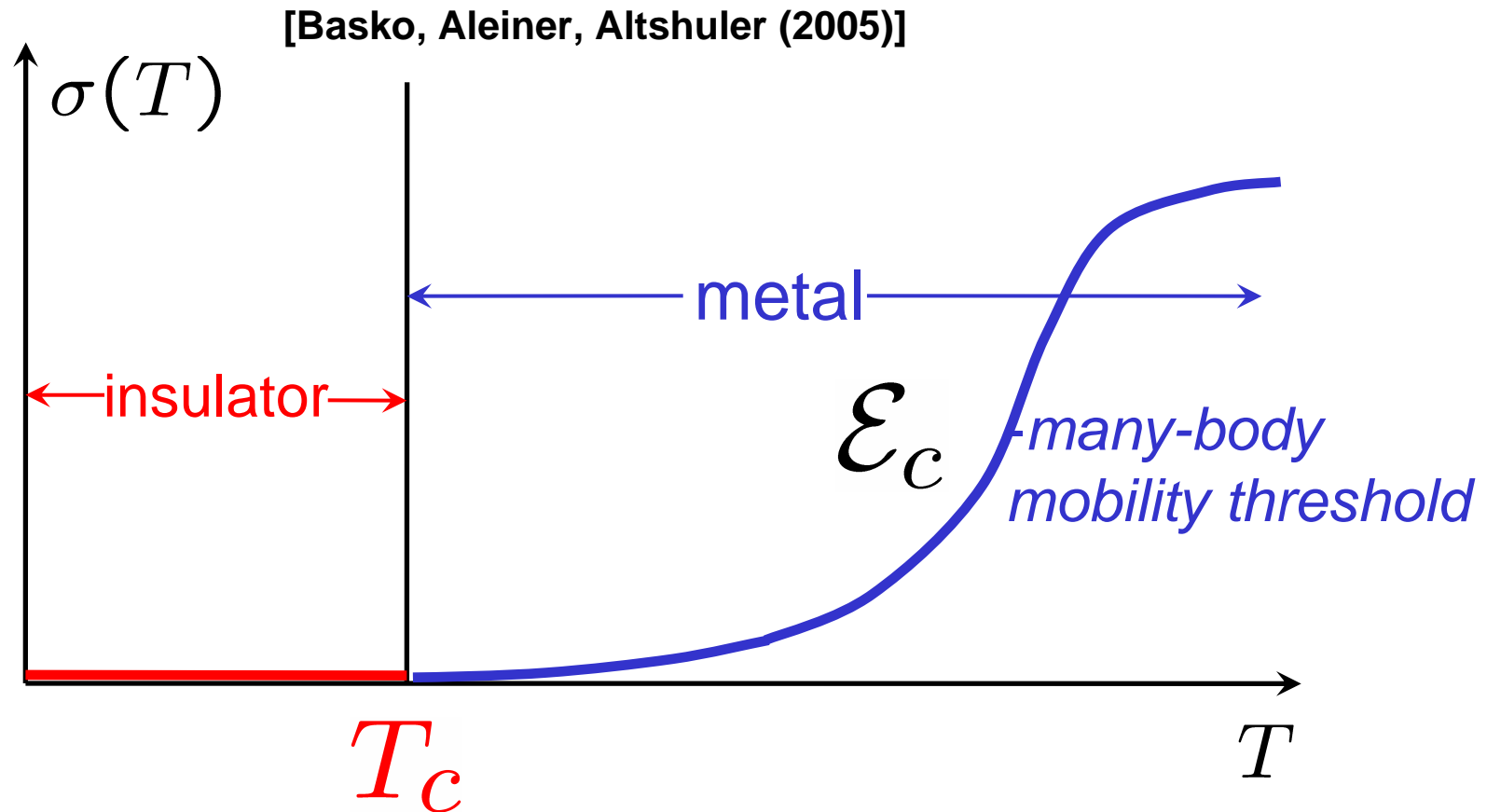






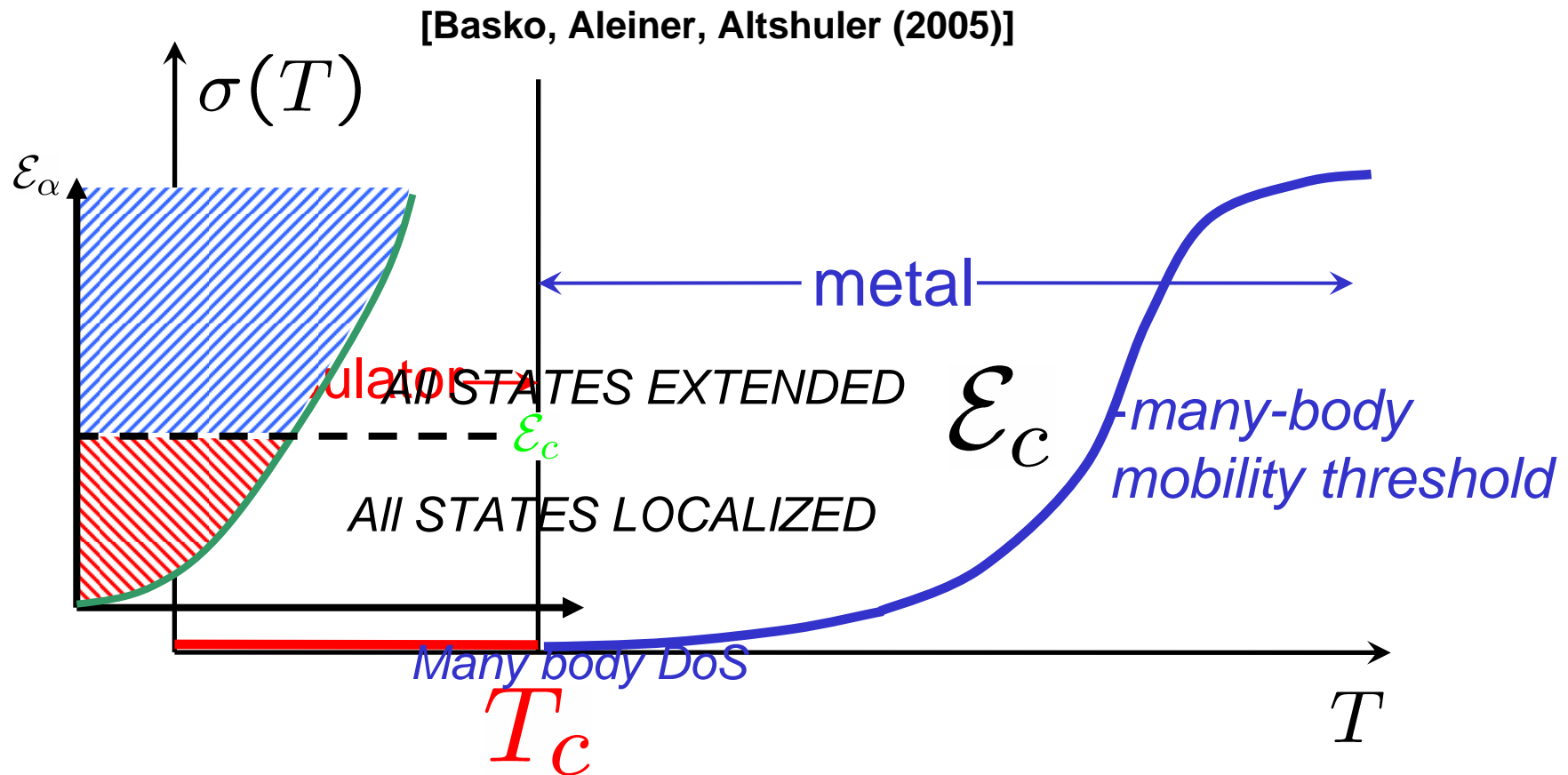
# Many-body mobility threshold

$$\left[ \hat{H}_1 + \hat{H}_{int} \right] \Psi_\alpha = \mathcal{E}_\alpha \Psi_\alpha$$



# Many-body mobility threshold

$$\left[ \hat{H}_1 + \hat{H}_{int} \right] \Psi_\alpha = \mathcal{E}_\alpha \Psi_\alpha$$



“All states are **localized**”

means that

Probability to find an extended state:

$$\mathcal{P}_{ext} \propto \exp \left( -\# \frac{\mathcal{V}}{\mathcal{V}_{loc}(\mathcal{E})} \right)$$

$$\lim_{\mathcal{E} \rightarrow \mathcal{E}_c - 0} \mathcal{V}_{loc}(\mathcal{E}) = \infty$$

System volume

# Main energy scale

$$\delta_{\xi} \equiv \frac{1}{\nu \xi^d}$$

Energy spacing  
between the states  
localized nearby

$\xi$

localization  
length

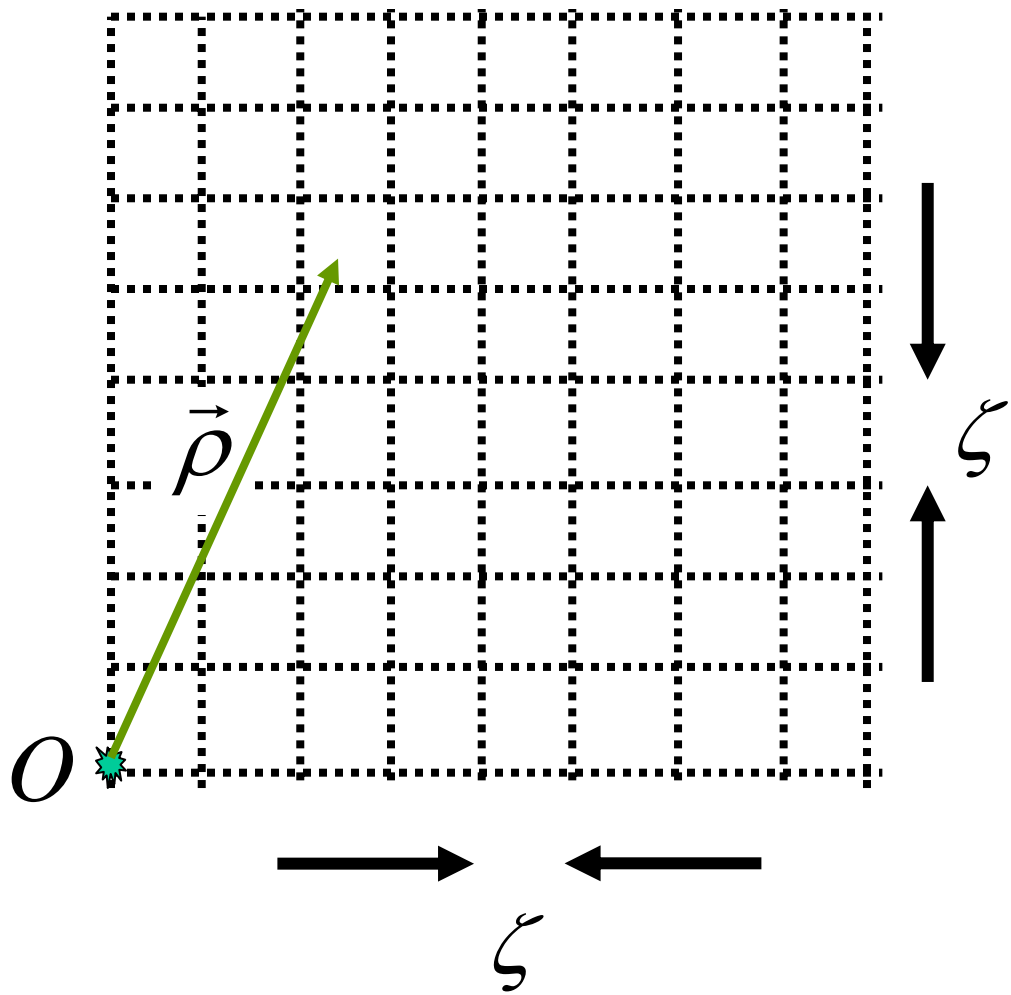
$\nu$

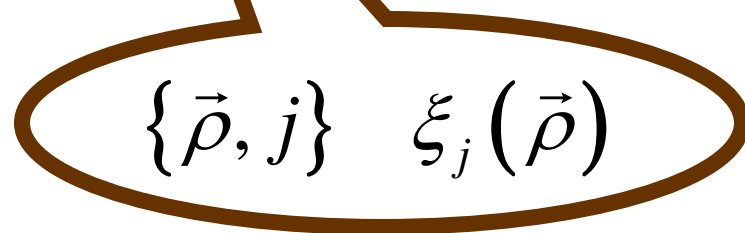
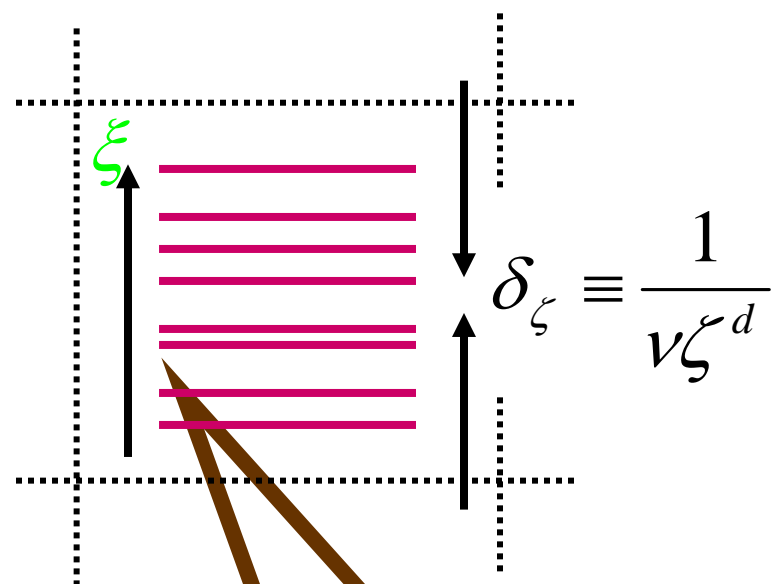
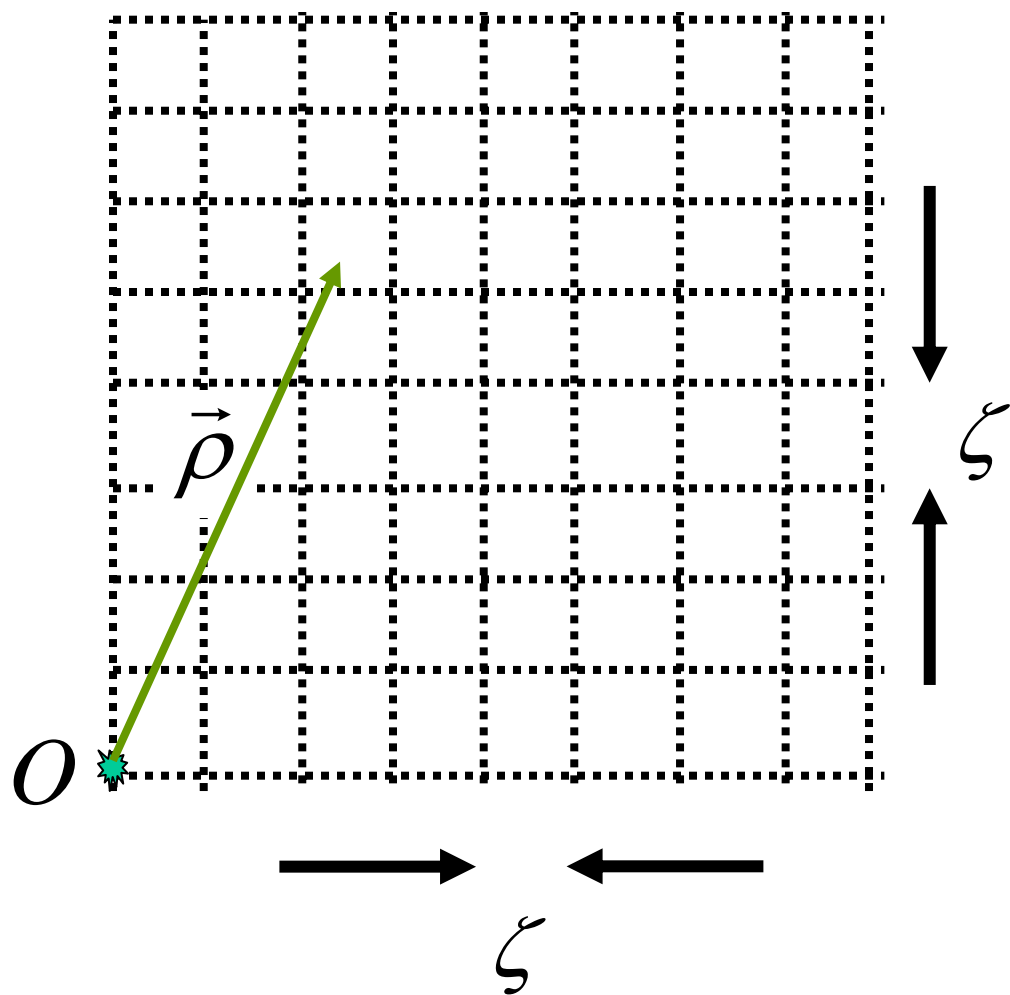
one-electron  
density of states

Need a model with small parameters

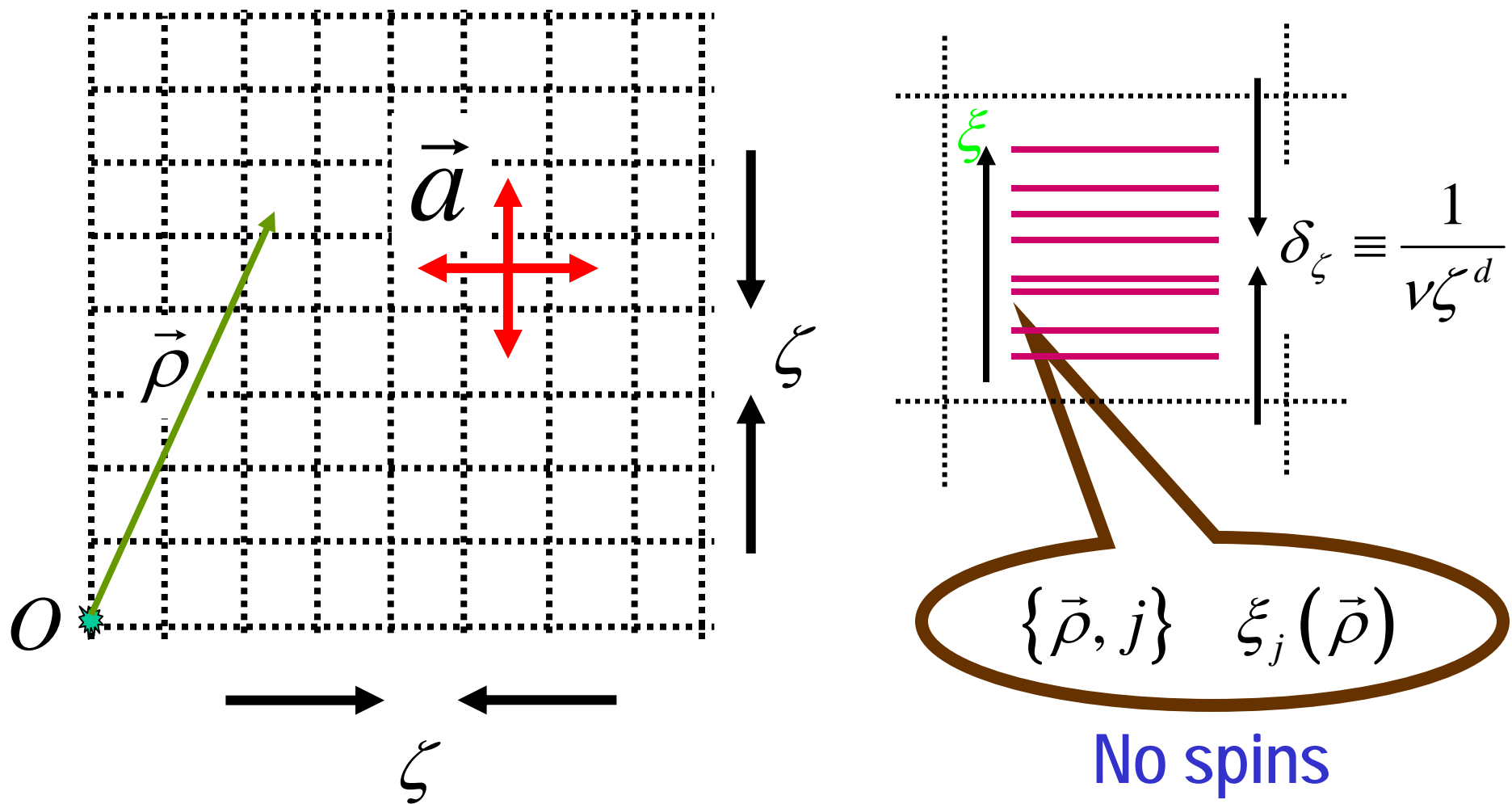
## We have to take into account that

1. A one-electron wave function decays exponentially as a function of the distance from its center.
2. There is level repulsion for the states localized nearby
3. Matrix elements of the interaction decay (probably as a power law) when differences between the energies of involved quasiparticles is increased.
4. These matrix elements have random sign.



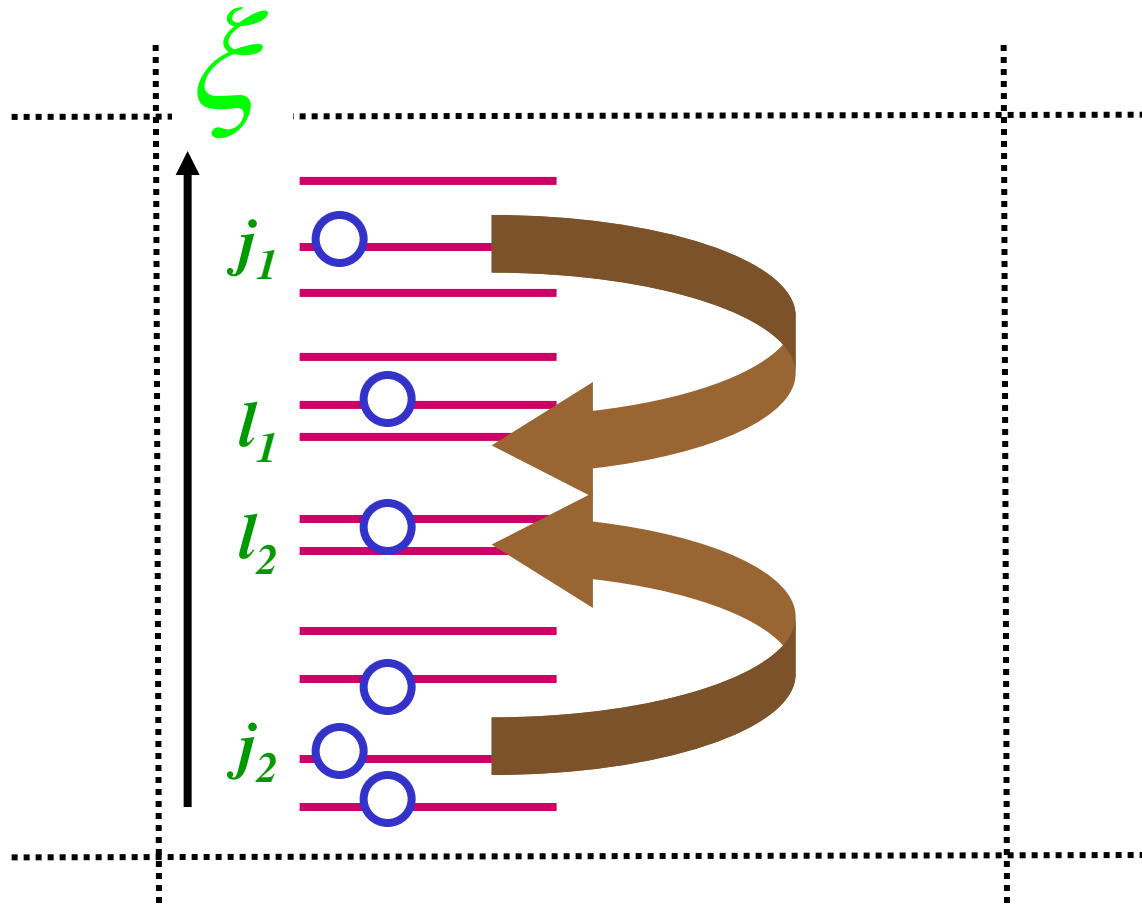


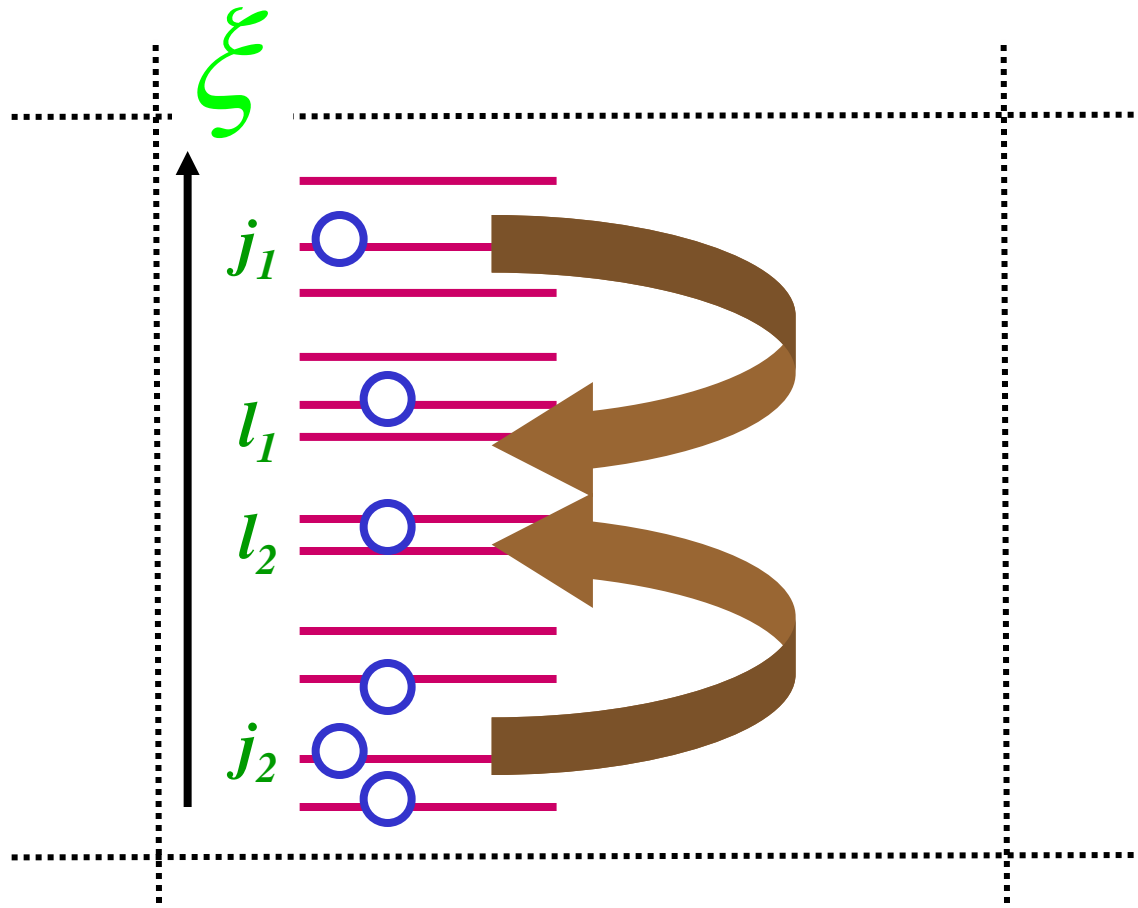
No spins



$$\hat{H}_0 = \sum_{\vec{\rho}, l} \hat{c}_l^\dagger(\vec{\rho}) \left[ \xi_l(\vec{\rho}) \hat{c}_l(\vec{\rho}) + I \delta_\zeta \sum_{\vec{a}, m} \hat{c}_m(\vec{\rho} + \vec{a}) \right]$$



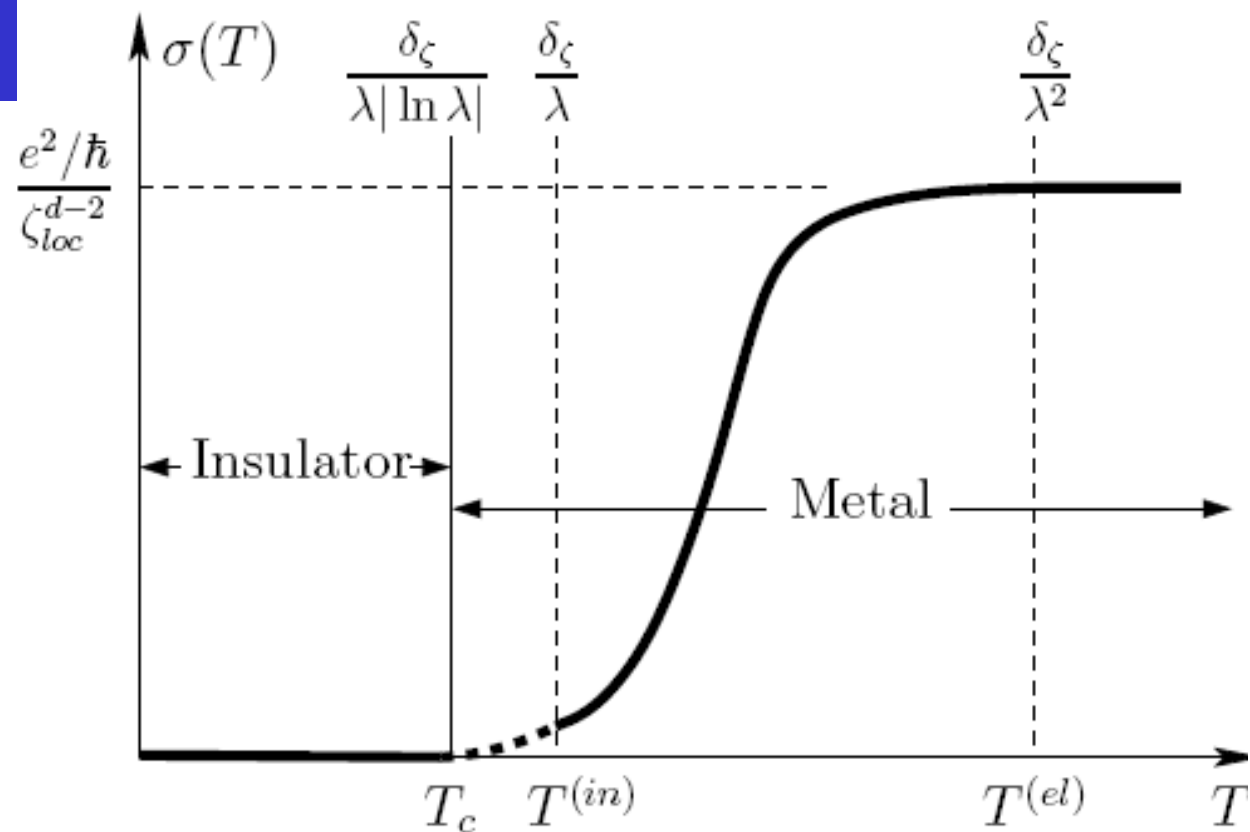




$$\hat{V}_{\text{int}} = \frac{1}{2} \sum_{\vec{\rho}; l_1, l_2; j_1, j_2} V_{l_1, l_2}^{j_1, j_2}(\vec{\rho}) \hat{c}_{l_1}^\dagger(\vec{\rho}) \hat{c}_{l_2}^\dagger(\vec{\rho}) \hat{c}_{j_1}(\vec{\rho}) \hat{c}_{j_2}(\vec{\rho})$$

Interaction only within the same cell;  
no diagonal matrix elements

# Technique



Self-  
 Consistent  
 Born  
 Approximation

Boltzmann  
 Equation

## Absence of Diffusion in Certain Random Lattices

P. W. ANDERSON

*Bell Telephone Laboratories, Murray Hill, New Jersey*

(Received October 10, 1957)

This paper presents a simple model for such processes as spin diffusion or conduction in the "impurity band." These processes involve transport in a lattice which is in some sense random, and in them diffusion is expected to take place via quantum jumps between localized sites. In this simple model the essential randomness is introduced by requiring the energy to vary randomly from site to site. It is shown that at low enough densities no diffusion at all can take place, and the criteria for transport to occur are given.

J. Phys. C: Solid State Phys., Vol. 6, 1973. Printed in Great Britain. © 1973

## A selfconsistent theory of localization

R Abou-Chacra†, P W Anderson‡§ and D J Thouless†

† Department of Mathematical Physics, University of Birmingham, Birmingham, B15 2TT

‡ Cavendish Laboratory, Cambridge, England and Bell Laboratories, Murray Hill, New Jersey, 07974, USA

Received 12 January 1973

# Idea of the calculation:

1. Start with some infinitesimal width  $\eta$  (*Im* part of the self-energy due to a bath) of each one-electron eigenstate
2. Consider *Im* part of the self-energy  $\Gamma$  in the presence of tunneling and *e-e* interaction.
3. Calculate the **probability distribution** function  $P(\Gamma)$
4. Consider the limit: 
$$\lim_{\eta \rightarrow 0, V \rightarrow \infty} P(\Gamma) \equiv P_0(\Gamma)$$

$V$  is the  
volume of  
the system

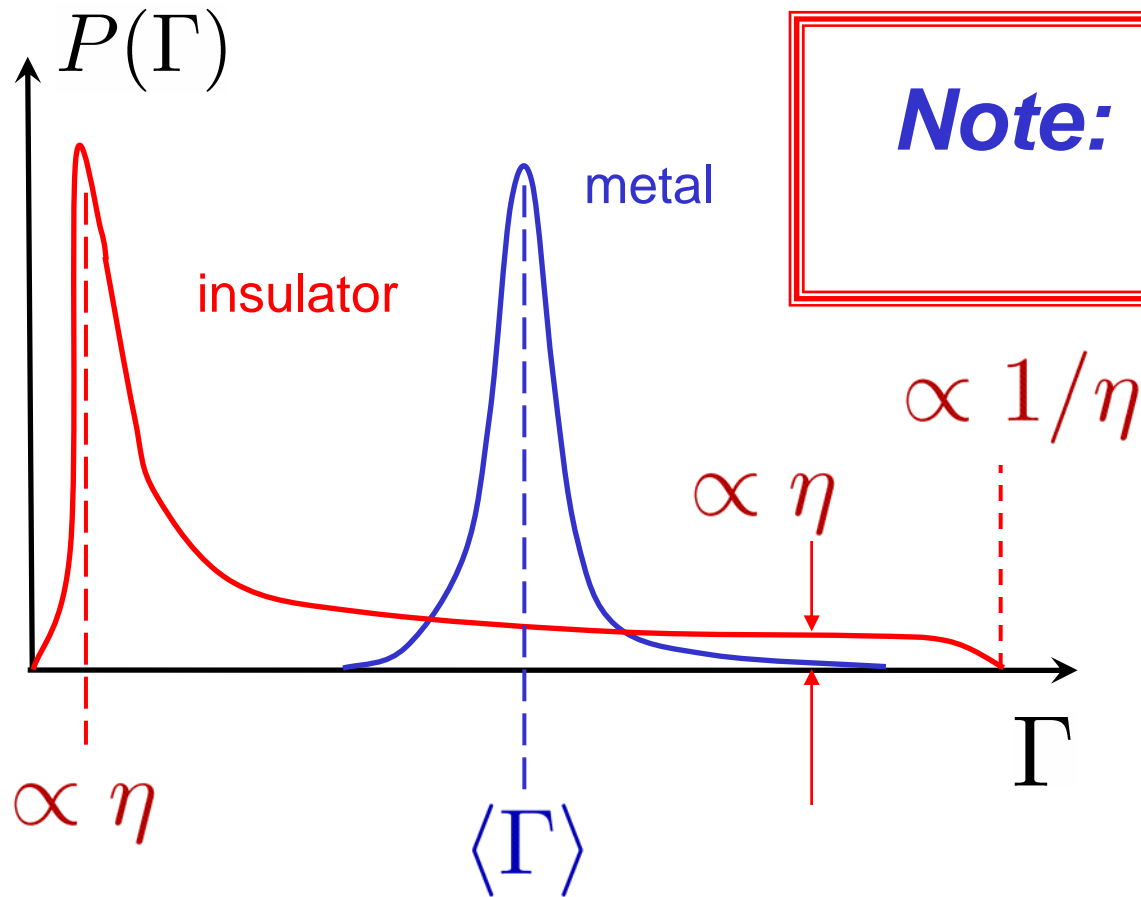
# Idea of the calculation:

1. Start with some infinitesimal width  $\eta$  (*Im* part of the self-energy due to a bath) of each one-electron eigenstate
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3. Calculate the probability distribution function  $P(\Gamma)$
4. Consider the limit:  $\lim_{\eta \rightarrow 0, V \rightarrow \infty} P(\Gamma) \equiv P_0(\Gamma)$

$V$  is the volume of the system

$$P_0(\Gamma) = \delta(\Gamma) \quad \text{- insulator}$$
$$\neq 0 \text{ for } \Gamma \neq 0 \quad \text{- metal}$$

# Probability Distribution



**Note:**  $\langle \Gamma \rangle = \langle \Gamma \rangle$

**Look for:**

$$\lim_{\eta \rightarrow +0} \lim_{\nu \rightarrow \infty} P(\Gamma > 0) = \begin{cases} > 0; & \text{metal} \\ 0; & \text{insulator} \end{cases}$$

# Stability of the insulating phase: **NO** spontaneous generation of broadening

$$\Gamma_\alpha(\varepsilon) = 0$$

is always a solution

$$\varepsilon \rightarrow \varepsilon + i\eta$$

linear stability analysis

$$\frac{\Gamma}{(\varepsilon - \xi_\alpha)^2 + \Gamma^2} \rightarrow \pi\delta(\varepsilon - \xi_\alpha) + \frac{\Gamma}{(\varepsilon - \xi_\alpha)^2}$$

After  $n$  iterations of  
 the equations of the  
**S**elf Consistent  
**B**orn **A**pproximation

$$P_n(\Gamma) \propto \frac{\eta}{\Gamma^{3/2}} \left( \text{const} \frac{\lambda T}{\delta_\zeta} \ln \frac{1}{\lambda} \right)^n$$

**first**  $n \rightarrow \infty$   
**then**  $\eta \rightarrow 0$

(...) < 1 – insulator is stable !



# Stability of the metallic phase: Finite broadening is self-consistent

- $$P(\Gamma) = \frac{1}{\sqrt{2\pi\langle\delta\Gamma^2\rangle}} \exp\left[-\frac{(\Gamma - \langle\Gamma\rangle)^2}{2\langle\delta\Gamma^2\rangle}\right]$$

$\sqrt{\langle\delta\Gamma^2\rangle} \ll \langle\Gamma\rangle$  as long as

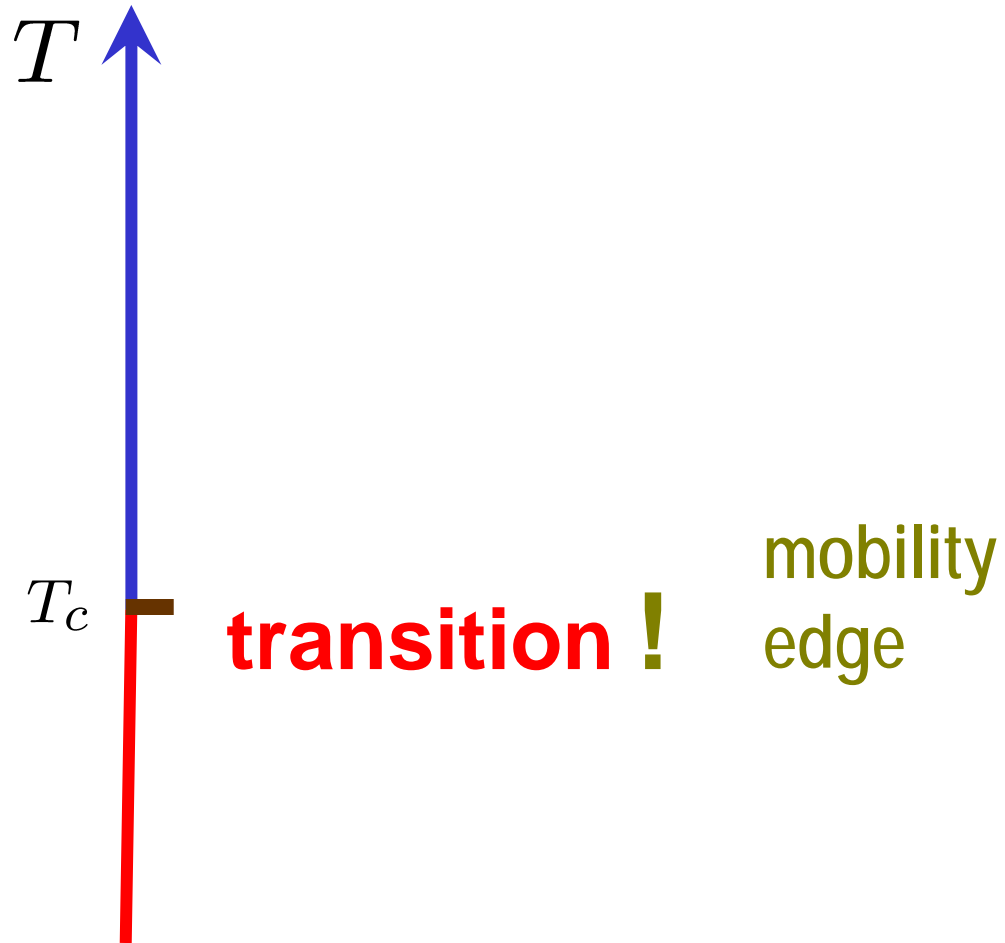
$$T \gg \frac{\delta_\zeta}{\lambda}$$

- $\langle\Gamma\rangle \ll \delta_\zeta$  (levels well resolved)

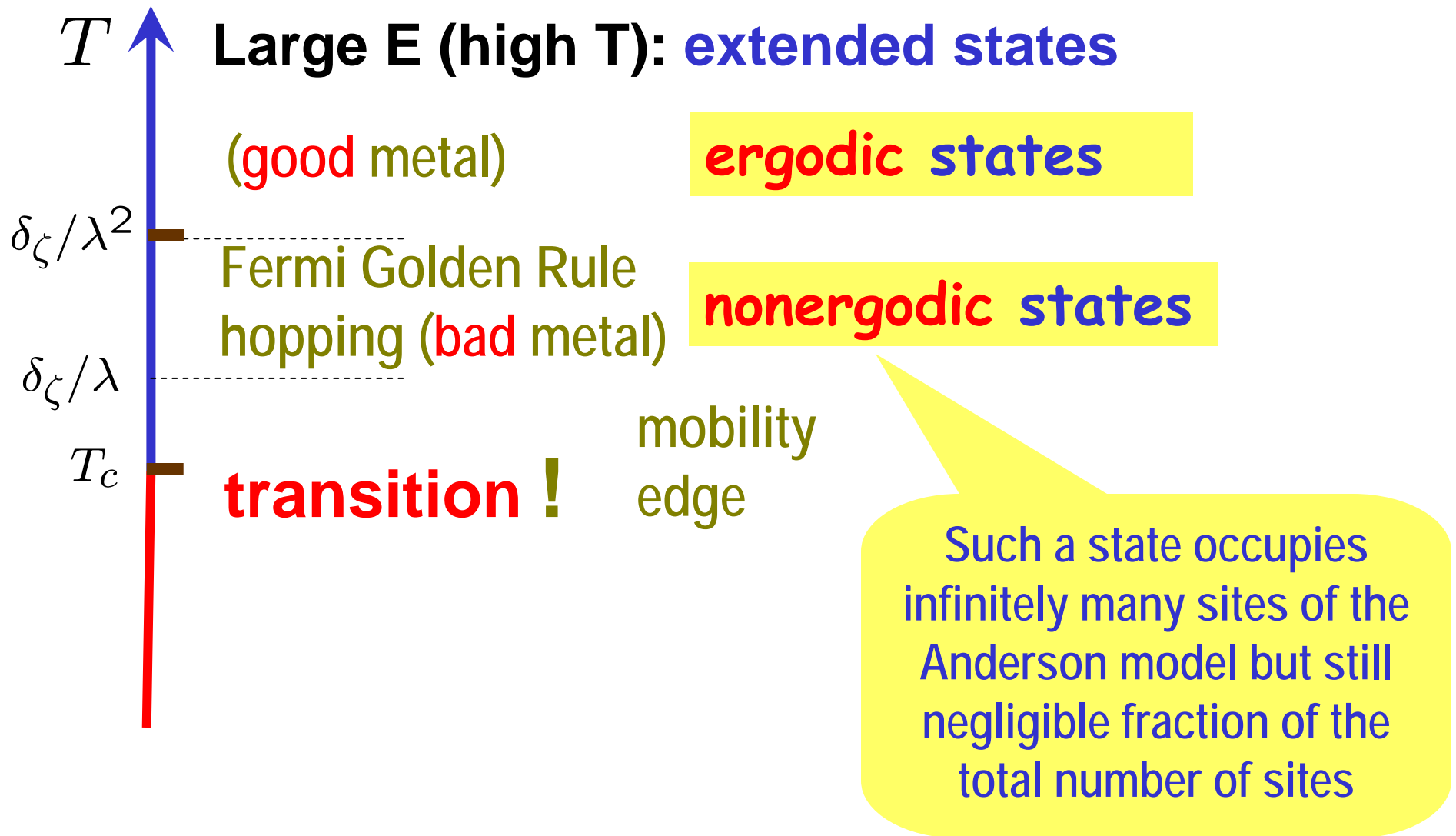
- quantum kinetic equation for transitions between localized states

$$\sigma(T) \propto \lambda^2 T^\alpha \quad (\text{model-dependent})$$

# Many-body mobility edge



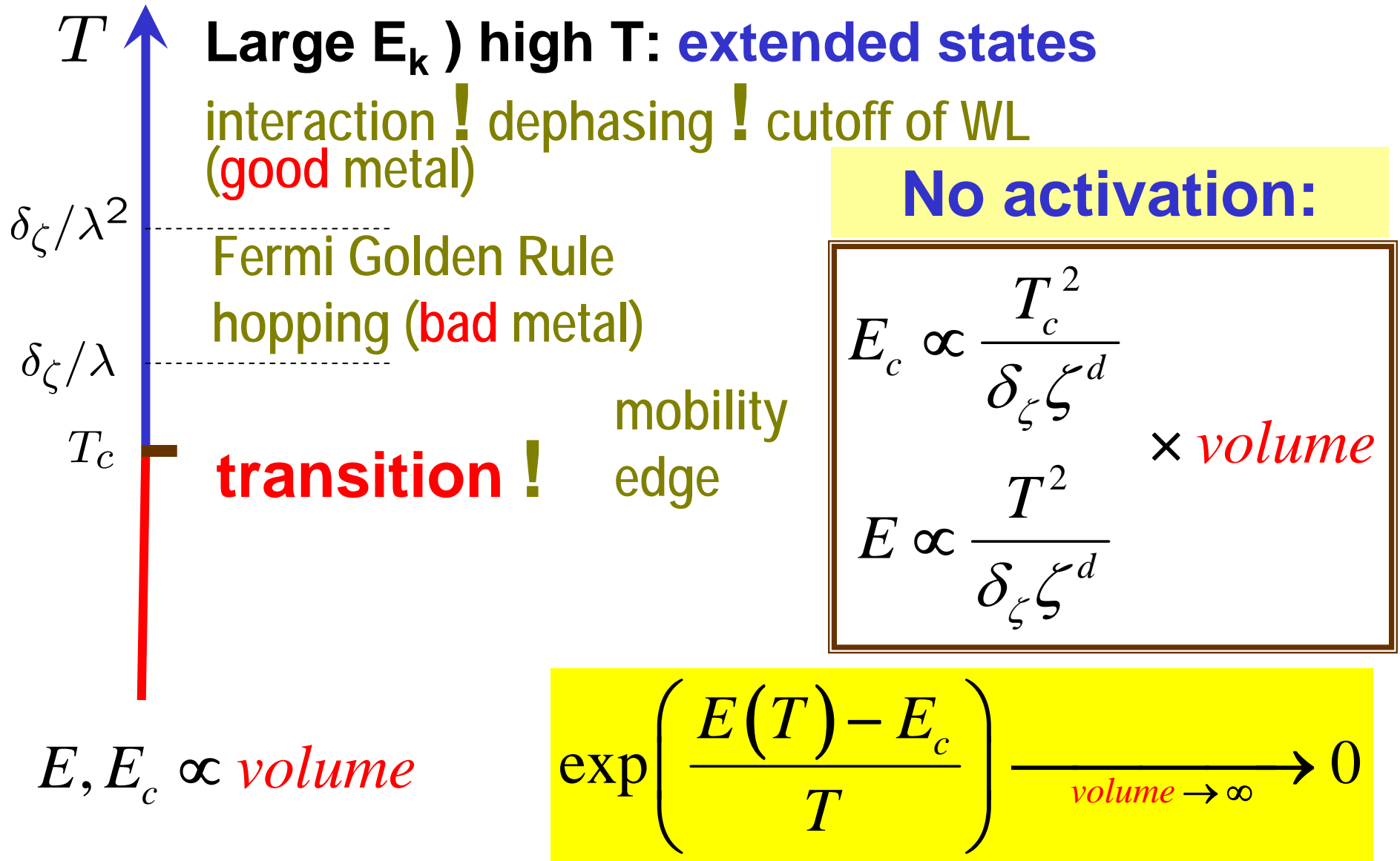
# Many-body mobility edge



# Many-body mobility edge



# Many-body mobility edge



## Conclusions & Some speculations

Conductivity exactly vanishes at finite temperature. **Finite temperature phase transition without any apparent symmetry change!**

Is it an ordinary thermodynamic phase transition or low temperature phase is a glass?

We considered weak interaction.

What about strong electron-electron interactions?

Melting of a pinned Wigner crystal?

What if we now turn on phonons?

Cascades.

Is conventional hopping conductivity picture ever correct?