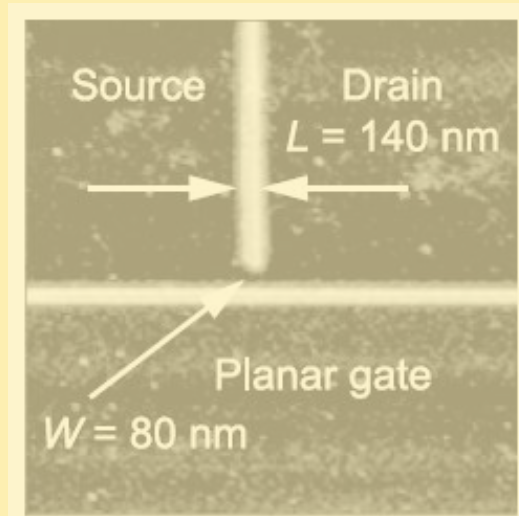
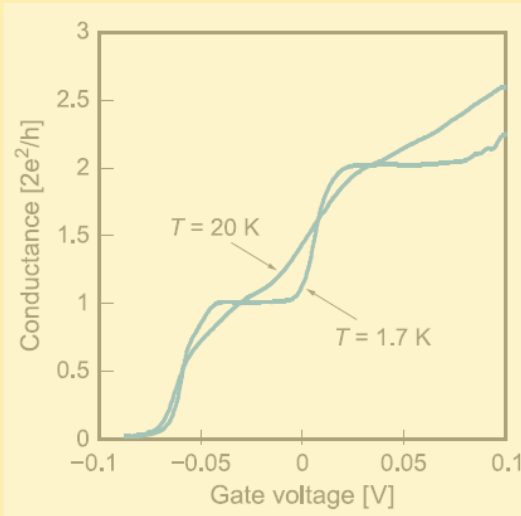


1 Mesoscopic electron transport



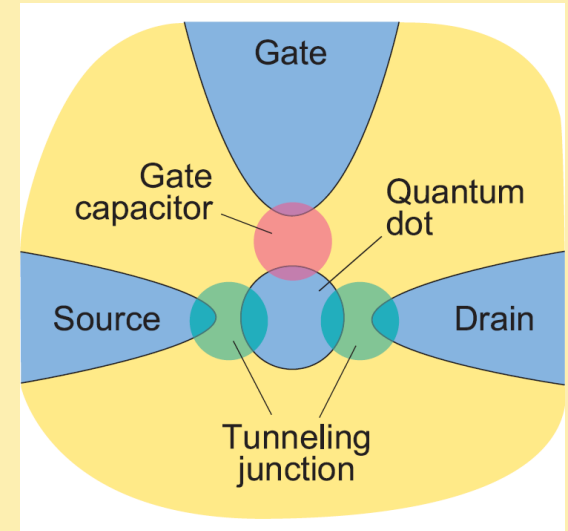
Transport in low dimensional structures

*When is a system low dimensional?
When do we observe size effects?*



Conductance quantization in ballistic 1D structures

Does the conductance diverge when there is no scattering?

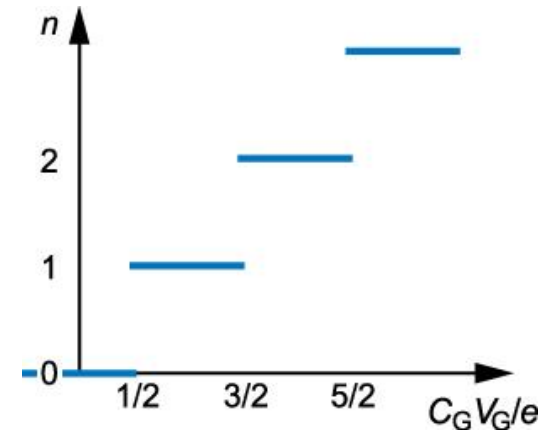
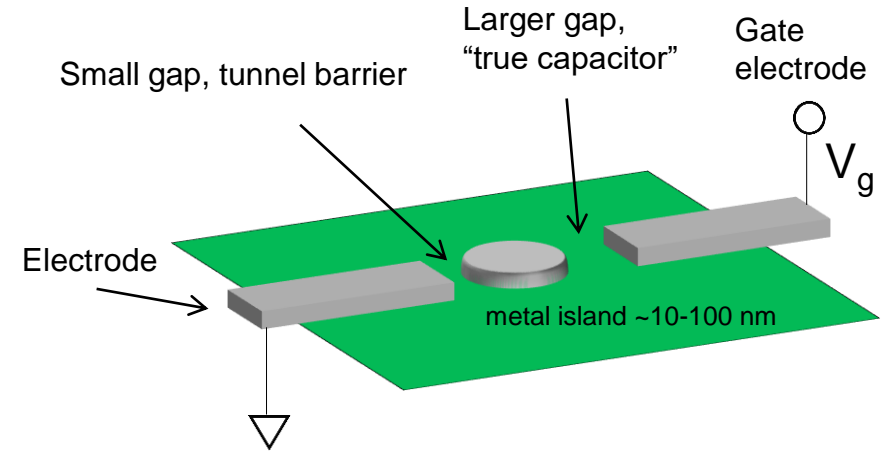
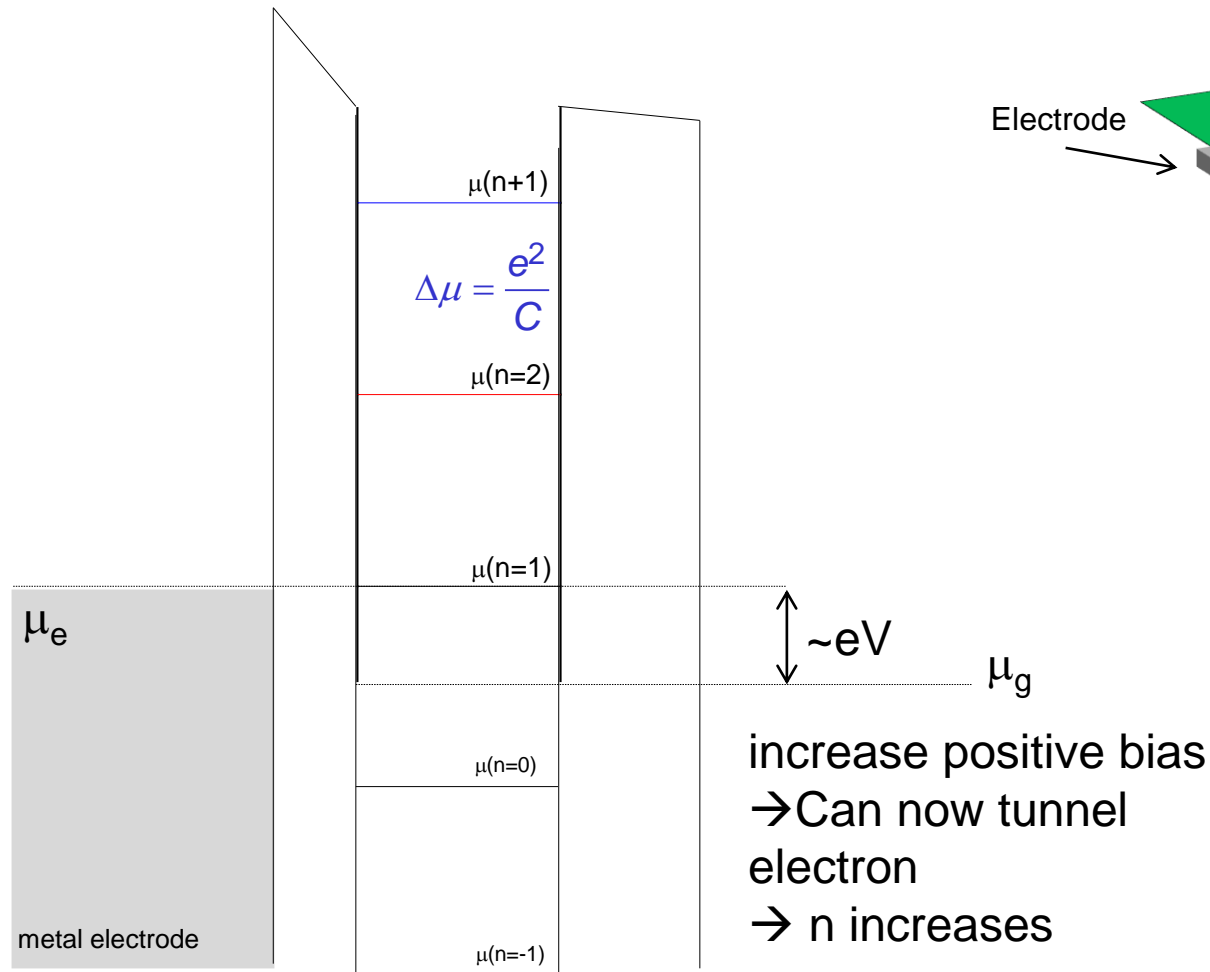


Coulomb blockade in quantum dot structures

Transferring single electrons across a quantum dot

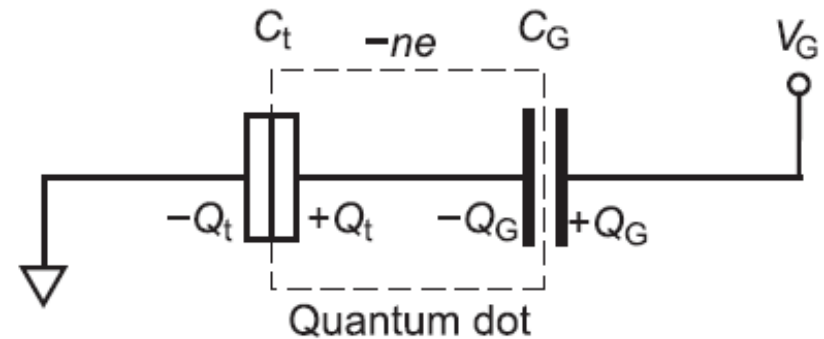
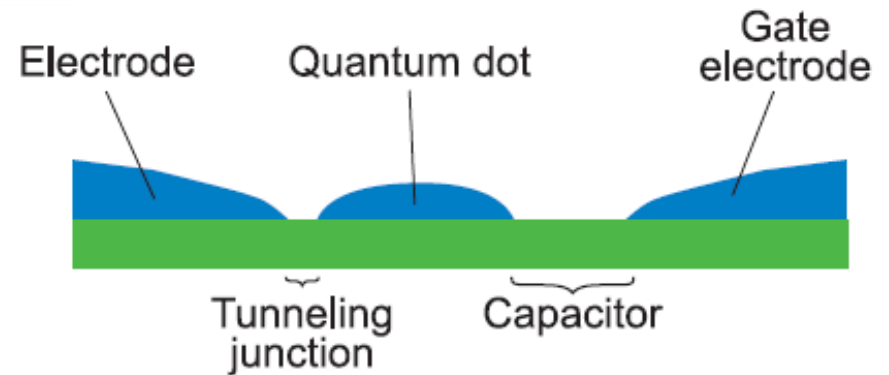
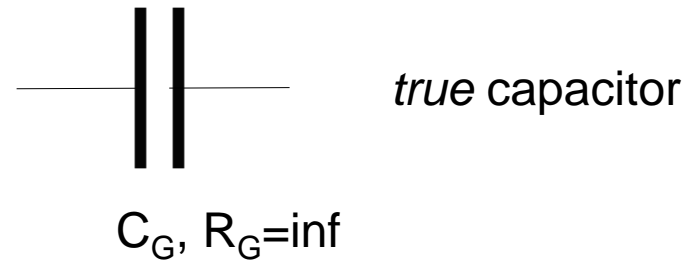
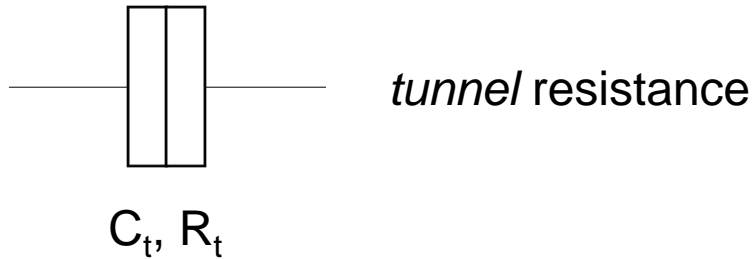
Recap: Single electron box

How can we charge the quantum dot?



Single electron box – quantitative model

Equivalent circuit: QD is *coupled* to the surrounding via



- applying a gate voltage will produce charge Q_G
- system will try to form counter charge
- Charge on QD is *quantized*, $Q_{\text{dot}} = -ne$, while charge on capacitors is continuous
 → A second (continuous) charge is formed at the tunnel resistance, Q_t

Single electron box – quantitative model

What is the free energy of the system when we apply a gate voltage?

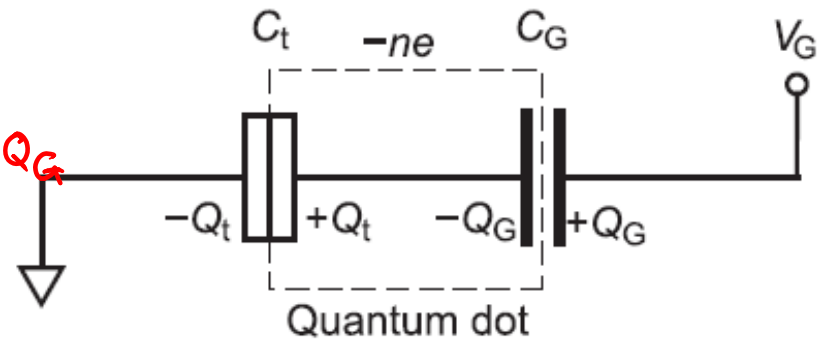
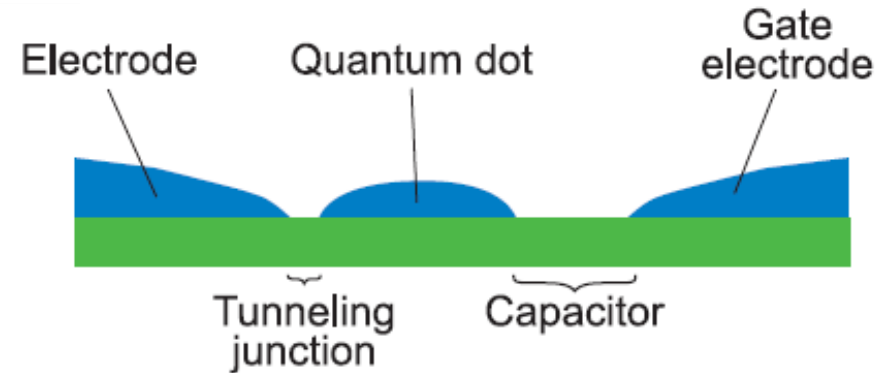
$$F(n) = W_Q(n) - A(n)$$

electrostatic energy stored
in the two capacitors

Work done by external
voltage source to get C_G
charged

$$W_Q(n) = \frac{Q_t^2}{2C_t} + \frac{Q_G^2}{2C_G} \quad \text{use } I = \frac{dQ}{dt}$$

$$A(n) = \int_0^\infty dt I(t) \cdot V_G = V_G \int_0^{Q_G} dQ = V_G \cdot Q_G$$



Single electron box – quantitative model

What is the free energy of the system when we apply a gate voltage?

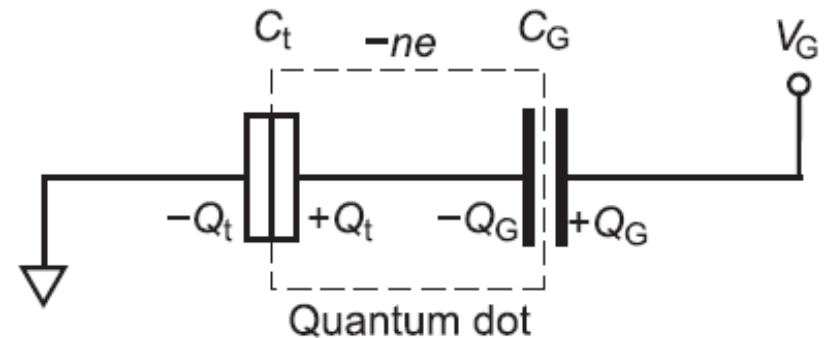
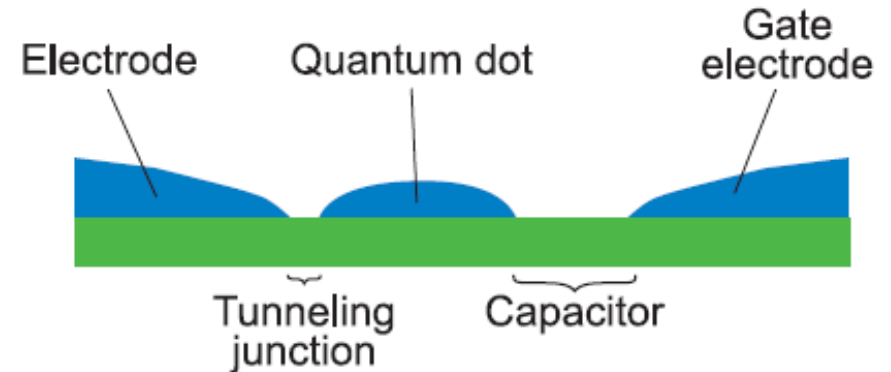
Boundary conditions

Charge balance

$$Q_t - Q_G = -en$$

Voltage drop

$$V_G = \frac{Q_t}{C_t} + \frac{Q_G}{C_G}$$



Single electron box – quantitative model

What is the free energy of the system when we apply a gate voltage?

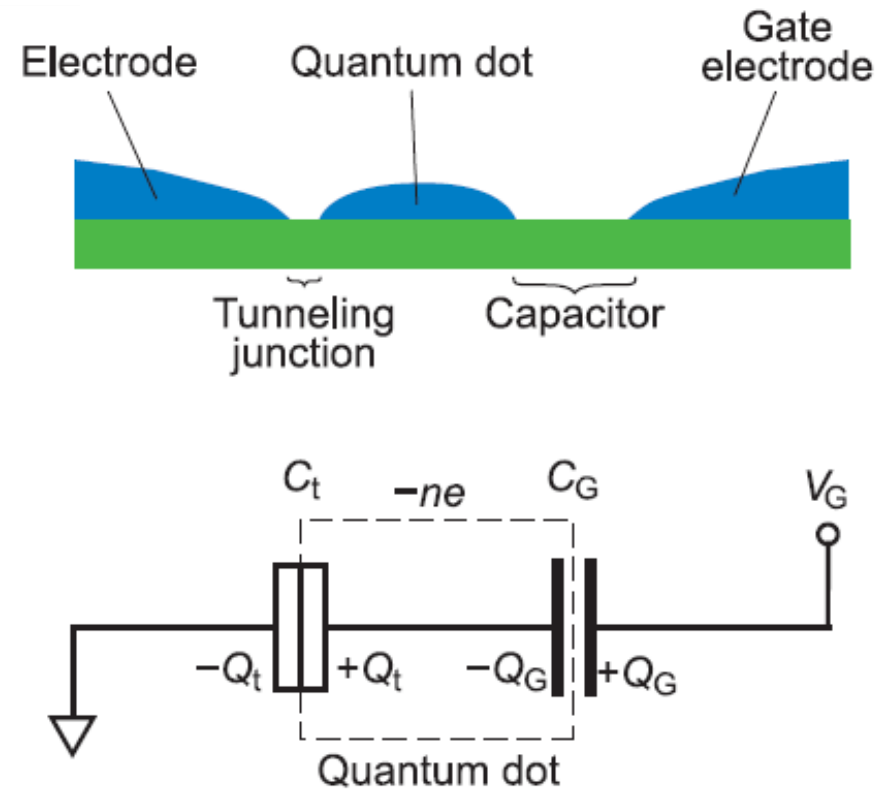
From boundary conditions it follows

$$Q_t = \frac{C_t C_G}{C_\Sigma} \left(V_G - \frac{en}{C_G} \right)$$

$$Q_G = \frac{C_t C_G}{C_\Sigma} \left(V_G + \frac{en}{C_t} \right)$$

with $C_\Sigma = C_t + C_G$

take Q_t and Q_G and insert into $F(n)$



Single electron box – quantitative model

What is the free energy of the system when we apply a gate voltage?

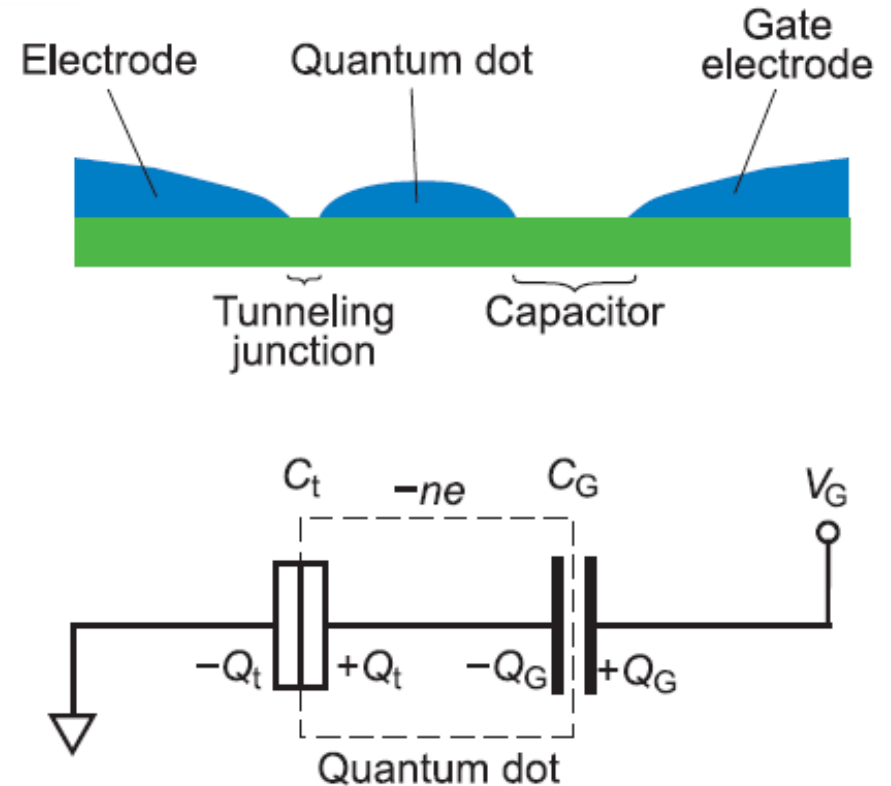
Plugging in Q_t and Q_G , it follows

$$W_Q(n) = \frac{1}{2C_\Sigma} (en)^2 + \frac{C_t C_G V_G^2}{2C_\Sigma}$$

$$= \frac{Q_t^2}{2C_t} + \frac{Q_G^2}{2C_G}$$

$$A(n) = \frac{C_t C_G}{C_\Sigma} V_G^2 + \frac{C_G}{C_\Sigma} V_G (en)$$

$$= V_G \cdot Q_G$$



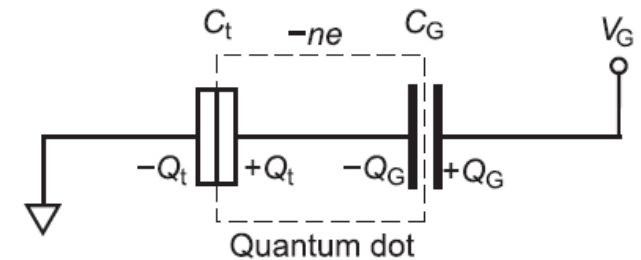
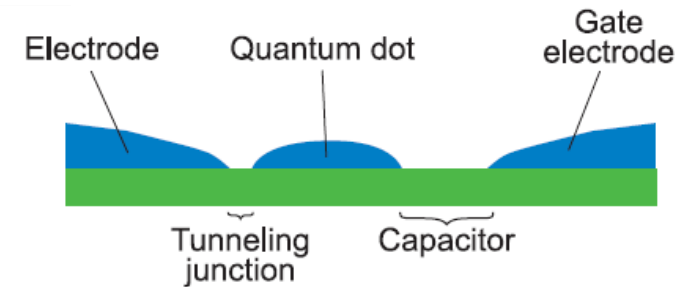
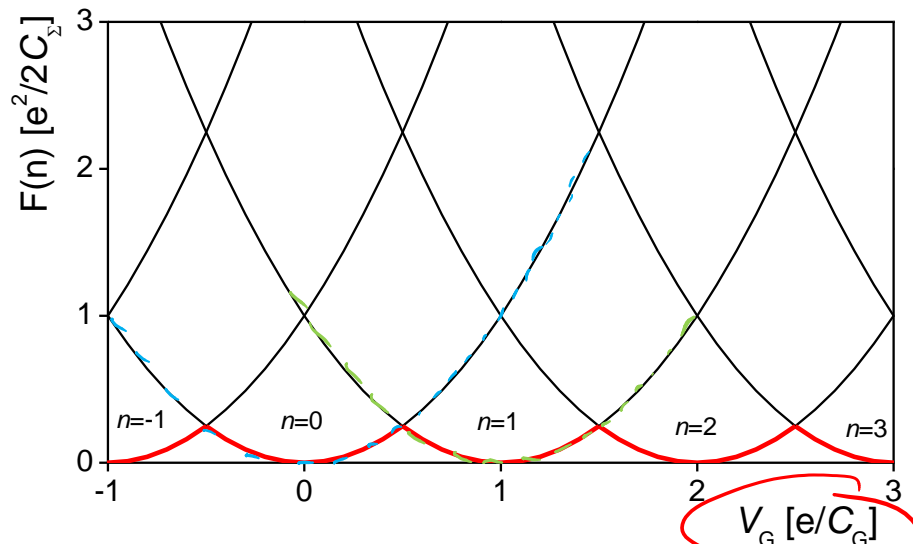
$$\Rightarrow F(n) = W_Q(n) - A(n) = \text{const.} (n - \text{const.})^2 + \text{const.}$$

Single electron box – quantitative model

How many electrons will sit on the QD when we apply a gate voltage?

$$F(n) = \underbrace{\frac{e^2}{2C_\Sigma} \left(n - \frac{C_G V_G}{e} \right)^2}_{\text{n-dependent}} + \underbrace{f(V_G)}_{\text{n-independent}}$$

graphical solution



With increasing bias voltage V_G the free energy of the system is minimum for a different charge number n !

→ Electrons can be transferred into/out of the dot **one by one**!

Weak and strong coupling

Requirements for the observation of *Coulomb blockade phenomena*

n should be a well-defined number, QD has to be reasonable decoupled from surrounding

1. thermal fluctuation of n has to be suppressed $\frac{e^2}{C} \gg k_B T$
2. Quantum fluctuations have to be suppressed: the electrons have to be reasonably localized on the QD,
i.e. spend enough time on the quantum dot before fluctuating back and forth

$$\Delta U \cdot \Delta t = \frac{e^2}{C_t} \cdot R_t \cdot C_t > h$$

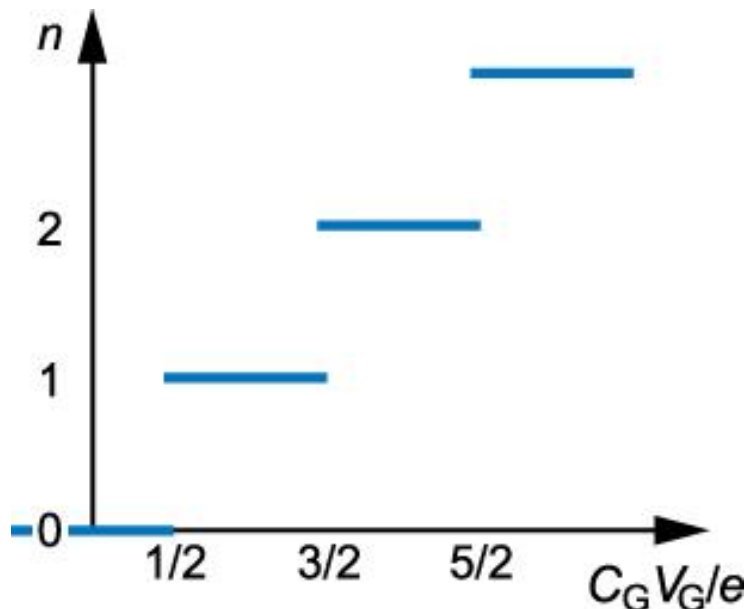
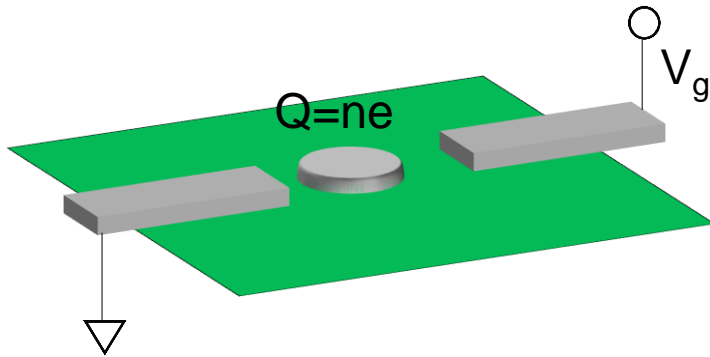
This timescale Δt is given by the $R_t C_\Sigma$ value of the coupling tunnel contact. Heisenberg's uncertainty relation gives

$$R_t > \frac{h}{e^2} \approx 25.8 \text{ k}\Omega$$

$R_t > \frac{h}{e^2}$ is called **weak coupling**

$R_t < \frac{h}{e^2}$ is called **strong coupling**

Coulomb blockade



The effect of an electron not being able to tunnel into a quantum dot due to an energy barrier is called **Coulomb blockade**

The Coulomb blockade occurs due to

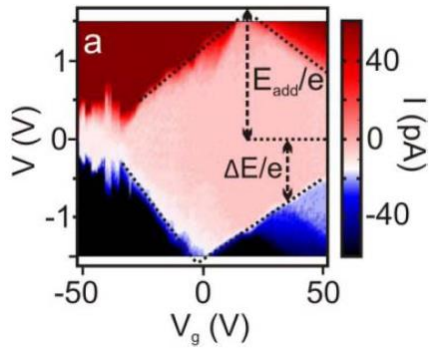
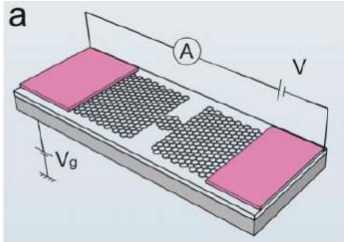
- low capacitance of quantum dots (metals) *# more 2 points*
- size quantization (semiconductors) *# additional effect*

Using the Coulomb blockade effect single electron charges can be transferred and controlled on a nanoscale quantum dot

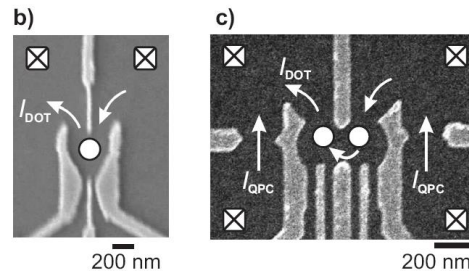
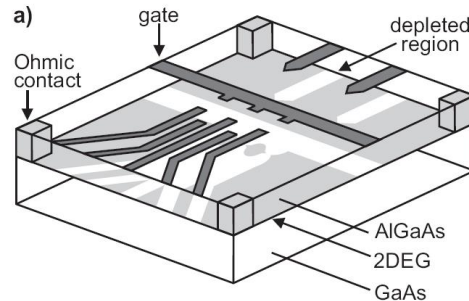
Requirements

$$R_t > \frac{h}{e^2} \quad \frac{e^2}{C} \gg k_B T$$

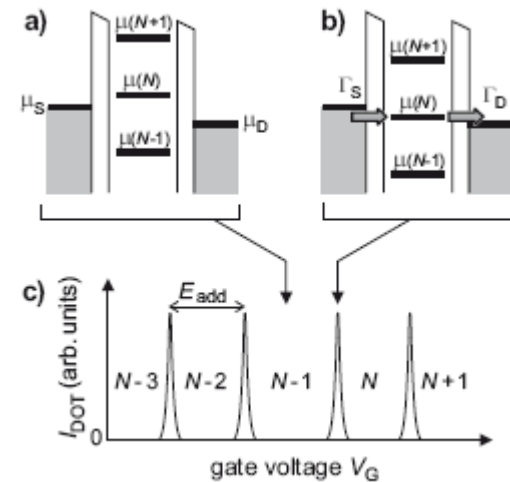
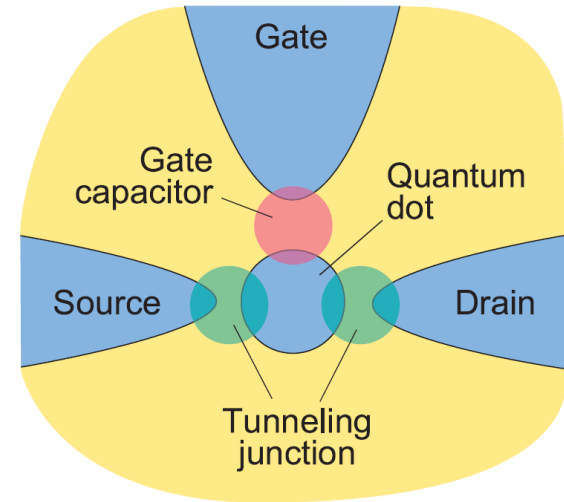
Coulomb blockade - examples



A. Barreiro, H.S.J. van der Zant, L.M.K. Vandersypen, "Quantum Dots at Room Temperature carved out from Few-Layer Graphene", arXiv:1211.4551 (21 Nov. 2012)



Hanson et al., Rev. Mod. Phys. 79, 1217 (2007)



Hanson et al., Rev. Mod. Phys. 79, 1217 (2007)

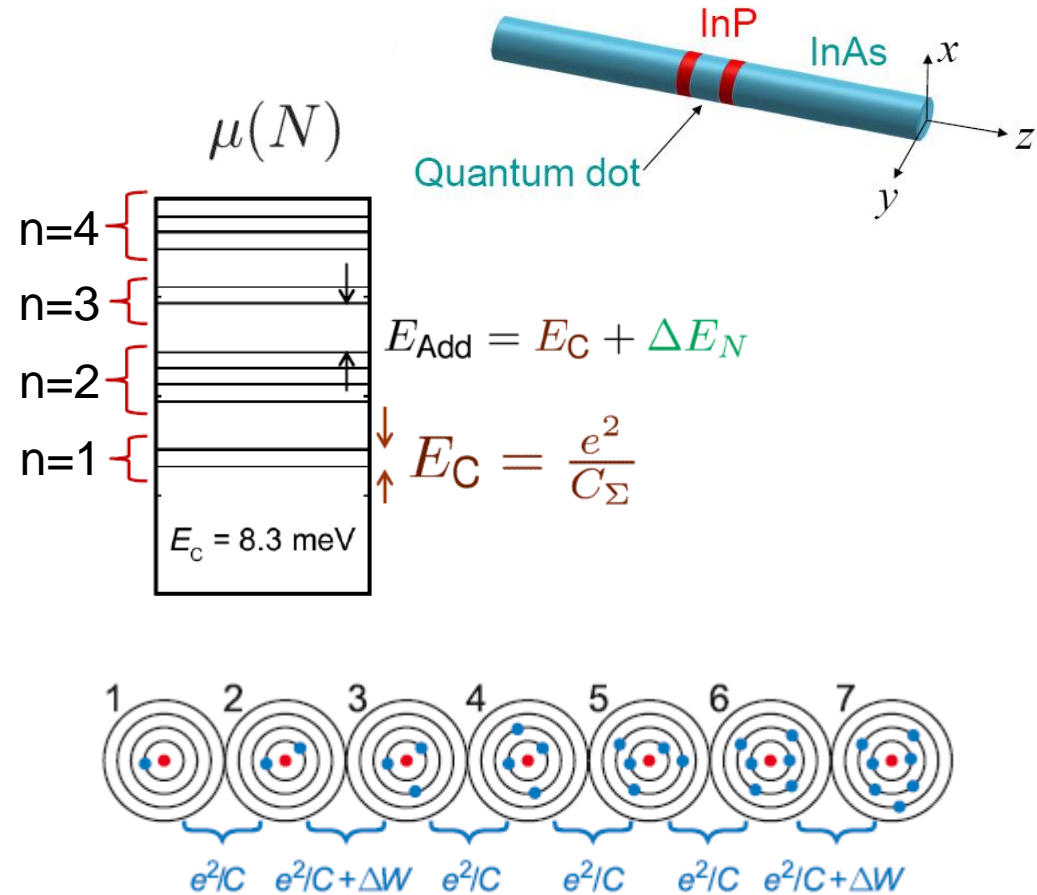
Artificial atoms: Semiconductor quantum dots

Charging energy

$$\mu(n) = W_C n + W_{k_n}$$

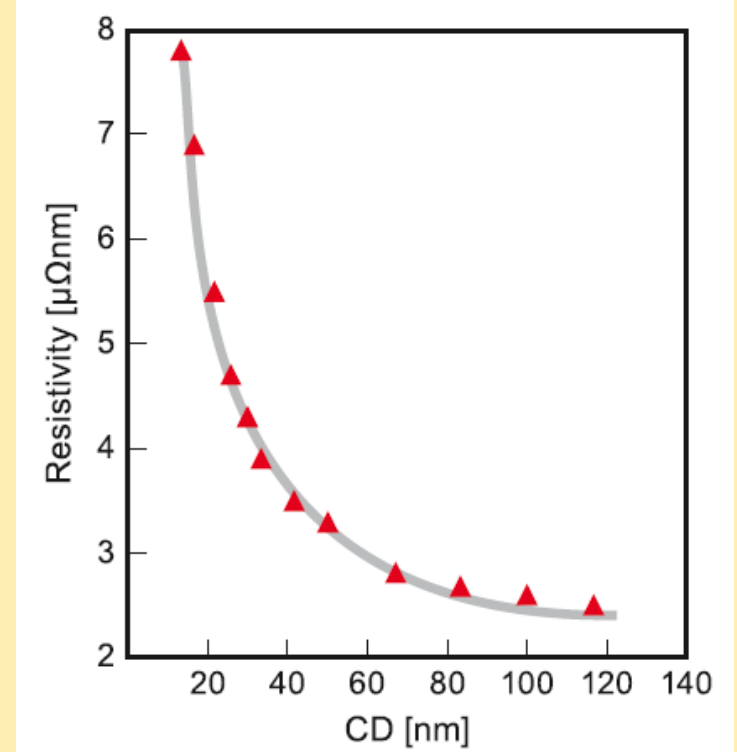
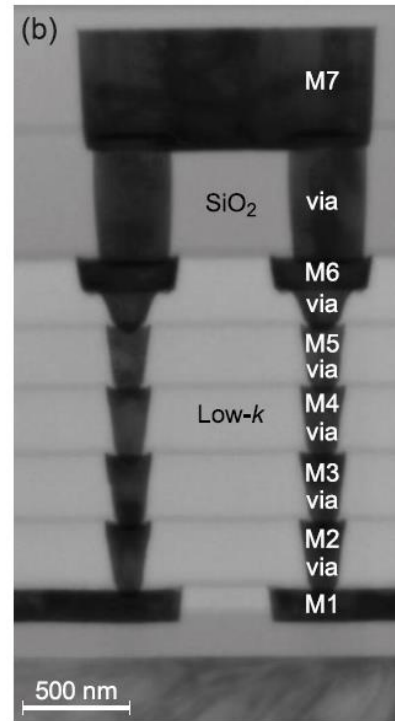
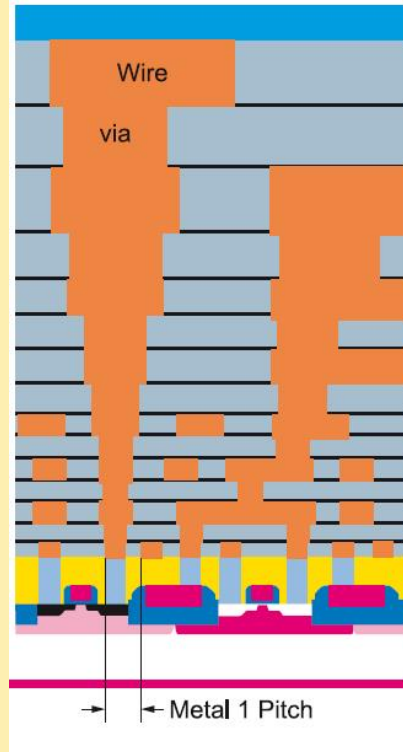
Size quantization energy

$$W_{k_n} = \frac{h^2}{2m^* L^2} (n_x^2 + n_y^2 + n_z^2)$$



Band-like clustering of states in semiconductor QD

2 Metal interconnects



Interconnects – basic principles

Delay time

Power consumption

Size-dependent resistivity of metal wires
Furcs-Sondheimer formalism

Resistivity of metal point contacts
Sharvin-Conductance

Metal interconnects - basics

Interconnects *interconnect* transistors and functional entities on chips

- provide power
- provide ground
- usually a stacked-layer wiring system/network

global

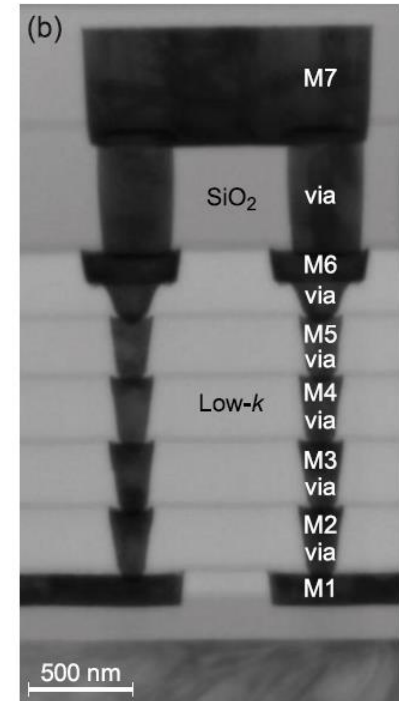
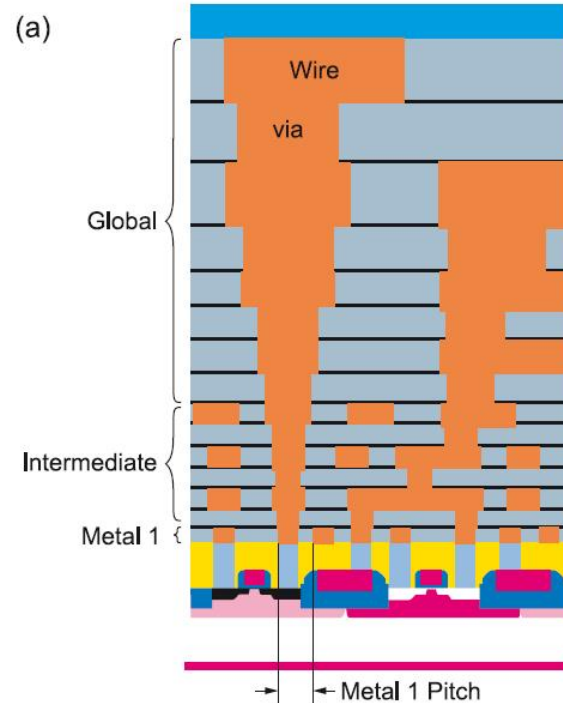
- ground, power supply, long distance communications
- 500 nm – 20 μm

intermediate

- interconnections between cells
- 100 nm – 500 nm

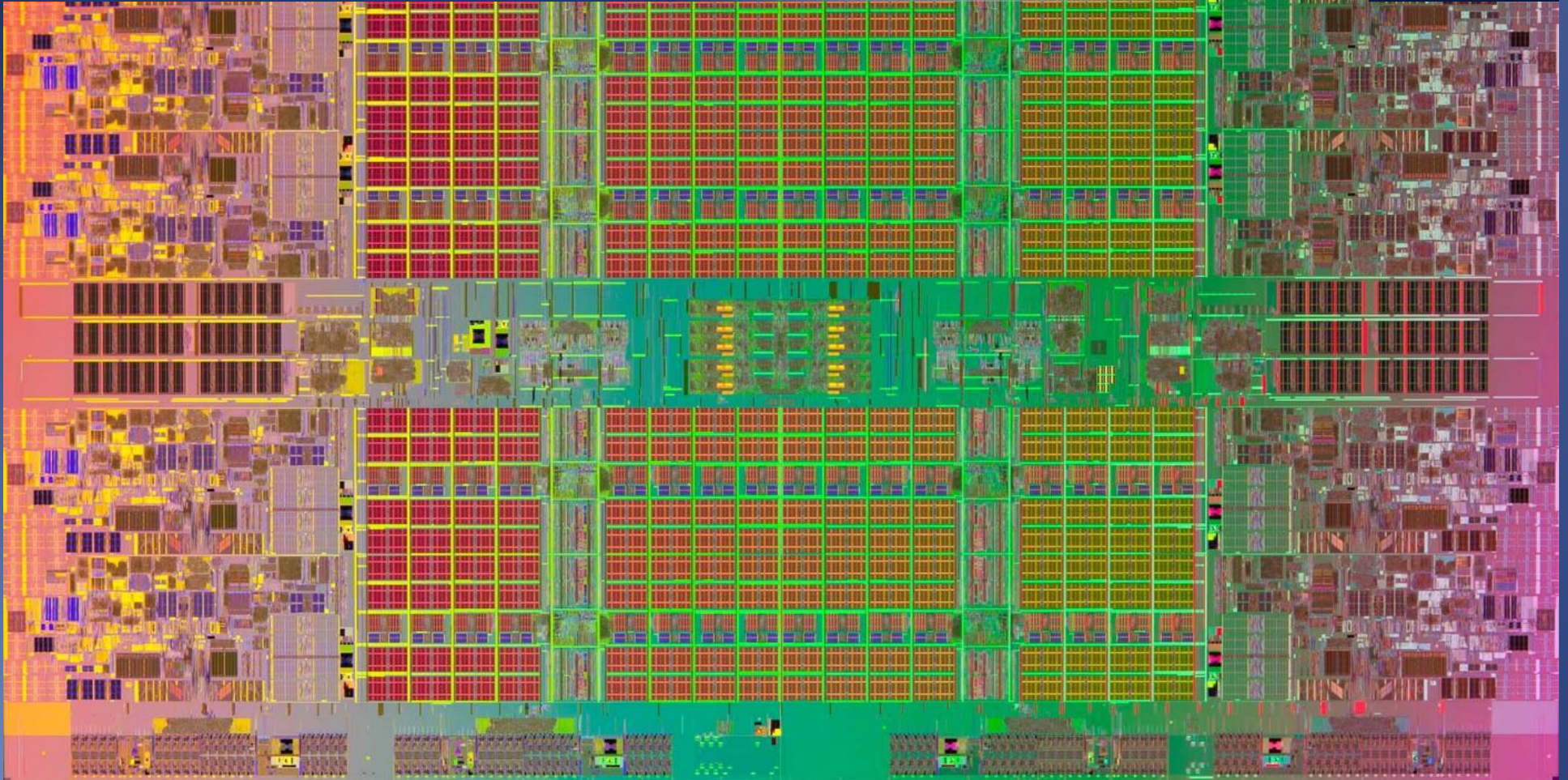
local

- Connecting single transistors
- < 100 nm

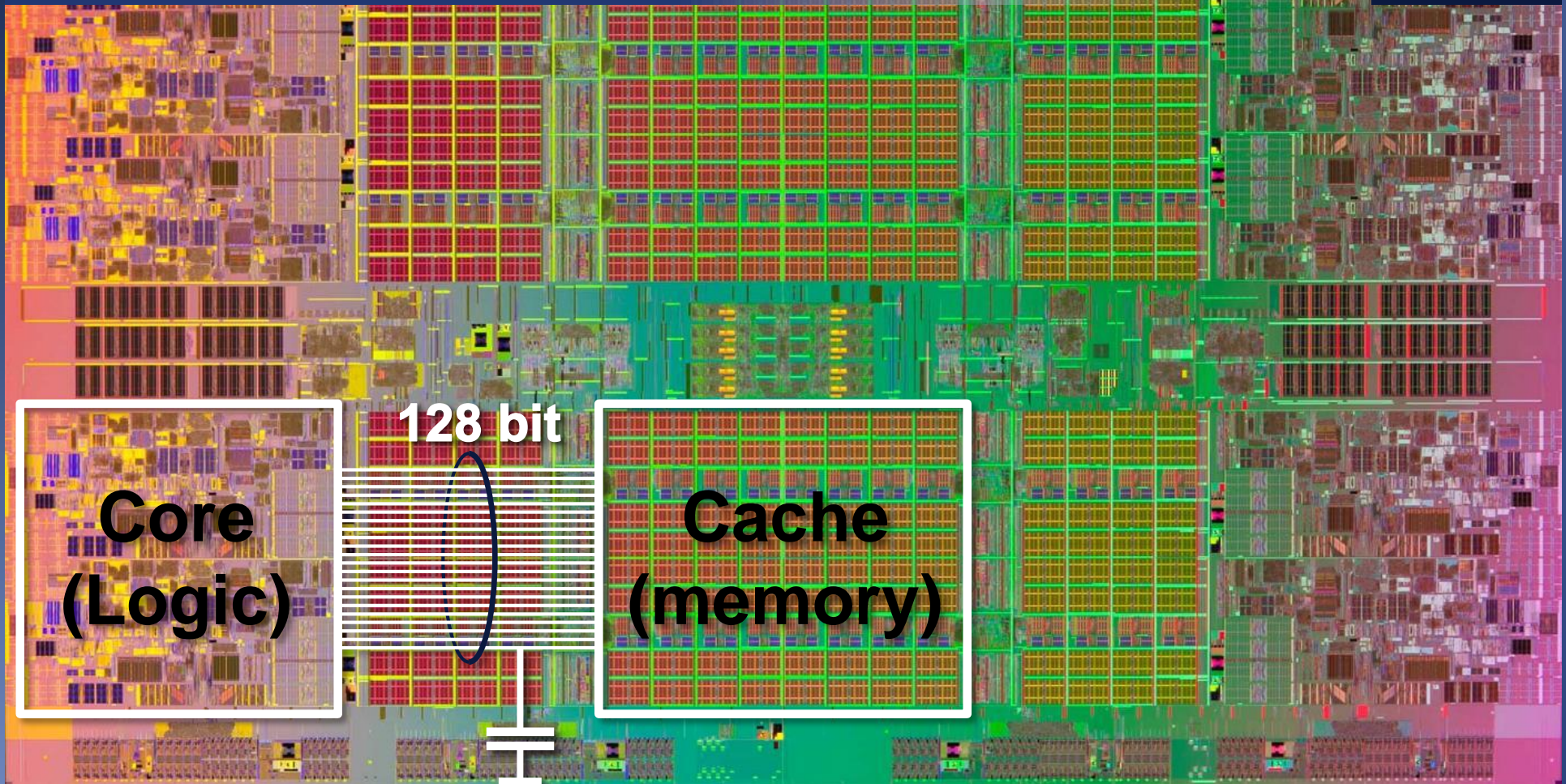


interconnects can cause >50% power consumption of a system
→ resistance of *thin* metal wires
→ capacitive coupling to surrounding

Data transmission in a modern processor



Data transmission in a modern processor



2.1 Delay time and power consumption

Delay time and power consumption

Delay time of signal propagation through wire system

$$\tau = R \cdot C$$

total capacitance of the system

$$C = 2(C_p + C_v) = 2\epsilon\epsilon_0 Ltw \left(\frac{1}{w^2} + \frac{1}{t^2} \right)$$

resistance of central wire

$$R = \rho \cdot \frac{L}{wt}$$

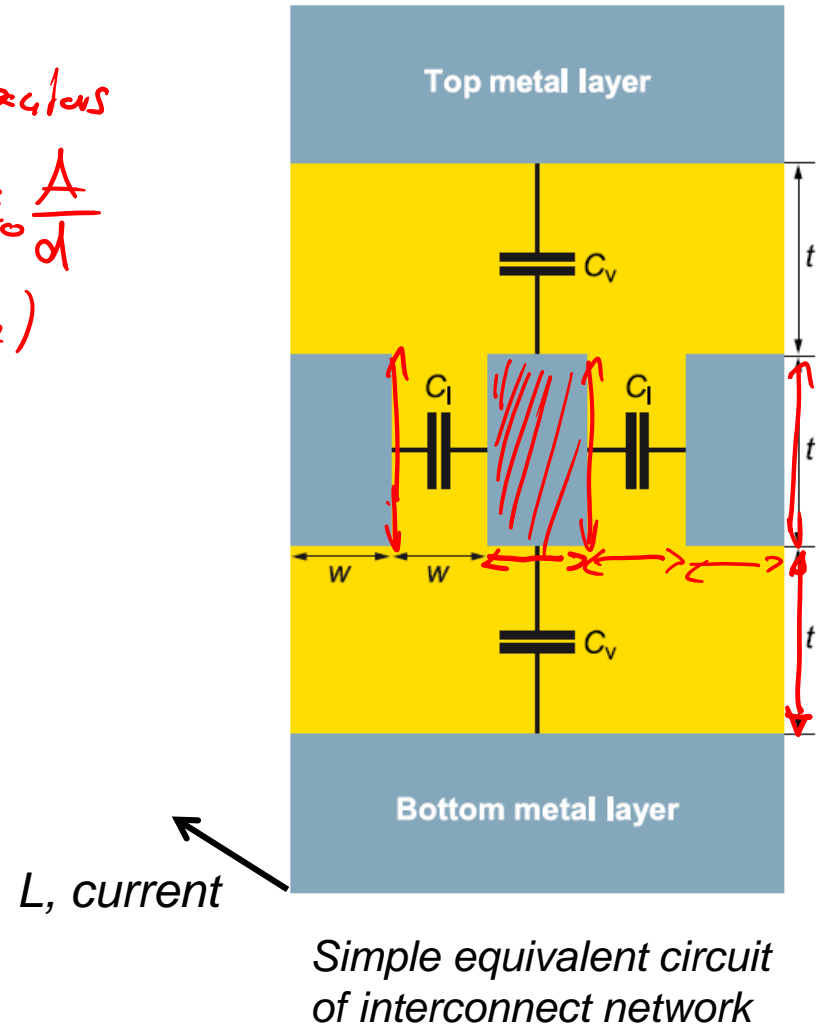
$$\tau_{delay} = 2\rho\epsilon_0\epsilon L^2 \left(\frac{1}{w^2} + \frac{1}{t^2} \right)$$

dynamic power dissipation

$$P = \frac{1}{2} C \cdot V^2 \cdot f$$

plate capacitors

$$C = \epsilon\epsilon_0 \frac{A}{d}$$



Strategies to reduce delay time and/or power consumption

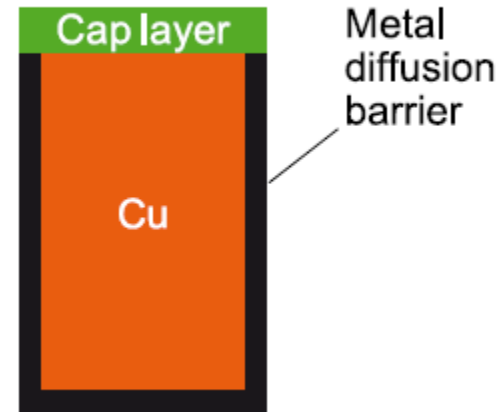
Reducing ρ

from W ($5.6 \mu\Omega\text{cm}$)

to Al alloys ($3.0 \mu\Omega\text{cm}$)

to Cu ($1.67 \mu\Omega\text{cm}$)

challenge: electromigration of Cu



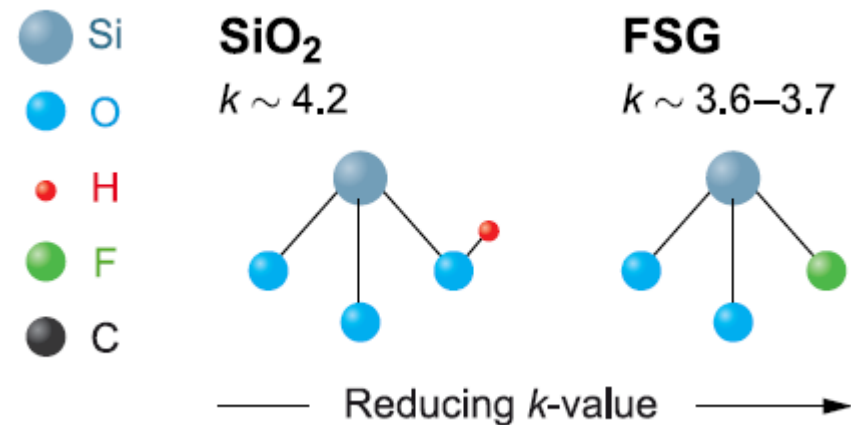
Reducing ϵ (or k)

chemical manipulation of SiO_2

→ reduces dipole moment

e.g. Si-O-bond replaced by Si-F-bond
(fluorosilicate glasses, FSG)

challenge: too low ϵ increases leakage



$$k \equiv \epsilon$$

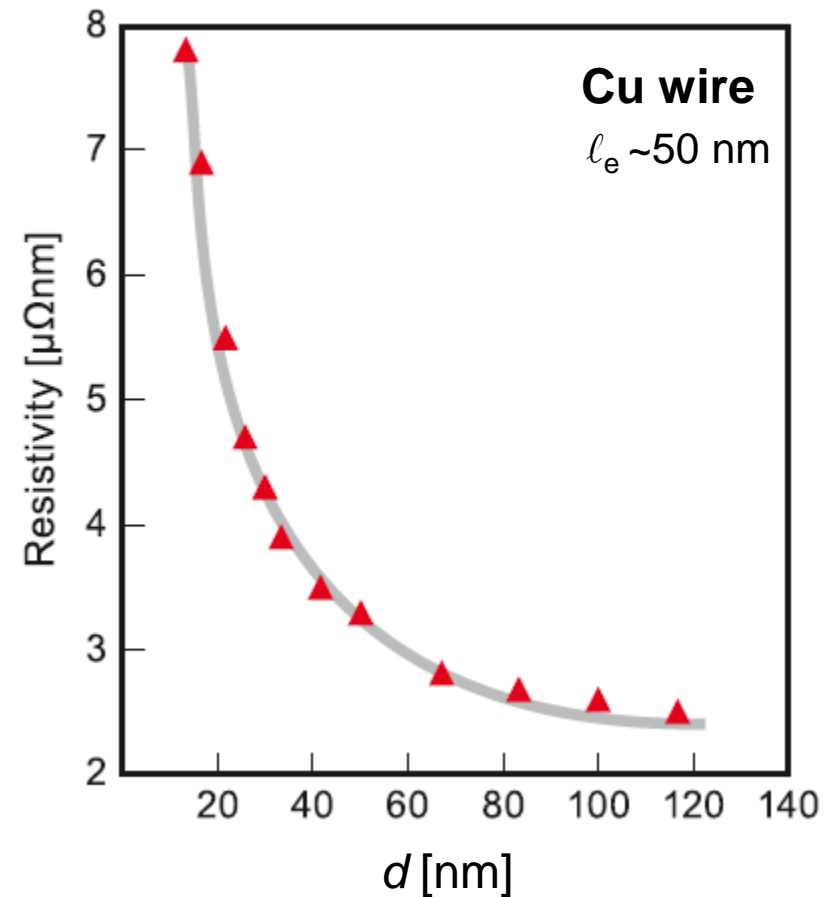
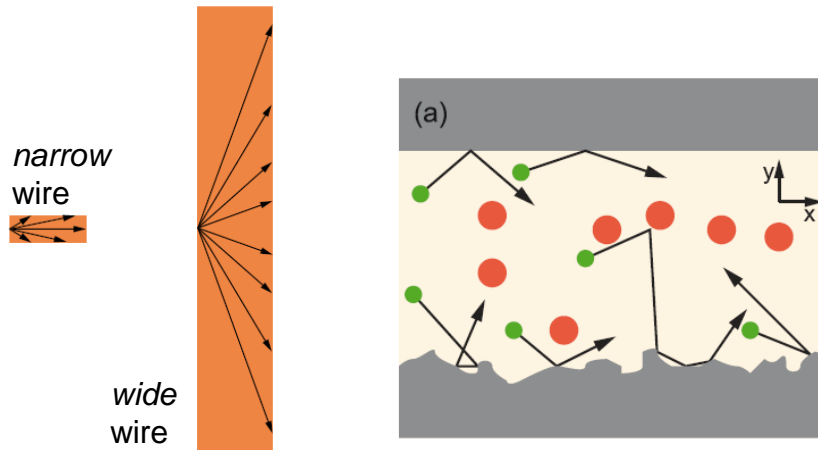
2.2 Size-dependent resistivity of metal wires

Size-dependent resistivity of metal wires

bulk resistivity: $\rho_{bulk} \neq \rho_{bulk}(L, w, t)$

→ independent of geometry

However, **experiments** show that ρ increases with decreasing feature size



Reminder. Boundary scattering!

For metals ($\lambda_F \sim 1$ nm), the boundary of the wire appear *rough* and limits ℓ_e

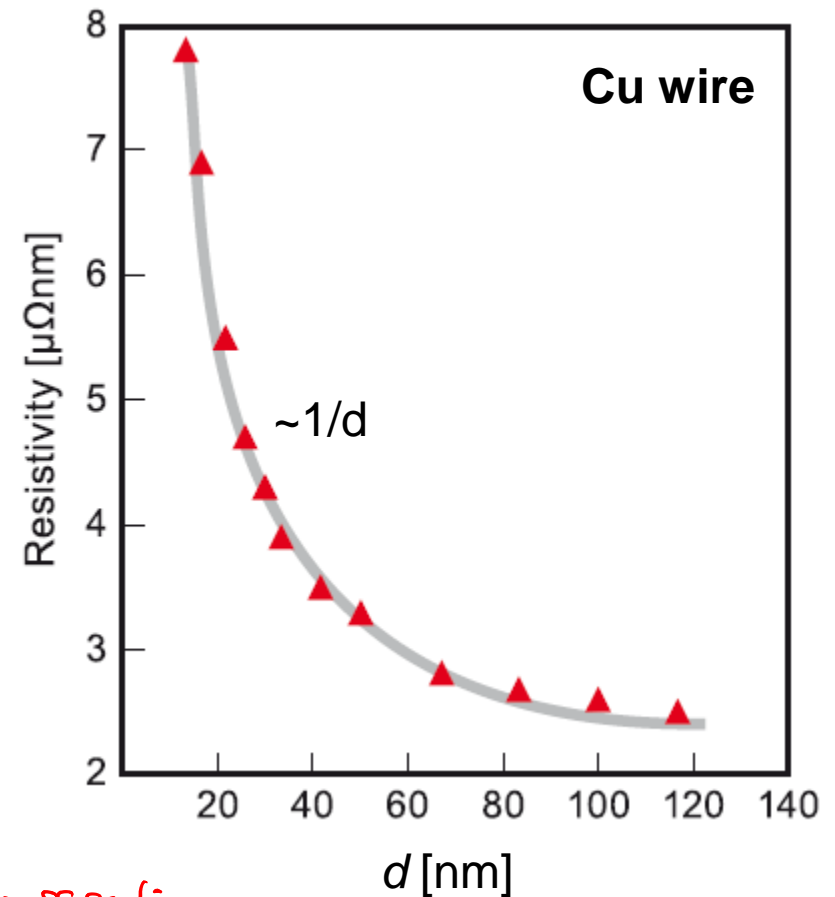
Fuchs-Sondheimer formalism

First quantitative model by Fuchs and Sondheimer considering **surface scattering**

$$\rho_{FS} = \frac{\rho_{bulk}}{1 - \frac{3}{2k}(1-p) \int_1^{\infty} \left(\frac{1}{t^3} - \frac{1}{t^5} \right) \frac{1-e^{-kt}}{1-pe^{-kt}} dt}$$

p = surface reflectivity

$k = \frac{d}{\ell_e}$ ℓ_e means "bulk" mean free path here



for $d \gg \ell_e$

$$\frac{3}{4} \frac{\ell_e}{d}$$

rectangular cross section

$$\rho_{FS} = \rho_{bulk} \left[1 + \frac{3\ell_e}{8d} (1-p) \right]$$

Fuchs-Sondheimer formula

→ circular cross section of the wire