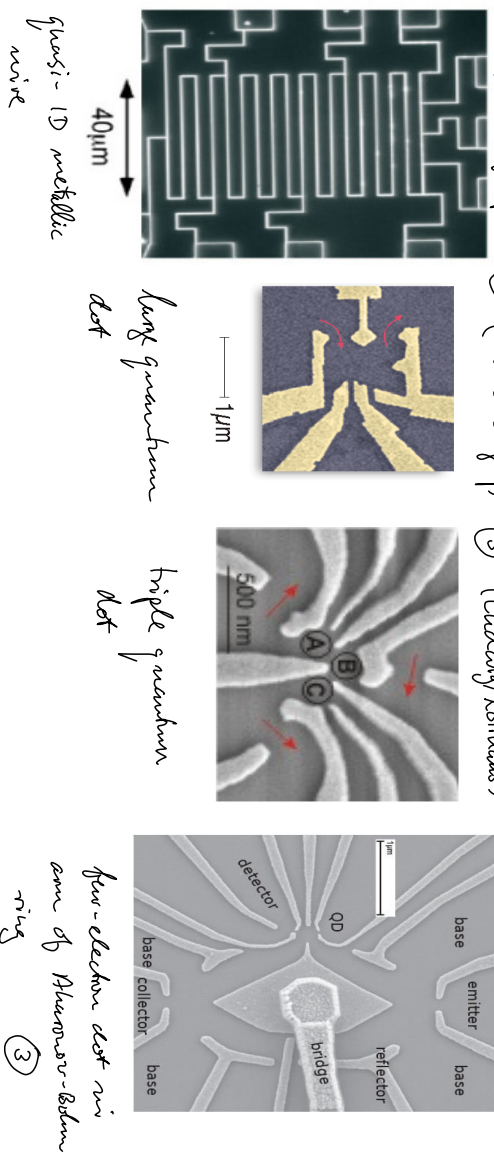


Textbook: Nazarov & Blanter (NB), "Quantum Transport",
Cambridge University Press, 2009.

- ① (Bisseg group) ② (Marcus group) ③ (Ludwig/Kottmanns) ④ (Weidmann group)



"Map of Quantum Transport"

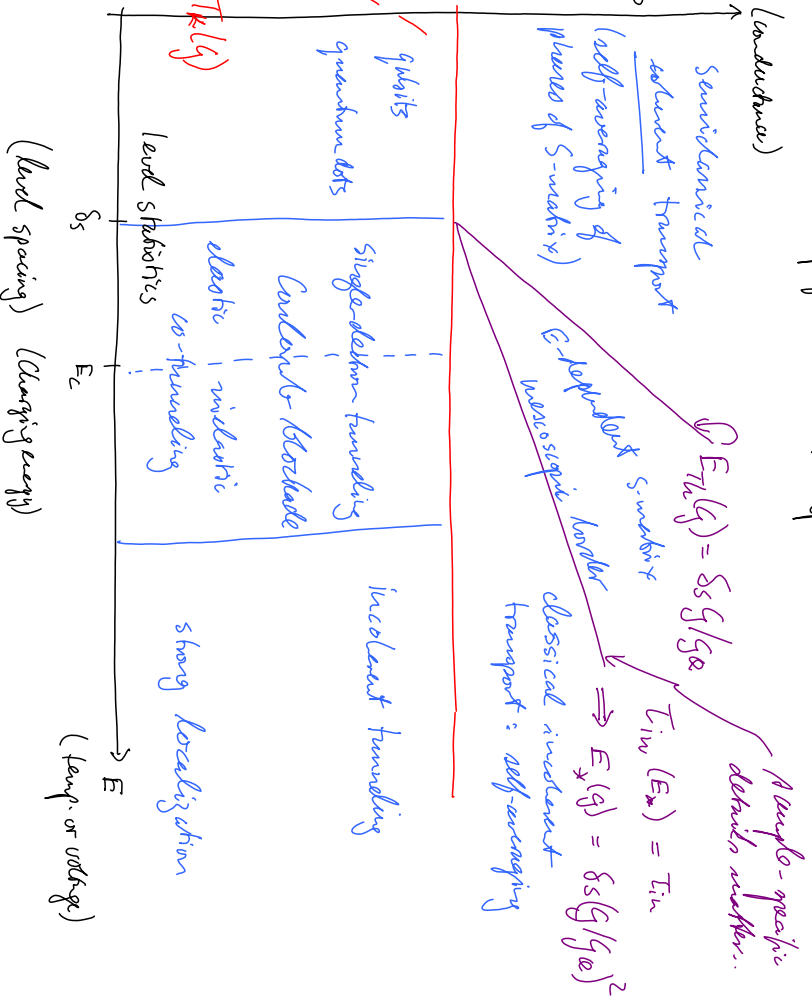
②

low- $\hbar$
plate

$\mathcal{G}$ (conductance)
scattering
semiclassical
coherent transport
(self-averaging of
phases of S-matrix)

$$g_0 = \frac{2e^2}{h}$$

(conductance
quantum)
 g_0
 $E = T(k, q)$

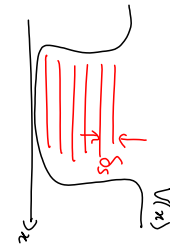


$$\text{Conductance} = G = \frac{\sigma A}{L} = \text{conductivity} \cdot \frac{\text{Area}}{\text{Length}} = \frac{1}{R} = \frac{1}{\text{resistance}} \quad (3)$$

Quantum conductance: $G_Q = \frac{e^2}{h} = \text{max. conductance of single quantum channel}$

$G \gg G_Q$: many "transport channels" (page 1, picture (1)(2))

$G \ll G_Q$: rare tunneling events (page 1, picture (3), (4))

$$\begin{aligned} \delta_S &= \text{mean level spacing} \\ &= \frac{1}{\nu} \quad (\nu = \text{s.p. density of states}) \end{aligned}$$


$E_C = \text{charging energy (cost of adding one electron)}$

for above: $\delta_S \approx 1 \text{ eV}$, $E_C \approx 10 \text{ eV}$

system size $\uparrow$: δ_S and $E_C \downarrow$, with $\delta_S \ll E_C$

For $G \gg G_Q$: electrons don't "bump" in nanostructure (4)

long enough to feel δ_S or E_C , relevant energies instead:

$$\text{Thomson Energy: } E_{Th} = \frac{\hbar}{\tau_{Th}} \quad \text{where } \tau_{Th} = \text{time to cross the nanosystem} \quad (1)$$

$$\text{Ballistic systems: } \tau_{Th} = L/v_F = \frac{\text{Length}}{\text{Fermi velocity}} \quad (2)$$



$$\text{Diffusive systems: } \tau_{Th} = L^2/D \quad (3)$$

$$D = \text{diffusion constant} = \frac{v_F^2 \ell}{3} \quad (\ell = \text{scattering length}) \quad (4)$$



$$\text{Can be shown: } E_{Th} \approx \delta_S \quad G/G_Q \quad (5)$$

$$\text{photochemical decay rate} = \gamma_{\text{in}}(E) = \frac{1}{\text{time between inelastic collisions}} = \frac{1}{\tau_{\text{in}}(E)} \quad (5)$$

$$\text{Classical transport happens if } \tau_{\text{Th}} \gg \tau_{\text{in}}(E) \quad (2)$$

$$(\text{system loses phase coherence before forming a sample}) \quad (3)$$

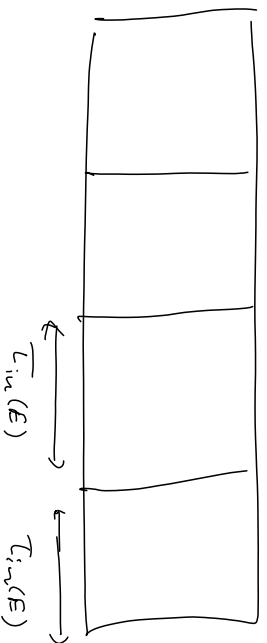
Boundary of classical region is where $\tau_{\text{in}}(E) = \tau_{\text{Th}}$

$$\text{can be shown that } E_* = \delta_S (g/g_0)^2 \quad (4)$$

$$= E_{\tau_{\text{Th}}} (g/g_0) \geq E_{\text{TH}}$$

(6)

Therefore: Think of big conductor divided in many small ones, where transport time are of order the inelastic time, $\tau_{\text{in}}(E)$



combination of these losses "self-averaging", i.e. sample-specific properties don't matter.

$$\text{When entire system becomes so small so one single loss, i.e. when } \tau_{\text{TH}} = \tau_{\text{in}}(E) \quad (\text{"mesoscopic boundary"}) \quad (1)$$

then there is no self-averaging, hence "mesoscopic"

(sample-specific) effects begin to show up.

Chapter 1: Scattering approach ($g \gg g_0$, $E < E_x$) ⑦

Conductance is a large number, described by scattering matrix S .

For $E < E_{TH}$, S -matrix is E -independent.

For $E_{TH} < E < E_x$, S -matrix becomes E -dependent.

Chapter 2: semi-classical coherent regime ($g \gg g_0$)
 more efficient description than scattering approach is available:
 self-averaging over phases of S -matrix can be done, to arrive
 at "quantum circuit theory" (similar in spirit to
 classical circuit theory)

Chapter 3: Coulomb blockade ($g \ll g_0$) ⑧

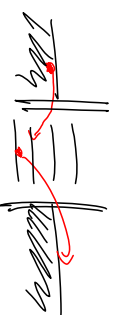
single-electron tunneling, charging energy E_c becomes important!

Quantum - connection to SET: cotunneling.

$E < \sqrt{g_s E_c}$: elastic cotunneling
 (through same level)



(scattering approach can be used!)



$\sqrt{g_s E_c} < E < E_c$: inelastic cotunneling

tunneling + superconductivity $\Rightarrow$ Josephson junction! Replaces
 coherence through sample!



Chapter 4: Randomness and interference (9)

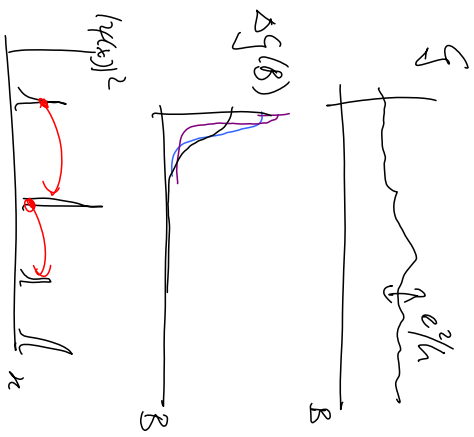
$\gamma \ll \gamma_0$ and $E \approx \delta_s$: "level statistics" (fluctuations of discrete energy levels)

$\gamma \gg \gamma_0$ and $E \approx E_x$:

fluctuations in transmission eigenvalues
"radiation fluctuations".

interference corrections: small localization:

$\gamma \ll \gamma_0$ and $E \gg E_x$: strong localization of wave functions, electron hopping



Chapter 5: Rubib and Quantum Dots (10)

need discrete energy levels $\Rightarrow \gamma \ll \gamma_0$, $E \ll \delta_s$
quantum information & manipulation.

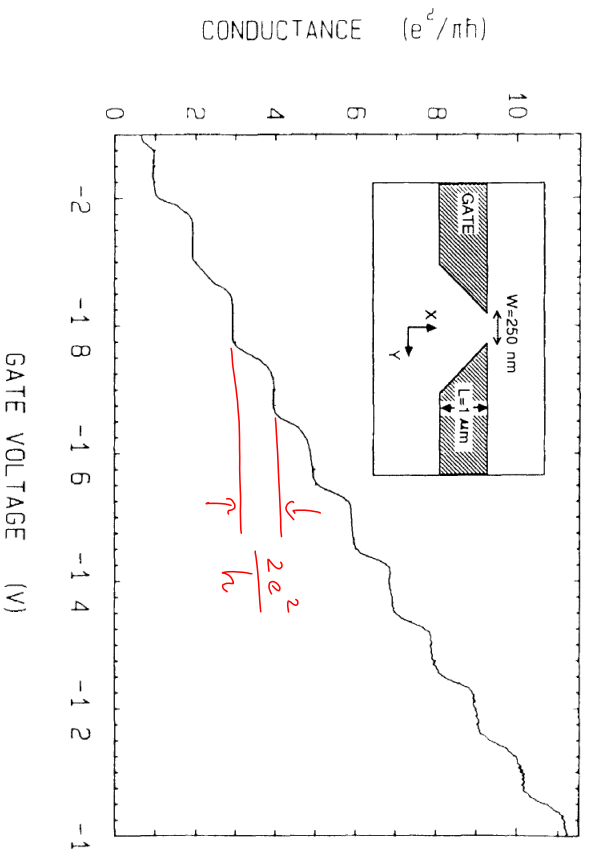
Chapter 6: Interaction, relaxation, decoherence (effects beyond Coulomb blockade)

- dissipative quantum mechanics
 - $\gamma \ll \gamma_0$: effect of environment on tunneling
 - $\gamma \approx \gamma_0$: interaction effects at higher conductance (broad effect, T_e depends exponentially on S).
 - departing for qubits and electrons (at mesoscopic boundaries)

1.2 Quantum Point Contact : Conductance quantization

(11)

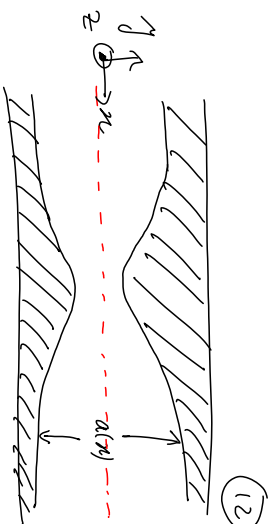
van Wees, B. J.; van Houten, H.; Beekker, C. W. J.; Williamson, J. G.; Kouwenhoven, L. P.; van der Marel, D. & Foxon, C. T. Quantized conductance of point contacts in a two-dimensional electron gas Phys. Rev. Lett., 60, 848-850 (1988)



Point contact $\approx$ adiabatic wave guide :

"adiabatic" = charges smoothly :

$$\left| \frac{da}{dx} \right| \ll 1, \quad a(x) \left| \frac{d^2a}{dx^2} \right| \ll 1 \quad (1)$$



$$\text{Potential: } V(x, y, z) = \begin{cases} 0 & \text{if } |y| < a(x)/2, |z| < b(x)/2 \\ \infty & \text{otherwise.} \end{cases} \quad (2)$$

$$\text{Schrödinger eq: } \left[-\frac{\hbar^2}{2m} \nabla^2 + V(x, y, z) - E \right] \psi(x, y, z) = 0 \quad (3)$$

$$\text{variables separate locally: } \psi_n(x, y, z) = \psi_n(x) \Phi_n(a(x), b(x); y, z) \quad (4)$$

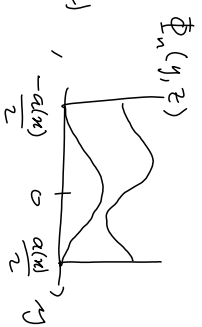
$$\left(-\frac{\hbar^2}{2m} \partial_x^2 + E_n(x) - E \right) \psi(x) = 0 \quad (5)$$

Effective potential for 1D motion in x-direction.

$$\left[-\frac{\hbar^2}{2m} (\partial_y^2 + \partial_z^2) + V(y, z) - E_n(\lambda) \right] \Phi_n(a(x), b(x), y, z) = 0 \quad (1)$$

transverse modes are standing waves:

$$\Phi_n(a, b; y, z) = \frac{2}{\sqrt{ab}} \sin(k_y^{(n)}(y - a/2)) \sin(k_z^{(n)}(z - b/2)) \quad (2)$$



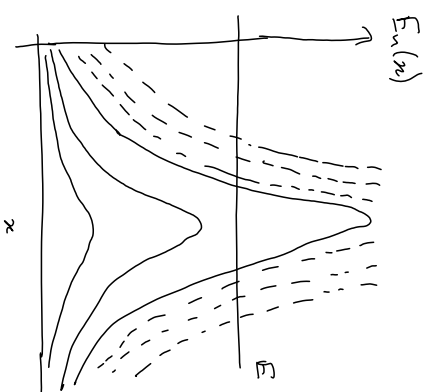
$$\text{mode index: } n \equiv (n_y, n_z) \quad (3)$$

transverse momenta:

$$k_y^{(n)} = \frac{\pi n_y}{a}, \quad k_z^{(n)} = \frac{\pi n_z}{b} \quad (4)$$

transverse energy:

$$E_n(a(x), b(x)) = \frac{\pi^2 \hbar^2}{2m} \left(\frac{n_y^2}{a^2(x)} + \frac{n_z^2}{b^2(x)} \right) \quad (5)$$

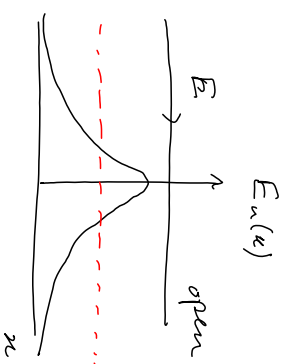


For given total energy E and given n , we have 1-d transmission problem: (1c)

If $E > E_n(x)$, channel is open
< lower energy (closed)

open: Transmission $T \simeq 1$
 closed: $T \simeq 0$

in general: $T_n(E) \in [0, 1]$ (tunneling)

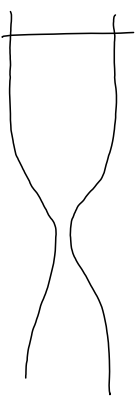


$\Rightarrow$ there are only finite number of open channels:

Calculating the current:

$$I = \int dy \int dz j(x = -\infty, y, z) \quad (1)$$

current density, midpt. of y, z at $x = -\infty$



(15)

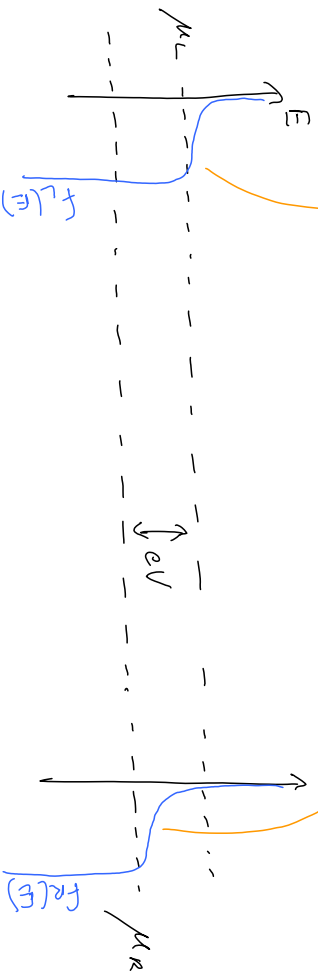
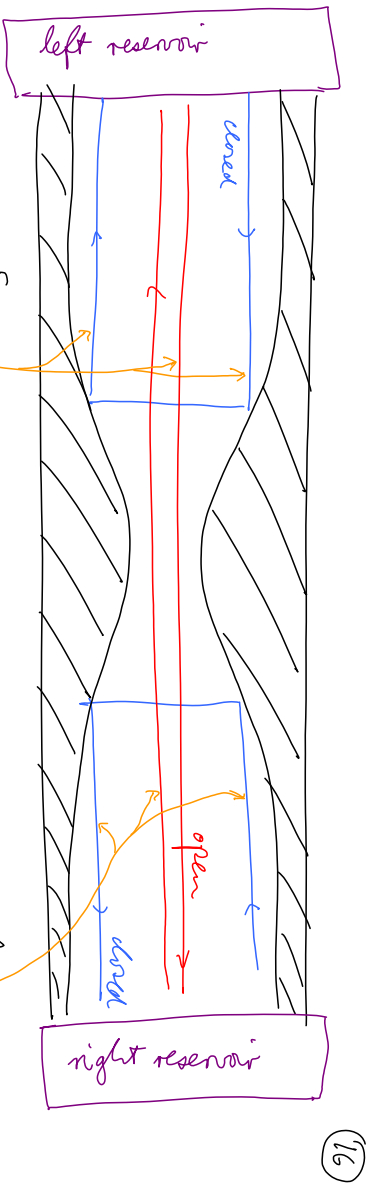
$$= ab j = ab \frac{2s}{\tau_{\text{spin}}} \int \frac{d^3 \hbar}{(2\pi)^3} e^{i\vec{r} \cdot \vec{v}(\vec{r})} f(\vec{r}) \quad (2)$$

↑ ↑
change relating occupation function

(13,4)

↑ transverse modes

$$= ab \int \frac{dk_x}{2\pi} \frac{1}{ab} \sum_n e^{i\psi_x(k_x)} f_n(k_x) \quad (3)$$



For closed channels;

$$f_n(k_x) = f_n(-k_x)$$

no contribution, arise:

$$\psi_x(k_x) = -\psi_x(-k_x)$$

(4)

For open channels:

(17)

$k_x > 0$ one from left reservoir, with
 $k_x < 0$ right

$$\left. \begin{aligned} f_L(E(k_x)) &= f_F(E(k_x) - \mu_L) \\ f_R(E(k_x)) &= f_F(E(k_x) - \mu_R) \end{aligned} \right\} \quad (1) \quad (2)$$

change variables: $\int dk_x \underbrace{v_x(k_x)}_{= \frac{1}{k} \frac{\partial E(k_x)}{\partial k_x}} = \frac{1}{k} \int dE \quad (3)$

$$\begin{aligned} (12.3) \mu_i (15.3) \quad I &= \frac{2s}{2\pi} \frac{E}{k} \sum_{k \in \text{open}} \underbrace{\int dE}_{N_{\text{open}}} \underbrace{[f_L(E) - f_R(E)]}_{eV = \mu_L - \mu_R} \\ k &= \frac{h}{2\pi} = \frac{2s}{N_{\text{open}}} \frac{e}{V} \end{aligned}$$

Conductance: $G = \frac{I}{V} = G_0 N_{\text{open}} \Rightarrow$ conductance quantization!
independent of material properties !!