

# Quantum Dots II

1. Open Dot Experiments

2. Kondo effect

3. Few Electron Dots

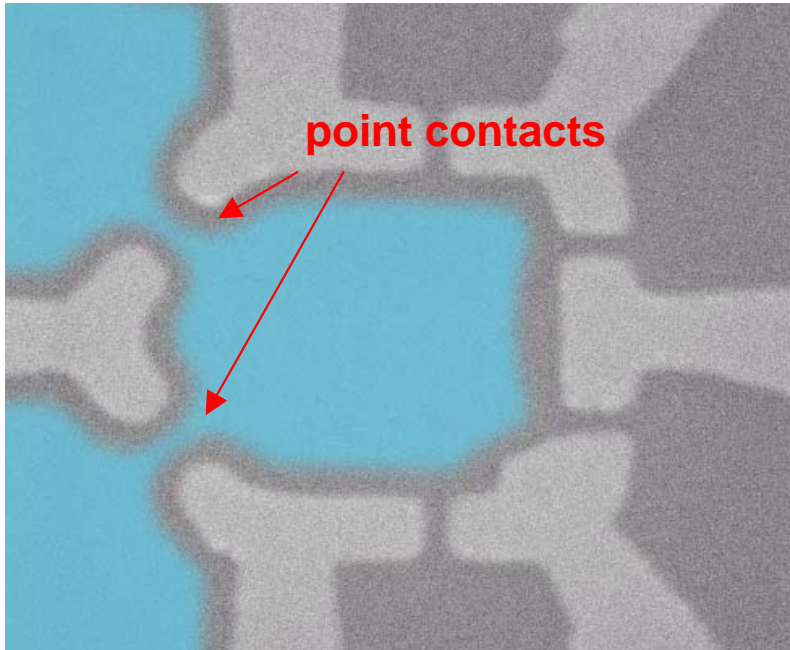
4. Double Quantum Dots

Huibers, Ph.D. Thesis (1999)

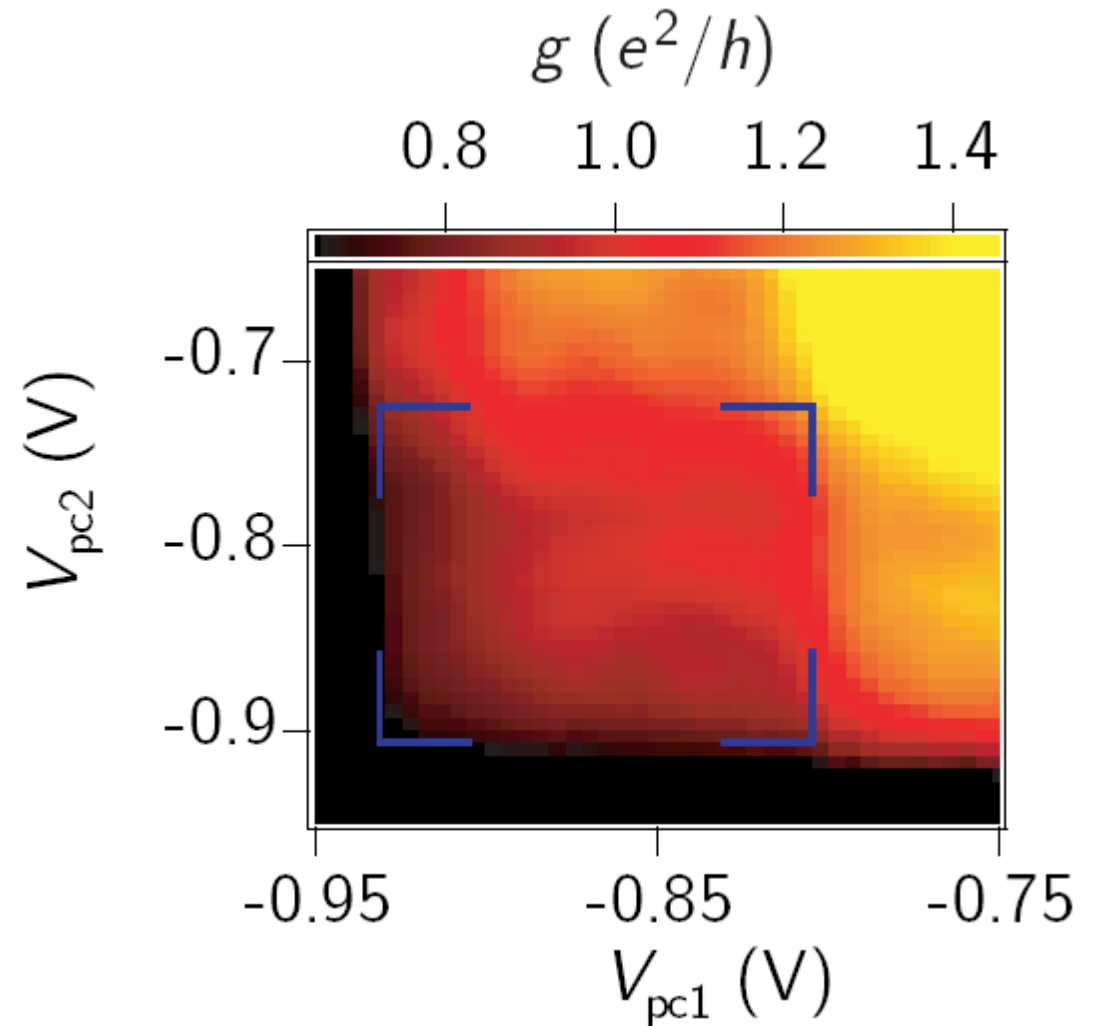
Huibers et al., PRL83, 5090 (1999)

# Open Dot Regime

## Open Dot



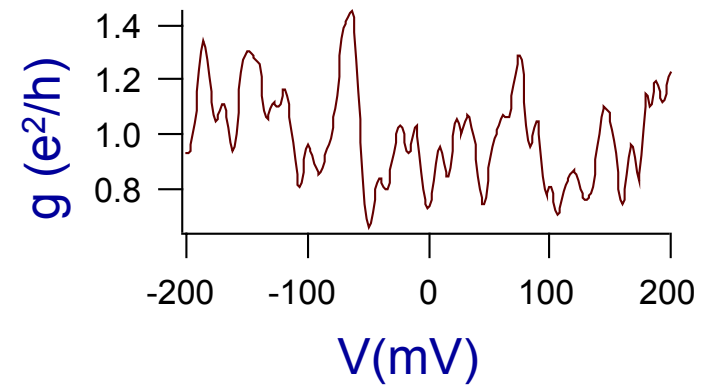
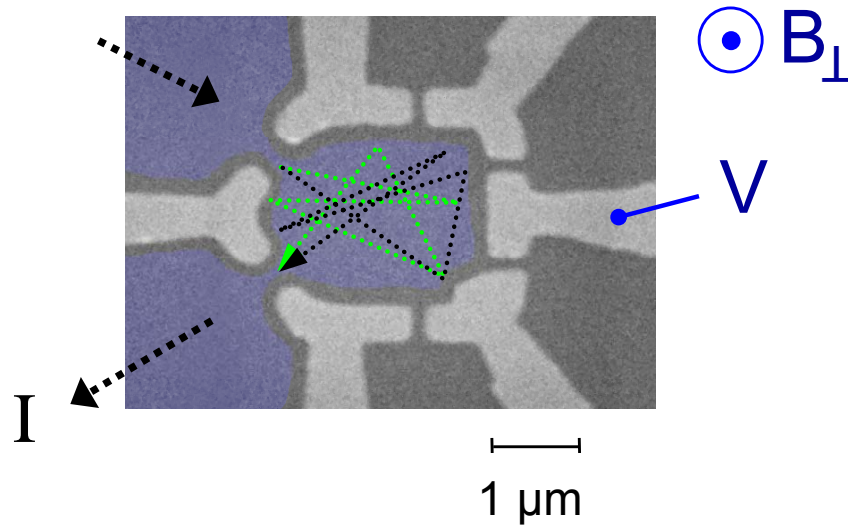
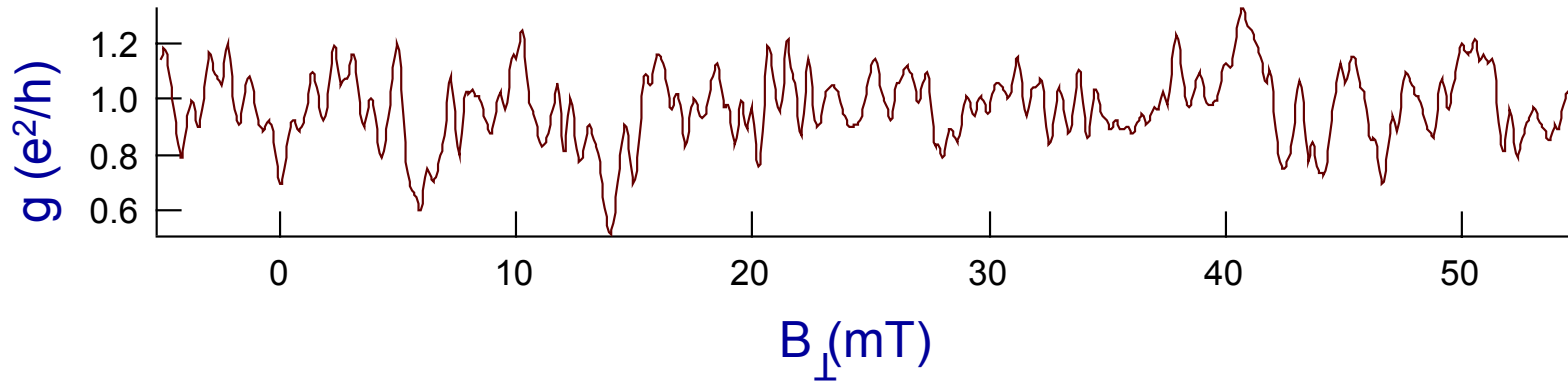
- $V_{\text{gate}}$  set to allow  $\geq 2e^2/h$  conductance through each point contact
- Dot is well-connected to reservoirs
- Transport measurements exhibit CF and Weak Localization



many open dot slides: A. Huibers and J. Folk

# Open Dot Regime: Conductance Fluctuations

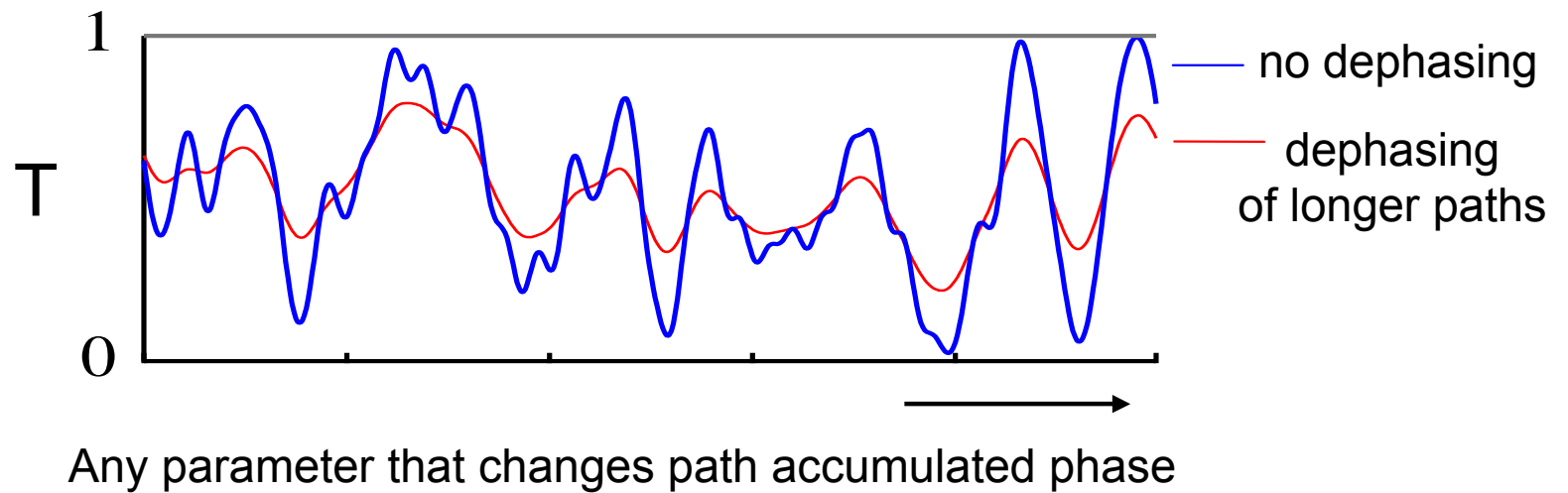
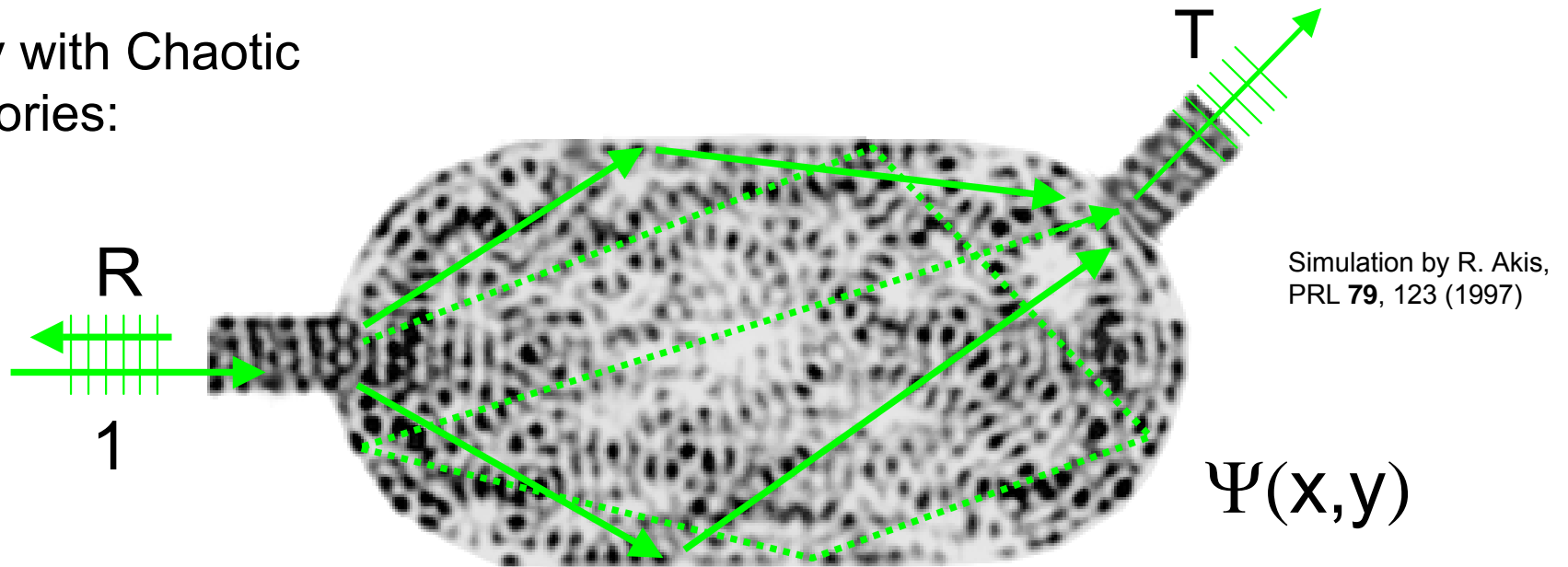
$$N_L = N_R = 1$$



Repeatable random  
interference fluctuations  
as function of dot parameters

# Two-Dimensional Quantum Dot

2D Cavity with Chaotic trajectories:

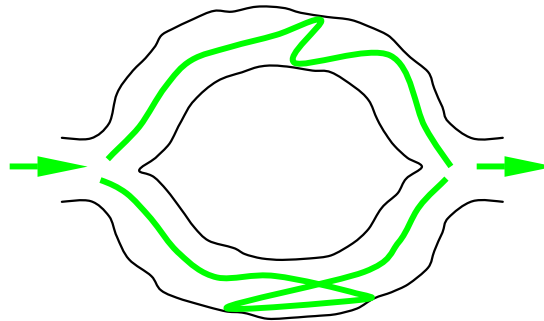


**Goal:** use quantum dot as a probe of quantum phase coherence

# Interferometers

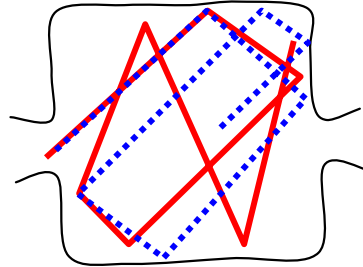
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Two-arm:



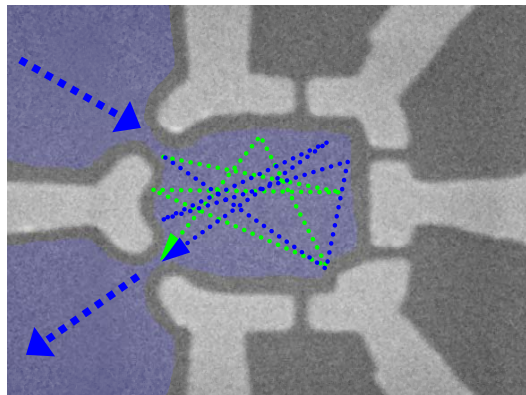
sometimes:  
reflections or  
small signal

Regular/  
Integrable:



Problem:  
partially chaotic

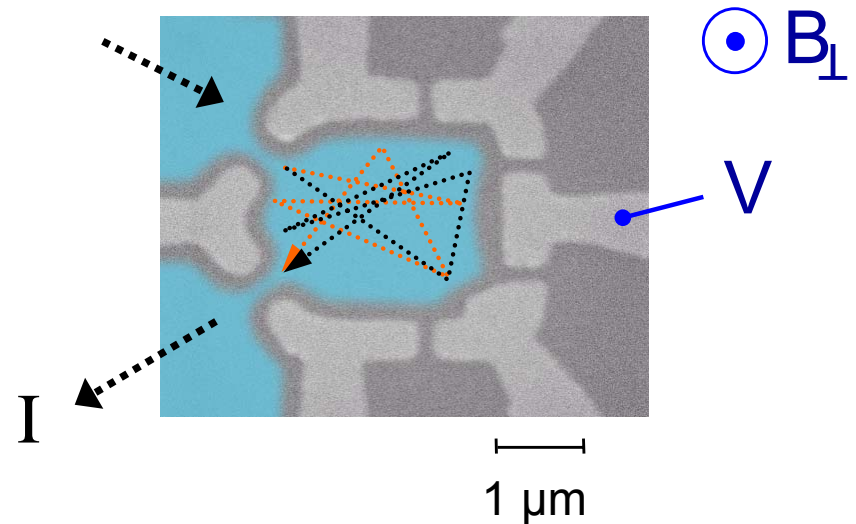
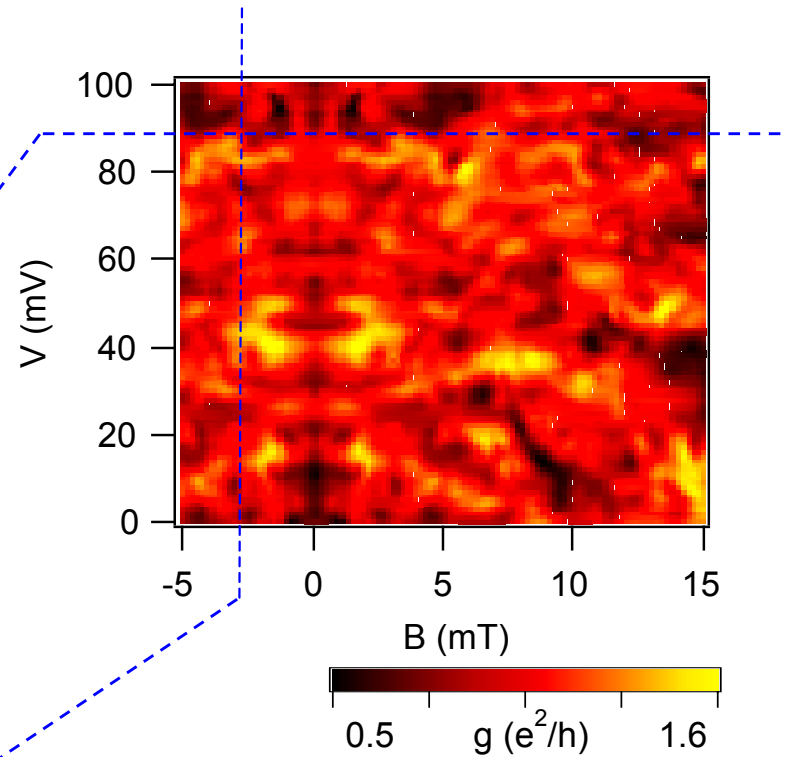
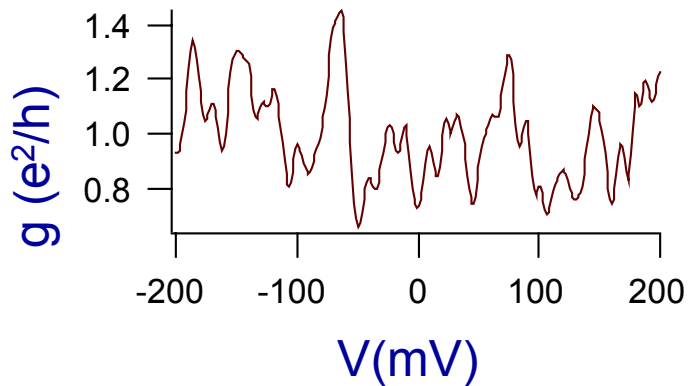
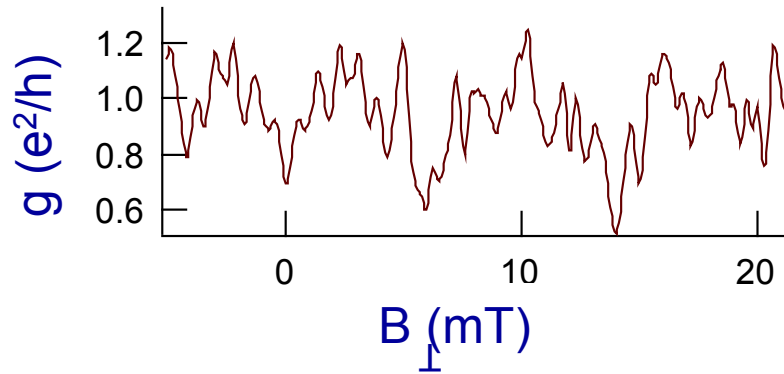
Chaotic:



1. Mostly chaotic/ergodic
2. Interesting physics & complete description

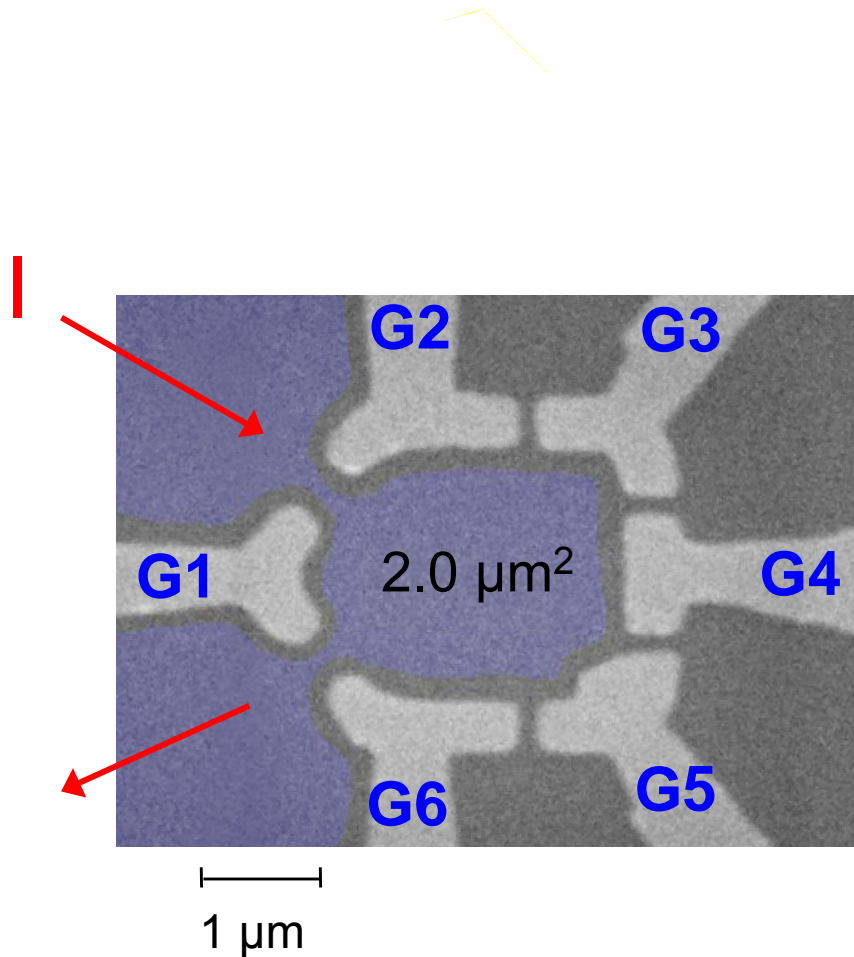
# Quantum Interference in Open Dots

Interference between all possible trajectories gives rise to repeatable random interference fluctuations as function of dot parameters



# Typical Quantum Dot

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2D conductor:

area =  $2.0 \mu\text{m}^2$

charge density =  $2 \cdot 10^{11} \text{ e/cm}^2$

$\lambda_F$  = Fermi wavelength = 50 nm

$v_F$  = Fermi velocity =  $200 \mu\text{m/ns}$

$E_F$  = Fermi energy = 7 meV

Dwell time in dot: 200 ps

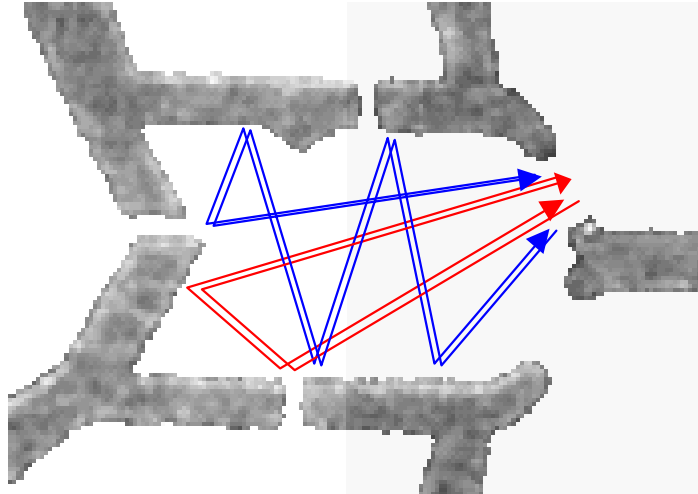
Crossing time: 7 ps

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30 bounces

bulk mean free path  $\ell_e \sim 2\text{-}10 \mu\text{m}$

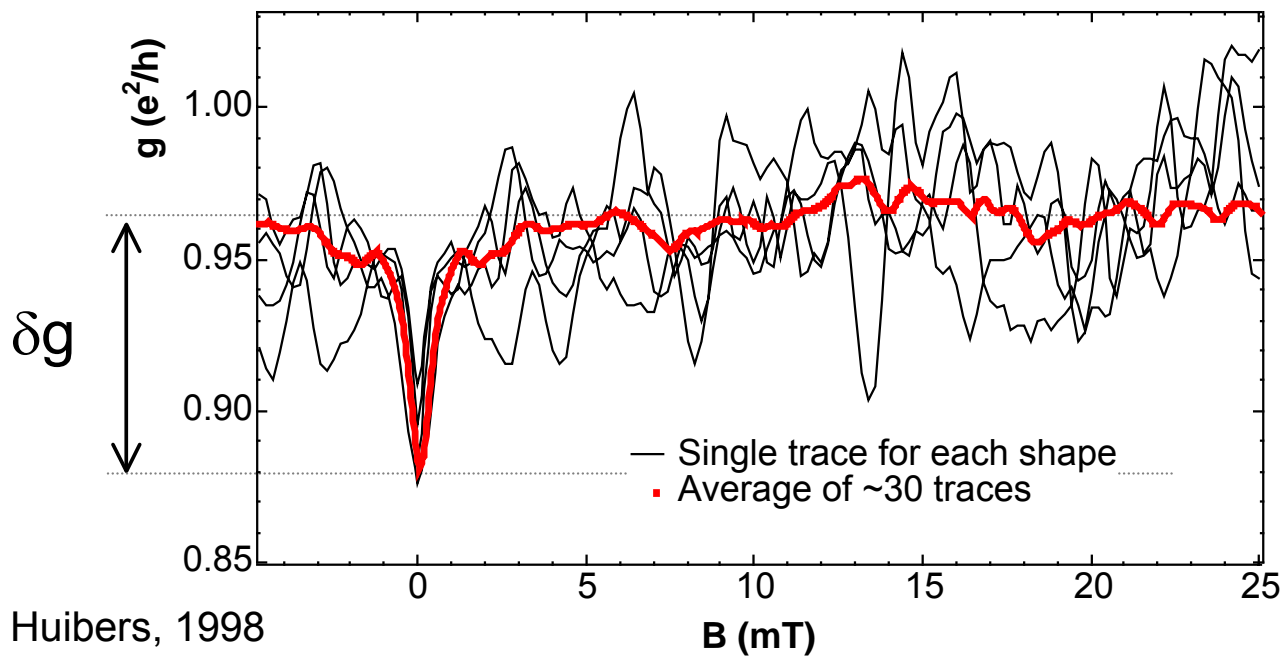
# Weak Localization



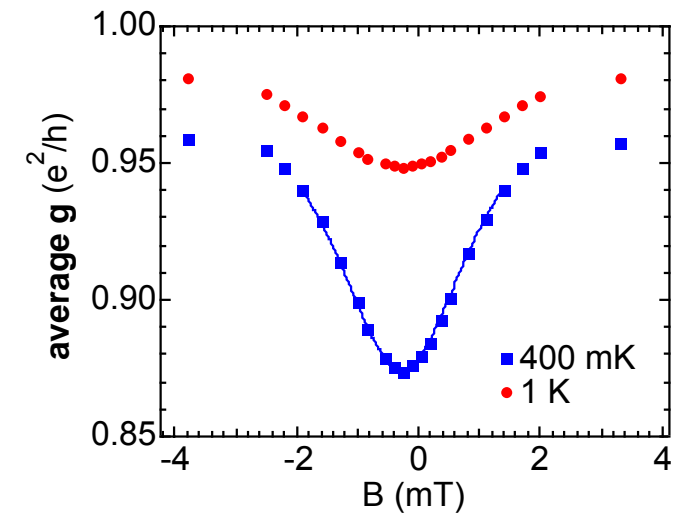
At  $B=0$ , phase-coherent backscattering results in “weak localization”



Conductance dip at  $B=0$



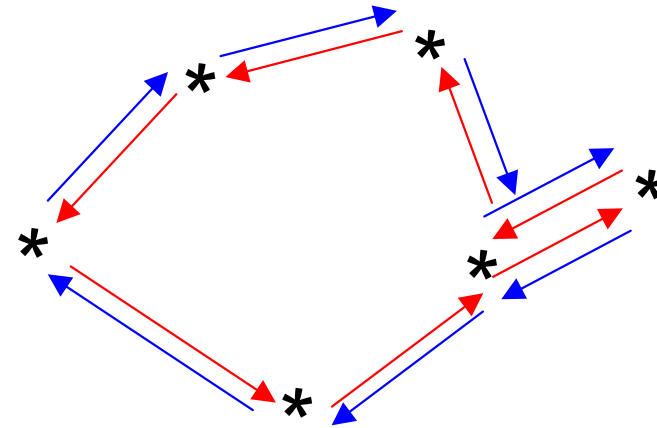
Huibers, 1998





# Quantum Correction: Weak Localization

constructive interference of coherently backscattered, time reversed trajectories decreases conductivity



$$\frac{\delta\sigma_{\text{loc}}}{\sigma} \propto -\frac{1}{k_{\text{F}}\ell} \ln\left(1 + \frac{\tau_{\phi}}{\tau}\right) \quad 2\text{D } (L_{\phi} \ll W)$$
$$\frac{\delta\sigma_{\text{loc}}}{\sigma} \propto -\frac{L_{\phi}}{W} \frac{1}{k_{\text{F}}\ell} \left(1 - \left(1 + \frac{\tau_{\phi}}{\tau}\right)^{-1/2}\right) \quad 1\text{D } (L_{\phi} \gg W)$$

magnetic field: AB-flux, cut off trajectories of area  $A > \phi_0 B$   
magnetoconductance

(assuming spinless electrons)

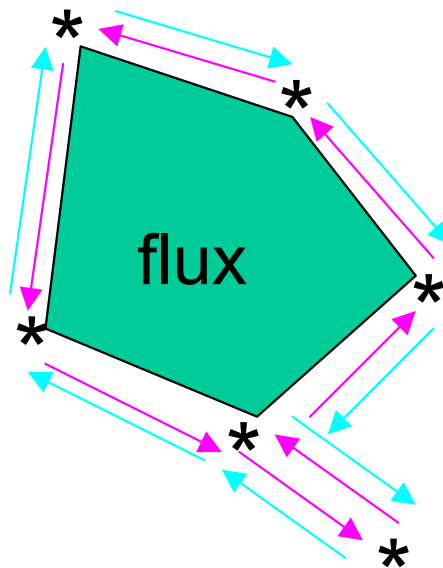
## Weak Localization in Magnetic Fields

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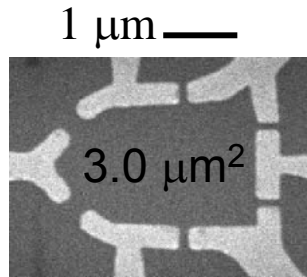
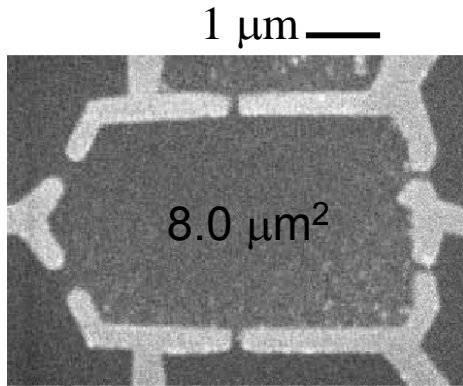
in a given magnetic field  $B$ , trajectories enclosing flux acquire additional Aharonov-Bohm phase:

$$\phi = \frac{2e}{\hbar} \int (\nabla \times \mathbf{A}) \cdot d\vec{S} = \frac{2eBS}{\hbar}$$

when summing over all trajectories, this  $\phi$  will effectively eliminate trajectories of area  $A \gg \phi_0/B$ . ( $\phi_0 = h/e$ )



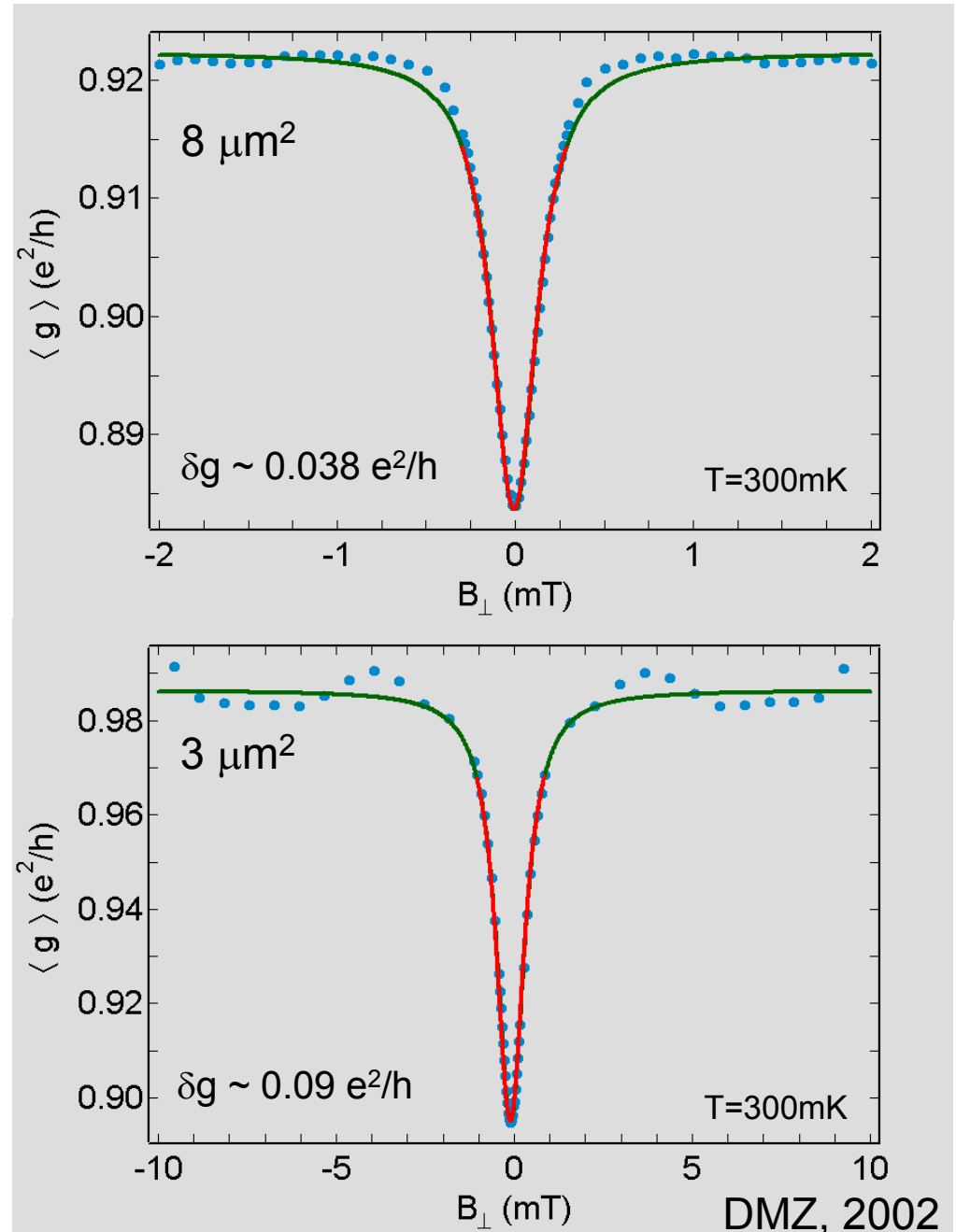
# Weak Localization: Measure of Dephasing



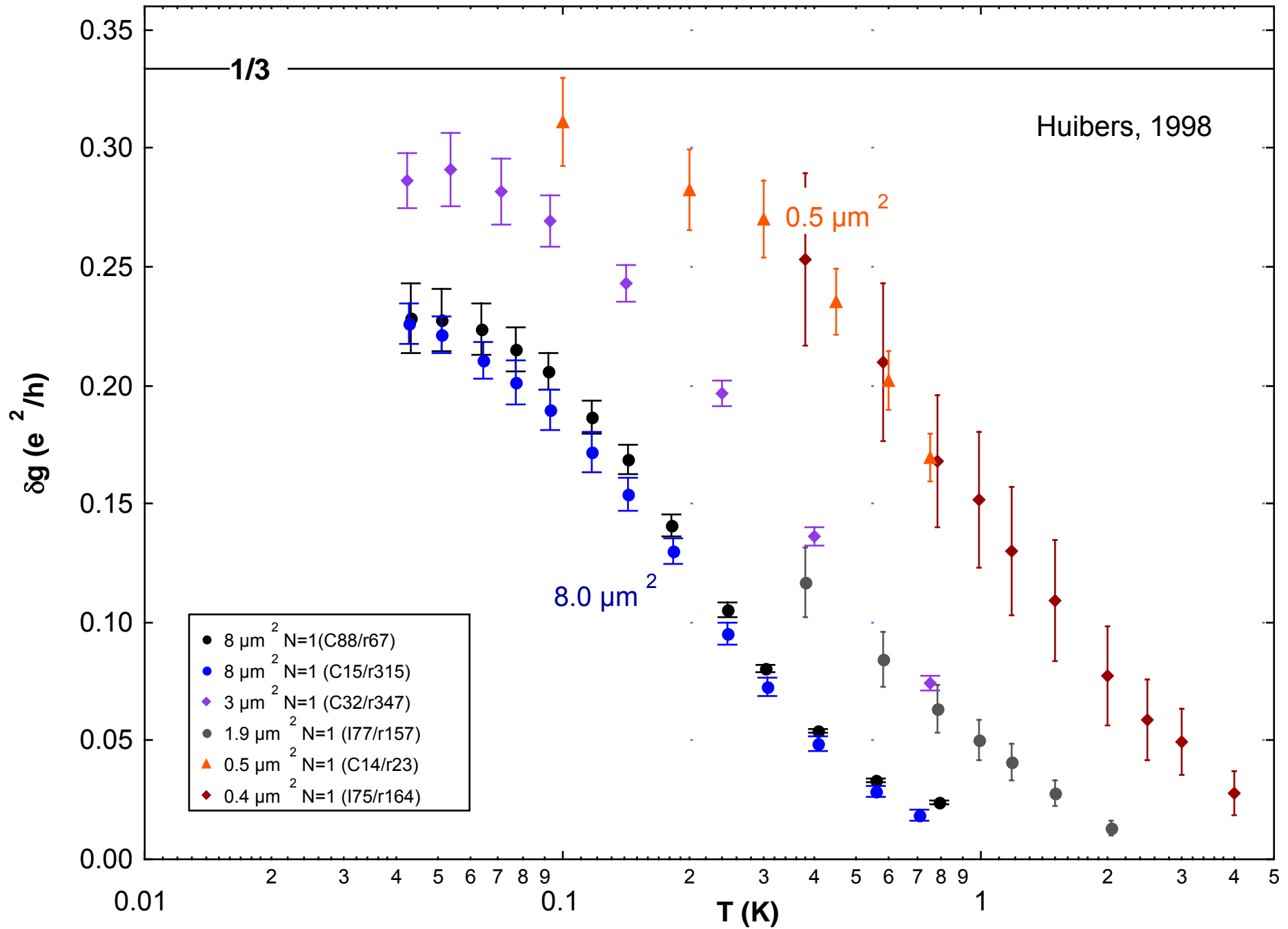
random matrix theory

$$\delta g = \frac{1}{2N + 1 + \gamma_\phi}$$

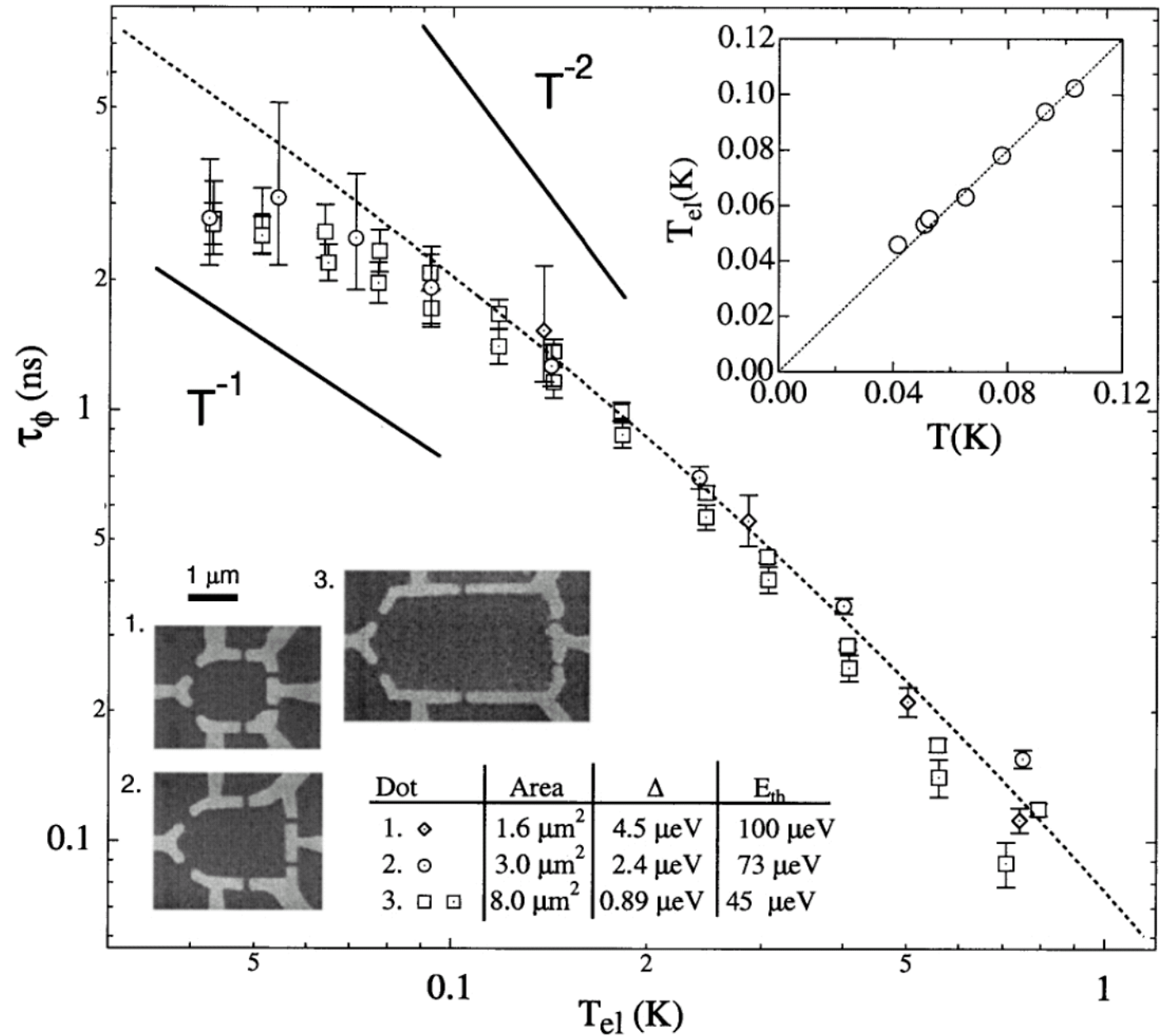
$$\tau_\phi^{-1} = \frac{\Delta}{h} \gamma_\phi$$



# Weak Localization vs T



# Low Temperature Saturation?



# Spin-Orbit Coupling

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electrons move with the Fermi velocity, electric fields in material appear as magnetic fields in the rest frame of the electron

these magnetic fields

- depend on magnitude of electron velocity (density dependence)
- couple to the electron spin via Zeeman coupling

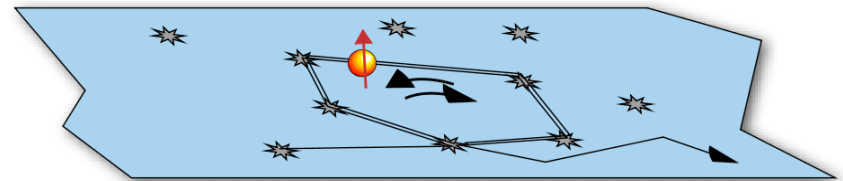
—————▶ spin-precessions

electric fields due to:

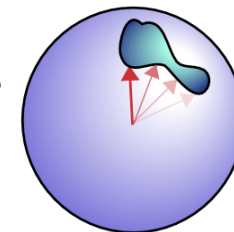
- heterointerface (*Rashba*)
- crystalline anisotropy in III-V zincblende crystal (*Dresselhaus*)

*spin precession affects phase interference  
( $2\pi$  in spin space gives -1 to phase)*

motion in real space



motion in spin space



# Spin-Orbit Coupling

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presence of electric fields  $\vec{E} = -\frac{1}{e}\vec{\nabla}V$

electrons are moving in these electric fields

rest frame of electrons: effective magnetic field

$$\vec{B}_{\text{so}} = -\frac{\vec{v}}{c} \times \vec{E}$$

magnetic moment  $\vec{\mu} = \frac{e\vec{S}}{mc}$  of electron couples to  $\vec{B}_{\text{so}}$

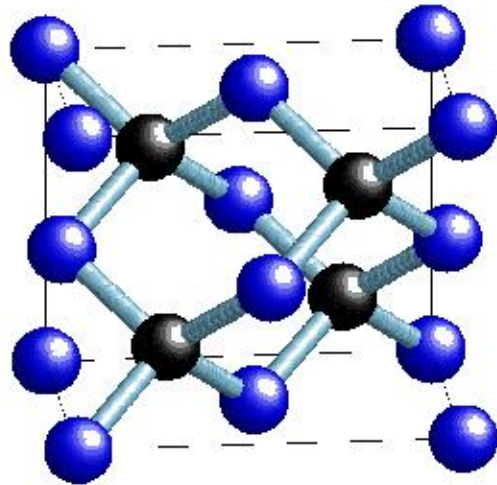
$$H_{\text{so}} = -\vec{\mu} \cdot \vec{B}_{\text{so}}$$

electrons precess around  $B_{\text{so}}$

$B_{\text{so}}$  depends on the electron momentum

spin rotation symmetry is broken,  
time reversal symmetry is NOT broken

# Spin-Orbit Coupling due to Crystal Anisotropy



Conventional cell

III-V Semiconductor

Zinkblende crystal structure:  
two interpenetrating fcc lattices  
with only Ga atoms on one lattice,  
only As on the other

absence of inversion symmetry

symmetry considerations:

$$H_{\text{SO}} = \gamma(\sigma_x k_x (k_y^2 - k_z^2) + \text{cycl.})$$

G. Dresselhaus,  
Phys. Rev. 100, 580 (1955)

after size quantization (2D):

$$\langle k_z \rangle = 0 \quad \alpha = \gamma \langle k_z^2 \rangle$$

$$H_D^{(1)} = \alpha(\sigma_x k_x - \sigma_y k_y)$$

linear Dresselhaus term

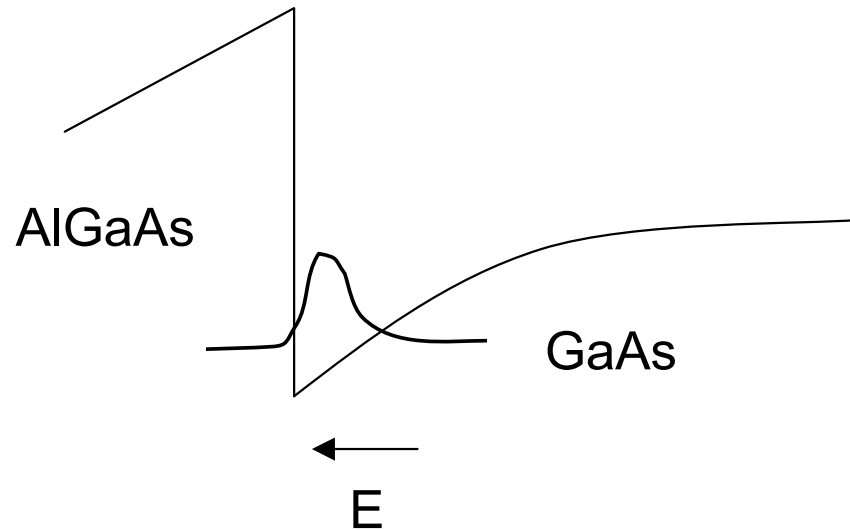
$$H_D^{(3)} = \gamma(\sigma_y k_y k_x^2 - \sigma_x k_x k_y^2)$$

cubic Dresselhaus term



# Spin-Orbit Coupling due to Heterointerface

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electric field at heterointerface  
perpendicular to 2D plane

$$\vec{B}_{so} \propto (k_y E, -k_x E, 0) \perp \vec{k}$$

$$H_R = \beta(\sigma_x k_y - \sigma_y k_x)$$

Rashba term (linear)

coupling strength parameters  $\beta$  and  $\gamma$  can be determined  
from Band structure, for example in k·p approximation

# Weak Antilocalization

initial state:  $|i\rangle$

final (forward):  $|f_f\rangle = R_N \dots R_2 R_1 |i\rangle = R|i\rangle$

final (backward):  $|f_b\rangle = R_1^{-1} R_2^{-1} \dots R_N^{-1} |i\rangle = R^{-1}|i\rangle$  (TRS)

$R_i$ : spin rotations

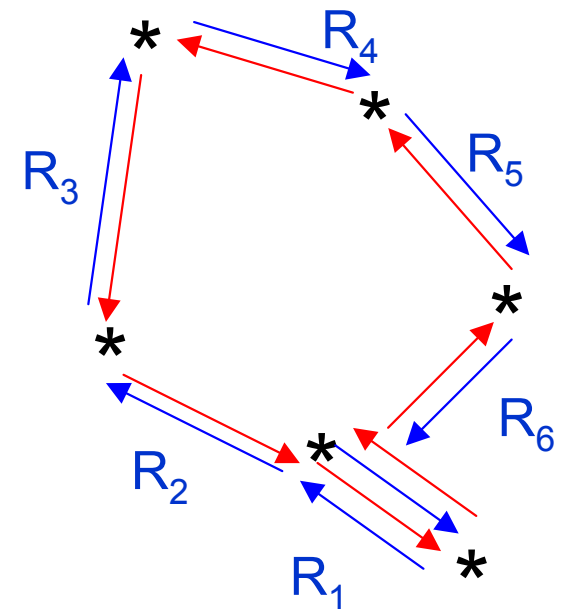
$$R = R_N \dots R_2 R_1$$

$$R^\dagger R = 1$$

$$R^{-1} = R^\dagger$$

interference term  $\langle f_b | f_f \rangle = \langle i | R^2 | i \rangle$

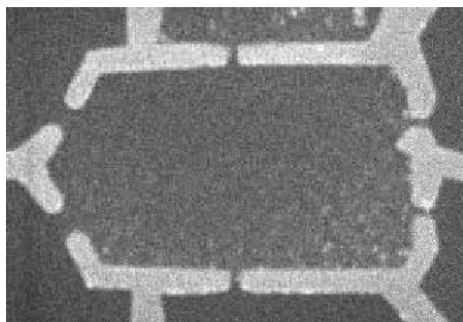
assuming strong spin-orbit coupling,  
summing over all trajectories is  
equivalent to averaging  $R^2$  over sphere



$$\overline{\langle f_f | f_b \rangle} = -\frac{1}{2}$$

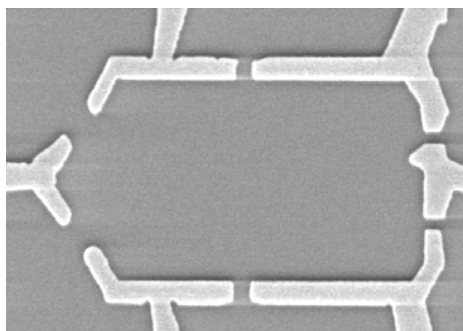
**destructive interference**  
**opposite sign for Magnetoconductance**

low density  
weaker SO coupling  
weak localization (WL)



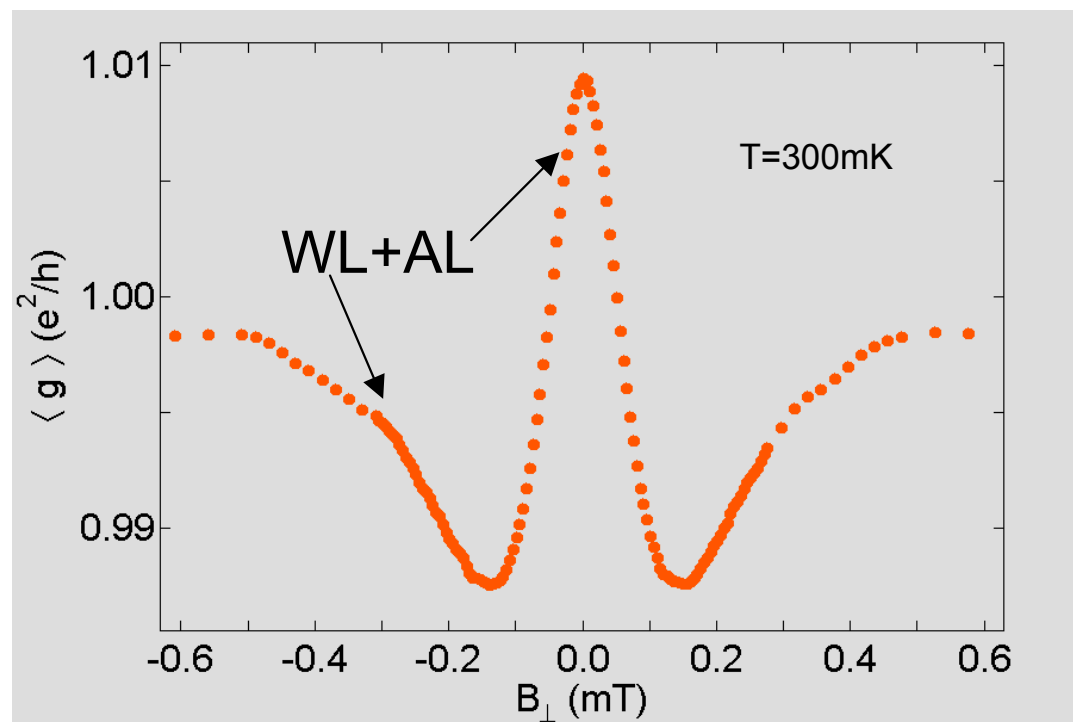
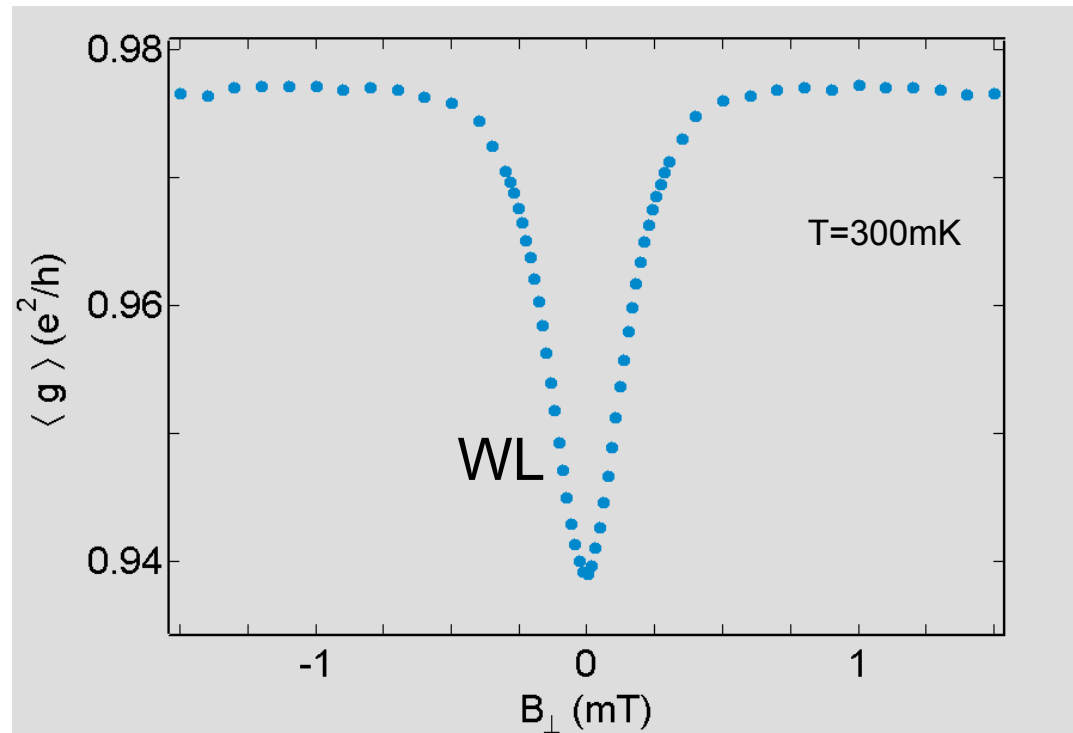
4 $\mu$ m

dots are on **different** wafers



4 $\mu$ m

high density  
stronger SO coupling  
antilocalization (AL)



1. Open Dot Experiments

**2. Kondo effect**

3. Few Electron Dots

4. Double Quantum Dots

Goldhaber-Gordon et al., Nature **391**, 156 (1998)

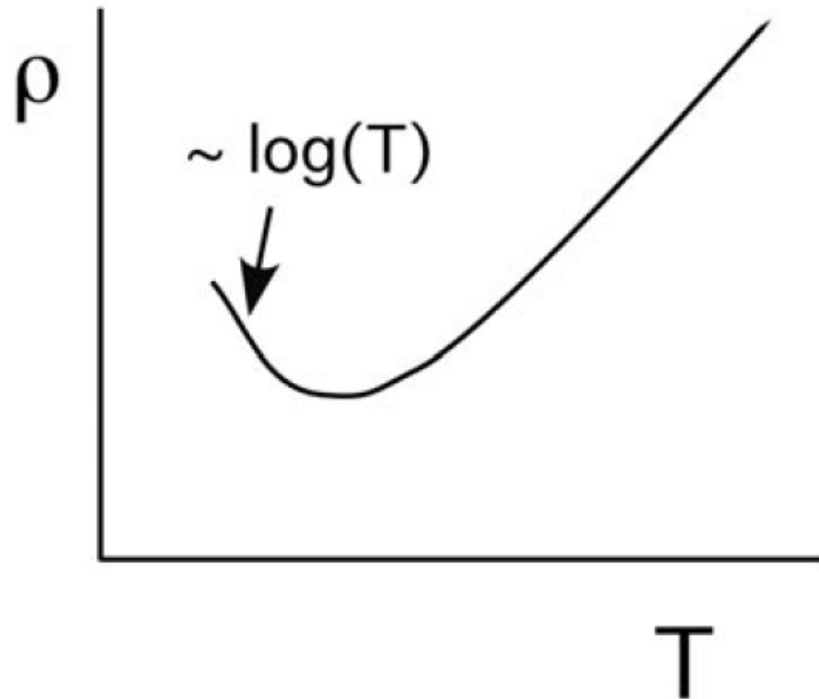
Cronenwett et al., Science **281**, 540 (1998)

S. Cronenwett, Ph. D. Thesis (2001)

# Kondo Effect in Metals

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1930s experiments:

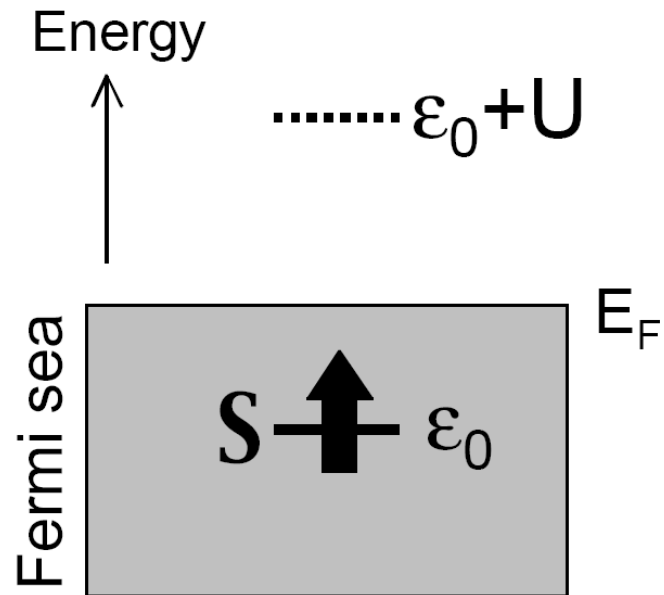


1960s: (exp) related to magnetic impurities

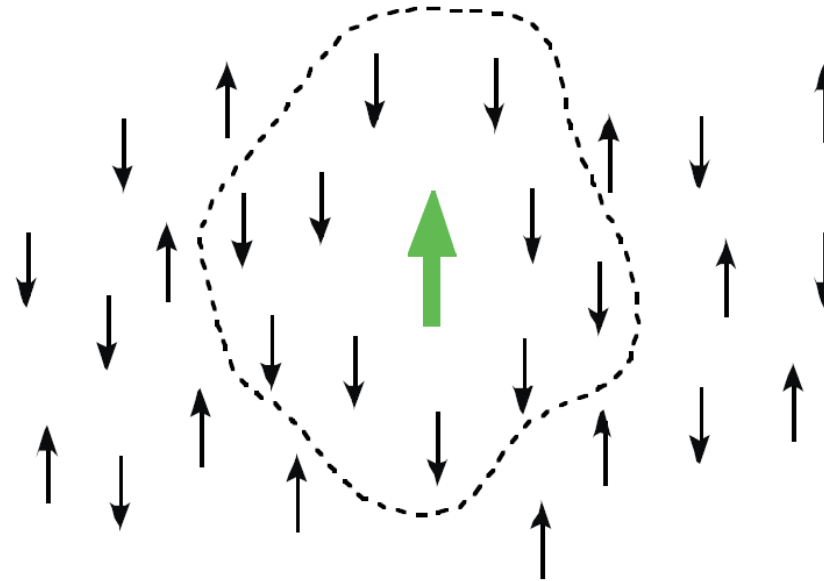
theoretical explanation by Jun Kondo  
spin-flip scattering on mag. impurities

$$\rho \sim \underbrace{\rho_0}_{\text{lattice}} + \underbrace{aT^5}_{\text{phonons}} - b \log(T)$$

# Kondo Effect in Metals: Model



new energy scale: Kondo temperature  $T_K$   
 formation of spin-singlet screening cloud



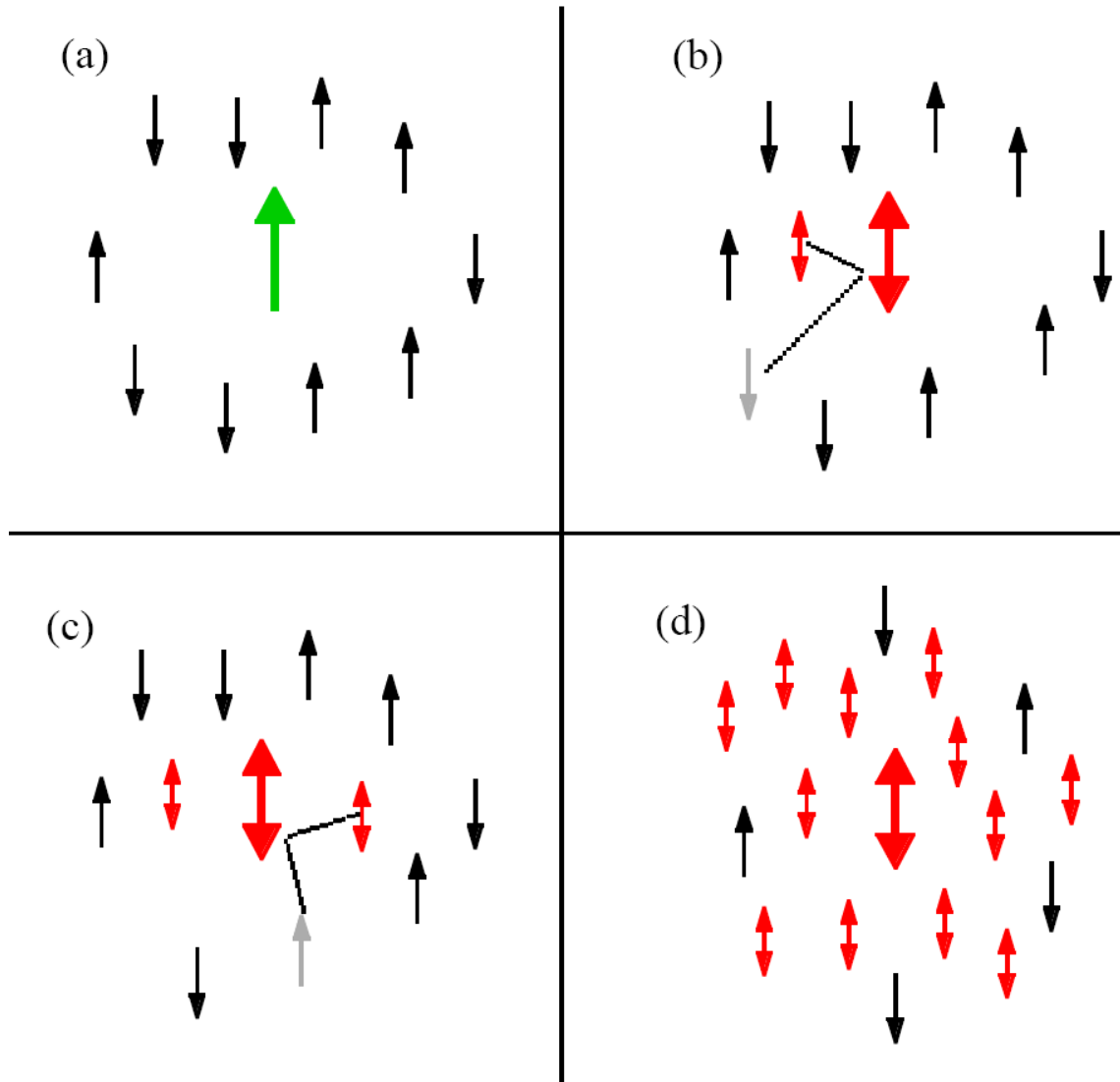
cloud: more effective scatterer  
 increase in resistance

Anderson Hamiltonian

$$H_A = \sum_{\sigma; k < k_f} \epsilon_{k\sigma} c_{k\sigma}^\dagger c_{k\sigma} + \sum_{\sigma} \epsilon_{\sigma} d_{\sigma}^\dagger d_{\sigma} + \frac{1}{2} U n_{\sigma} n_{\sigma'} + \sum_{\sigma; k < k_f} t_{k\sigma} c_{k\sigma}^\dagger d_{\sigma} + H.c..$$

free electrons                      localized electrons                      on site charging                      coupling between localized and free ele.

# Kondo Effect in Metals: spin flip scattering

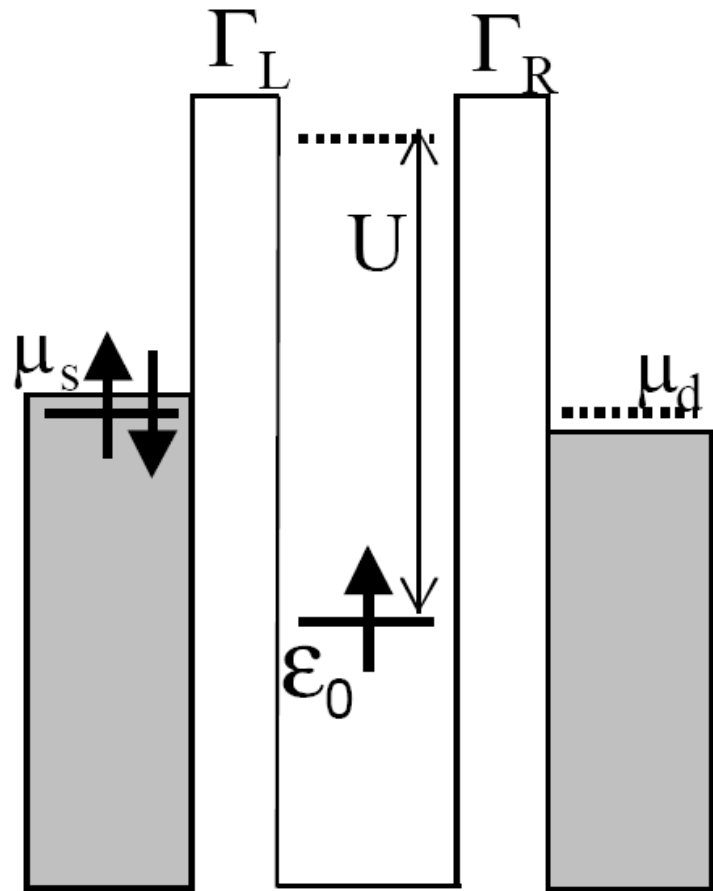


each scattering event  
entanglements impurity with  
conduction electron

singlet cloud formation

temperature scale  $T_K$

# Kondo Effect in Quantum Dots

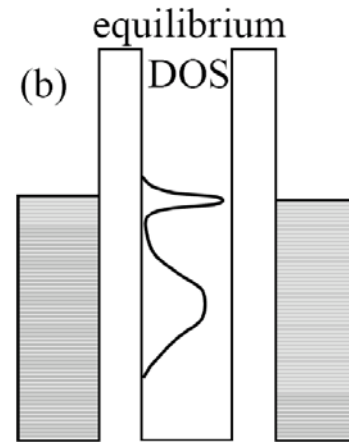
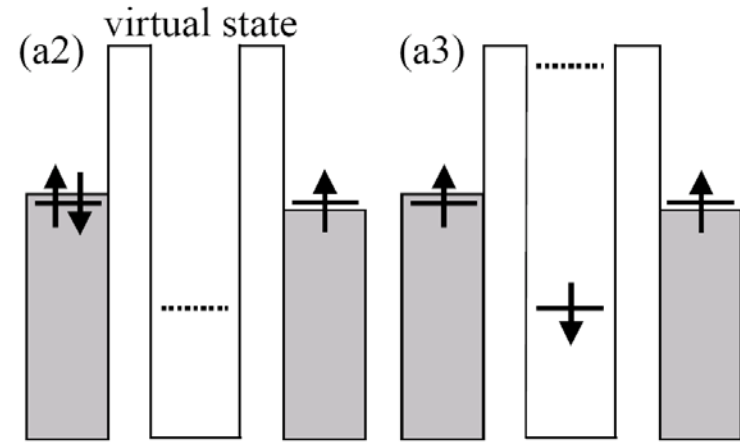


unpaired spin

$$T_K \propto \sqrt{U\Gamma} \exp(\pi\epsilon_0/2\Gamma)$$

$$\Gamma = \Gamma_L + \Gamma_R$$

spin-flip cotunneling (elastic)



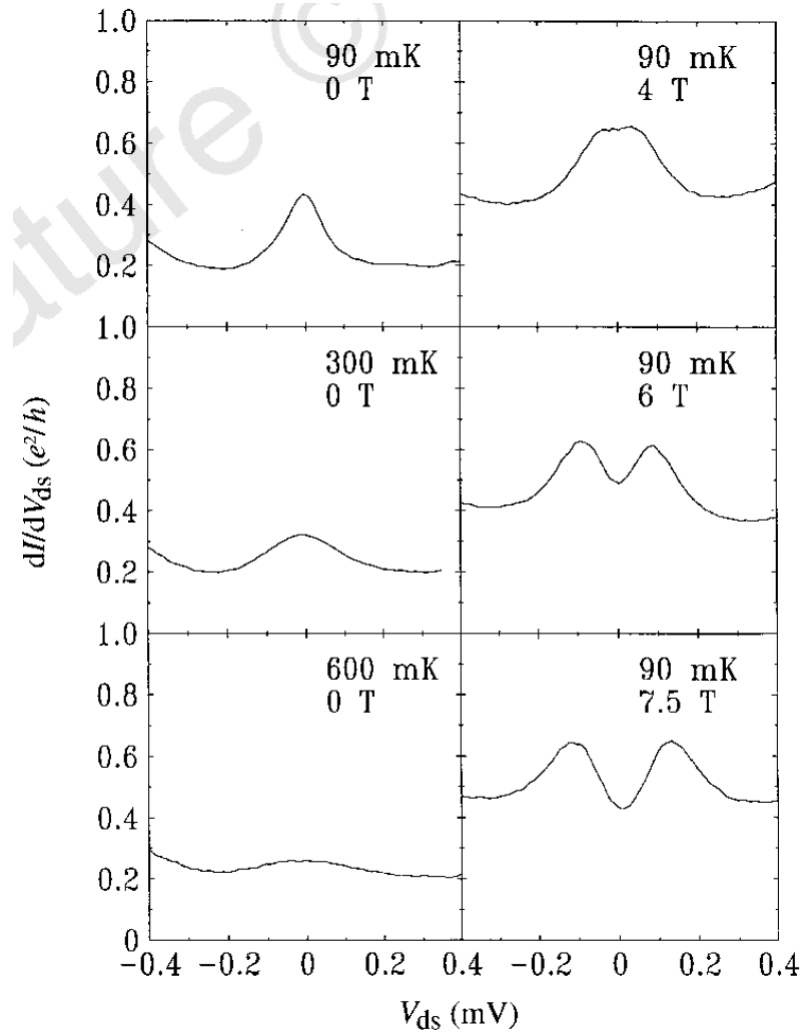
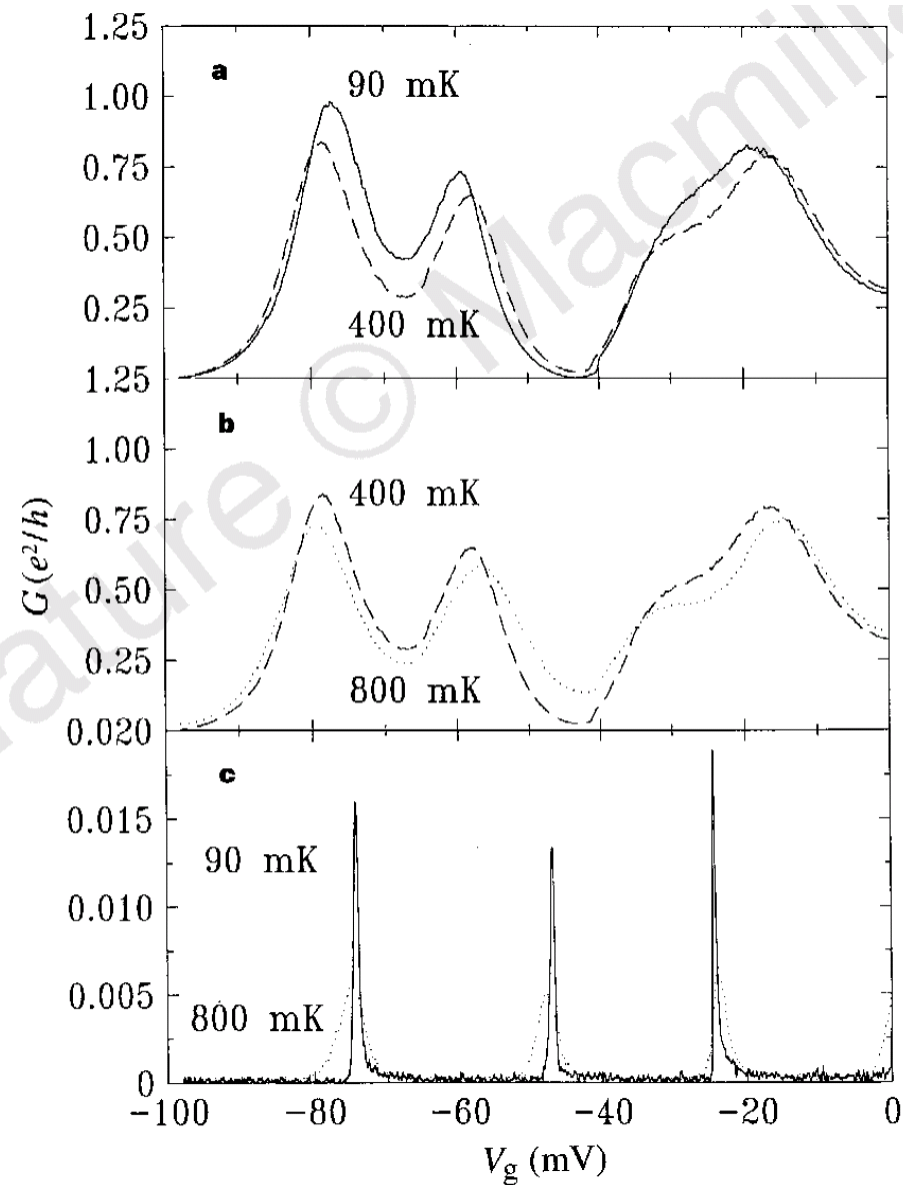
for  $T < T_K$  : DOS at  $\mu_{S,D}$  enhanced zero bias conductance!!

for  $T \gg T_K$  DOS peak suppressed

dots: parameters tunable  
SINGLE impurity

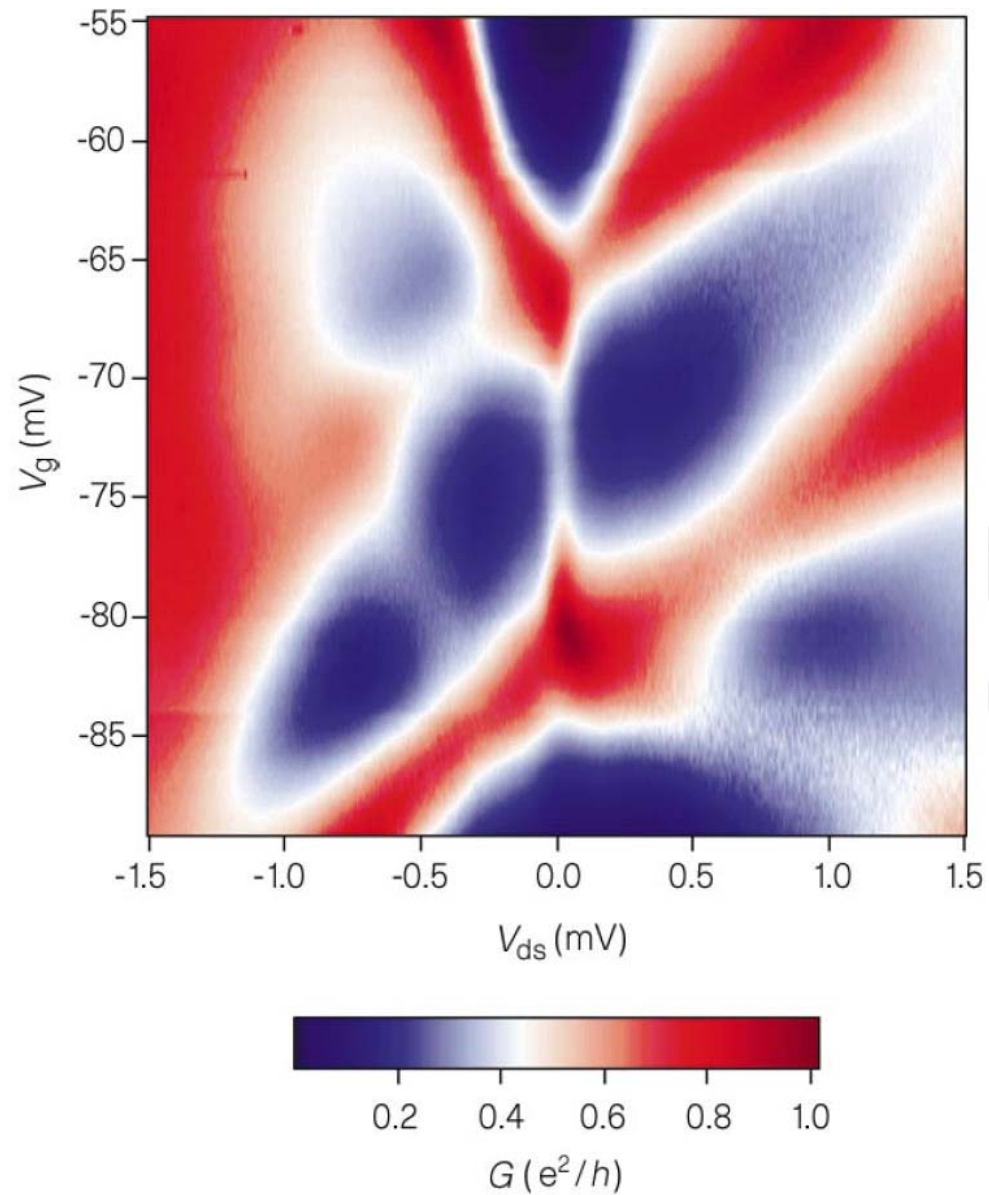


# Kondo Effect in Quantum Dots: Experiment



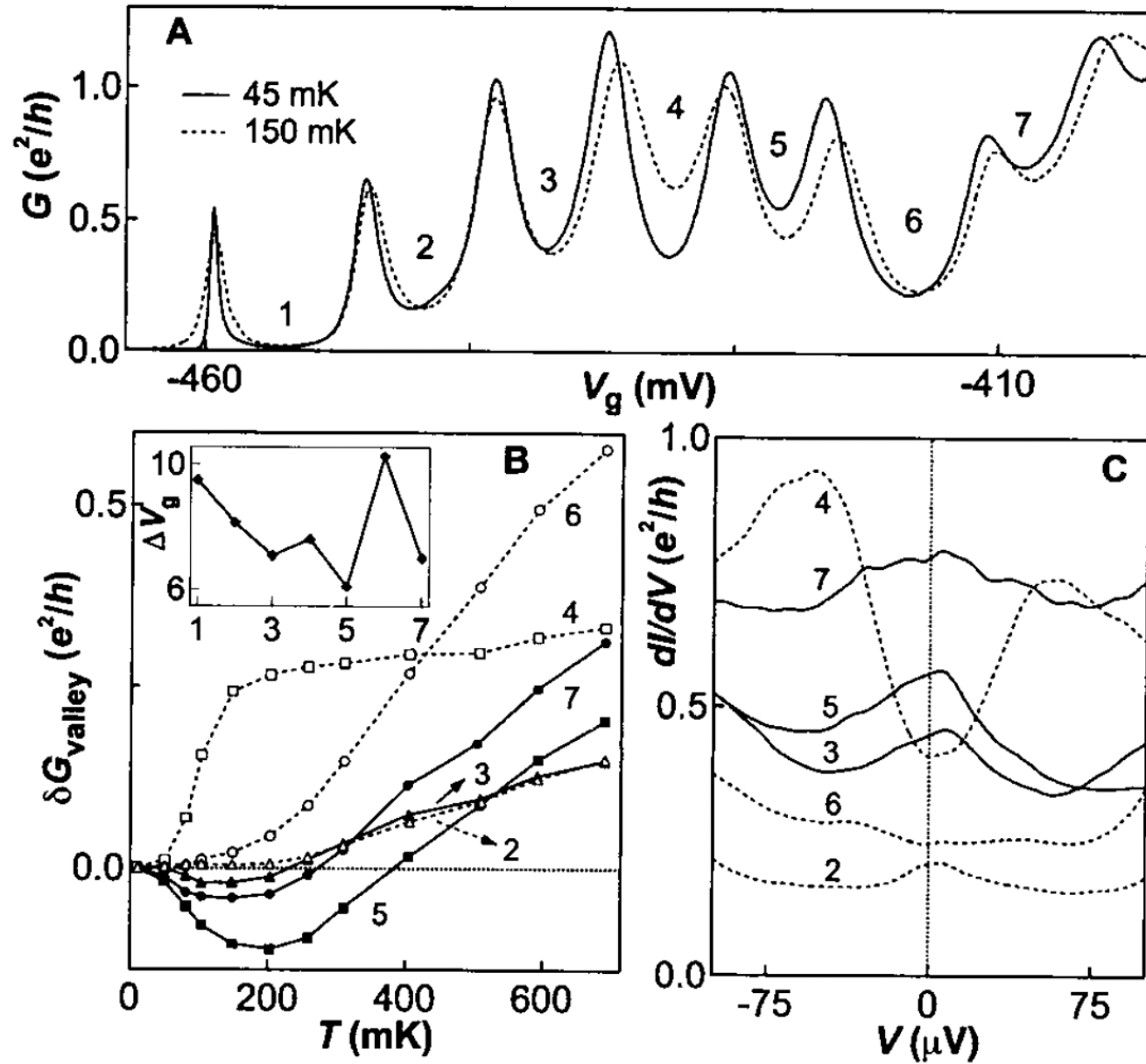
# Kondo Effect in Quantum Dots: Experiments

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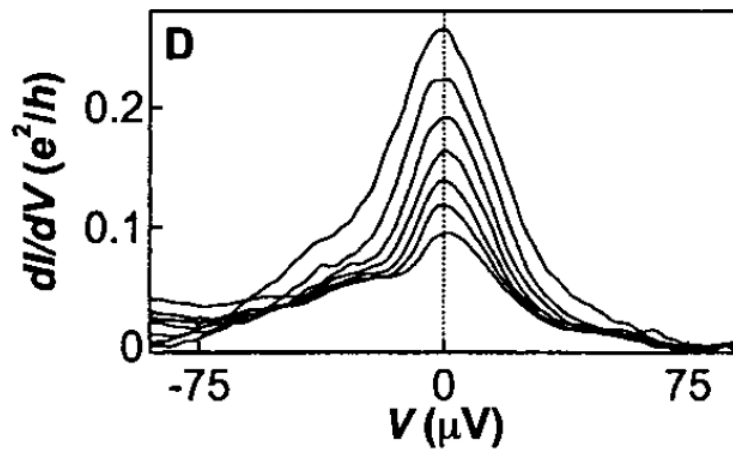
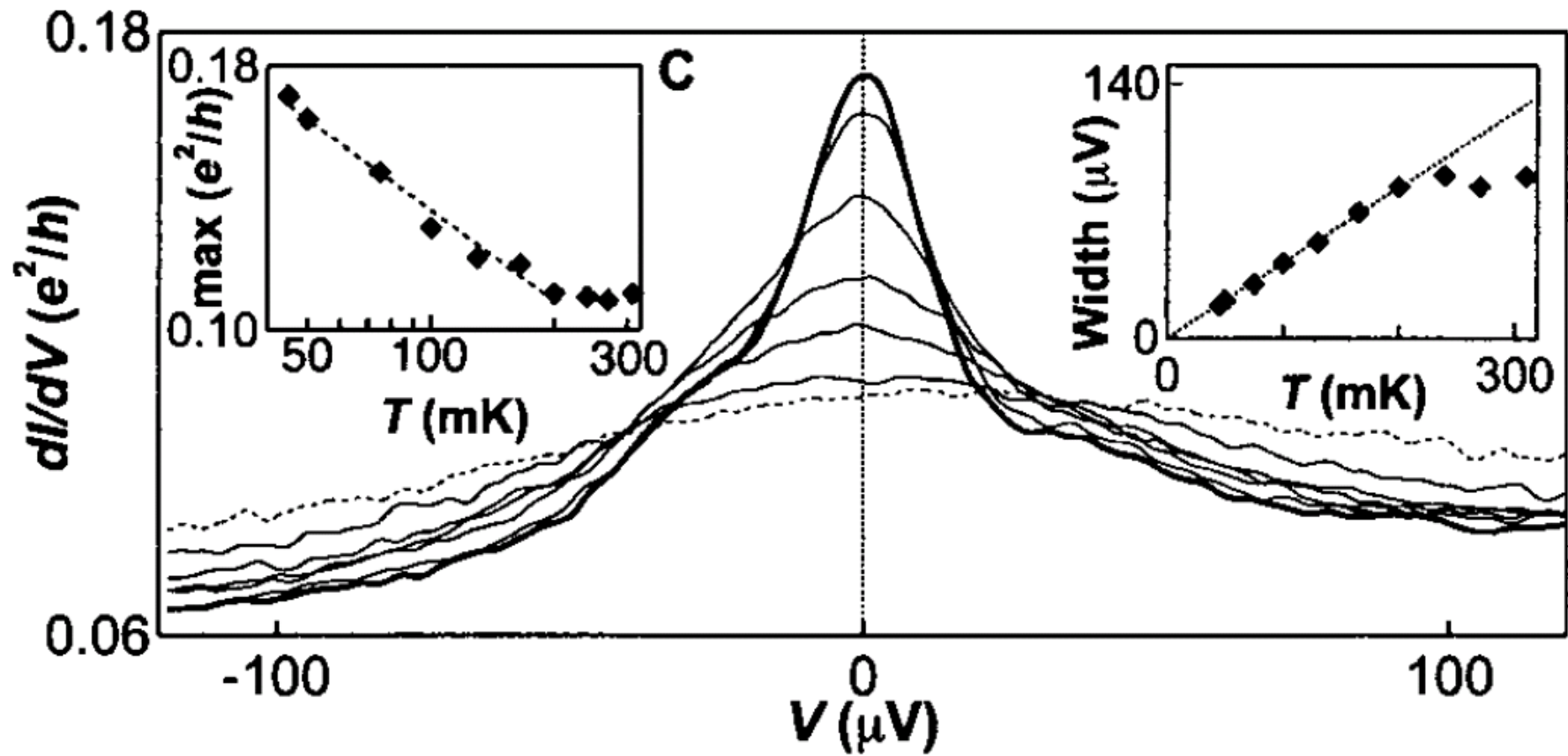


Goldhaber-Gordon et al., Nature **391**, 156 (1998)

# Kondo Effect in Quantum Dots: Experiments

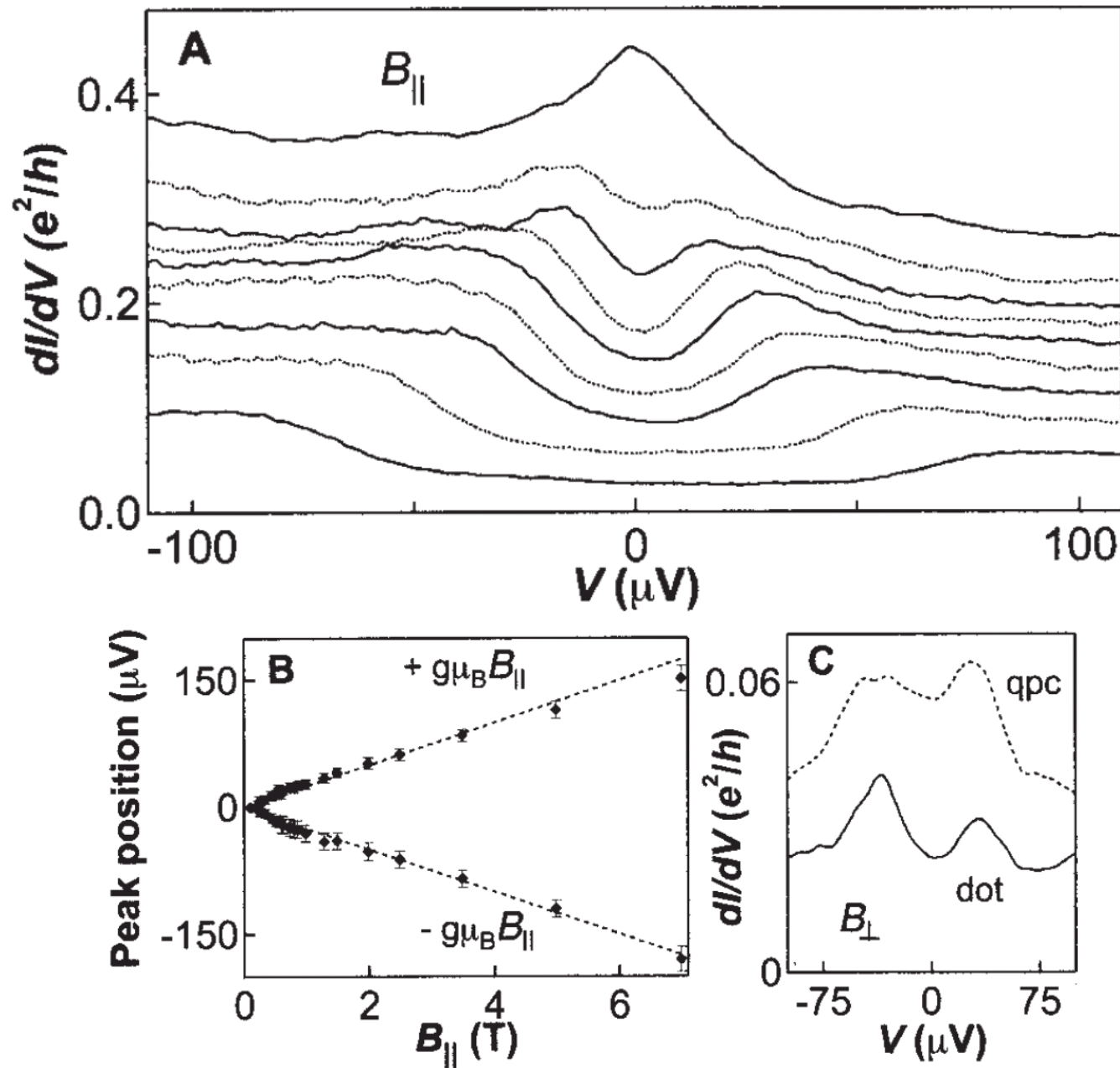


# Kondo Effect in Quantum Dots: Experiments



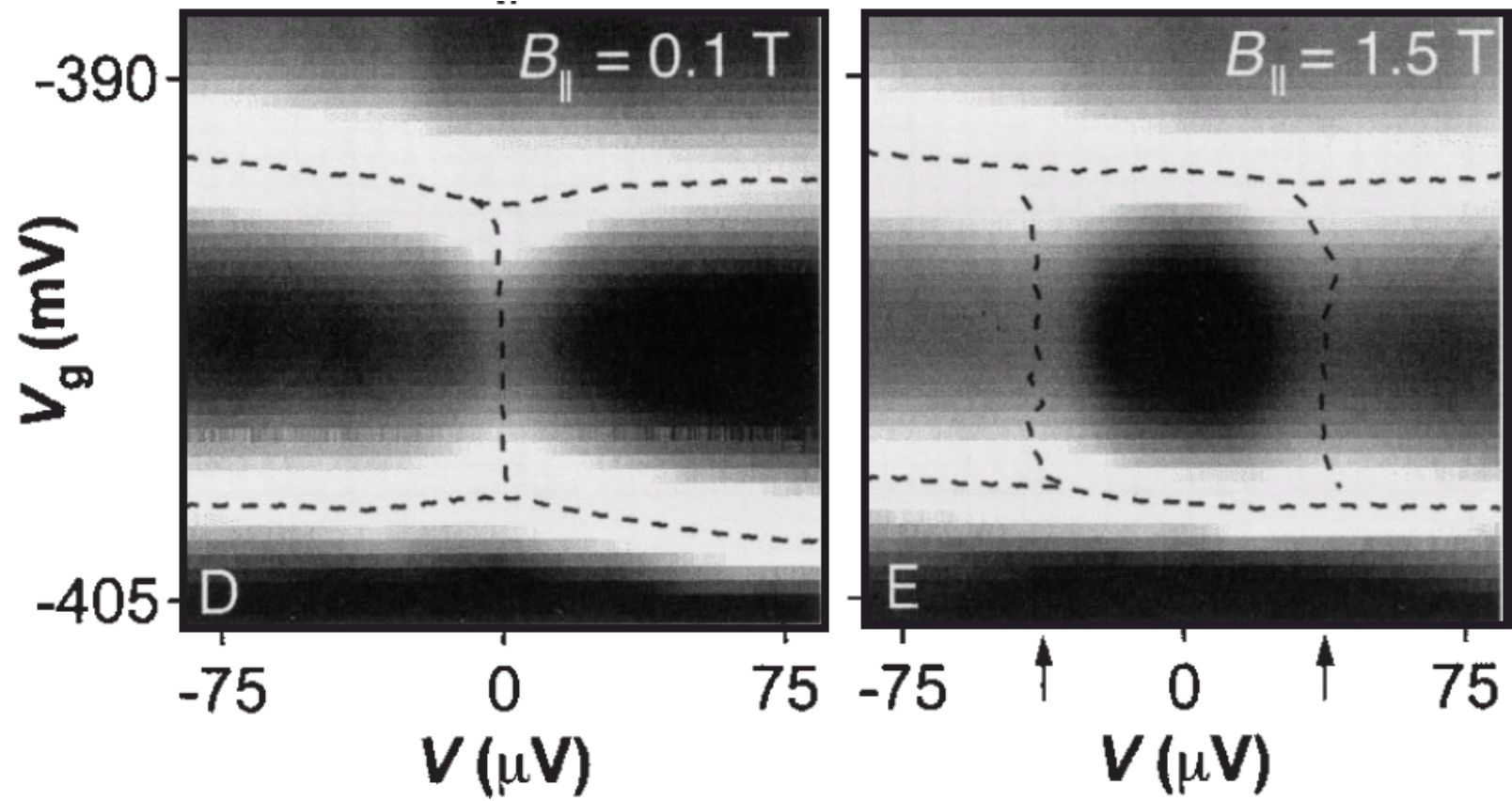
gate voltages  
into odd valley

# Kondo Effect in Quantum Dots: Experiments



# Kondo Effect in Quantum Dots: Experiments

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**3. Few Electron Dots**

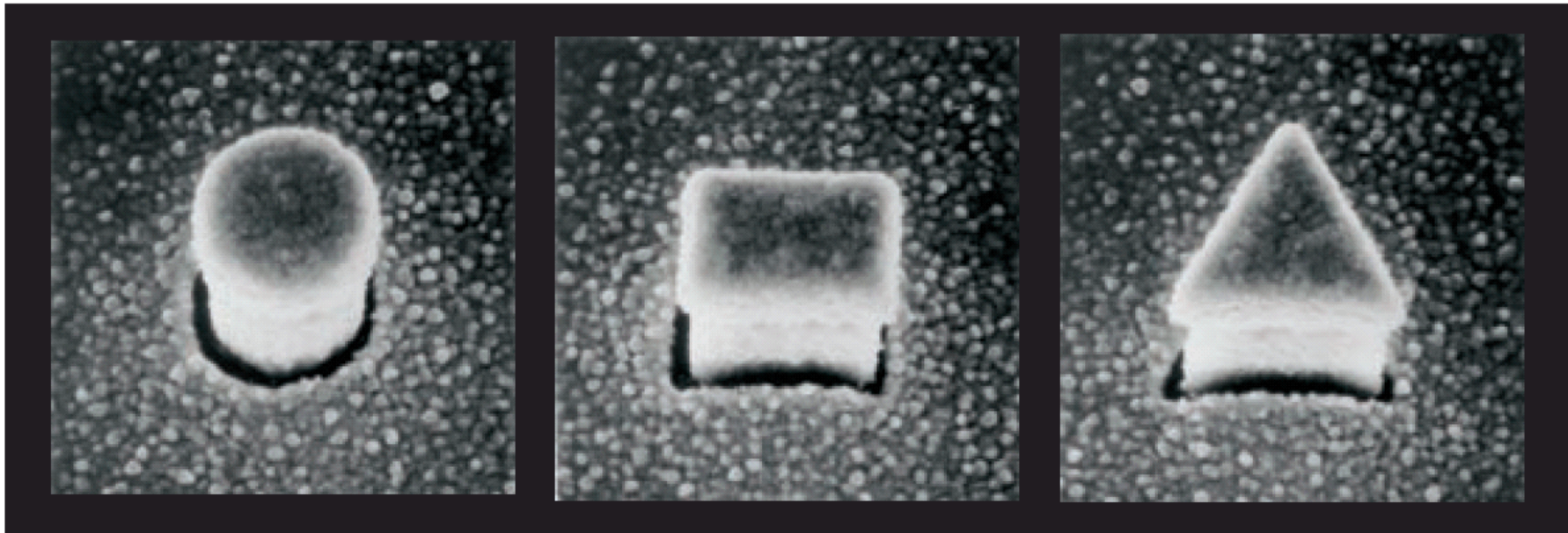
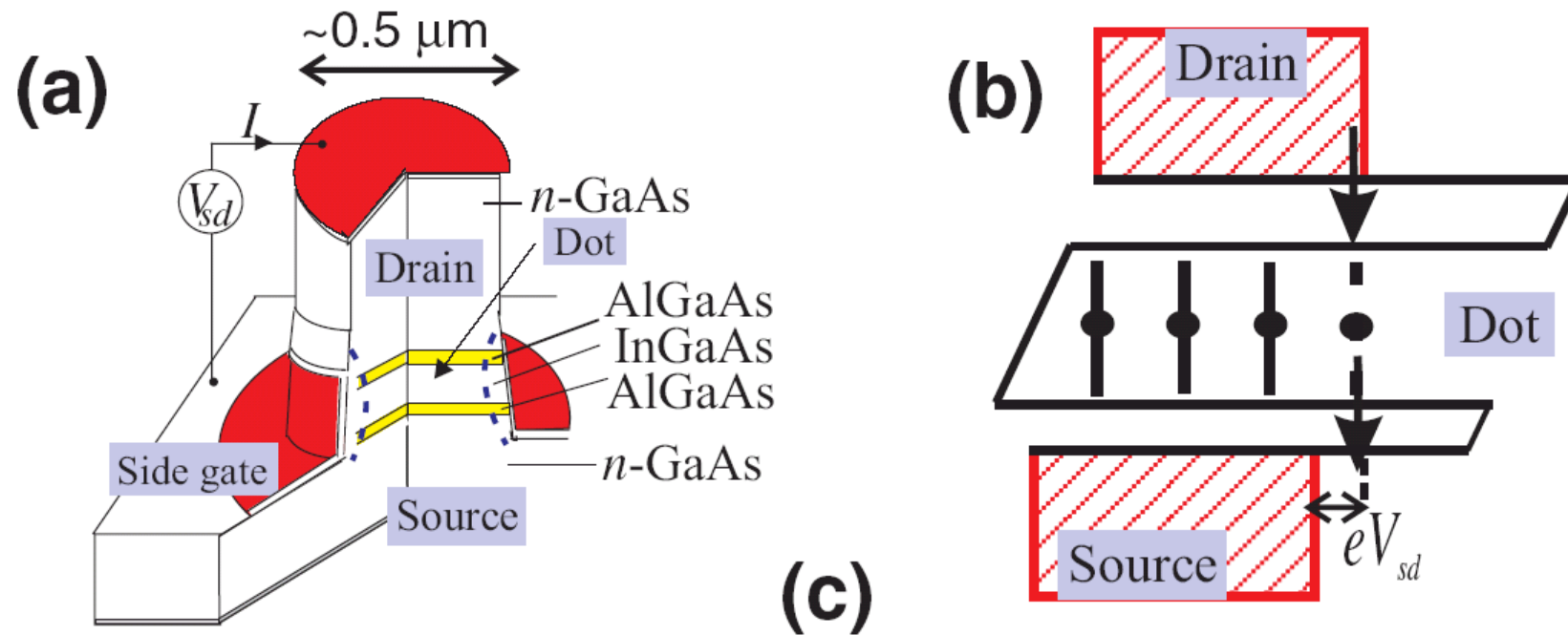
4. Double Quantum Dots

Kouwenhoven, Austing and Tarucha, RPP 64, 701 (2002)

Tarucha et al., PRL77, 3613 (1996)

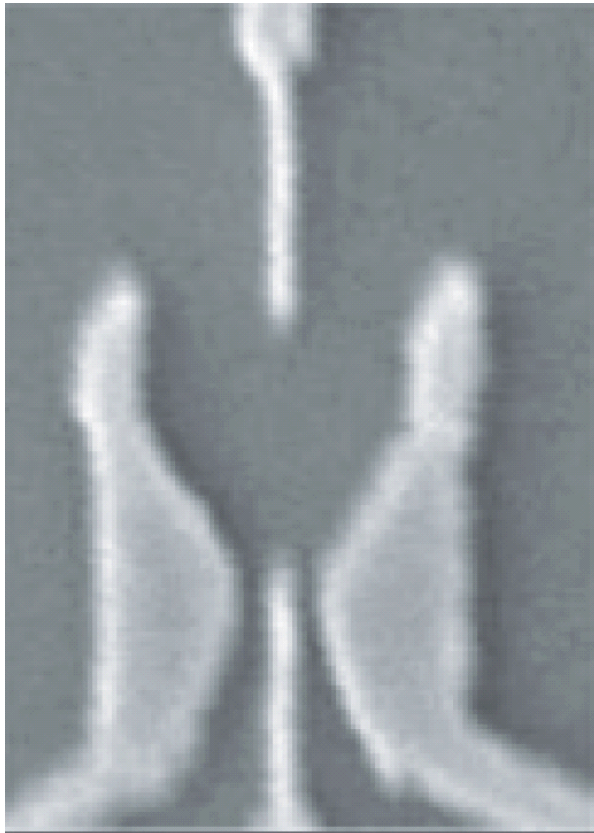
Kouwenhoven et al., Science 278, 1788 (1997)

# Few Electron Quantum Dots: Vertical

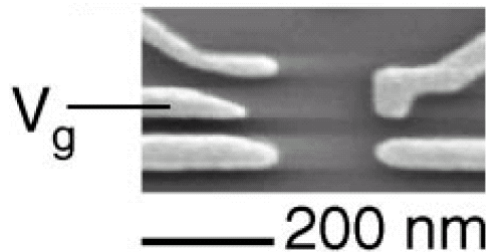




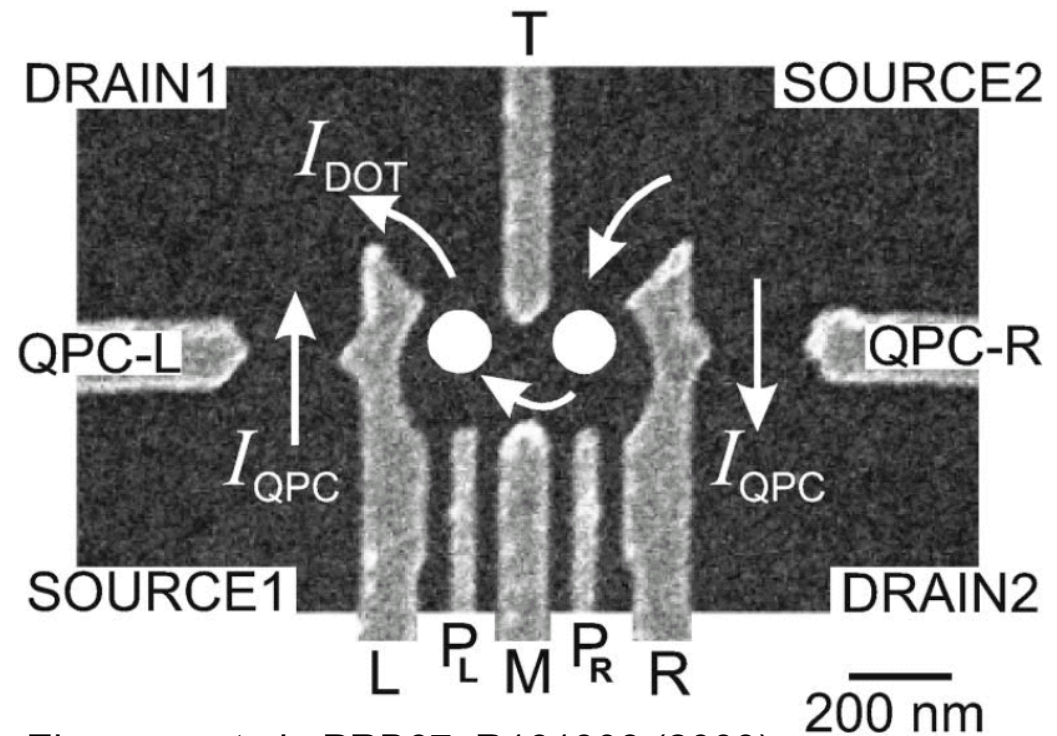
# Few Electron Quantum Dots: Lateral



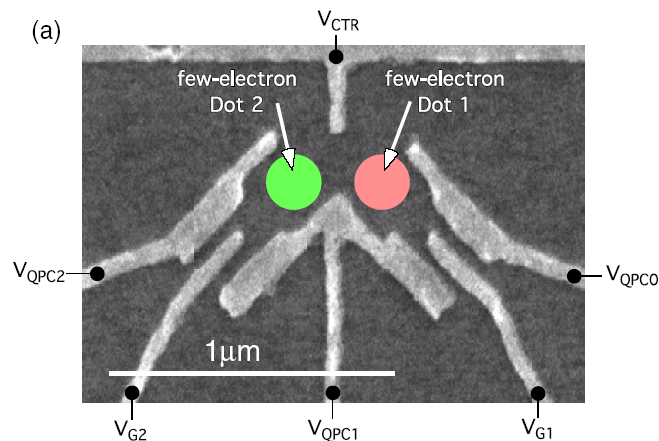
Ciorga et al., PRB61, R16315 (2000)



Zumbuhl et al., PRL93, 256801 (2004)



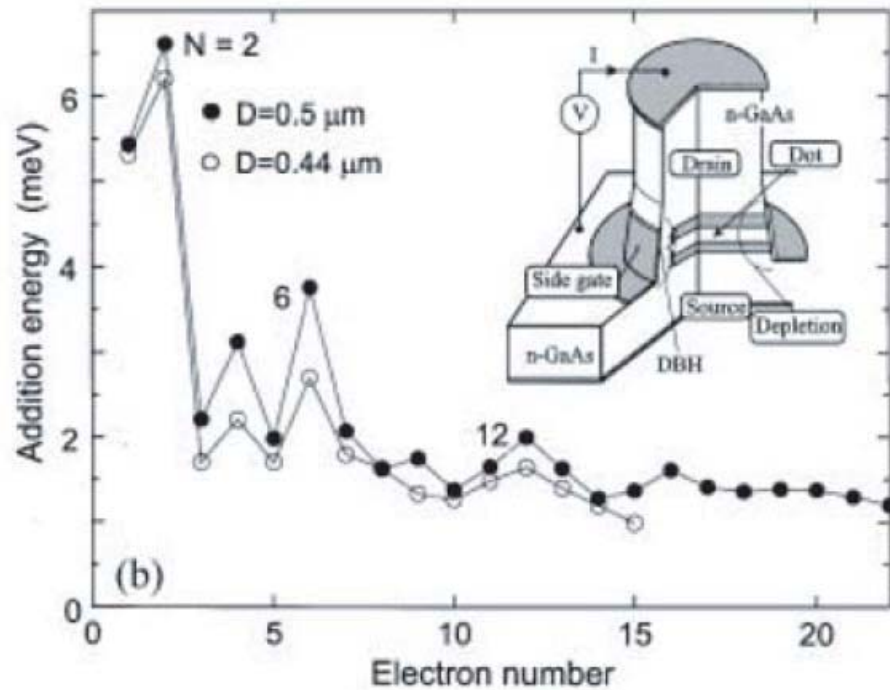
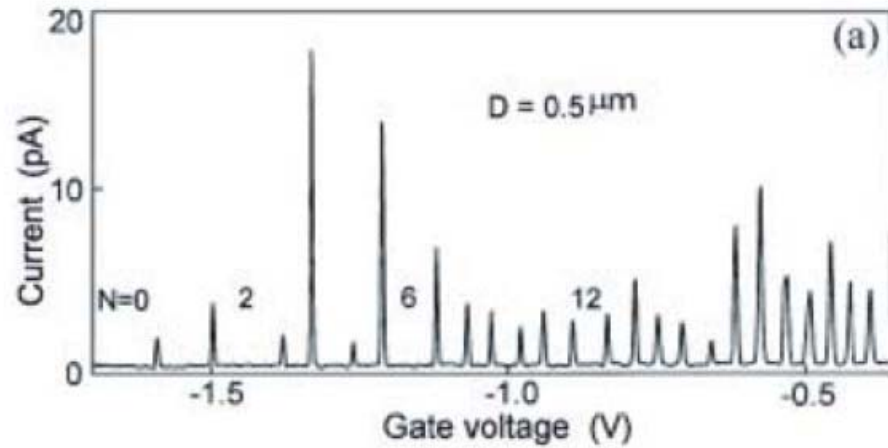
Elzerman et al., PRB67, R161308 (2003)  
Petta et al., PRL93, 186802 (2004)



Chan et al., Nanotech. 15, 609 (2004)

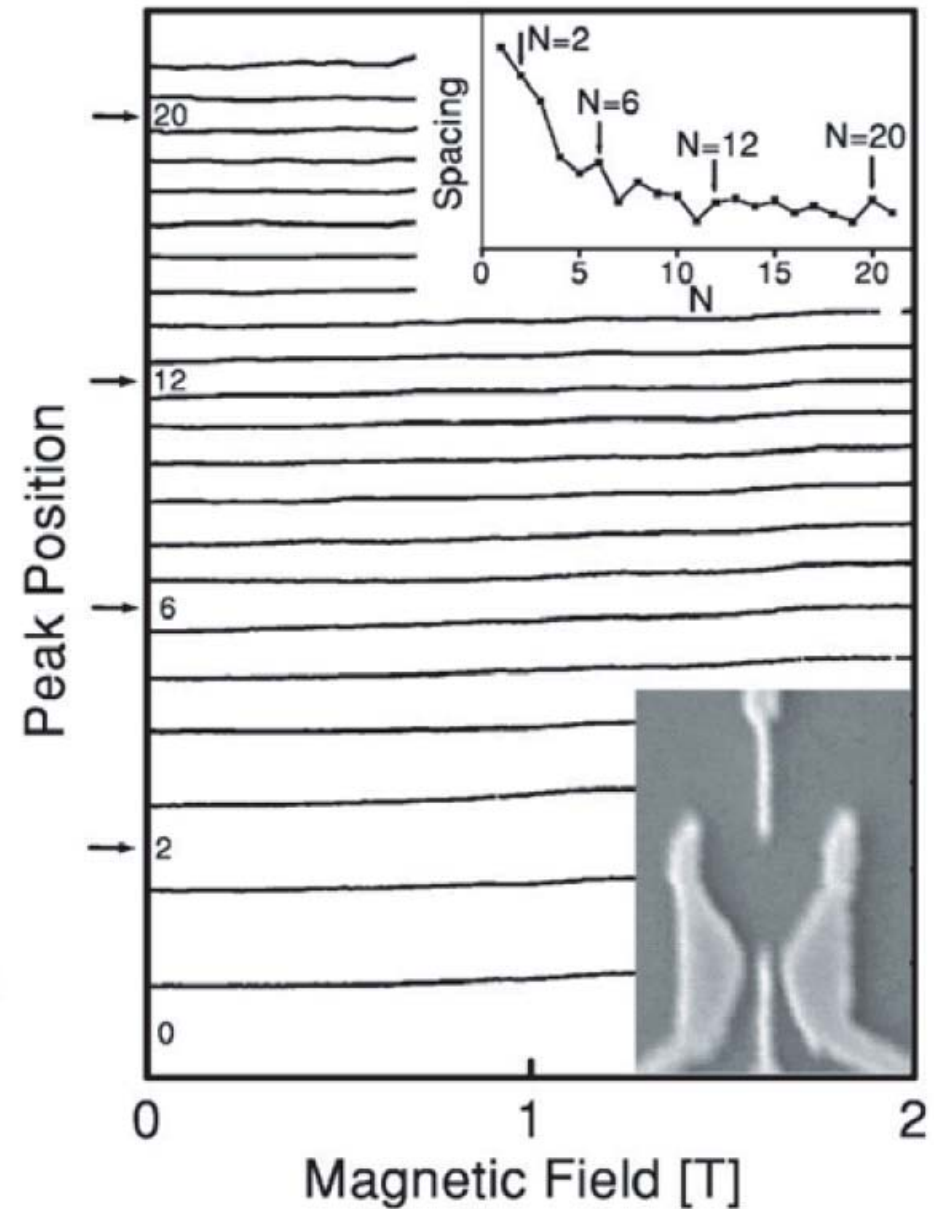
# Rotation Symmetry and Angular Momentum

circular symmetry: 2D shell filling



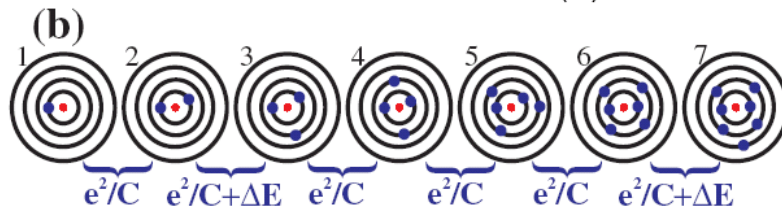
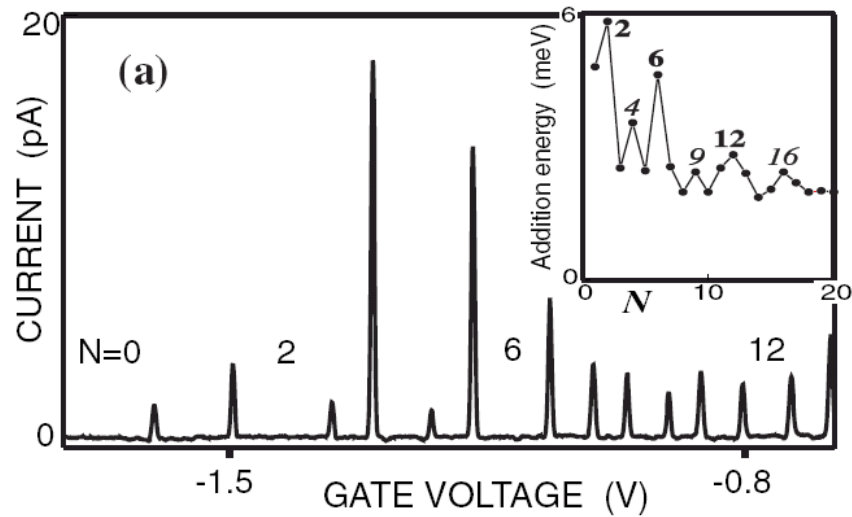
Tarucha et al., PRL77, 3613 (1996)

circular symmetry broken



Ciorga et al., PRB61, R16315 (2000)

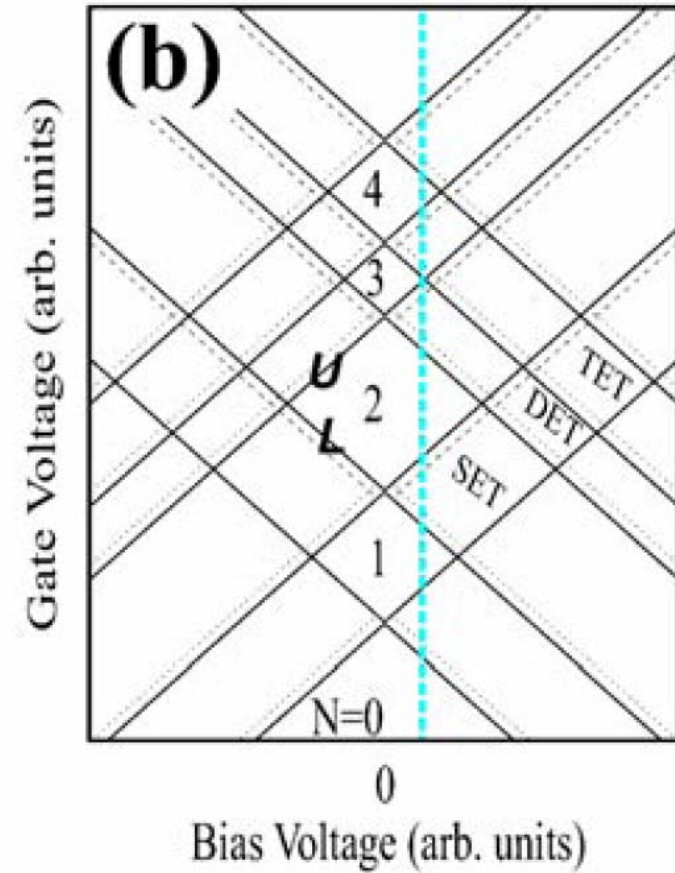
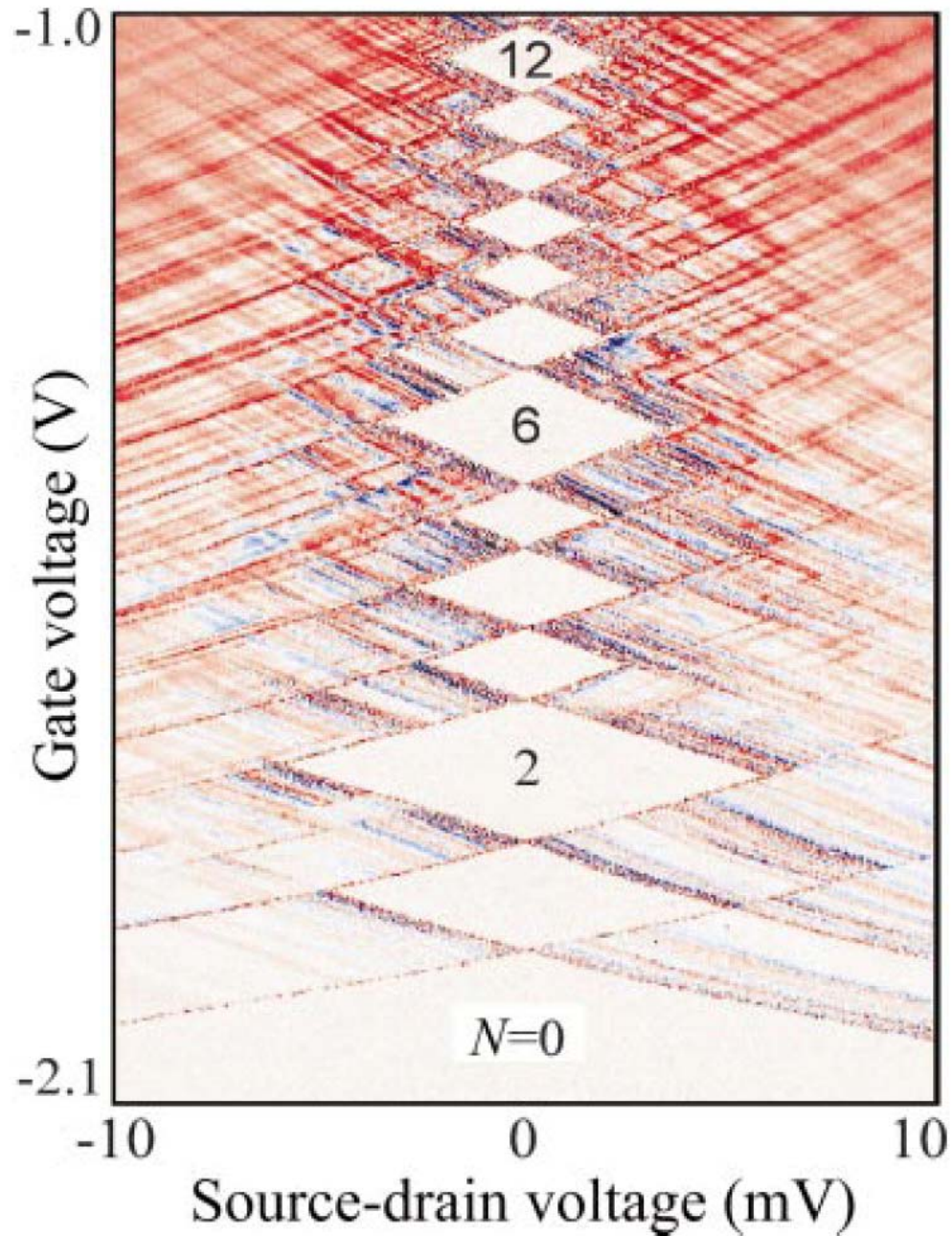
# 2D Periodic Table of Elements



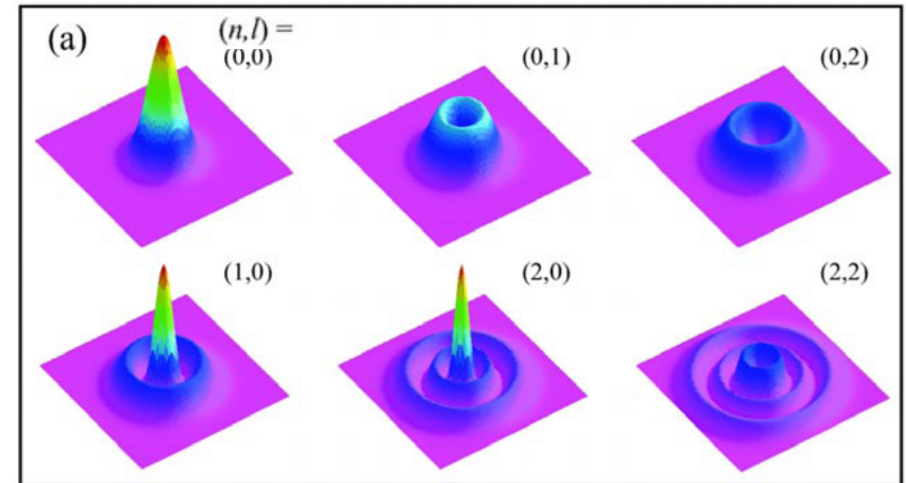
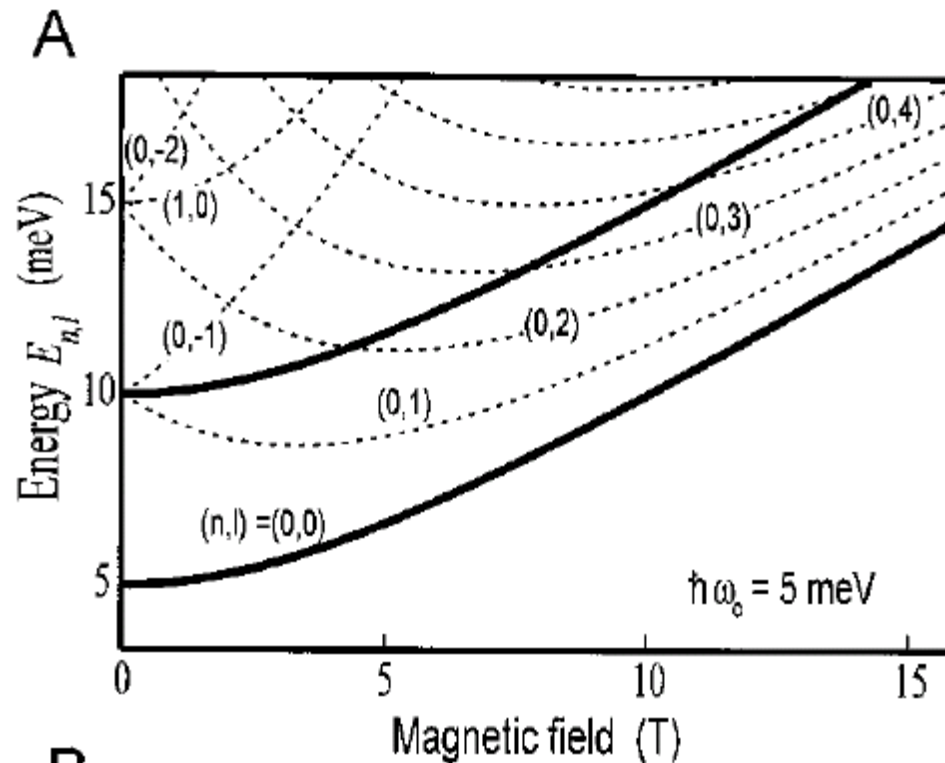
(c) **Periodic Table of 2D Artificial Atoms**

1 Ta						2 Ha
3 Et	4 Au				5 Ko	6 Oo
7 Sa	8 To	9 Ho		10 Mi	11 Cr	12 Ja
13	14	15	16 Wi	17 Fr	18 El	19
						20 Da

# Excitation Spectra of Circular, Few Electron Dots



# Fock-Darwin States: Single Particle Levels



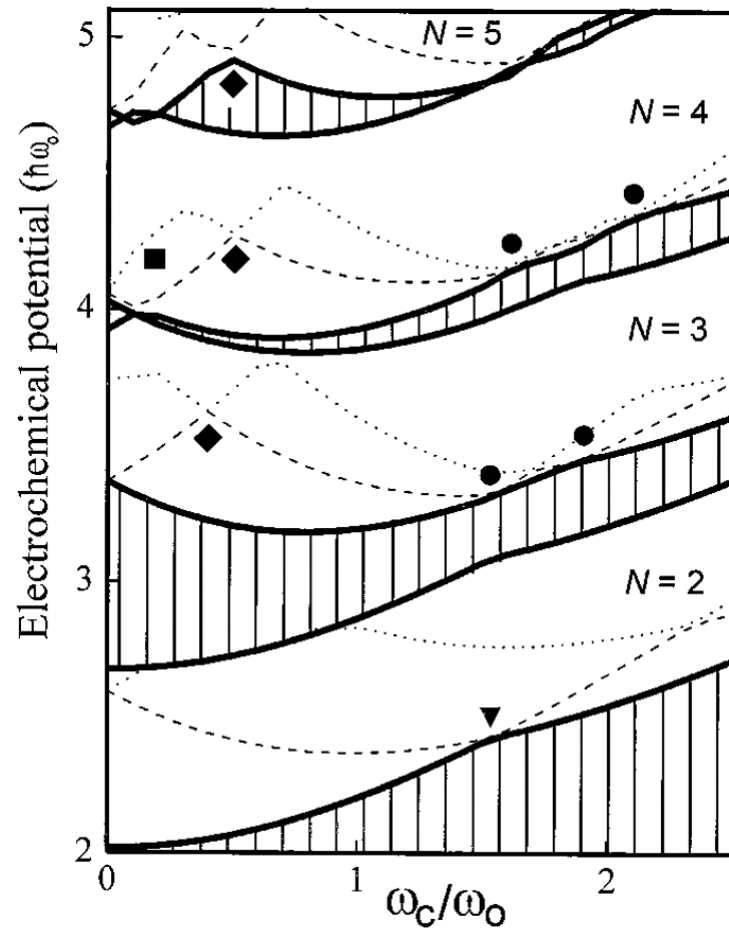
Fock-Darwin Energies

$$E_{n,\ell} = (2n + |\ell| + 1)\hbar \sqrt{\left(\frac{1}{4}\omega_c^2 + \omega_o^2\right)} - \frac{1}{2}\ell\hbar\omega_c$$

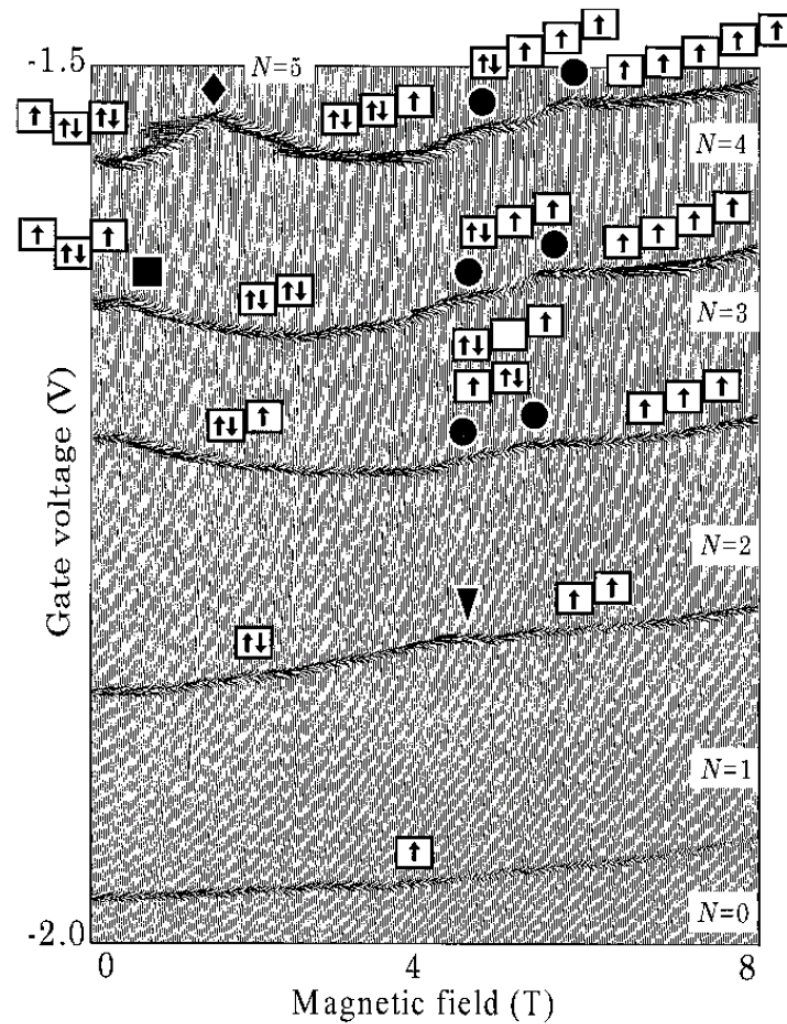
$n = 0, 1, 2, \dots$  radial

$l = 0, \pm 1, \pm 2, \dots$  angular momentum

# Magnetic Field Transitions



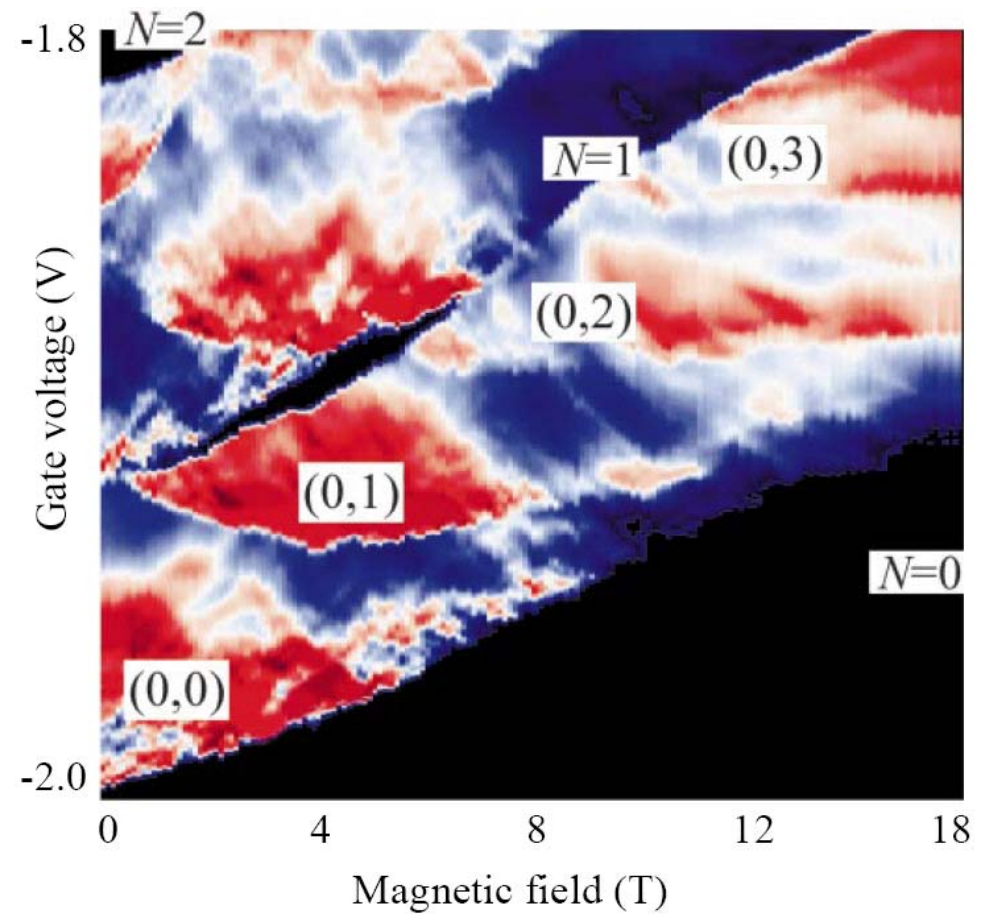
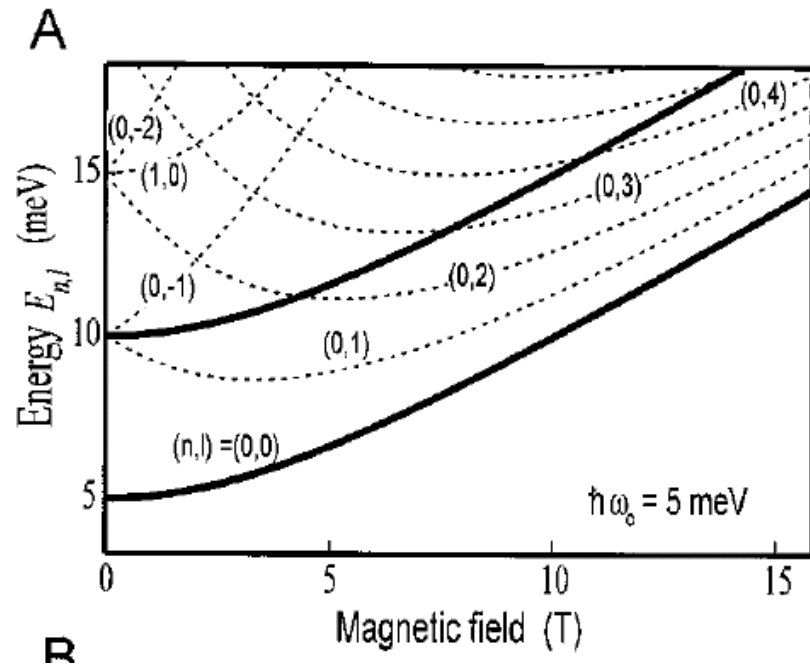
exact calculation



experiment

“atomic physics” like experiments not accessible in real atoms!!

# Zero to One Electron Transition



# Higher Transitions

