

# Electron transport in graphene

## **“Graphene = a material for future electronics”**

- Conductivity & mobility in graphene
  - Drude picture
  - Effective mass
  - Boltzmann equation
- Scattering mechanisms, limitation of mobility
- Present status
- Applications

## **“Dirac physics” in transport**

- Klein tunneling
- Reflectionless transmission in p-n junction
- Evidence of Klein backscattering in interference pattern of n-p-n junction

## **Graphene based Hybrid Quantum Devices**

Sources:

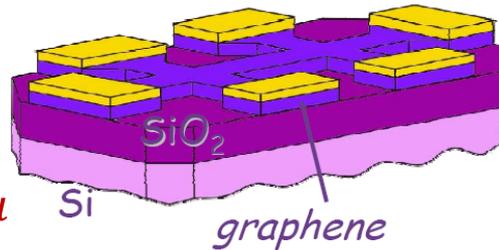
[http://www.tntconf.org/2010/Presentaciones/TNT2010\\_Geim.pdf](http://www.tntconf.org/2010/Presentaciones/TNT2010_Geim.pdf)

# Conductivity, mobility

## Mobility, conductance:

$$v_d \equiv \mu E$$

$$j \equiv env_d = en\mu E = \sigma E \quad \sigma = en\mu$$



$$E(\vec{k}) = \hbar v_F |\vec{k}|$$

## Conductivity (Drude model):

$$\sigma = \frac{e^2 n \tau}{m} \quad \mu = \frac{e \tau}{m}$$

How to separate mobility ( $\mu$ ) and e density ( $n$ )?

Measure:  $\rho$  + Hall resistance  $R_H = \frac{E_y}{j_x B} = -\frac{1}{ne}$

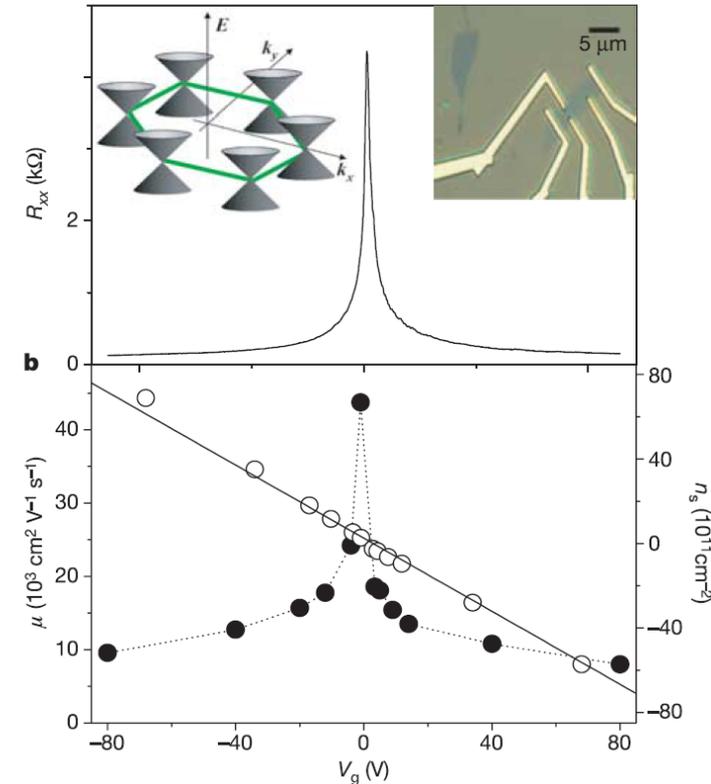
Effect of gate voltage,  $V_g$ ?

$$N = C_g V_g / e \quad \rightarrow \quad n \sim V_g \quad \rightarrow \quad k_F \sim \sqrt{V_g}$$

## Measurement:

- At  $V_g$  zero,  $R_H$  (and  $n$ ) changes sign  $\rightarrow$  boarder between e and h bands
- mobility largest at Dirac point ( $V_g = 0$ ).

## R vs. $V_g$ Transport characteristics



# Conductivity, mobility

How to calculate conductivity?

$$\sigma = e^2 \tau \frac{n}{m}$$

What is  $m$ , effective mass?

$$\frac{1}{m} = \frac{1}{m_{xx}} = \frac{1}{\hbar^2} \frac{\partial^2 E}{\partial k_x^2}$$

For quadratic dispersion:  $E = \frac{\hbar^2 k^2}{2m_{eff}}$ ,  $m = m_{eff}$

For Dirac electrons, where  $E(\vec{k}) = \hbar v_F |\vec{k}|$  ?

Naively  $1/m = 0$ , but NOT.

To calculate  $1/m$ :

$$\frac{\partial^2 |k|}{\partial k_x^2} = \dots = \frac{k_y^2}{|k|^3} \rightarrow \frac{1}{m_{xx}} = \frac{1}{\hbar} v_F \frac{k_y^2}{k^3}$$

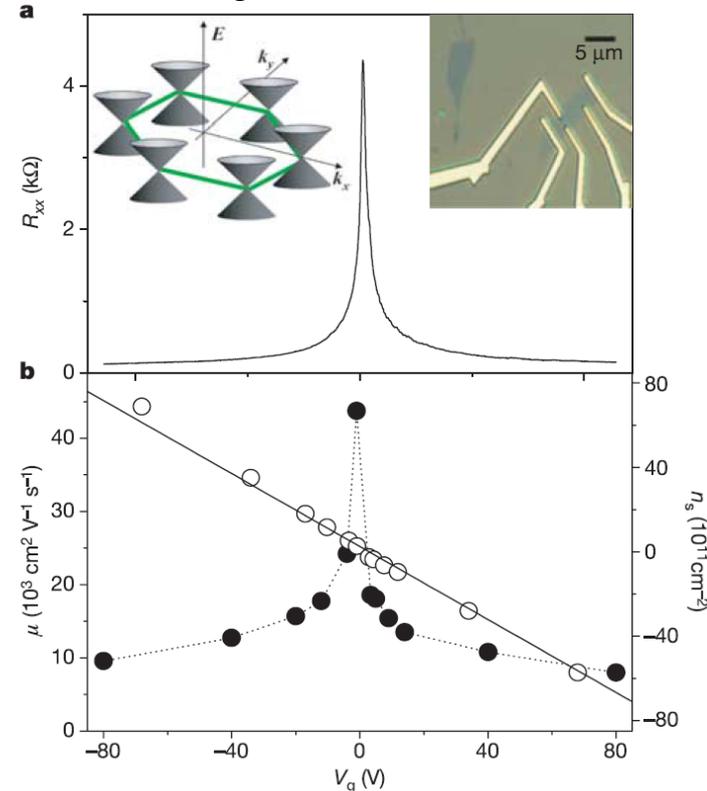
$$\frac{\partial |k|}{\partial k_x} = \frac{1}{2} \frac{2k_x}{|k|}$$

→ Effective mass depends on  $k$

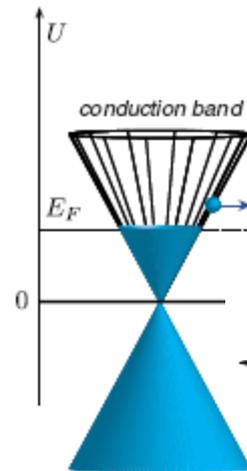
One has to average  $1/m$  for all filled states:

$$E(\vec{k}) = \hbar v_F |\vec{k}|$$

**R vs.  $V_g$  Transport characteristics**



Nature 438, 201 (2005)



# Conductivity, mobility

Accurate calculation of  $\sigma$ ?

From Boltzmann equation (see Solyom 24.3.39.):

$$\sigma = e^2 \tau \frac{n}{m} = e^2 \tau \cdot 2 \cdot 2 \cdot \int_{\text{filled } k \text{ states}} \frac{d^2 k}{(2\pi)^2} \frac{1}{m_{xx}}$$

**HOMEWORK** Calculate  $\sigma$

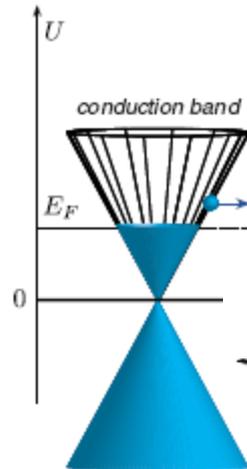
Result: 
$$\sigma = e^2 \tau \frac{v_F}{\hbar \pi} k_F$$

with relaxation length  $l \equiv v_F \tau$

$$\sigma = \frac{2e^2}{h} l k_F$$

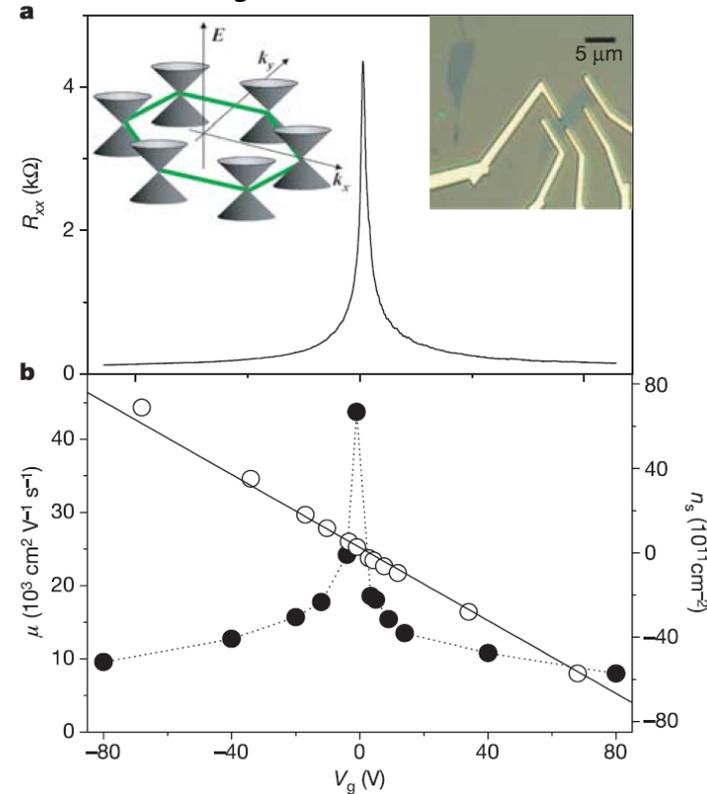
Mean free path: 
$$l = \frac{\hbar}{e} \mu \sqrt{n\pi}$$

E.g. for mobility = 600.000,  $l$  is only  $\approx 3\mu\text{m}$



$$E(\vec{k}) = \hbar v_F |\vec{k}|$$

**R vs.  $V_g$  Transport characteristics**



Nature 438, 201 (2005)

# Scattering mechanisms in graphene

## What limits the mobility at room T?

Source of  $1/\tau$  ?

### Scattering mechanisms resulting resistivity:

- potential scattering: impurities, defects, vacancies
- Electron – phonon scattering
- Etc.

Usual terms: (see Solyom II.)

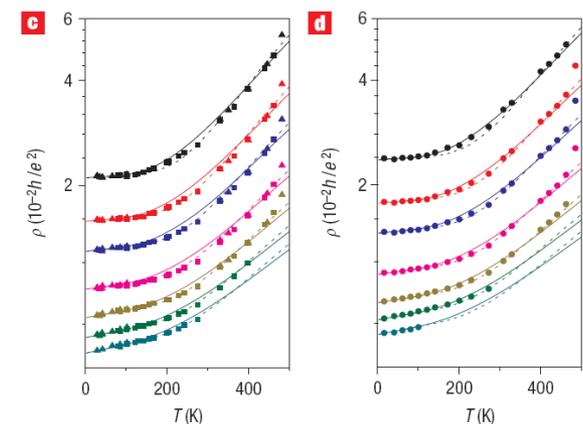
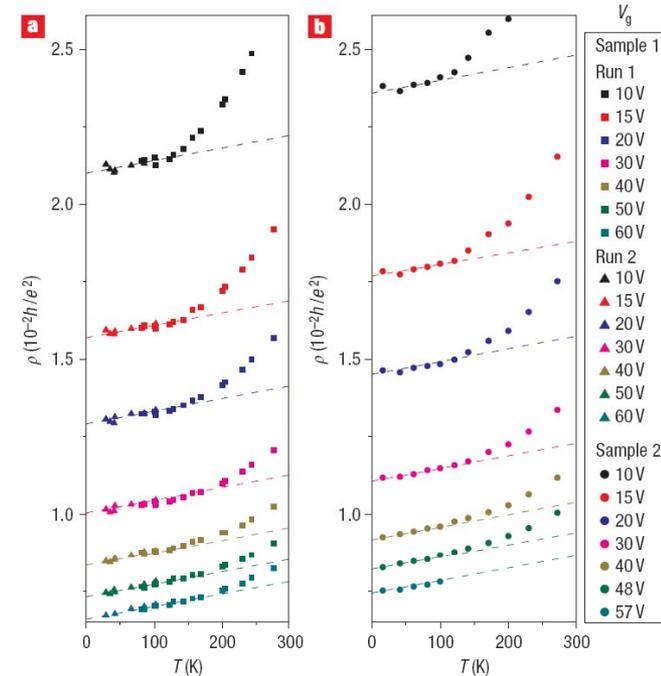
- Residual resistivity ( $\rho_0$ ): T independent
- Longitudinal acoustic phonons ( $\rho_A$ ): linear in T

$$\rho(V_g, T) = \rho_0(V_g) + \rho_A(T); \quad \rho_A(T) = \left(\frac{h}{e^2}\right) \frac{\pi^2 D_A^2 k_B T}{2h^2 \rho_s v_s^2 v_F^2}$$

### Measurements (see Fig. a,b)

- At higher T, strong deviation from linear T dependence
- Dependence also on  $V_g$
- It suggests scattering on high energy phonon modes

Sample on  $\text{SiO}_2$  substrate, UHV  
 $T=[20\text{K}-500\text{K}]$ , 4-point



# Scattering mechanisms in graphene

$$\rho(V_g, T) = \rho_0(V_g) + \rho_A(T) + \rho_B(V_g, T);$$

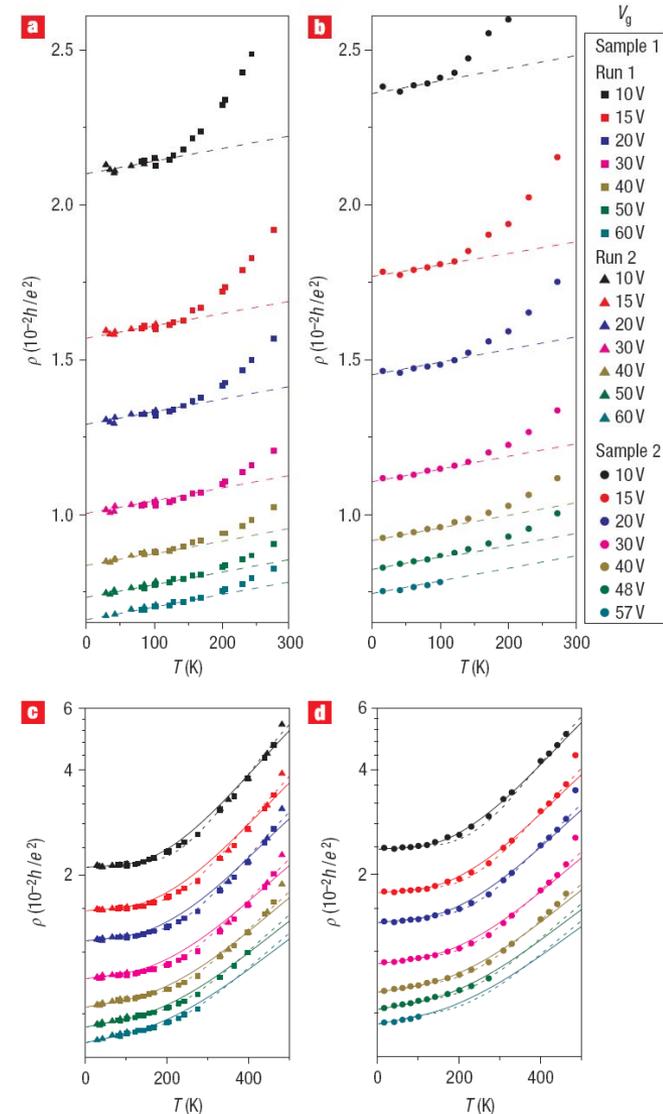
$$\rho_B(V_g, T) = B_1 V_g^{-\alpha_1} \left( \frac{1}{e^{(59\text{meV})/k_B T} - 1} + \frac{6.5}{e^{(155\text{meV})/k_B T} - 1} \right)$$

$\rho_B$ : additional term to fit the measurements (see Fig. c,d)  
 Bose-Einstein distribution  $\sim$  population of high energy phonon modes, e.g. optical phonons  
 Very good fit of the measured curves with  $\alpha_1=1.04$

Optical phonons of graphene?

- Strong  $V_g$  dependence is not expected
- Mainly out of plane phonons at this energy. It is not expected to give strong contribution

**Interfacial phonon scattering:** Surface optical phonon modes in  $\text{SiO}_2$  couples to e-s in graphene  
 The expected phonon energies and coupling strength (1:6.5) are inserted into  $\rho_B$   
 Strong  $V_g$  dependence also expected



# Scattering mechanisms in graphene

## What limits the mobility at room T?

Different T dependence of  $\rho_0$ ,  $\rho_A$ ,  $\rho_B$  allows to separate the three contributions. ( $\rho_B = \rho - \rho_0 - \rho_A$ )

Fig. a

$\rho_A$   $V_g$  independent

$\rho_B \sim V_g^{i-1.04}$  relation confirmed

→ **Residual resistivity dominates**

Fig. b

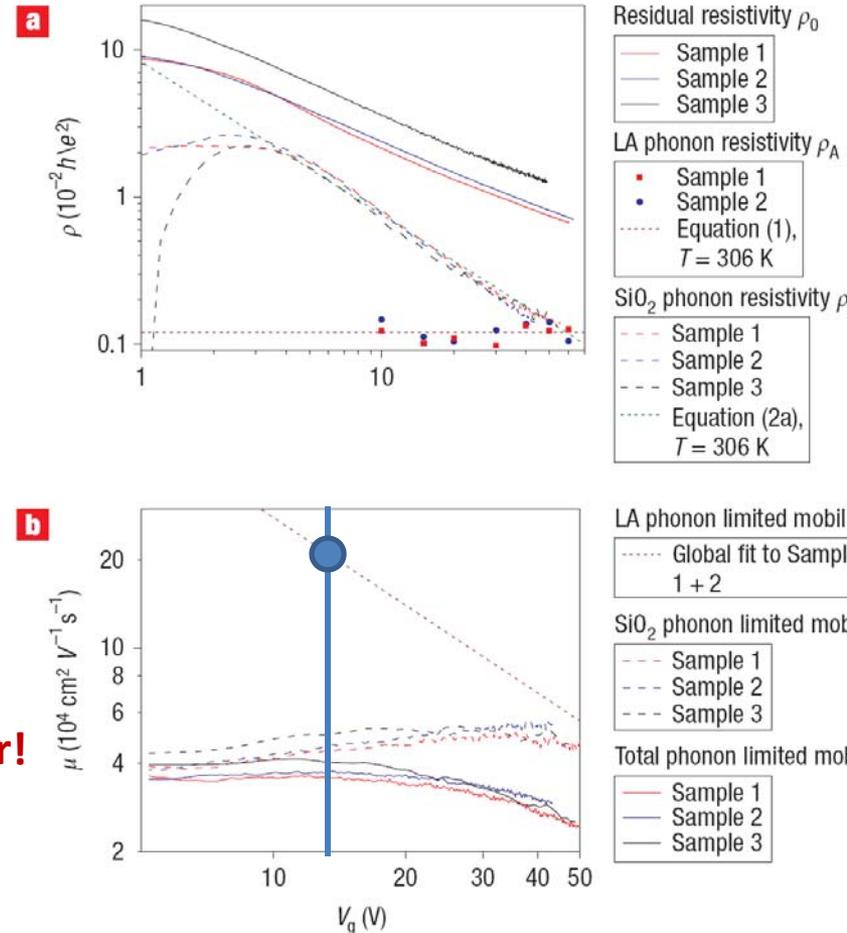
Derive the mobility related to two e-p processes :

$\mu = 1/nep = 1/c_g V_g e \rho$

→ SiO<sub>2</sub> contribution (c) dominates

→ **The intrinsic, LA phonon scattering mobility** at  $n = 10^{12} \text{cm}^{-2}$  (technologically relevant) :  $\mu \approx 200\,000$  (see blue dot) **Higher than any known semiconductor!** (E.g. InSb  $\approx 77\,000$  and carbon nanotubes  $\approx 100\,000$ ).

## Contributions at Room T



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Fig. c

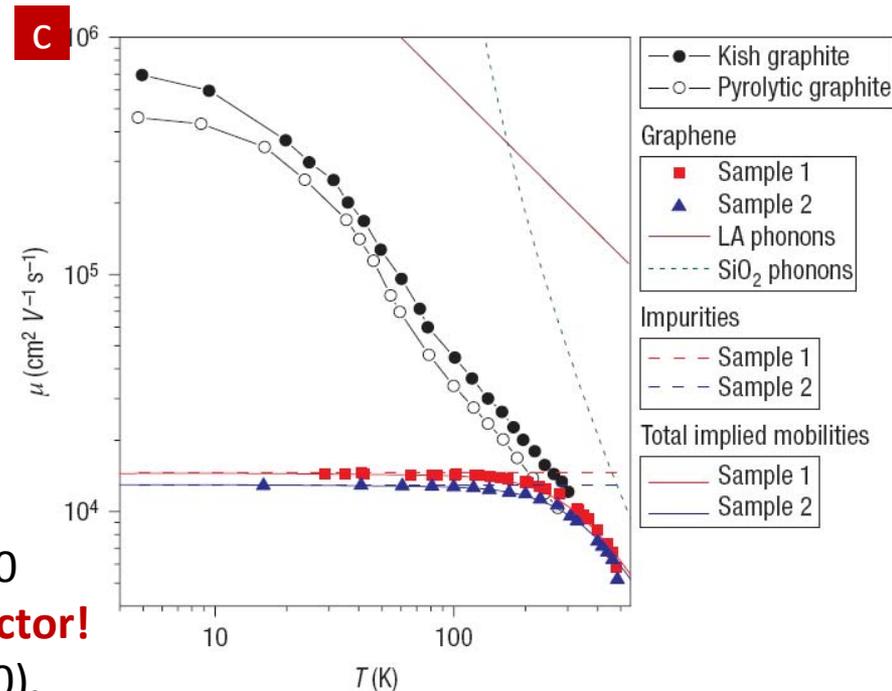
Comparison with graphites, sources of exfoliated graphene

Mobility is much smaller than for graphites. It is impurity dominated.

→ Residual res. not due to point defects

but due to **charge impurities in SiO<sub>2</sub> substrate**

## T dependence



# Scattering mechanisms in graphene

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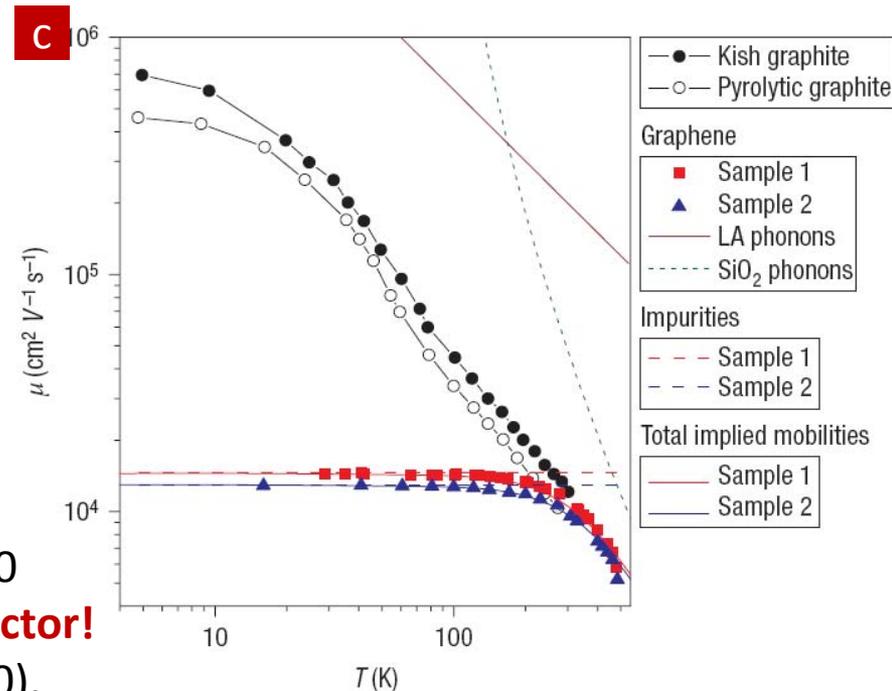
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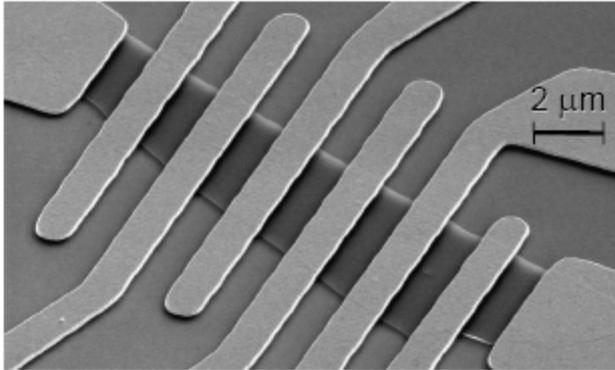
## T dependence



**$\Sigma$ : Problem, SiO<sub>2</sub> is bad substrate**

# Suspended flakes

SiO is etched by BHF

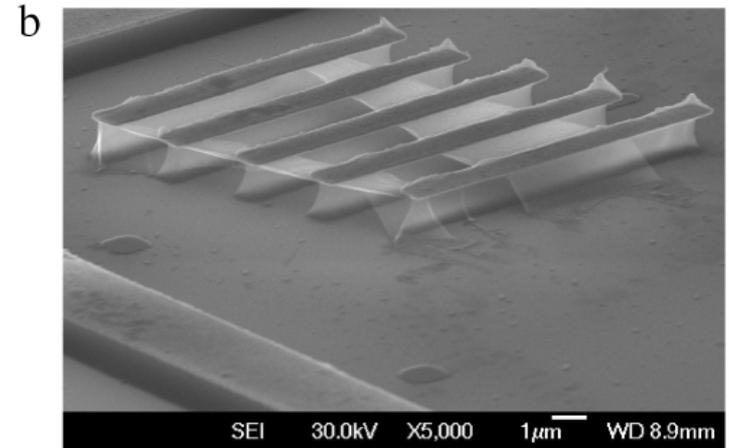


mobility up to  $200,000 \text{ cm}^2/\text{V}\cdot\text{s}$   
Mean free path  $L \sim \mu\text{m}$

low-T mobilities  
few million  $\text{cm}^2/\text{V}\cdot\text{s}$   
Manchester, arxiv 2010

→ Demonstration of FQHE

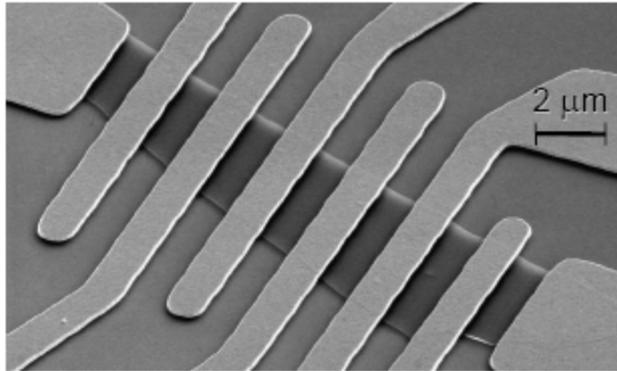
PMGI based organic polymer  
N. Tombros [arXiv:1009.4213](https://arxiv.org/abs/1009.4213)



Possible with any metals ->  
spin physics, superconductivity  
 $600,000 \text{ cm}^2/\text{Vs}$  at  $n = 5.0 \text{ E}9 \text{ cm}^{-2}$  at 77K.  
 $L \sim 3 \mu\text{m}$

# Suspended flakes

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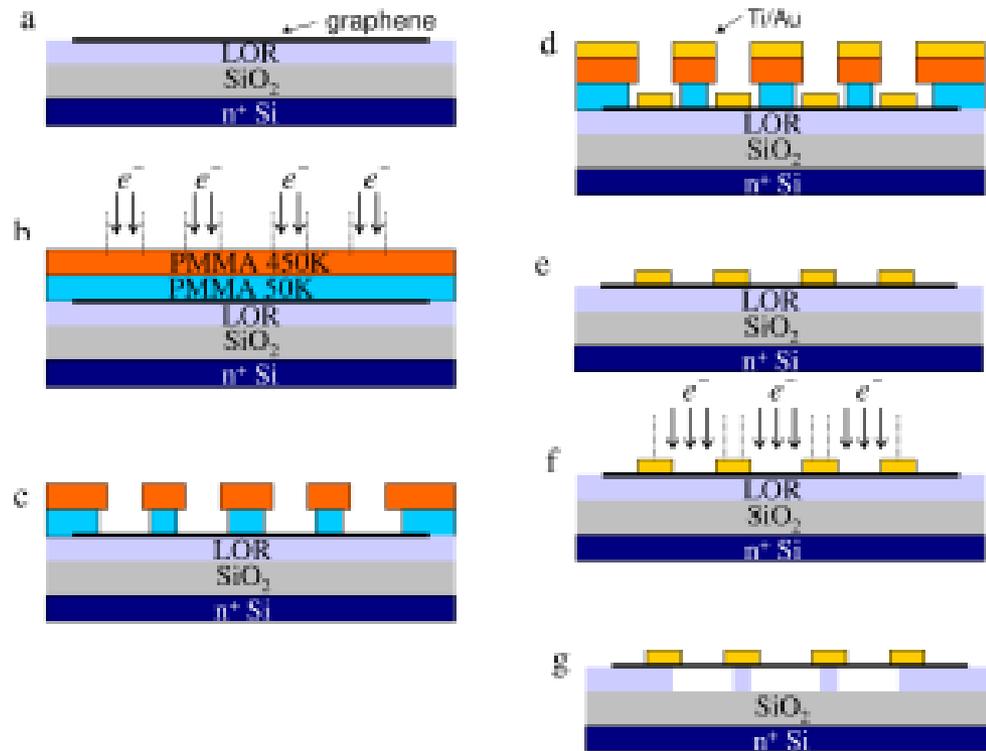


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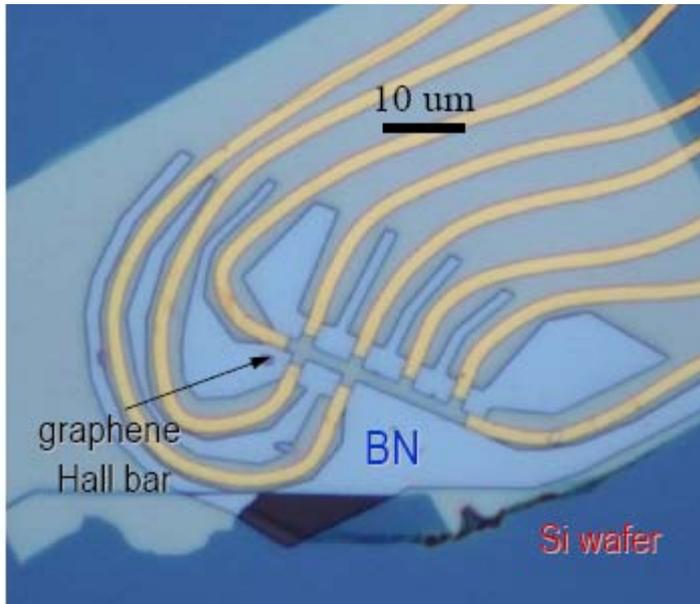
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→ Demonstration of FQHE

PMGI based organic polymer  
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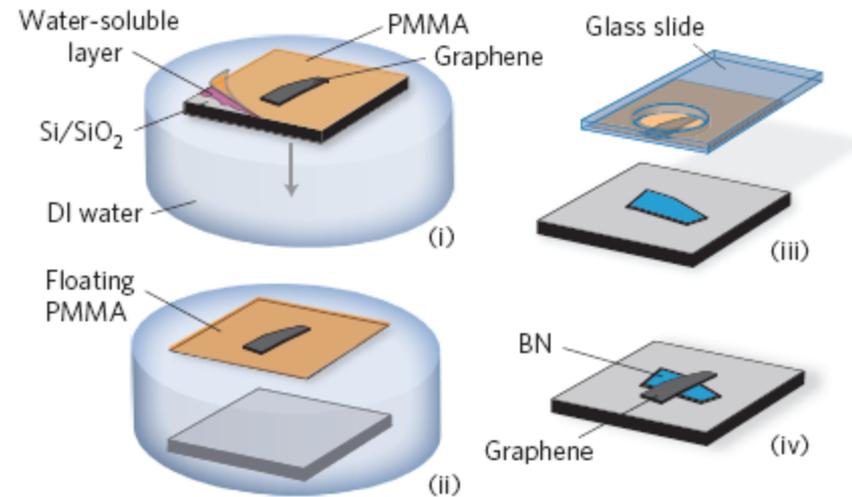
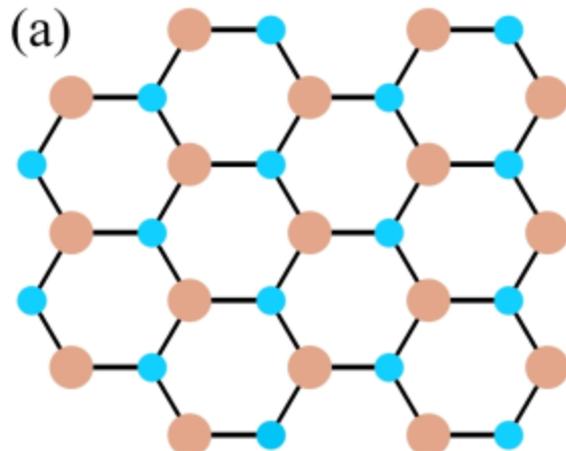


# Better substrate – Boron Nitride



room-T mobility  
close to 100,000 cm<sup>2</sup>/V·s

because it has an atomically smooth surface that is relatively free of dangling bonds and charge traps. It also has a lattice constant similar to that of graphite, and has large optical phonon modes and a large electrical bandgap.



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## “Dirac physics” in transport

- Klein tunneling
- Reflectionless transmission in p-n junction
- Evidence of Klein backscattering in interference pattern of n-p-n junction

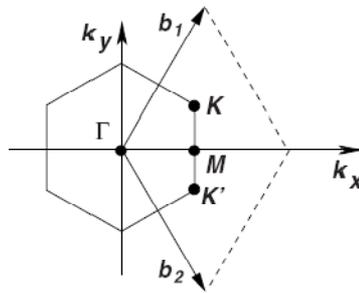
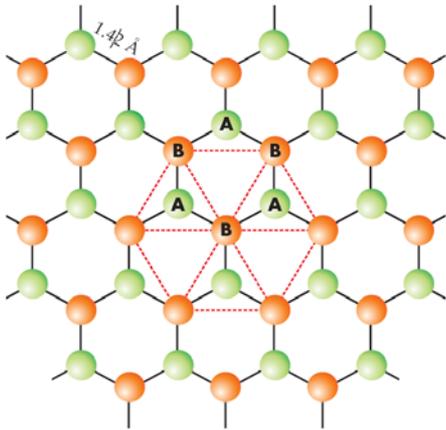
## Graphene based Hybrid Quantum Devices

Sources:

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# Band structure

Lattice, reciproce lattice  
Unit cell with two atoms  
A and B sublattice



Grafén tight binding sávszerkezeti leírása:

Első szomszéd hopping közelítésben (csak másik alrácra ugorhat az elektron)

Hamilton 2\*2 mátrix, A és B alrác komponensekre:

$$H = v_F \boldsymbol{\sigma} \cdot \mathbf{p}$$

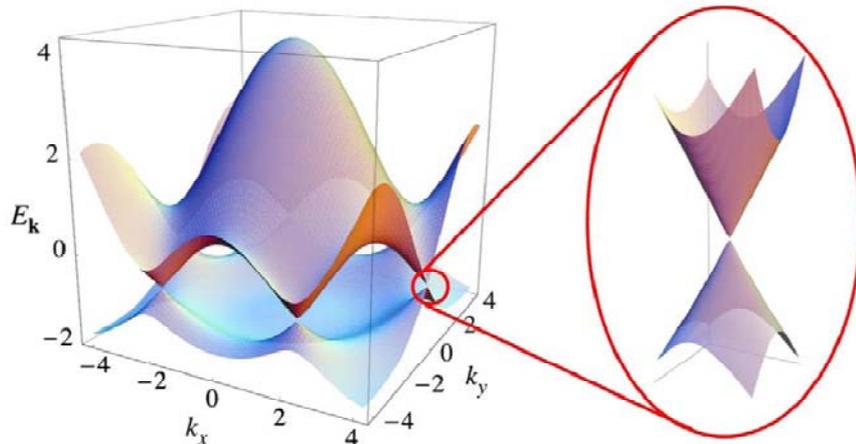
ahol  $\sigma$  a Pauli mátrixok,  $\mathbf{p}$  az e. impulzusa,  $v_F \approx 10^6 \text{m/s}$

A két alrác pseudospinként viselkedik:

$|\uparrow\rangle$  : A alrácson tartózkodás (zöld)

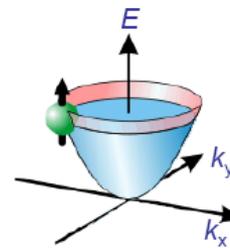
$|\downarrow\rangle$  : B alrácson tartózkodás (piros)

→ Formailag a Dirac egyenlettel megegyező leírást, ahol spin szerepét átveszi a pseudospin



“Schrödinger fermions”

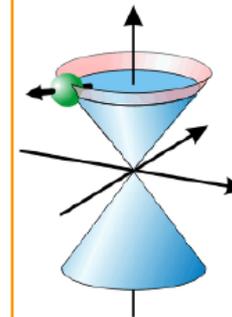
$$\hat{H} = \hat{p}^2 / 2m^*$$



metals and semiconductors

ultra-relativistic particles

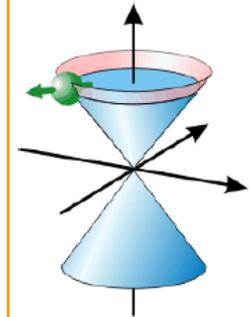
$$\hat{H} = c \bar{\sigma} \cdot \hat{p}$$



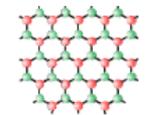
neutron stars and accelerators

massless Dirac fermions

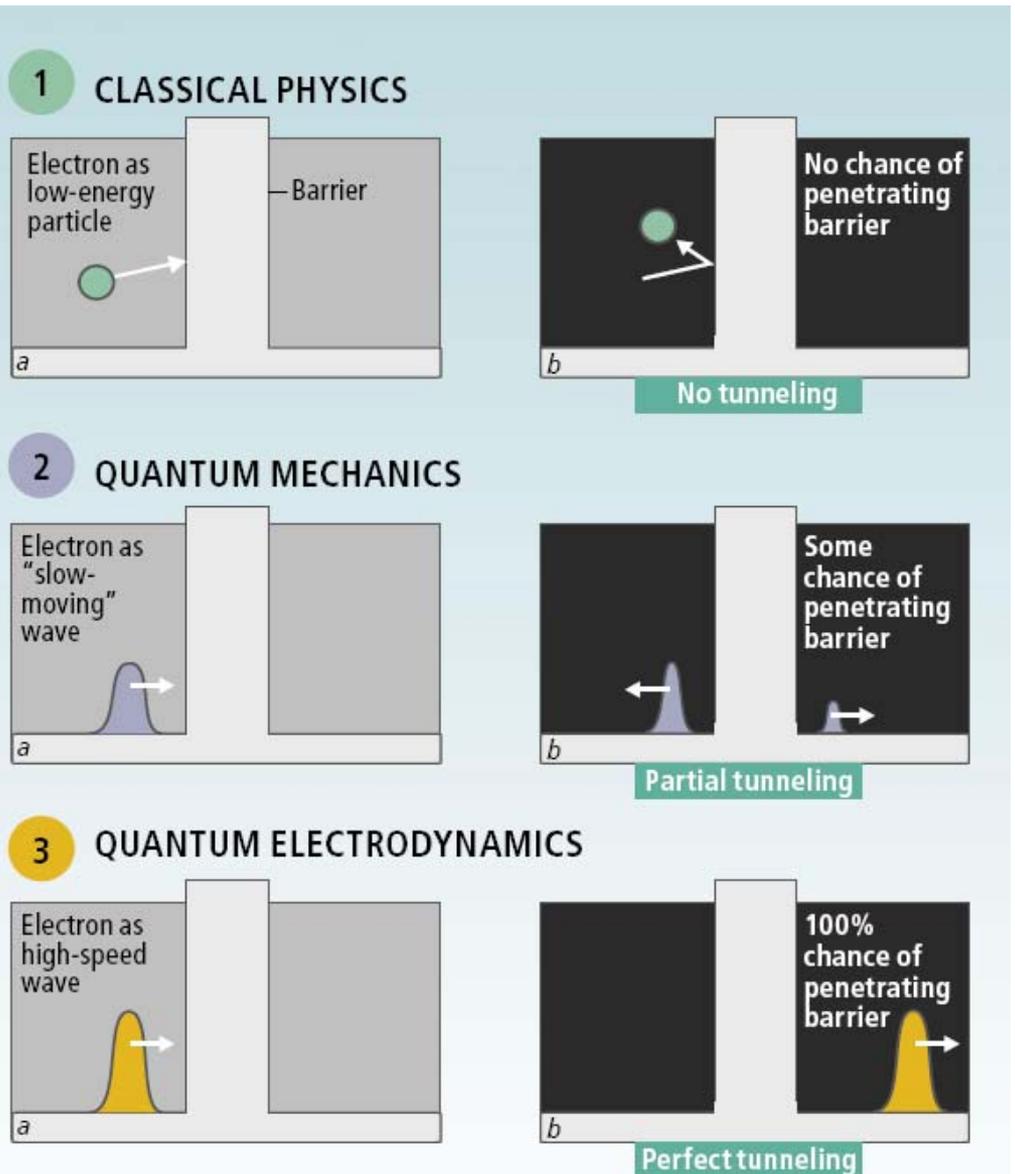
$$\hat{H} = v_F \bar{\sigma} \cdot \hat{p}$$



monolayer graphene



# Tunneling, Klein tunneling



# Tunneling, Klein Backscattering

$$\frac{1}{m_{xx}} = \frac{v_F}{\hbar} \frac{k_y^2}{|k|^3}$$

Evolution of group velocity:

$$\frac{dv_x}{dt} \equiv \frac{1}{m_{xx}} F_x = \frac{1}{m_{xx}} (-e) E_0 \quad (*)$$

In linear electrostatic potential:

$$V = E_0 x, \quad E_x = E_0, \quad F_x = -e E_0$$

At normal incident:  $k_y = 0 \rightarrow \frac{dv_x}{dt} = 0 \rightarrow$

backscattering is avoided

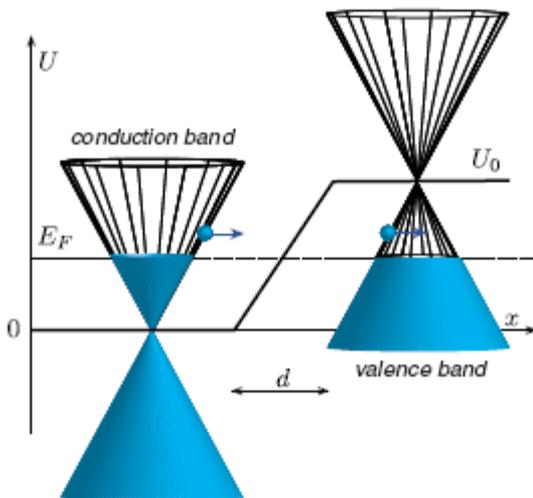
Electron can propagate through an infinite high potential barrier.

**Klein scattering:**

*perfect transmission at normal incident*

N-P junction:

*Potential profile with a step of  $U_0$  at a distance  $d$*



# Tunneling, Klein tunneling

## More precisely, quasi classical dynamics

Two Dirac cones: Conduction band  $E = \hbar v_F |k|$ ,

Valance band  $E = -\hbar v_F |k|$

$$\vec{v} \equiv \frac{1}{\hbar} \frac{\partial E}{\partial \vec{k}} = \frac{1}{\hbar} \hbar v_F \frac{\vec{k}}{|k|} = v_F \vec{e}_k = v_F^2 \frac{\vec{k}}{E},$$

$$\text{thus } |v| = v_F, \vec{v} \parallel \vec{k}$$

$$\hbar \dot{\vec{k}} \equiv \vec{F} = -e E_0 \vec{e}_x$$

Effect of the potential profile,  $U$  (see figure):

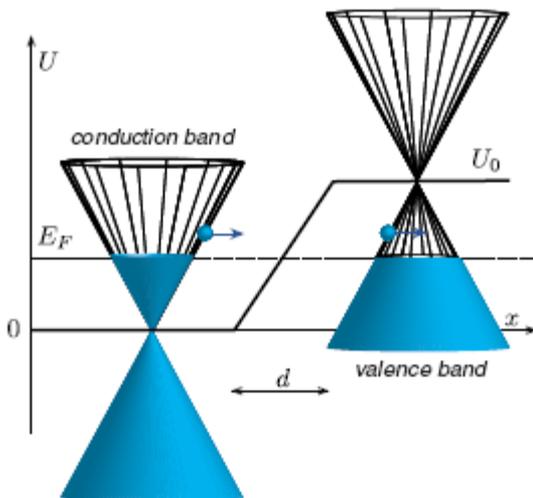
- $k$  decreases and changes sign
- based on (\*),  $\vec{v}$  stays constant, i.e.  $\vec{v} = v_F \vec{e}_x$ .
- e ends up in the valence band

## Klein scattering:

*perfect transmission at normal incident*

N-P junction:

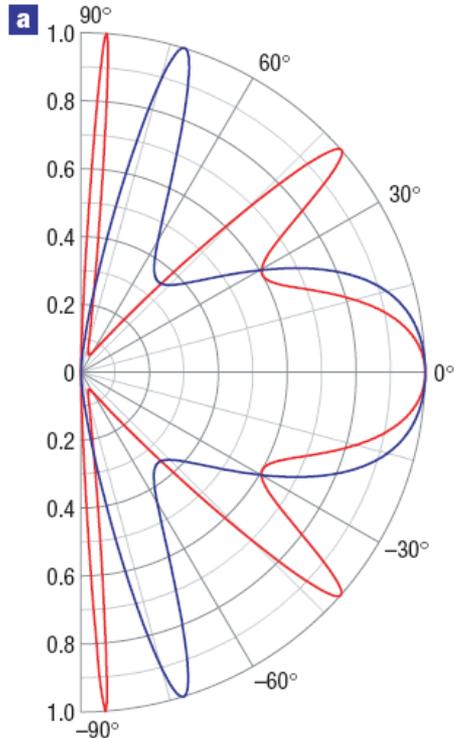
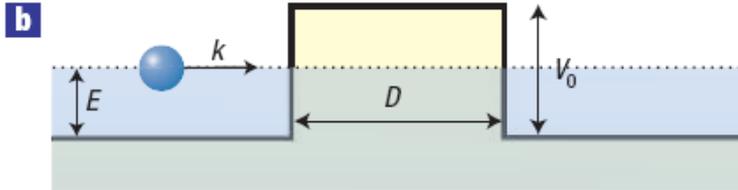
*Potential profile with  
a step of  $U_0$  at a distance  $d$*



# Tunneling, Klein tunneling

## Result of proper calculation

Wave function matching

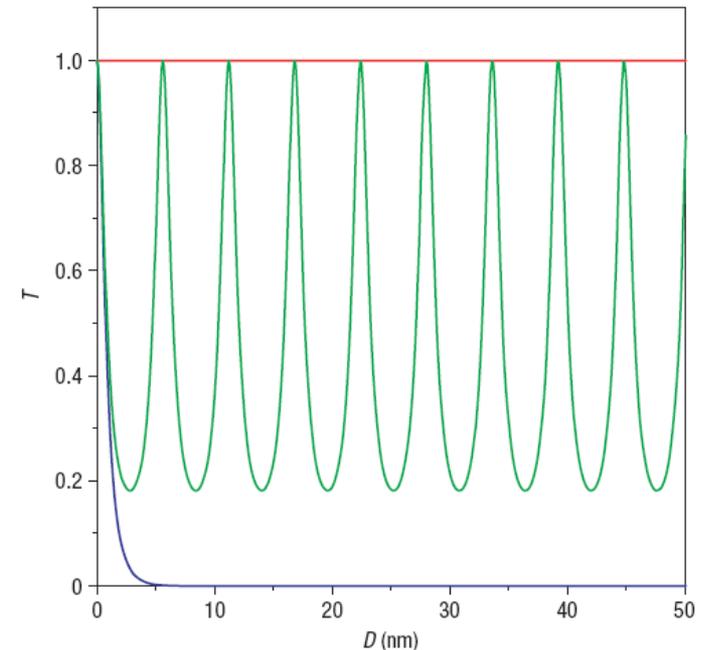


**Transmission probability T** through a 100-nm-wide barrier as a function of the incident angle, two different barrier height

## Transmission probability vs. D

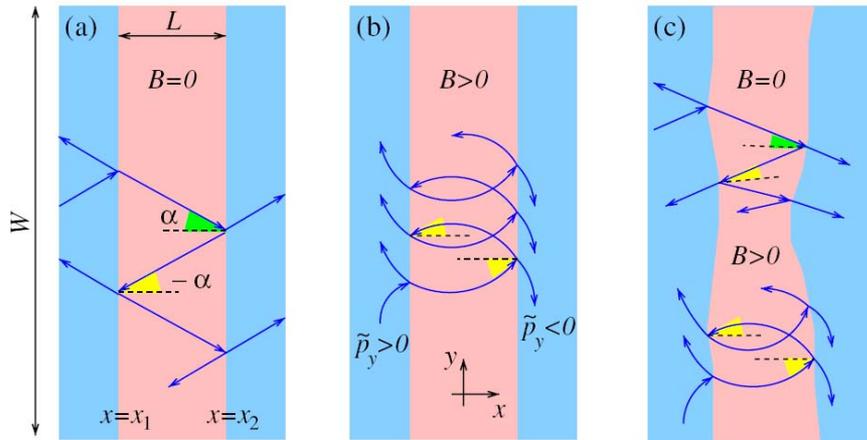
of normally incident electrons

- in single- and bi-layer graphene (red and blue curves, respectively) and in a non-chiral zero-gap semiconductor (green curve)



→ Difficult to measure since e-s out of normal incident also arrive

# Klein backscattering & Fabry-Perot Interferences



## Backscattering on P-N-P junction

When incident angle,  $\alpha$  is varied from positive to negative, phase of the reflection amplitude ( $R$ ) jumps  $\pi$ . Its sign changes. (At  $\alpha=0$ ,  $R=0$ ).

If  $\alpha \neq 0 \rightarrow R > 0$ , several scatterings in P-N-P  $\rightarrow$  interference pattern

Accumulated phase in one circle:

$$\Delta\theta = 2\theta_{\text{WBK}} + \Delta\theta_1 + \Delta\theta_2$$

where  $\theta_{\text{WBK}}$  phase from travelling in N

$\Delta\theta_1, \Delta\theta_2$  Klein backreflection phase of the interfaces

At  $B=0$  (see Fig. a) the incident angles

$\Delta\theta_{1(2)}$  at P-N and N-P have opposite signs  $\rightarrow$  jumps in  $\Delta\theta_1, \Delta\theta_2$  cancels

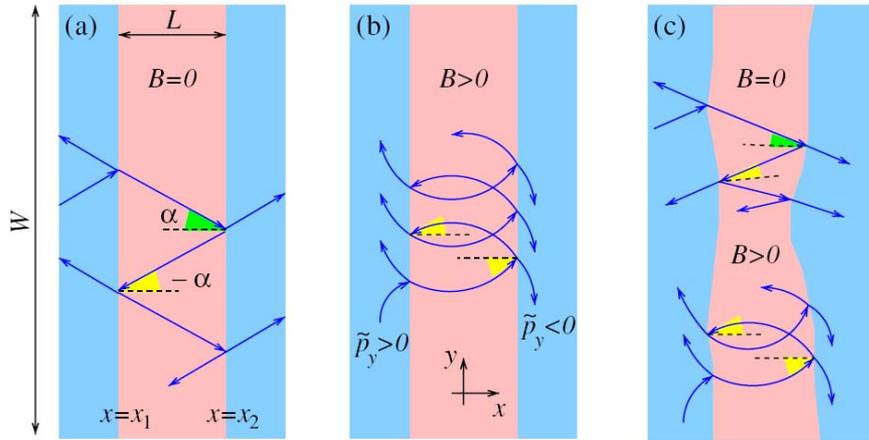
At  $B>0$  (see Fig. b), trajectories are curved,  $\rightarrow$  incident angles at P-N and N-P can be equal

In this case one can show that  $\Delta\theta_1 + \Delta\theta_2 = \pi$  (It is a Berry phase!)

Thus for  $B=0$   $\nearrow$  and trajectories with small  $p_y$   $\pi$  shift is expected (i.e. sign change) transmission amplitude

(Fig.c) one can show, it is robust against barrier roughness

# Klein backscattering & Fabry-Perot Interferences



## Remark (Berry-phase):

Trajectory in Fig.a corresponds to **1**

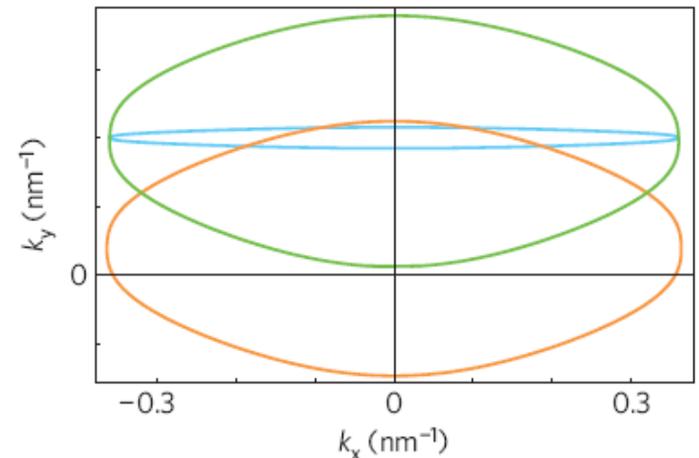
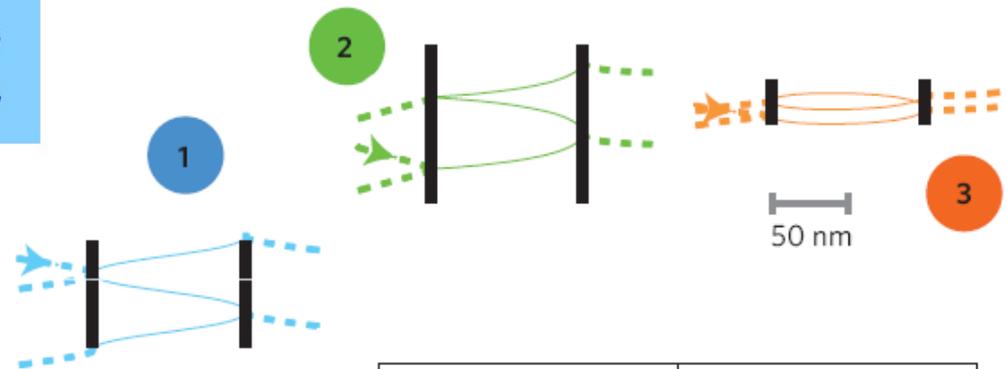
Trajectory in Fig.b corresponds to **3**

The main difference that during one circle between P-N and N-P:

the  $k$  vector of **3** goes around  $k=0$  while for **1** NOT.

This generates the Berry phase:

Due to the chiral symmetry, topological singularity at degeneracy point of the band structure  $k=0$ .

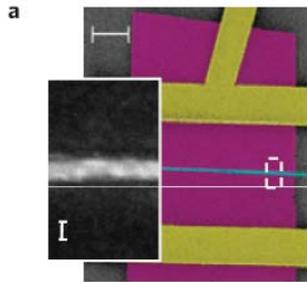


# Klein backscattering & Fabry-Perot Interferences

## N-P-N device

Separate gating by backgate and topgate

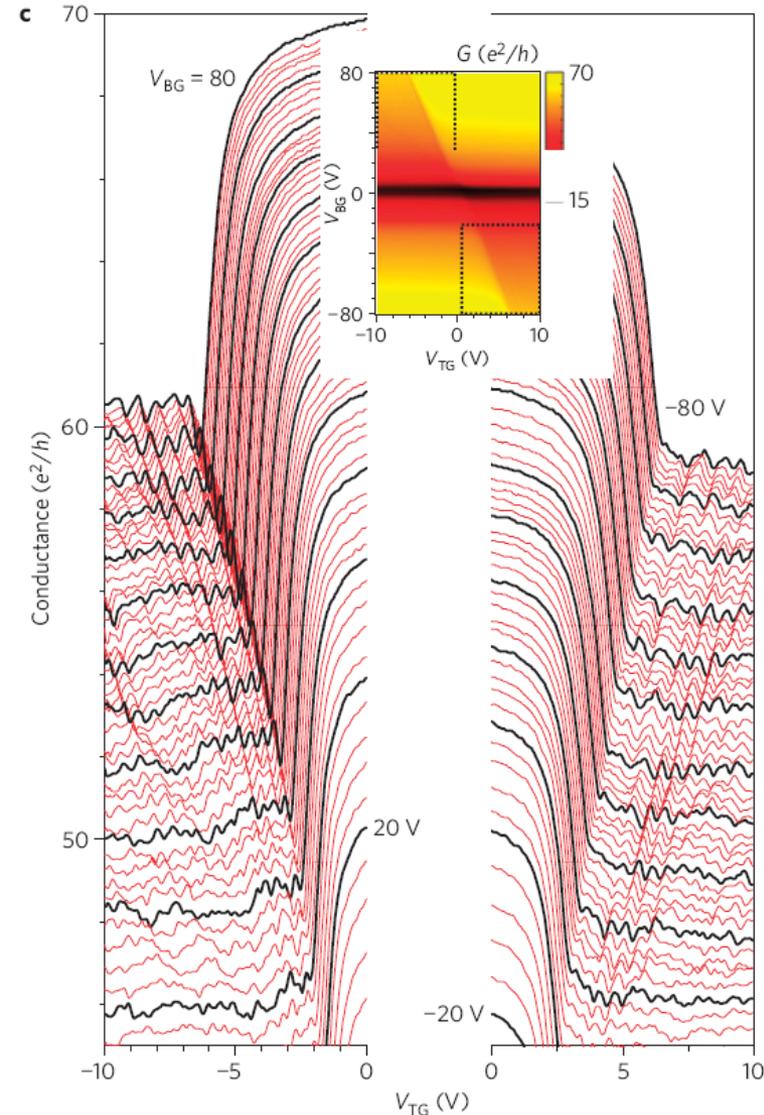
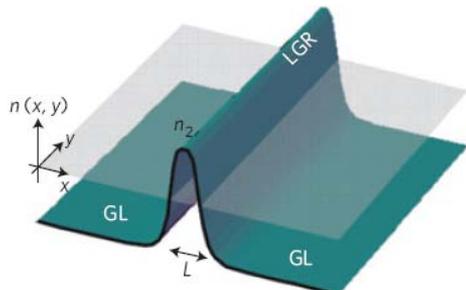
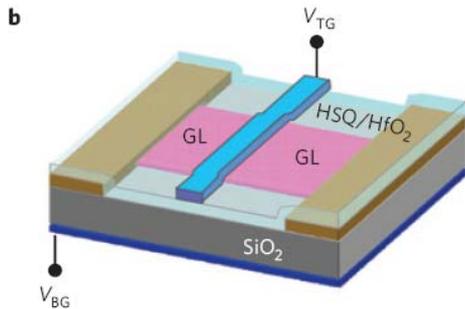
Topgate width=20nm!  $\rightarrow$  ballistic



### G vs. $V_{TG}$ vs. $V_{BG}$

- Conductance is lower when N-P-N setting instead of N-N-N
- Oscillations at N-P-N configuration:

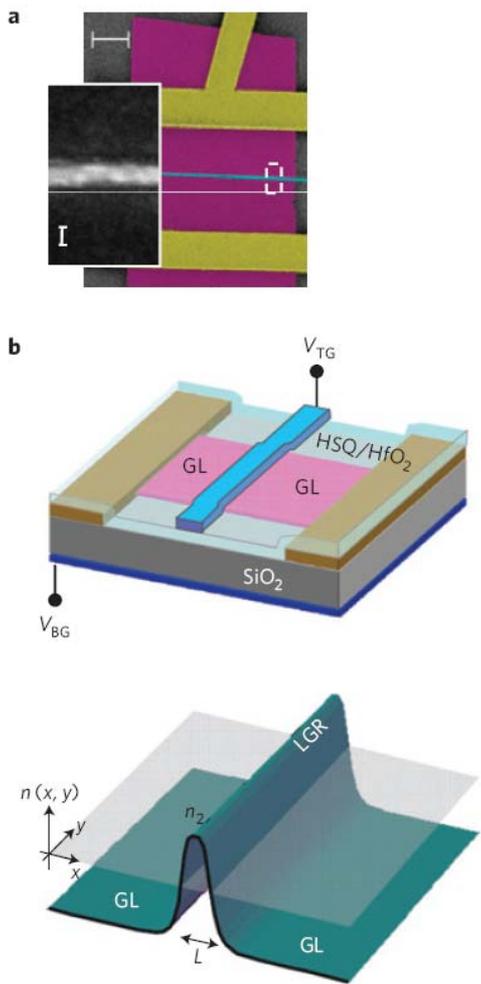
- $V_{TG}$  varies pot. barrier  $\rightarrow \theta_{WBK} \rightarrow$  oscillations
- Oscillatory G is induced by trajectories with incident angle where neither T, nor R is large (i.e.  $\alpha$  not too small)



# Klein backscattering & Fabry-Perot Interferences

## N-P-N device

Separate gating by backgate and topgate  
 Topgate width=20nm!  $\rightarrow$  ballistic

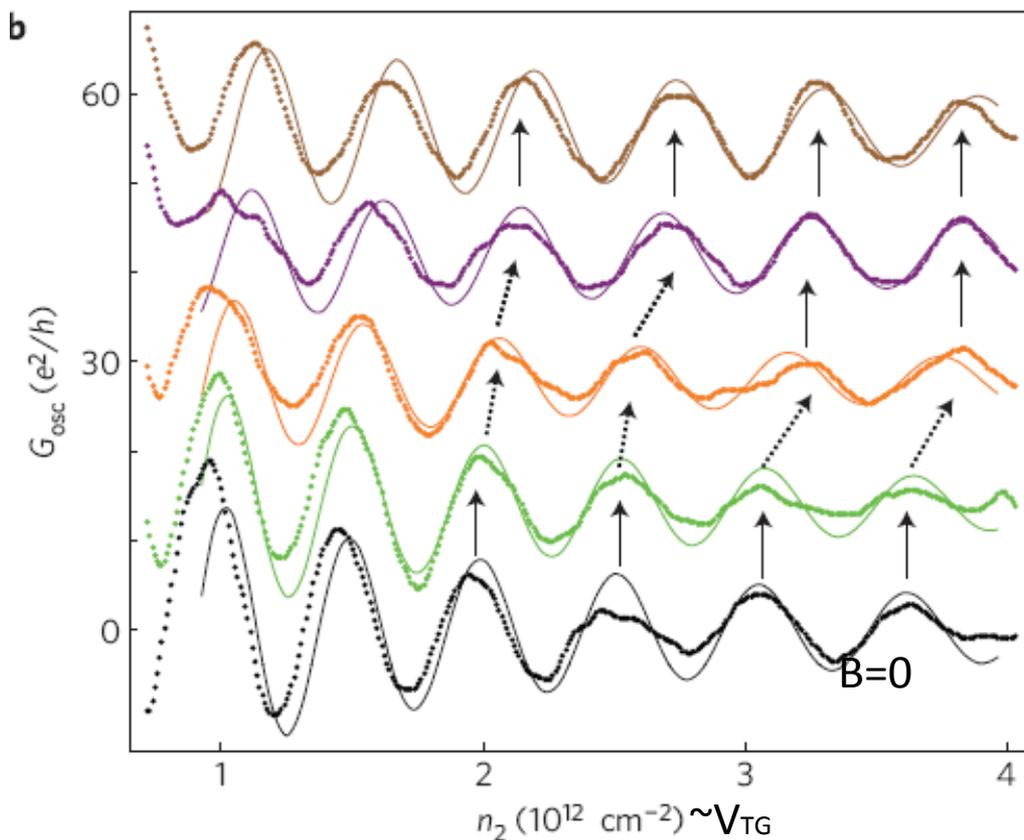


Young et al. Nature Physics 5, 222 (2009)

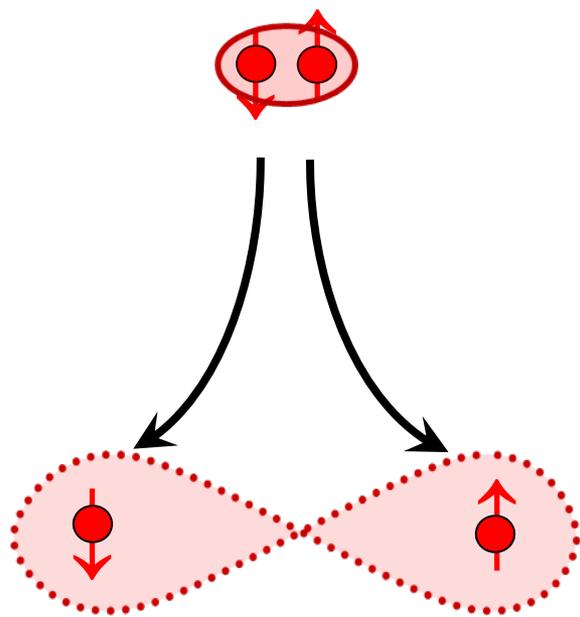
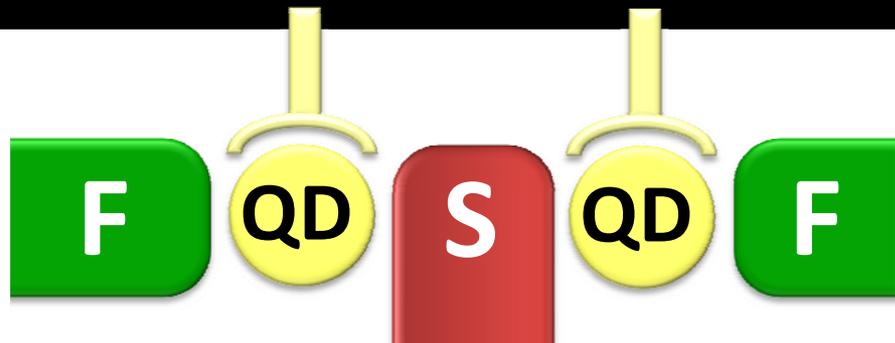
## G oscillations vs. B (Dots experiment, line theory)

At different B fields (B=0, 200, 400, 600, 800mT) the oscillations of G.

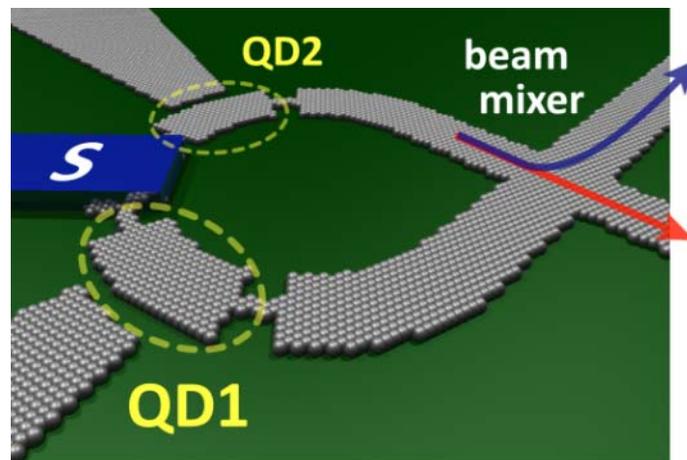
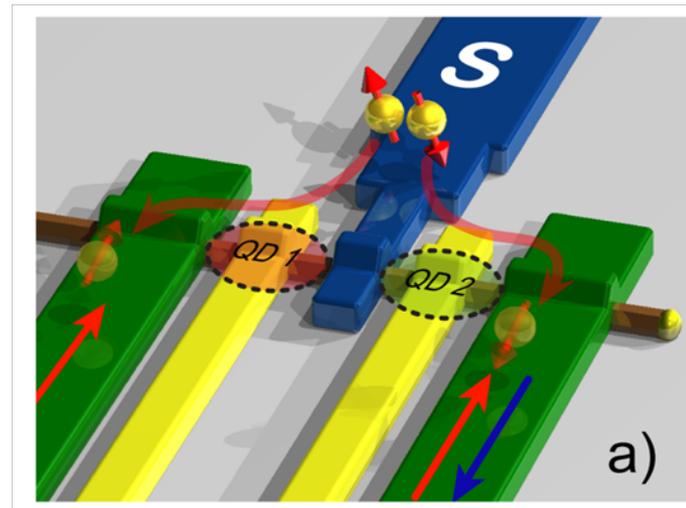
In this B range  $\approx \pi$  shift is induced in the interference pattern.



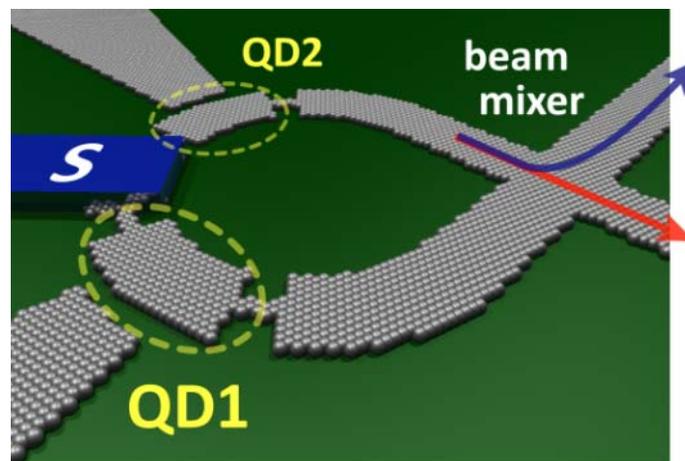
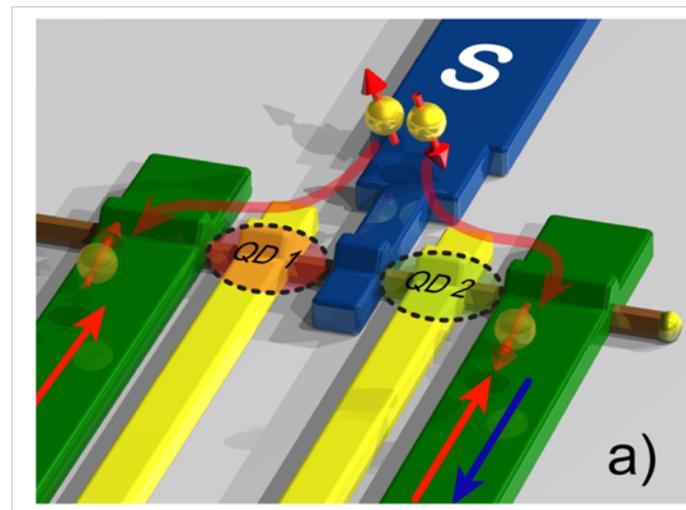
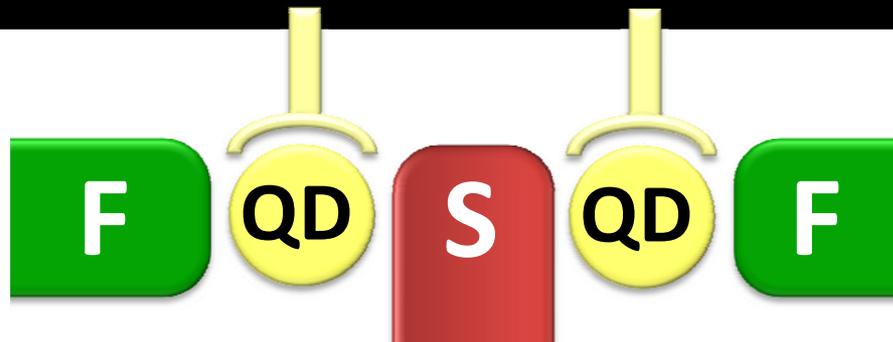
# Hybrid Graphene Devices



Entanglement?



# Hybrid Graphene Devices



# Infrastructure

## BME:

**Low T transport lab:** He liquefier, 4 cryostats, He4, He3 systems, electronics, MCBJ, Kerr ...

## BME & MFA Joint lab:

**E-lithography:** JEOL 848 + Raith Elphy;  
LeoXBeam SEM/FIB, AFM, STM,  
Clean room (300m<sup>2</sup>), Raman, ...

Collaboration with L.P. Biro MFA



BME

MFA

