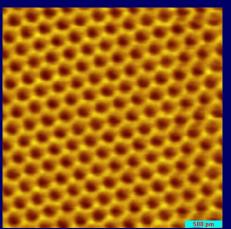
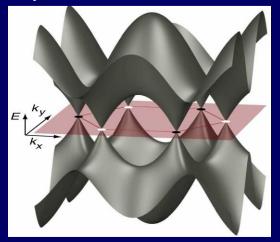
Electronic properties of Graphene and 2-D materials

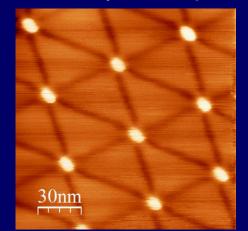
Graphene



Graphene band structure



Twisted Bilayer Graphene



- 2D materials
- ❖ Graphene
- STM and transport measurements
- Flat bands and correlations
- Twisted bilayer graphene

Eva Y. Andrei Rutgers University



2D crystals do not exist in nature



Herbert Wagner

...continuous symmetries cannot be spontaneously broken at finite temperature in systems with sufficiently short-range interactions in dimensions $D \le 2$.

ABSENCE OF FERROMAGNETISM OR ANTIFERROMAGNETISM
IN ONE- OR TWO-DIMENSIONAL ISOTROPIC HEISENBERG MODELS*

N. D. Mermin[†] and H. Wagner[‡] Laboratory of Atomic and Solid State Physics, Cornell University, Ithaca, New York (Received 17 October 1966)

long-range fluctuations can be created with little energy cost and since they increase the entropy they are favored.

No long range order in 2D

HYSICAL REVIEV

OLUME 158. NUMBER

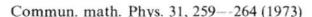
10 TUNE 19

Existence of Long-Range Order in One and Two Dimensions

P. C. Hohenberg

Bell Telephone Laboratories, Murray Hill, New Jersey (Received 24 October 1966)

It is pointed out that a rigorous inequality first proved by Bogoliubov may be used to rule out the existence of quasi-averages (or long-range order) in Bose and Fermi systems for one and two dimensions and Tatlo



There are no Goldstone Bosons in Two Dimensions*

Sidney Coleman **



J. Phys. C: Solid State Phys., Vol. 6, 1973. Brinted in Great Britain. @ 1973

Ordering, metastability and phase transitions in two-dimensional systems

J M Kosterlitz and D J Thouless

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





... No Magnets
....No superfluids
...No superconductors

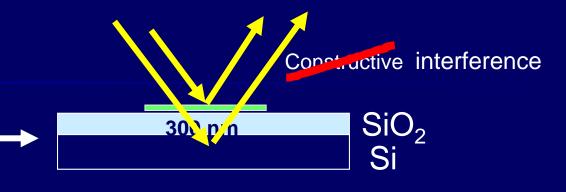
... No 2D crystals

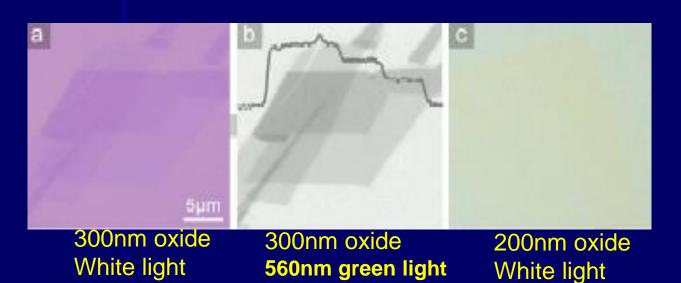


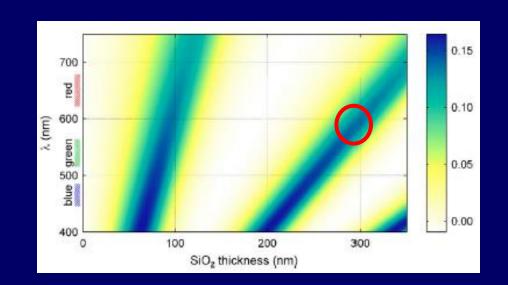
How to see an atomically thin layer?







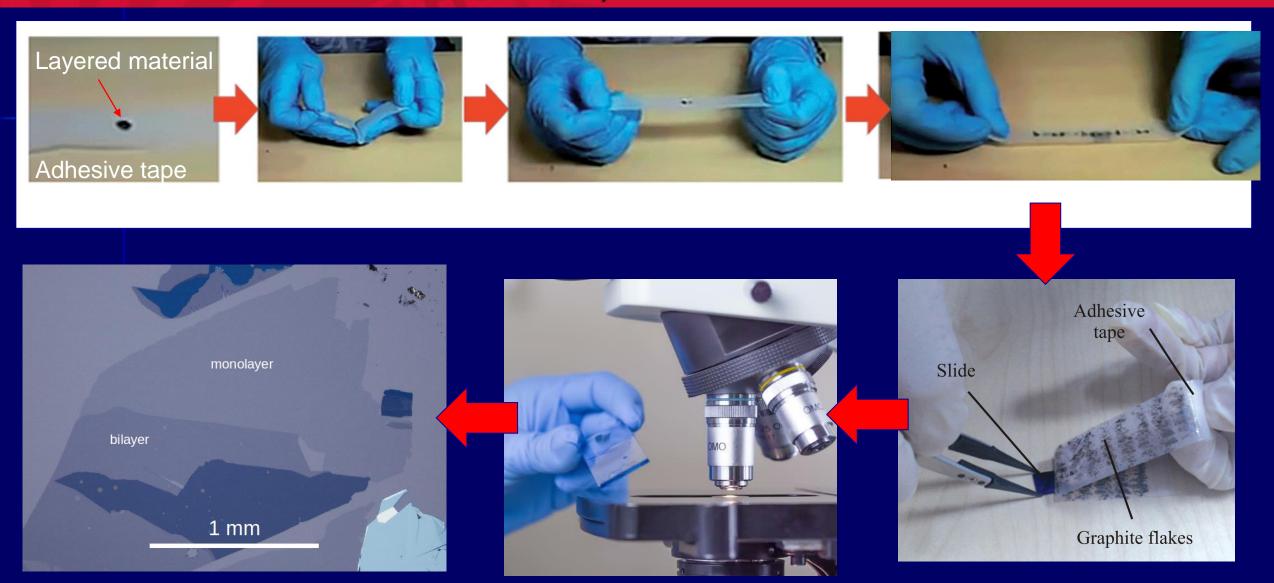




P. Blake et.al, APL, 91, 063124 (2007)



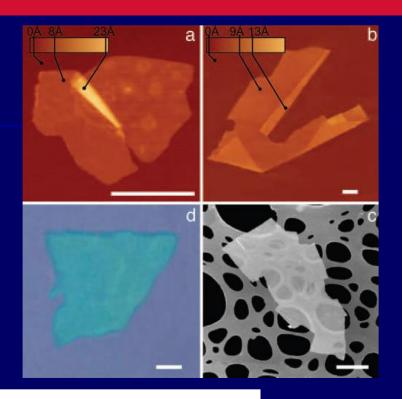
Scotch tape exfoliation



First 2D materials



Graphene, hBN, MoS₂, NbSe₂, Bi₂Sr₂CaCu₂O_x,



Two-dimensional atomic crystals

K. S. Novoselov*, D. Jiang*, F. Schedin*, T. J. Booth*, V. V. Khotkevich*, S. V. Morozov[†], and A. K. Geim*[‡]

*Centre for Mesoscience and Nanotechnology and School of Physics and Astronomy, University of Manchester, Manchester M13 9PL, United Kingdom; and †Institute for Microelectronics Technology, Chernogolovka 142432, Russia

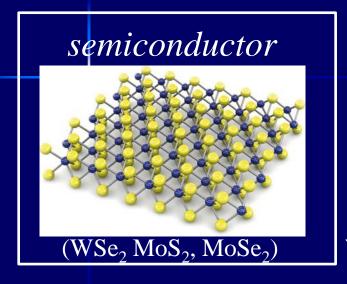
Edited by T. Maurice Rice, Swiss Federal Institute of Technology, Zurich, Switzerland, and approved June 7, 2005 (received for review April 6, 2005)

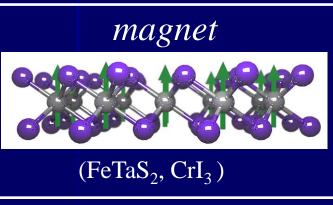
We report free-standing atomic crystals that are strictly 2D and can be viewed as individual atomic planes pulled out of bulk crystals or as unrolled single-wall nanotubes. By using micromechanical cleavage, we have prepared and studied a variety of 2D crystals including single layers of boron nitride, graphite, several dichalcogenides, and complex oxides. These atomically thin sheets (essentially gigantic 2D molecules unprotected from the immediate environment) are stable under ambient conditions, exhibit high crystal quality, and are continuous on a macroscopic scale.

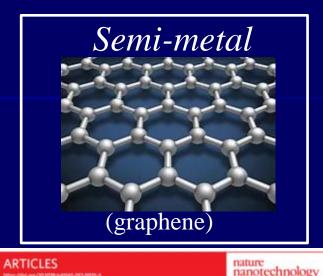
wafer (Fig. 1d), because even a monolayer adds up sufficiently to the optical path of reflected light so that the interference color changes with respect to the one of an empty substrate (phase contrast). The whole procedure takes literally half an hour to implement and identify probable 2D crystallites. Their further analysis was done by atomic force microscopy (AFM), for which single-layer crystals were selected as those exhibiting an apparent (12) thickness of approximately the interlayer distance in the corresponding 3D crystals.



Atomically thin crystals

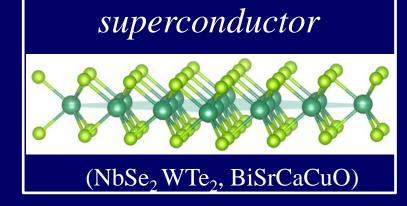


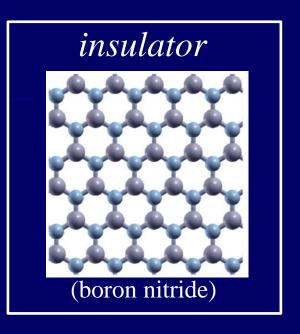


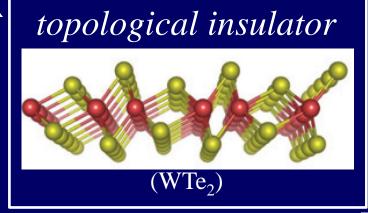


Two-dimensional materials from high-throughput computational exfoliation of experimentally known compounds

Nicolas Mounet[®], Marco Gibertini[®], Philippe Schwaller[®], Davide Campi^{*}, Andrius Merkys[®], Antimo Marrazzo[®], Thibault Sohier[®], Ivano Eligio Castelli[®], Andrea Cepellotti^{*}, Giovanni Pizzi[®] and Nicola Marzari^{**}





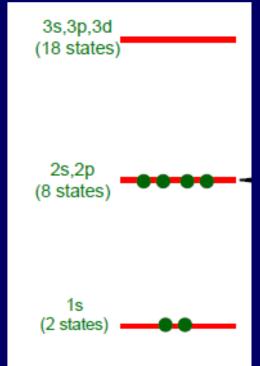


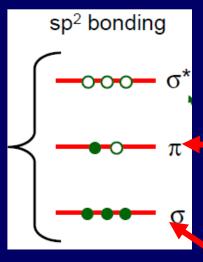


Graphitic bond



Carbon: Z=6 4 valence electrons 2s²2p²

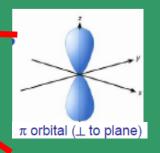




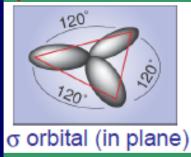
 sp^2

 $2S + 2p_x + 2p_y$ hybridize

3 planar σ orbitals: tetragon 1 Out of plane $2p_{\tau}$ [" π " orbital]



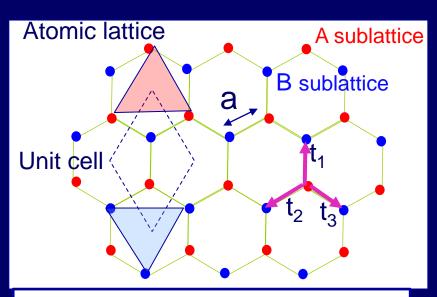
 π electrons allow conduction



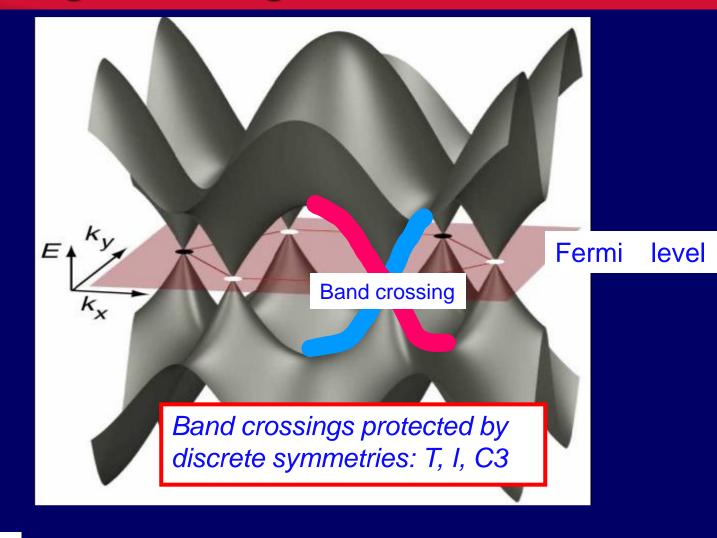
σ bonds: 2D and exceptional rigidity



Graphene: Tight binding model



- 1. 2D
- 2. Honeycomb structure (non-Bravais)
- 3. 2 identical atoms/cell
- 4. Identical NN hopping



Phys 769: Selected Topics in Condensed Matter Physics

Summer 2010

Lecture 5: Graphene: Electronic band structure and Dirac fermions

Lecturer: Anthony J. Leggett TA: Bill Coish

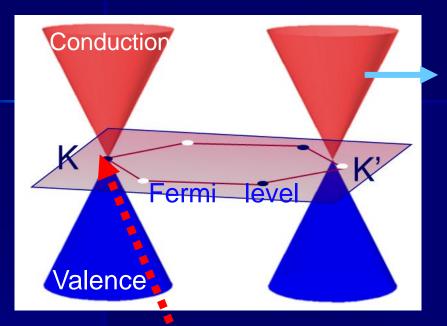
Good general references:

CN: A. H. Castro Neto et al., Rev. Mod. Phys., 81, 109 (2009).

Geim et al., Nature Materials 6 183 (2007).



Graphene



$$v_F = \frac{3}{2\hbar} at \sim 10^6 m/s$$

Dirac Weyl Hamiltonian

$$H = v_F \begin{pmatrix} \vec{\sigma} \cdot \vec{p} & 0 \\ 0 & -\vec{\sigma}^* \cdot \vec{p} \end{pmatrix} \mathbf{K}'$$

p momentum relative to K point σ Pauli matrices operate on sublattice degree of freedom

Dirac point

6 Dirac cones, but only 2 independent at K and K' valleys



No Backscattering



 $\vec{p} = \hbar(\vec{k} - \vec{K})$ Dirac Weyl Hamiltonian

$$H = v_F \begin{pmatrix} \vec{\sigma} \cdot \vec{p} \\ 0 \end{pmatrix} \begin{pmatrix} 0 \\ -\vec{\sigma}^* \cdot \vec{p} \end{pmatrix}$$

Helicity

$$H_{K} \propto \vec{\sigma}.\vec{p}$$

→ projection of pseudospin along momentum conserved



Two massless Dirac fermions with opposite chirality (helicity)

4 flavors: 2 Spin, 2 valley



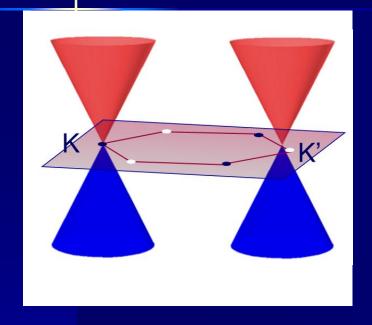
High carrier mobility

No backscattering between cones

No backscattering within cones

Density of states

Band structure

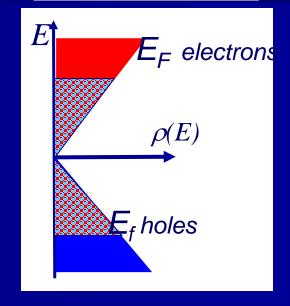


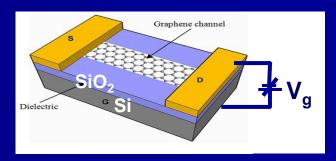


 $1V \longrightarrow 7x10^{10} \text{ cm}^{-2}$ $\Delta E \sim 80 \text{meV}$

Density of states

$$\rho(E) = 3^{3/2} a^2 \frac{|E|}{\pi (\hbar v_F)^2}$$

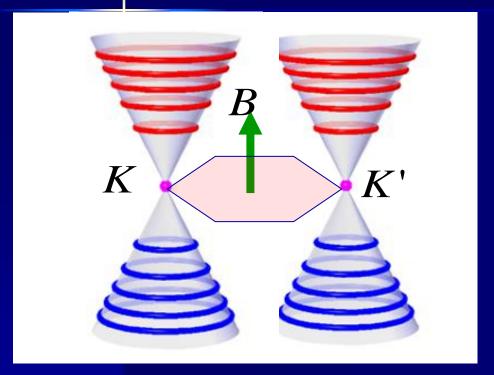




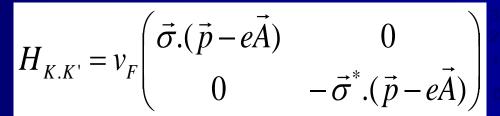


Landau levels

Band structure



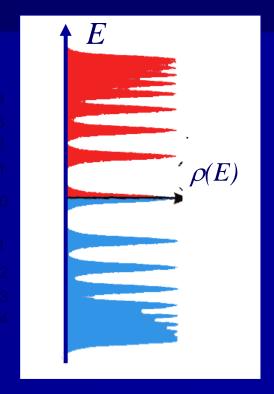
$ec{B} = \vec{ abla} imes \vec{A}$



Landau Levels

$$E_N = \pm v_F \sqrt{2e\hbar B|N|}$$
 $N = 0,\pm 1,...$

Density of states



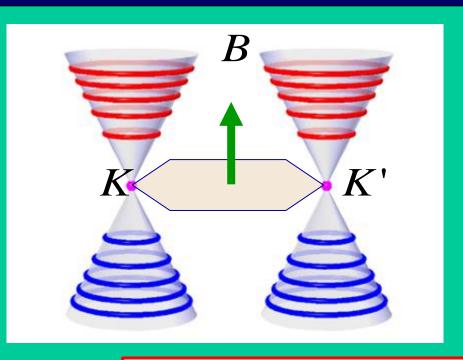


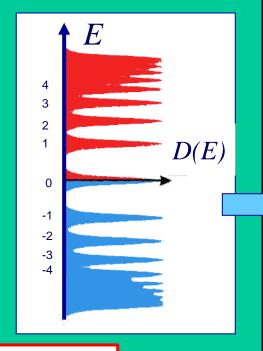
Sublattice Valley locking

Band structure

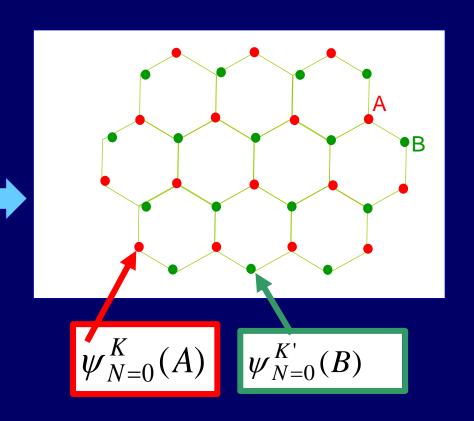
DOS

N=0 Wave function





$$E_N = \pm v_F \sqrt{2e\hbar B|N|} N = 0,\pm 1,...$$





Energy and length scales

$$E_N = \pm v_F \sqrt{2e\hbar B|N|} = \pm \varepsilon_0 \sqrt{2|N|}; \quad N = 0,\pm 1,...$$

$$\varepsilon_0 = \hbar v_F / l_B \approx 35 \sqrt{\mathrm{B}} \,\mathrm{meV}$$

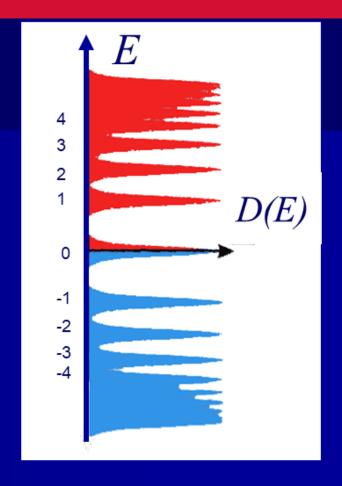
Orbital Degeneracy
$$\mathbf{g}_0 = B/\phi_0 = 2.5 \times 10^{14} \, m^{-2} B[T]$$

Total *Degeneracy*:
$$g = 4g_0$$
; $\phi_0 = h/e$



$$E_N = (N + 1/2)\hbar\omega_c$$

$$\hbar\omega_c \approx (0.1B)meV$$





Exquisite sensitivity of 2Dmaterials

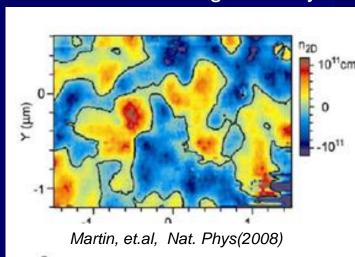
- 2D materials
- Graphene
- Transport and STM
- Flat bands and correlations
- Twisted bilayer graphene

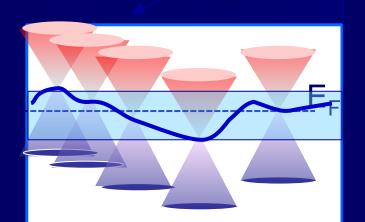


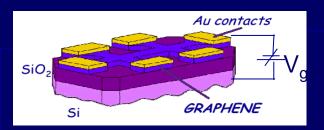
Transport measurements: Substrate trouble

e-h puddles smeared Dirac point

SET: Local charge density



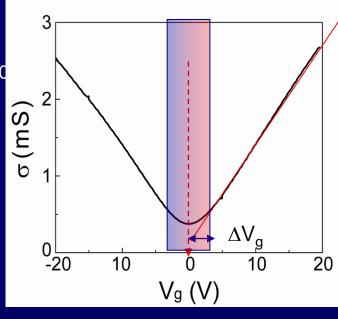




STM

- Ishigami et al Nano letters 200
- Stolyarova et al PNAS 2007
- Geringer et al PRL 2009
- Zhang et al Nat Phys 2009
- Luican et al PRB R 2011
- S. Jung et al Nat Phys 2011

Conductivity

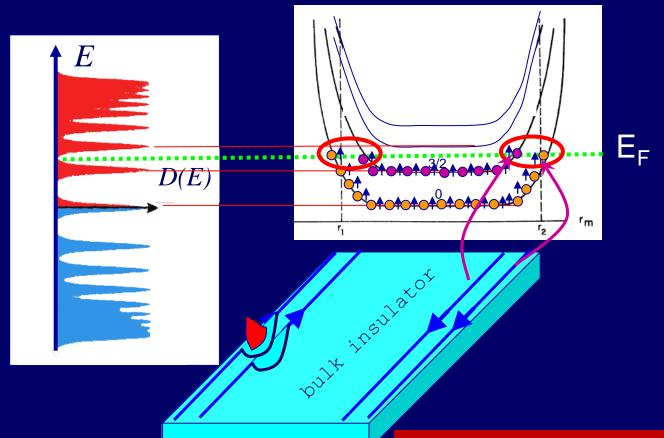


 $V_{gmin} \sim 1-10V$ $n_{min} \sim 10^{11} - 10^{12} \text{ cm}^{-2}$ $(\Delta E_{RP} \sim 30-100 \text{meV})$



Quantum Hall Effect

For $E_F \neq E_N$ and sharp edge each filled Landau level contributes one (g-fold degenerate) edge state



- Ballistic channelsOne dimensional edge states
- No backscattering:as long as channels far apart
- -Robust against disorder

Prerequisites for QHE in graphene

IQHE

Decoupled edges

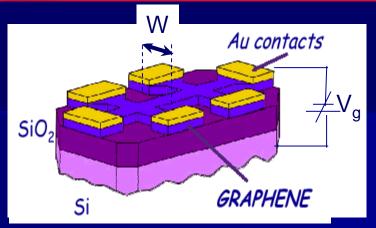
$$l_c = 25nm/(B[T])^{1/2} << W$$

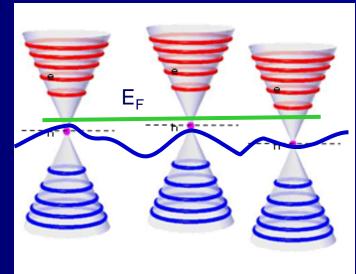
2. Well defined Landau levels

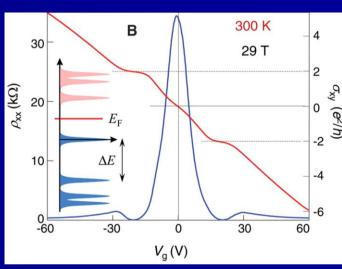
$$E_1 - E_0 = 35 meV \cdot (B[T])^{1/2} > \Delta E_{RP}, k_B T$$

 $\Delta E_{RP} \sim 100 \text{meV} \rightarrow \text{IQHE for B} > 10 \text{T}$

random potential







Room-Temperature Quantum Hall Effect in Graphene







Unteger and fractional QHE in graphene

IQHE

1. Decoupled edges

$$l_c = 25nm/(B[T])^{1/2} << W$$

2. Well defined Landau levels

$$E_1 - E_0 = 35 meV \cdot (B[T])^{1/2} > \Delta E_{RP}, k_B T$$

$$\Delta E_{RP} \sim 100 \text{meV} \rightarrow \text{IQHE for B} > 10 \text{T}$$

random potential

IQHE easily observed in graphene even at room temperature!

FQHE

1. Decoupled edges

$$l_c = 25nm/(B[T])^{1/2} << W$$

2. Well defined "Λ" levels

$$E_{\Lambda=1/3} = v^2 \frac{e^2}{\kappa l_c} \approx 5meV \frac{\left(B[T]\right)^{1/2}}{\kappa} > \Delta E_{RP}$$

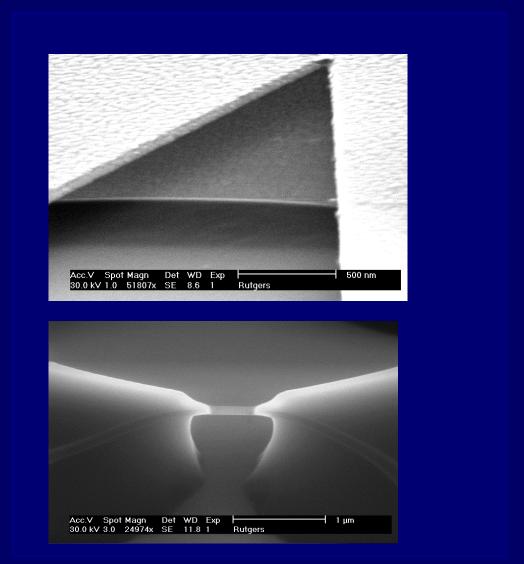
$$\Delta E_{RP} \sim 100meV \Longrightarrow B > 6400 T$$

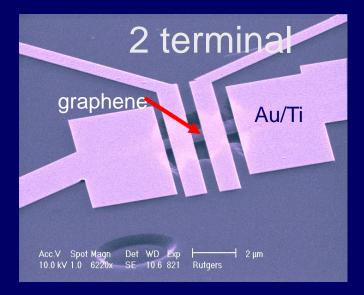
FQHE not observable in graphene on SiO₂ substrates



Suspended Graphene

- •X. Du, I. Skachako, A. Barker, E. Y. A. Nature Nanotech. 3, 491 (2008)
- Bolotin et al , Solid State Communications (2008)

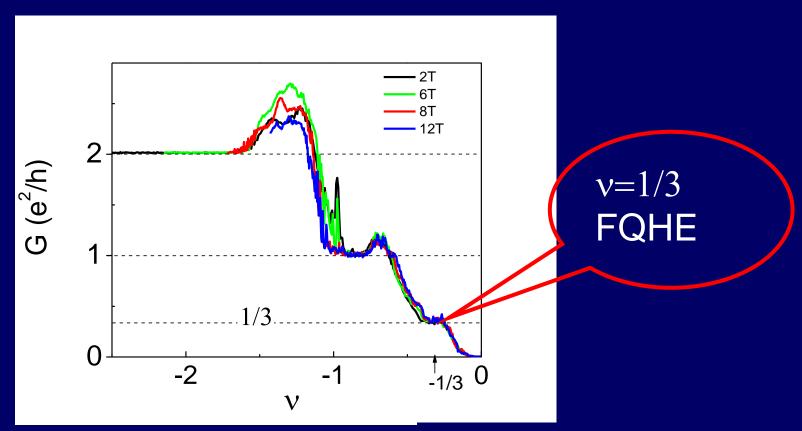








X. Du, I. Skachko, F. Duerr, A. Luican, EYA, Nature **462**, 192 (2009)



FQHE in graphene
Seen already at 2T at 0.3K
Persists up to 20K (in 12T)

- Bolotin et al (Columbia) Nature 462 (2009)
- Geim & Novoselov (Manchester)

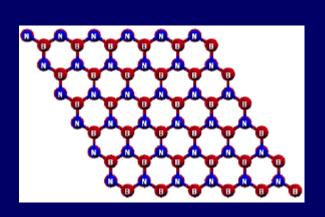
G/BN C. R. Dean et. al arxiv1010.1179

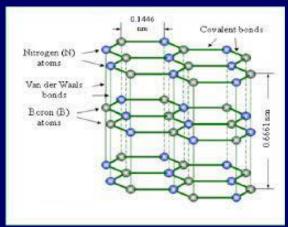


The magic of hexagonal Boron Nitride (hBN)

Hexagonal Boron Nitride - hBN

- Honeycomb structure no dangling bonds, chemically inert.
- Protects graphene from electrical random potential, and chemical contamination

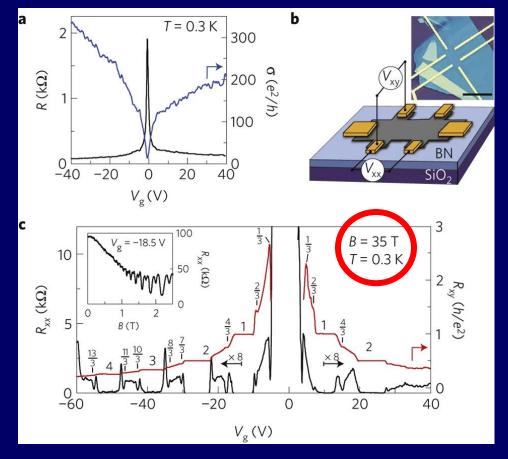




Multicomponent fractional quantum Hall effect in graphene

C. R. Dean, A. F. Young, P. Cadden-Zimansky, L. Wang, H. Ren, K. Watanabe, T. Taniguchi, P. Kim □, J. Hone & K. L. Shepard

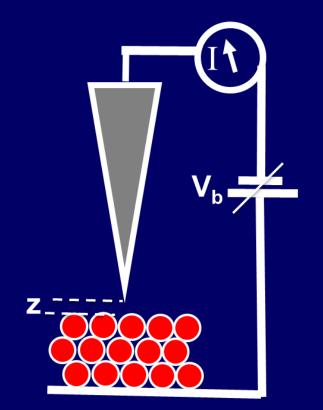
Nature Physics 7, 693–696 (2011) Cite this article



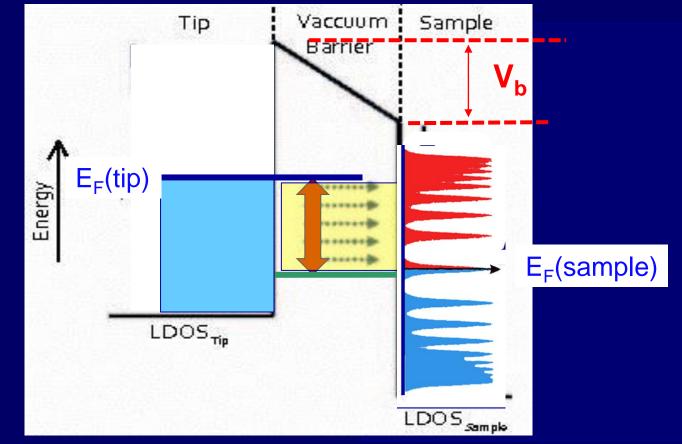


Scanning Tunneling Microscopy (STM

STM measures tunnel current across a vacuum barrier



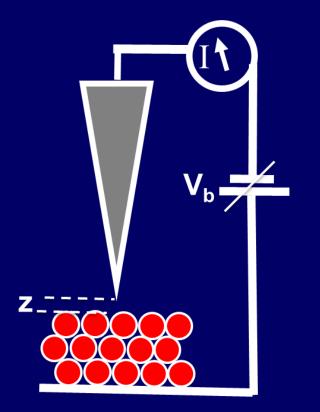
$$I(r, z, V) \propto \left[\int_{0}^{eV_b} \rho(r, \varepsilon) d\varepsilon\right] \exp^{-z(r)\kappa}$$

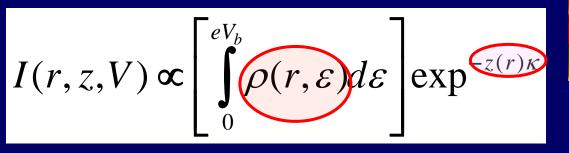


STM probes all states between $E_F(tip)$ and $E_F(Sample)$



Graphene: STM and STS



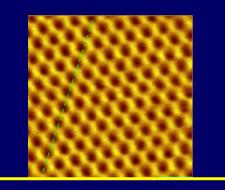




I, V_b constant

Topography: Imaging Atoms

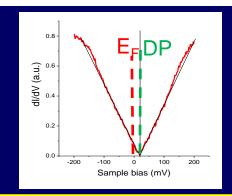
$$I(r,z) \propto \exp^{-z(r)\kappa}$$



z constant

Spectroscopy: Density of states

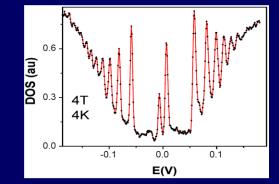
$$dI/dV_b \propto \rho(r, V_b)$$



Landau level Spectroscopy:

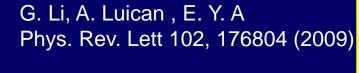
- Local Fermi velocity
- Quasiparticle lifetime
- Coupling to substrate

$$\left| E_N = \pm v_F \sqrt{2e\hbar B} |N| \right|$$



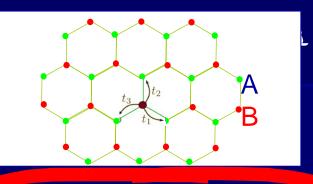
E.Y. Andrei



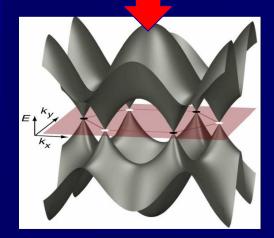


Graphene on graphite

Controlling the band structure?



- 1. 2D
- 2. Honeycomb structure
- 3. 2 identical atoms/cell
- 4. $t_1 = t_2 = t_3 = t$



Can one control/manipulate the band structure?

- \triangleright Relax the condition: $t_1 = t_2 = t_3$
 - Strain
- Impose external potential
 - Substrate
- Point Defects
- Electrostatic potential
- Magnetic field



2-terminal measurement of QHE

Conformal invariance and shape-dependent conductance of graphene samples

Dmitry A. Abanin and Leonid S. Levitov Phys. Rev. B **78**, 035416 – Published 10 July 2008

Fractional quantum Hall effect in suspended graphene: Transport coefficients and electron interaction strength

D. A. Abanin, I. Skachko, X. Du, E. Y. Andrei, and L. S. Levitov Phys. Rev. B **81**, 115410 – Published 8 March 2010

Berry phase of π

Electronic properties of graphene: a perspective from scanning tunneling microscopy and magnetotransport

Eva Y Andrei¹, Guohong Li¹ and Xu Du²

Published 19 April 2012 • © 2012 IOP Publishing Ltd

Reports on Progress in Physics, Volume 75, Number 5

PHILOSOPHICAL TRANSACTIONS OF THE ROYAL SOCIETY A

MATHEMATICAL, PHYSICAL AND ENGINEERING SCIENCES

Fractional quantum Hall effect in suspended graphene probed with two-terminal measurements

I. Skachko, X. Du, F. Duerr, A. Luican, D. A. Abanin ⊡, L. S. Levitov and E. Y. Andrei

Published: 13 December 2010 https://doi.org/10.1098/rsta.2010.0226

Onsager: Area of N'th orbit in k space

$$S(k_N) = \pi k_N^2 l_B^2 = 2\pi (N + 1/2 - \gamma/2\pi)$$

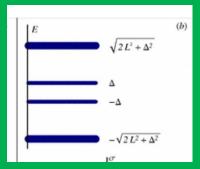
$$\gamma = \text{Berry Phase}$$

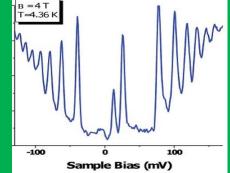
 $\oint_C d\lambda \langle \psi(\lambda) | i(\partial/\partial \lambda) | \psi(\lambda) \rangle$

Excitonic gap (Magnetic catalysis)

Excitonic gap, phase transition, and quantum Hall effect in graphene

V. P. Gusynin, V. A. Miransky, S. G. Sharapov, and I. A. Shovkovy Phys. Rev. B **74**, 195429 – Published 22 November 2006





More references

Graphene monolayer and bilayer tutorials

REVIEW ARTICLE

The electronic properties of bilayer graphene

Edward McCann¹ and Mikito Koshino²

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