

Localised Collective Excitations in Doped Graphene in a Strong Magnetic Field

[arXiv:0902.4176]

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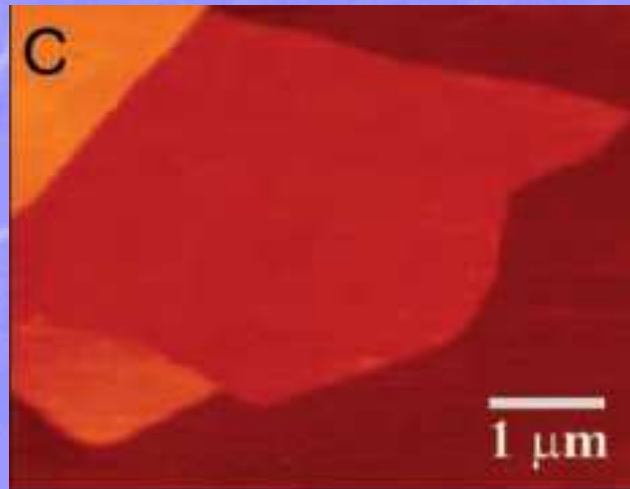
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What is Graphene?

- “*Graphene is the name given to a single layer of carbon atoms densely packed into a benzene-ring structure*”
- First isolated in 2004 by physicists at the University of Manchester, UK and the Institute for Microelectronics Technology, Chernogolovka, Russia



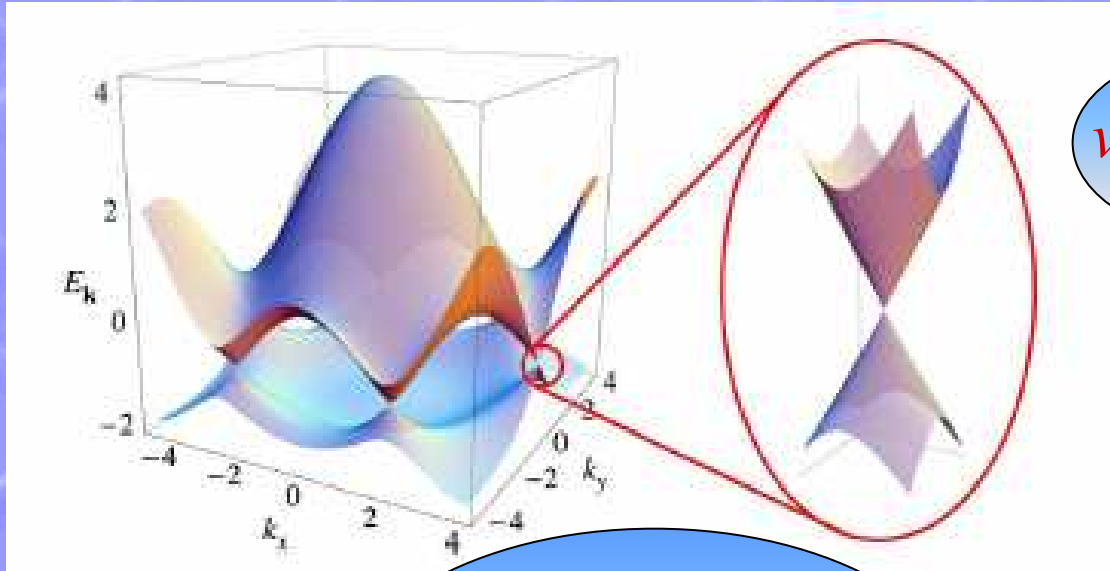
Andre Geim



K.S. Novoselov *et al.*,
Science **306**, 666
(2004)

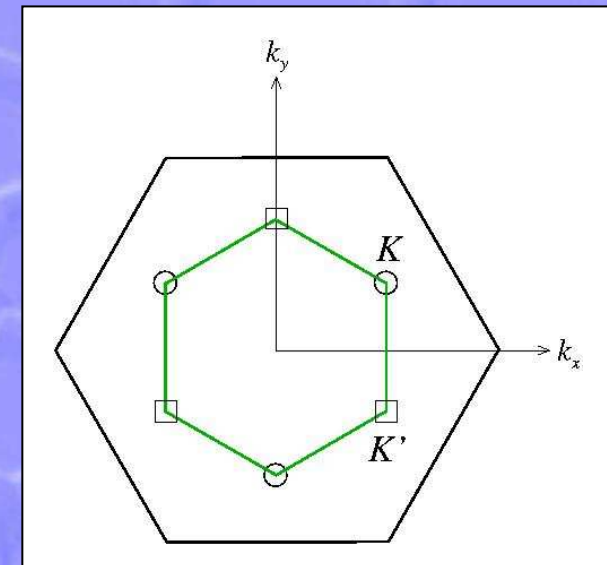
Dispersion Relation

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



$$v_F \sim 10^6 \text{ ms}^{-1}$$

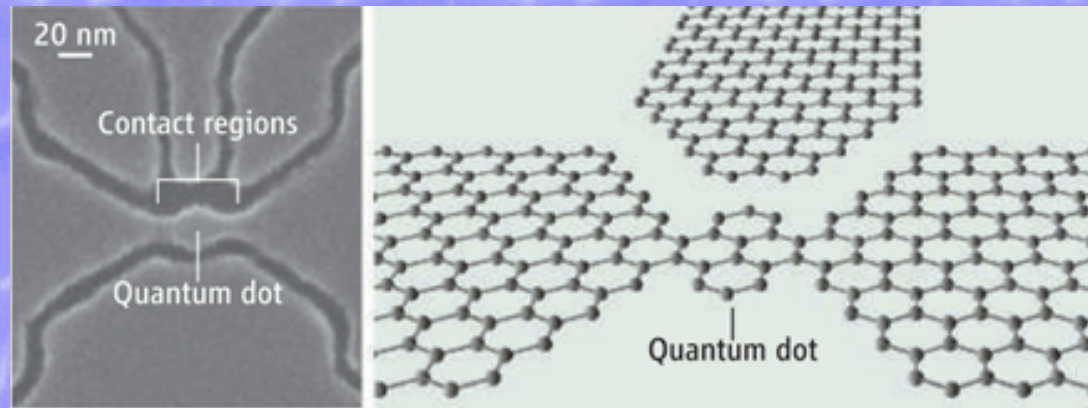
linear dispersion relation
at low energies



P.R. Wallace, Phys. Rev., **71**, 622 (1947)

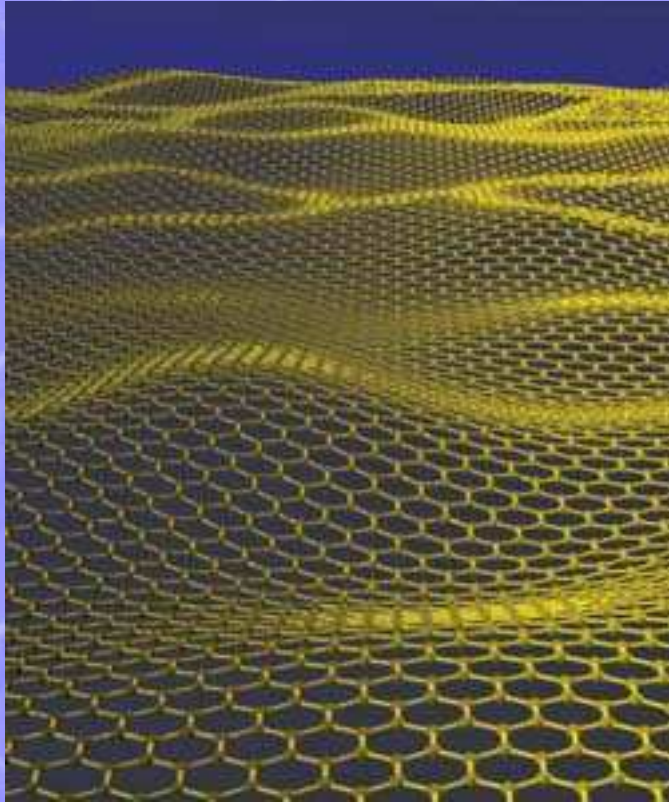
What's all the fuss?

- Fundamentally interesting physics and useful physical properties
- QED in a mesoscopic system – e.g. room temperature anomalous integer quantum Hall effect
K.S. Novoselov et al, Nature **438**, 197
- Important applications i.e. nanoelectronics – is graphene the new silicon?



Credit: Andre Geim

Unanswered Questions

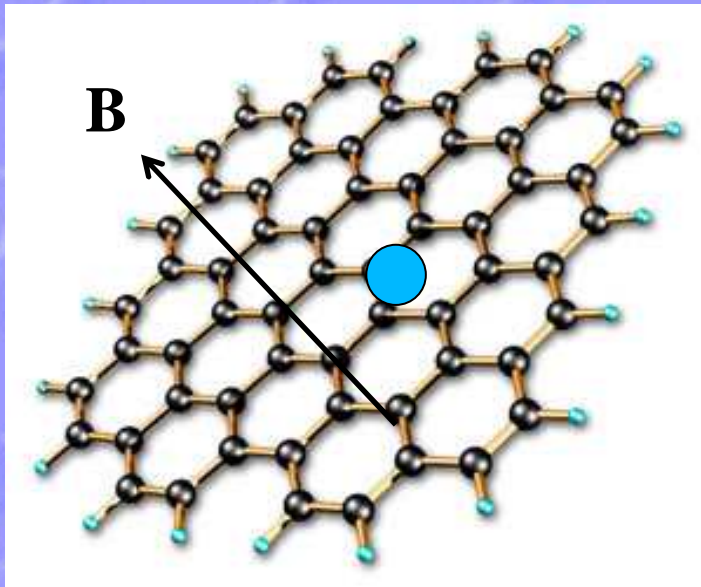


Credit: Andre Geim

- How does substrate affect graphene properties?
- What is the nature of disorder in graphene?
- How could an energy gap be opened at the K points?
 - constriction i.e. nanoribbons
 - application of an electric field

Aim of Project

- Single graphene sheet in a strong perpendicular magnetic field and single axially symmetric charged impurity - corresponds to a low impurity density



High magnetic field approximation – reduce disorder broadening, so LLs well defined

- Want to find collective excitations localised on impurity

Single Particle Picture

- Single impurity at origin – symmetric gauge: $\mathbf{A} = 1/2 \mathbf{B} \times \mathbf{r}$
- No intervalley scattering

$$H_D = v_F \sigma \cdot \mathbf{p} = \hbar \frac{\sqrt{2} v_F}{\ell_B} \begin{pmatrix} 0 & -\frac{i}{\sqrt{2}} \left(\frac{z^*}{2\ell_B} + 2\ell_B \partial_Z \right) \\ \frac{i}{\sqrt{2}} \left(\frac{z}{2\ell_B} - 2\ell_B \partial_{Z^*} \right) & 0 \end{pmatrix} = \hbar \omega_c \begin{pmatrix} 0 & a \\ a^\dagger & 0 \end{pmatrix}$$

- Corresponding eigensystem:

$$\Phi_{ns\uparrow m}(\mathbf{r}) = 2^{\frac{1}{2}(\delta_{n,0}-1)} \begin{pmatrix} s_n \phi_{|n|-1 m}(\mathbf{r}) \\ \phi_{|n| m}(\mathbf{r}) \end{pmatrix} \chi_s, \quad \epsilon_n = \text{sign}(n) \hbar \omega_c \sqrt{|n|}$$

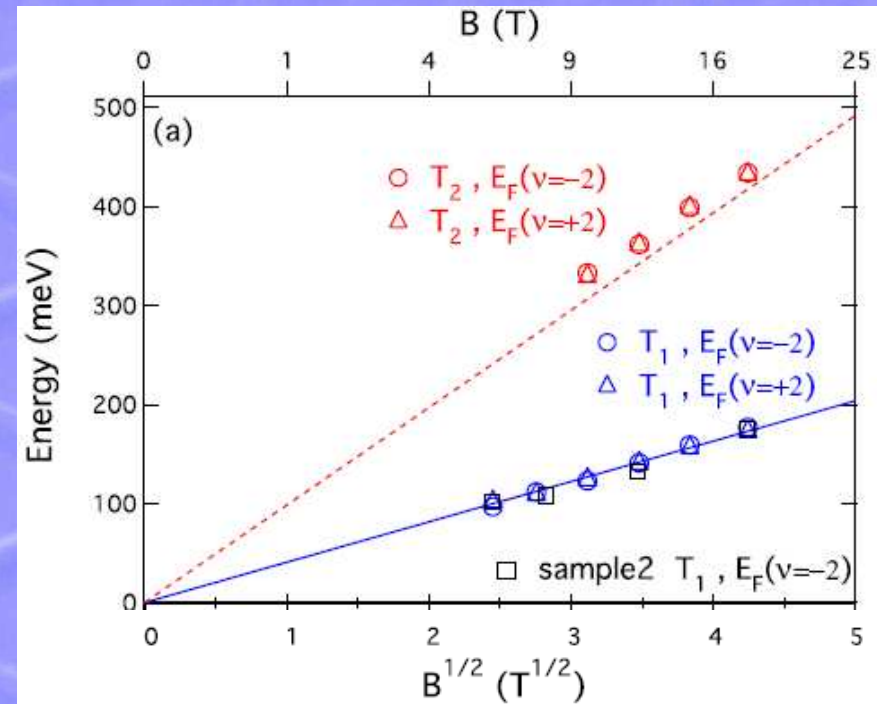
Quantum numbers: $n = \dots, -1, 0, 1, \dots$, $m = 0, 1, \dots$, $\sigma = \uparrow, \downarrow$, $s = \uparrow, \downarrow$

- C.f. 2DEG

$$\epsilon_n = \hbar \frac{eB}{mc} \left(n + \frac{1}{2} \right)$$

Self Energy Corrections to Cyclotron Resonance

- In 2DEG Kohn's theorem applies -CR independent of e-e interactions
- Doesn't hold for graphene – e-e interactions very important
- LL energy renormalised by self energy corrections due to exchange with electrons below Fermi level
- Corrections to CR diverge logarithmically with the cut off



Seen experimentally: Z. Jiang *et al.*, PRL **98**, 197403 (2007)

Collective Excitations

$$Q_{N_1 N_2 M_z}^\dagger = \sum_{m_1, m_2=0}^{\infty} A_{N_1 N_2 M_z}(m_1, m_2) c_{N_1 m_1}^\dagger d_{N_2 m_2}^\dagger, \quad A_{N_1 N_2 M_z}(m_1, m_2) \sim \delta_{M_z, |n_1| - m_1 - |n_2| + m_2}$$

c^\dagger creates electron
 d^\dagger creates hole

Renormalised single
particle energies

Impurity interaction

$$\begin{aligned} \hat{H}_{N_1 N_2}^{N'_1 N'_2} = & \sum_{m=0}^{\infty} (\tilde{\epsilon}_{N_1} + V_{N_1 m}) c_{N_1 m}^\dagger c_{N_1 m} \\ & - \sum_{m=0}^{\infty} (\tilde{\epsilon}_{N_2} + V_{N_2 m}) d_{N_2 m}^\dagger d_{N_2 m} \\ & + \sum_{\substack{m_1, m_2 \\ m'_1, m'_2}} \left(W_{N_1 m_1 N'_2 m'_2}^{N_2 m_2 N'_1 m'_1} - W_{N_1 m_1 N'_1 m'_1}^{N'_2 m'_2 N_2 m_2} \right) c_{N'_1 m'_1}^\dagger d_{N'_2 m'_2}^\dagger d_{N_2 m_2} c_{N_1 m_1} \end{aligned}$$

Dynamical exchange interaction

Direct electron-hole attraction

Connection to 2DEG

$$W_{N_1 m_1' N_2 m_2'}^{N_1 m_1' N_2 m_2'} \equiv \langle \Phi_{N_1 m_1'} \Phi_{N_2 m_2'} | U_{ee} | \Phi_{N_1 m_1} \Phi_{N_2 m_2} \rangle = \delta_{s_1, s_1'} \delta_{\sigma_1, \sigma_1'} \delta_{s_2, s_2'} \delta_{\sigma_2, \sigma_2'} \mathcal{U}_{n_1 m_1 n_2 m_2}^{n_1' m_1' n_2' m_2'}$$

$$\begin{aligned} \mathcal{U}_{n_1 m_1 n_2 m_2}^{n_1' m_1' n_2' m_2'} &= 2^{\frac{1}{2}} (\delta_{n_1 0} + \delta_{n_1' 0} + \delta_{n_2 0} + \delta_{n_2' 0} - 4) \left[U_{|n_1| m_1 |n_2| m_2}^{|n_1'| m_1'| |n_2'| m_2'} + s_{n_1} s_{n_1'} U_{|n_1| -1 m_1 |n_2| m_2}^{|n_1'| -1 m_1' |n_2'| m_2'} \right. \\ &\quad \left. + s_{n_2} s_{n_2'} U_{|n_1| m_1 |n_2| -1 m_2}^{|n_1'| m_1'| |n_2'| -1 m_2'} + s_{n_1} s_{n_2} s_{n_1'} s_{n_2'} U_{|n_1| -1 m_1 |n_2| -1 m_2}^{|n_1'| -1 m_1' |n_2'| -1 m_2'} \right] \end{aligned}$$

$$V_{nm} = 2^{\frac{1}{2}} (\delta_{n0} - 1) (V_{|n|m} + s_n V_{|n|-1m})$$

$$V_{0m} = V_{0m} = -(2m - 1)!! / 2^m m!$$

$$V_{1m} = \frac{1}{2} (V_{0m} + V_{1m}) = (8m - 3) / (8m - 4) V_{0m}$$

Energy units

$$E_0 = \sqrt{\frac{\pi}{2}} \frac{e^2}{\epsilon \ell_B}$$

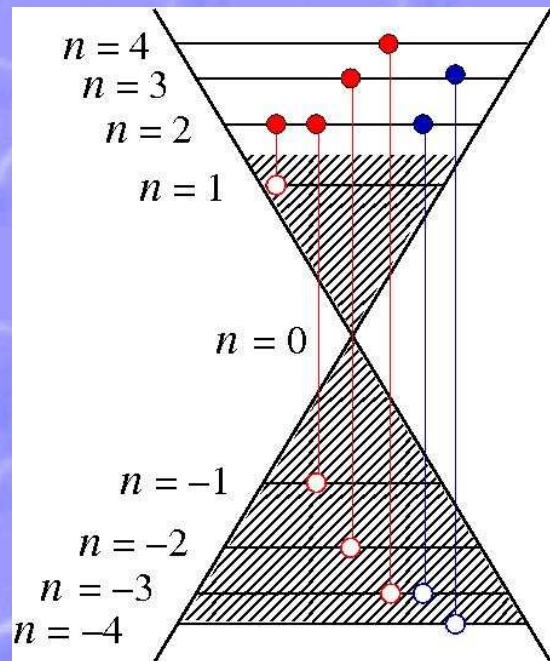
Examples

Which transitions are allowed?

In the dipole approximation, contribution to Hamiltonian from incoming circularly polarised light is

$$\delta H_{\pm} = \frac{ev_F \mathcal{E}}{i\omega c} e^{-i\omega t} \begin{pmatrix} \sigma_{\pm} & 0 \\ 0 & \sigma_{\mp} \end{pmatrix}$$

$\sigma_{\pm} = \sigma_x \pm i\sigma_y$ corresponds to left and right circularly polarised light



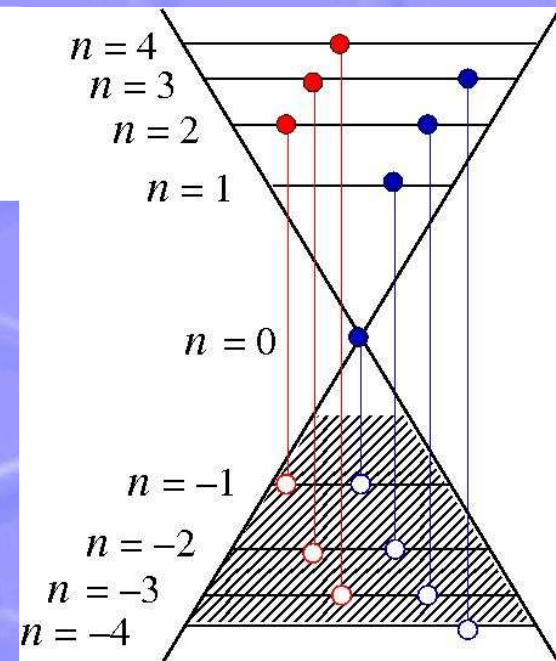
σ^+ σ^-

$$|n_{\text{final}}| - |n_{\text{initial}}| = \pm 1$$

$$m_{\text{final}} = m_{\text{initial}}$$

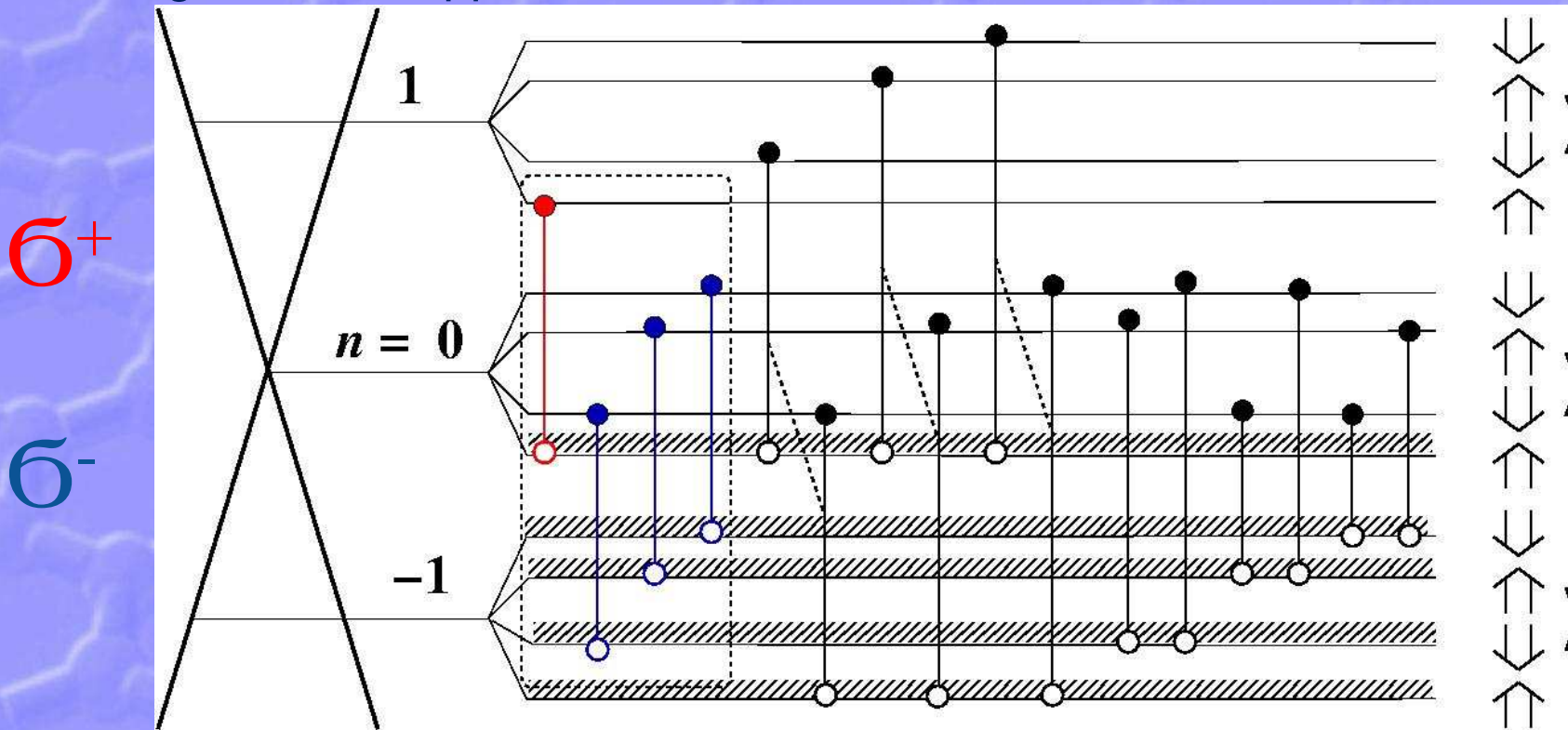
$$\sigma_{\text{final}} = \sigma_{\text{initial}}$$

$$S_{\text{final}} = S_{\text{initial}}$$



Which transitions should we take into account?

- Do calculations for sublevel filling factors of the zeroth LL.
- Infinitely many transitions with same M_z are mixed by Coulomb interactions
- Only consider mixing for those with the same energies – strong magnetic field approx



Numerical Considerations

- Matrix comes in 4x4 blocks
- Need to truncate at finite m – ok, because only seek localised states
- Size limit imposed by calculating matrix elements with high order Laguerre polynomials, not by diagonalisation
- Use $4 \times 50 = 200$ basis elements
- Results stable wrt changing matrix size

In the Absence of an Impurity

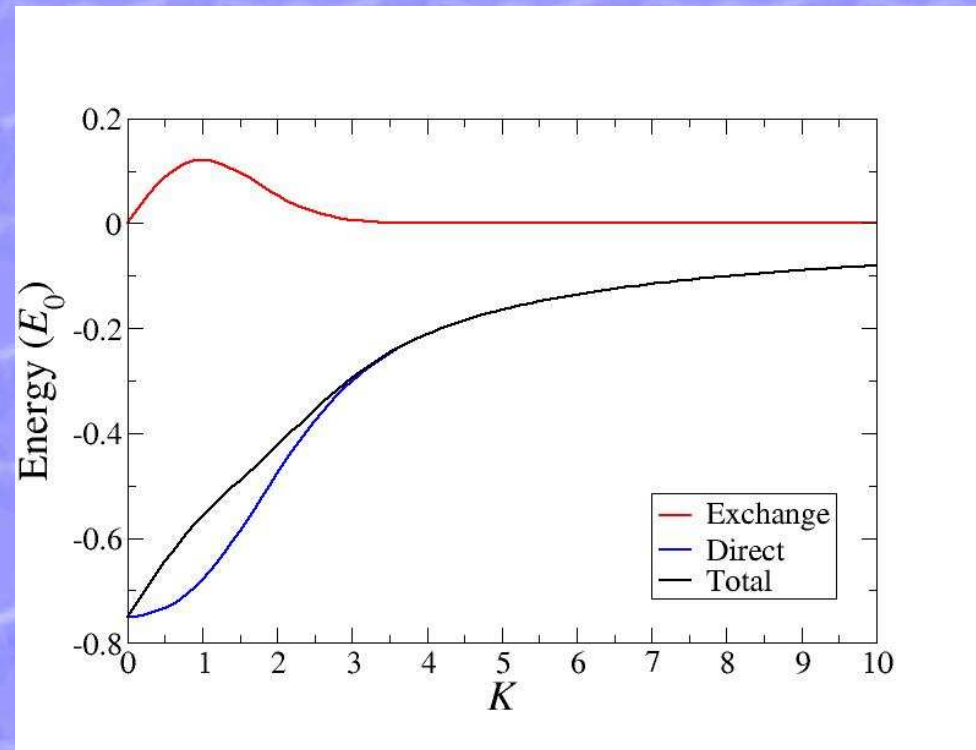
A. Iyengar *et al.* PRB **75**, 125430 (2007); Yu. A. Bychkov and G. Martinez, PRB **77**, 125417 (2008)

- Use Landau gauge

$$\mathbf{A} = -By\hat{\mathbf{x}}$$

- $m \rightarrow k_y$
- Thickness of band is $0.75E_0$ with

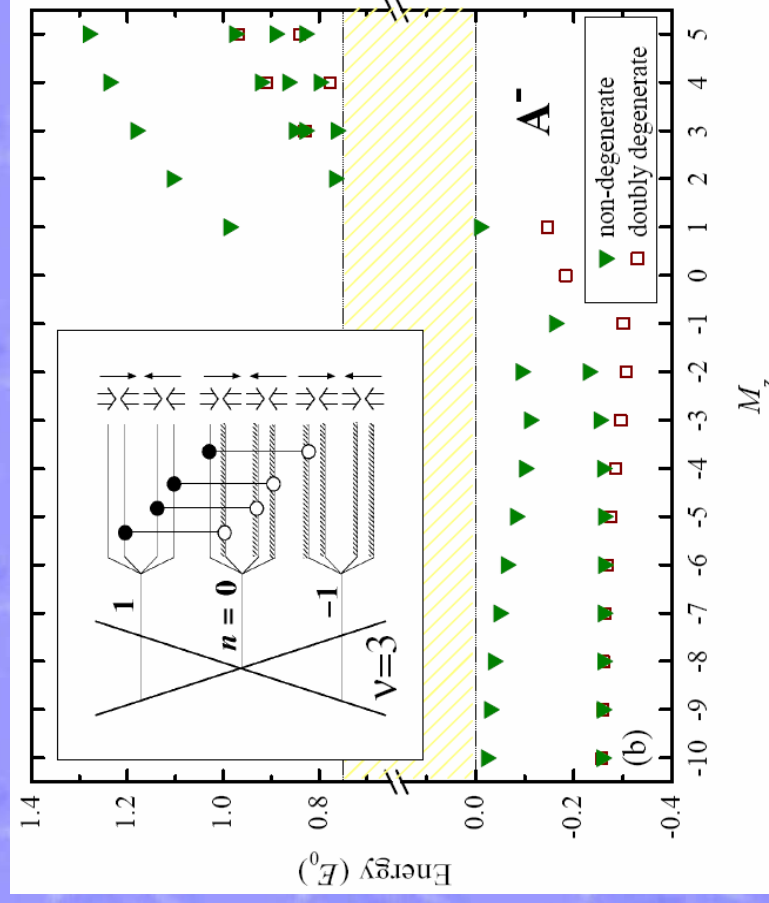
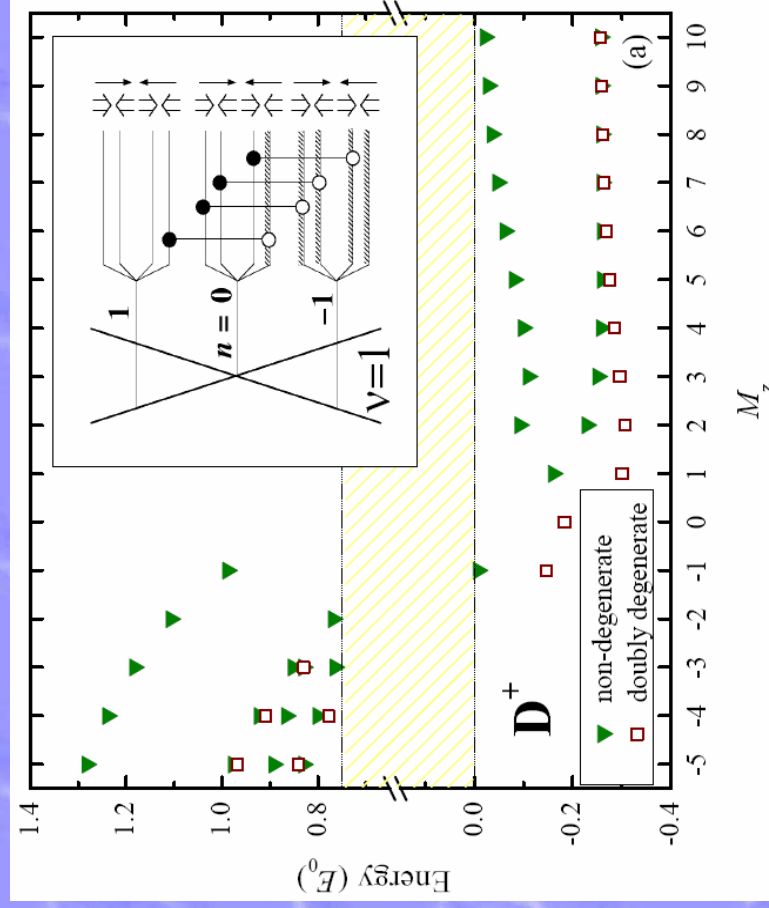
$$E_0 = \sqrt{\frac{\pi}{2}} \frac{e^2}{\epsilon \ell_B}$$



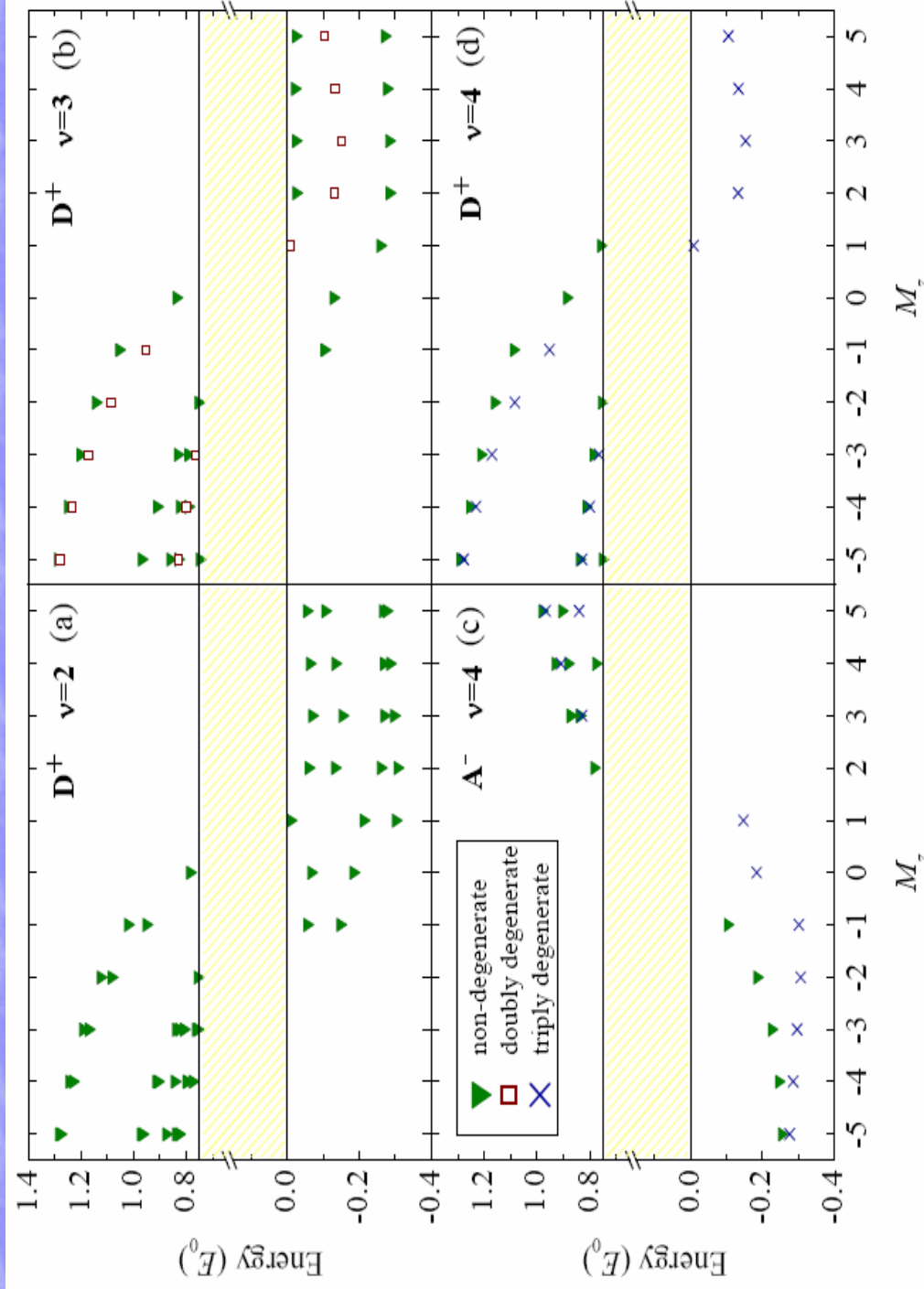
Results

Symmetry relation: Results same under

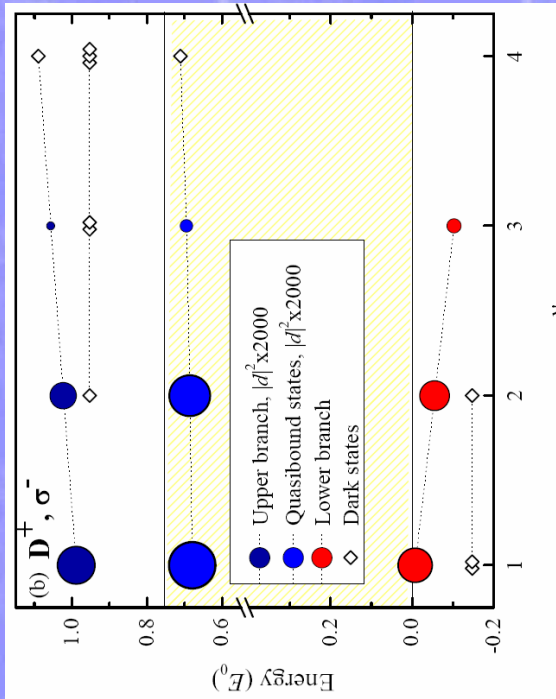
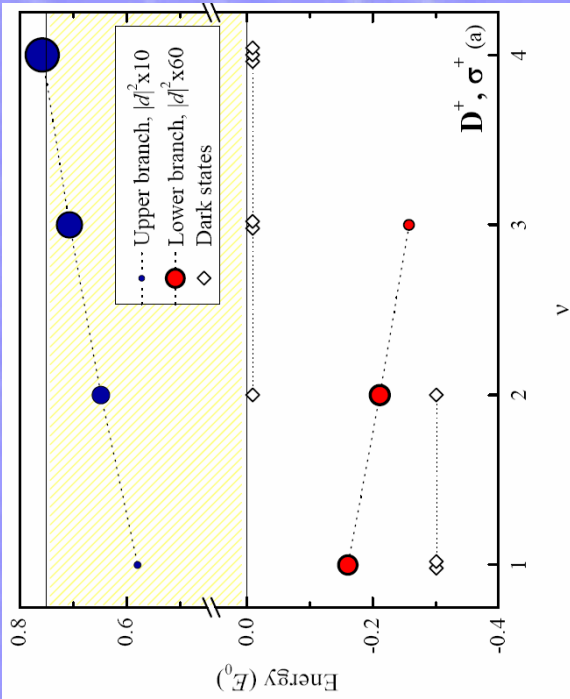
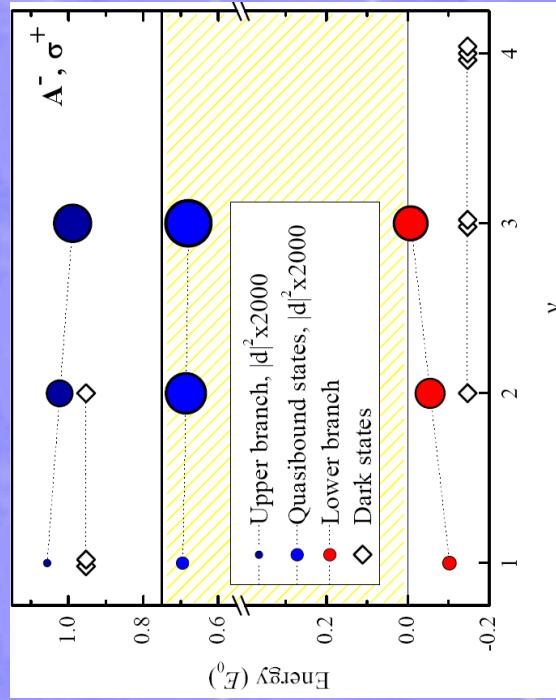
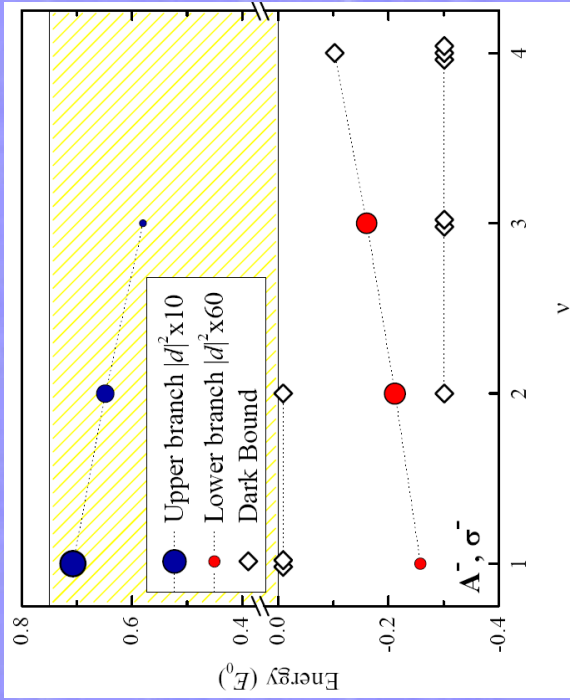
$$M_z \rightarrow -M_z, v \rightarrow v-4, D^+ \rightarrow A^-$$



Evolution with Filling Factor



SO(4)



Summary/Outlook

- Used a secondary quantised approach to determine collective excitations bound on an impurity in a single sheet of graphene in the presence of a strong perpendicular magnetic field [[arXiv:0902.4176](https://arxiv.org/abs/0902.4176)]
- Could our results be experimentally detected?
Only know relative strengths
Could think about:
 - rippling in graphene layer
 - intervalley scattering
 - nature of impurity
- First step towards combining work on magnetoplasma in pristine graphene and disordered graphene
- Could they enhance understanding of the nature of disorder in graphene?
- Further work:
Non integer filling factors: consider opposite limit of low electron density, so D- states are formed.

Thank you for listening



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Other work:

- Anderson transition in BCC and FCC lattices using transfer-matrix method [PRB **77**, 245117-8 (2008)]
- Excitonic storage in an Aharonov-Bohm nanoring with applied in-plane electric field [PRL **102**, 096405 (2009)]