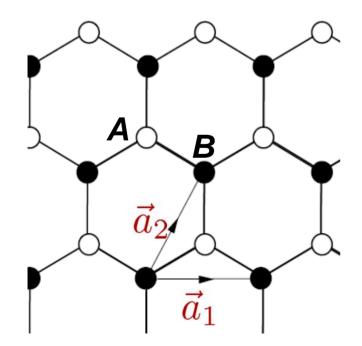
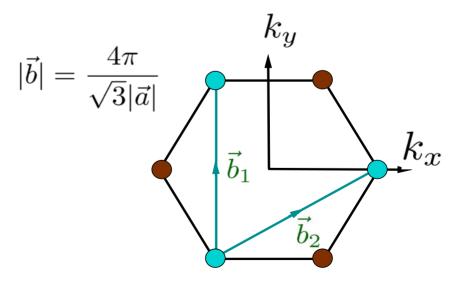
Transport properties of graphene.

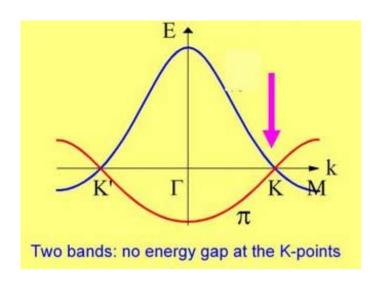
- 1. Effect of disorder on transport in graphene.
 - I.L. Aleiner and K.B. Efetov, PRL, 97, 236801, (2006)
- 2. Quantum dots in graphene.
- P.G. Silvestrov and K.B. Efetov, PRL, 98, 016802 (2007)

Lattice



Inverse lattice.





Hamiltonian H_0 for free electrons

$$H_0 = -iv\vec{\Sigma}\vec{\nabla}$$
 v-velocity

$$\Sigma_{x,y} = 1^{K,K'} \otimes \tau_{x,y}^{A,B}, \qquad \tau_{x,y,z}$$
 -Pauli matrices

A,B-sublattice space, K,K' -valley space

Full Hamiltonian (no spin): $H = H_0 + \hat{V}$

The external potential \hat{V} is generally a 4x4 matrix. It is proportional to unity matrix only if it varies slowly on the lattice period.

Two problems:

- 1. \hat{V} is disorder potential (full matrix)
- 2. \hat{V} is a barrier across the graphene strip (unity matrix).

New results for 1).

- a) Most general structure of \hat{V} is identified (using most general symmetries).
- b) Going beyond the self-consistent Born approximation (SCBA) (derivation and solution of RG equations).
- c) derivation of a supermatrix non-linear σ -model (ultimate localization but a complicated crossover).

Symmetries of the model:

1. Time reversal (TR) symmetry (exact)

$$\varphi^{T}(r) = ((\varphi_{A}, \varphi_{B})_{AB}; (\varphi_{B}^{*}, -\varphi_{A}^{*})_{AB})_{K,K'}$$

$$\varphi^*(r) = \hat{z}\varphi(r), \quad \hat{z} = \tau_y^{A,B} \otimes \tau_y^{K,K'}$$

$$H = \hat{z}H^T\hat{z}$$

Disorder may break all the symmetries except the time reversal one!

The most general form:

$$\hat{V}(r) = u_0(r) + \sum_{\{m,i\} = \{x,y,z\}} \hat{G}_{m,i} u_{m,i}(r), \qquad \hat{G}_{m,i} = \tau_m^{K,K'} \otimes \tau_i^{A,B}$$

 $u_0(r), u_m(r)$ are real independent random functions

However, averaging must restore rotation, reflection,

translation and C_{6v} symmetries!



$$\begin{split} \langle \hat{V}(\mathbf{r}_{1}) \otimes \hat{V}(\mathbf{r}_{2}) \rangle &= \delta(|\mathbf{r}_{1} - \mathbf{r}_{2}|) \,\, \gamma_{0} \,\, \mathbb{1} \otimes \mathbb{1} \\ &+ \delta(|\mathbf{r}_{1} - \mathbf{r}_{2}|) \,\, \gamma_{\parallel} \,\, \hat{G}_{z,z} \otimes \hat{G}_{z,z} \\ &+ \delta(|\mathbf{r}_{1} - \mathbf{r}_{2}|) \,\, \gamma_{\perp} \, \sum_{j=x,y} \hat{G}_{z,j} \otimes \hat{G}_{z,j} \\ &+ \delta(|\mathbf{r}_{1} - \mathbf{r}_{2}|) \,\, \beta_{\parallel} \, \sum_{m=x,y} \hat{G}_{m,z} \otimes \hat{G}_{m,z} \\ &+ \delta(|\mathbf{r}_{1} - \mathbf{r}_{2}|) \,\, \beta_{\perp} \, \sum_{j,m=x,y} \hat{G}_{m,j} \otimes \hat{G}_{m,j} \end{split}$$

Five independent real constants for the disorder.

 γ_0 is due to long range impurities (with respect to the lattice period)

 β -intervalley scattering

4

Impurities conserving the chirality.

$$\hat{V}(\mathbf{r}) = -\hat{\Sigma}_z \hat{V}(\mathbf{r}) \hat{\Sigma}_z \longrightarrow \gamma_0 = \gamma_{\parallel} = \beta_{\parallel} = 0$$

Consequence: One delocalized state exist at E=0! Minimum metallic conductivity at the center.

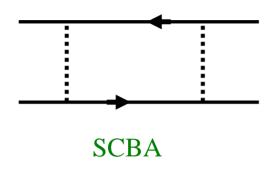
(See review: Altland, Simons, Zirnbauer, Phys. Rep. (2002)

However: no reason for this symmetry!

Moreover, all weak coupling constants grow in the process of the renormalization!

One should go beyond SCBA.

Scattering amplitudes in perturbation theory.

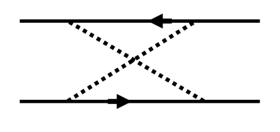


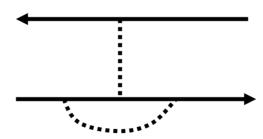
$$G_{\varepsilon}(p) = \frac{1}{\varepsilon - p\Sigma + i\delta}$$

$$\int G_{\varepsilon}(p)G_{\varepsilon}(p)d^{2}p \propto \ln(J/\varepsilon)$$

Logarithmic contributions to the amplitudes!

However, other logarithmic contributions exist!





SCBA fails: RG treatment.

Physical quantities as integrals over supervectors ψ (averaging has been performed)

$$\langle ... \rangle = \int ... \exp(-L[\psi])D\psi$$
, $L[\psi] = L_0[\psi] + L_{int}[\psi]$

$$L_{0} [\psi] = i \int \bar{\psi} \left[\varepsilon - \hat{\mathbb{H}}_{0} - \hat{\Lambda} \left(\frac{\omega}{2} + i0 \right) \right] \psi d\mathbf{r},$$

$$L_{int} [\psi] = \frac{1}{2} \int \left[\gamma_{0} \left(\bar{\psi} \psi \right)^{2} + \Gamma_{m}^{i} \left(\bar{\psi} \hat{\mathbb{G}}_{m,i} \psi \right)^{2} \right]$$

$$\Gamma_{z}^{z} = \gamma_{z}, \ \Gamma_{z}^{\{x,y\}} = \gamma_{\perp}, \ \Gamma_{\{x,y\}}^{z} = \beta_{z}, \ \Gamma_{\{x,y\}}^{\{x,y\}} = \beta_{\perp}$$

$$\Gamma_z^z = \gamma_z, \ \Gamma_z^{\{x,y\}} = \gamma_\perp, \ \Gamma_{\{x,y\}}^z = \beta_z, \ \Gamma_{\{x,y\}}^{\{x,y\}} = \beta_\perp$$

RG treatment: $\psi = \psi_0 + \widetilde{\psi}$

Slow **Fast**

has 16 components for the density of states

Integration over the fast field.

Reproducing the form of the Lagrangian: RG equations

Assumption:
$$\gamma_0 \gtrsim \gamma_{\parallel,\perp}, \beta_{\parallel,\perp}$$

New coupling constants:
$$egin{array}{ll} g_{\parallel} &= \gamma_{\parallel} + 2\gamma_{\perp}; & g_{\perp} &= \beta_{\parallel} + 2\beta_{\perp} \\ \delta g_{\parallel} &= \gamma_{\parallel} - \gamma_{\perp}; & \delta g_{\perp} &= \beta_{\parallel} - \beta_{\perp} \end{array}$$

Equations:

$$2\pi v \partial_t v = -\left(\gamma_0 + g_{\parallel} + 2g_{\perp}\right);$$

$$9\pi v^2 \partial_t \gamma_0 \approx 2\left(g_{\parallel}^2 + 2g_{\perp}^2\right);$$

$$9\pi v^2 \partial_t g_{\parallel} \approx -8g_{\parallel}^2 - 20g_{\parallel}g_{\perp} + 14g_{\perp}^2;$$

$$9\pi v^2 \partial_t g_{\perp} \approx \underline{4g_{\parallel}g_{\perp}} - 18g_{\perp}^2.$$

$$\pi v^2 \partial_t \delta g_{\parallel,\perp} \approx -3\gamma_0 \delta g_{\parallel,\perp};$$

Absence of intervalley scattering is not stable

Solution of the RG equations:

$$v\left(\varepsilon\right) = \left(\frac{\gamma_{0}}{\pi} \ln \frac{|\tilde{\varepsilon}|}{\varepsilon_{0}}\right)^{1/2}, \quad \gamma_{0}\left(\varepsilon\right) = \gamma_{0} + \mathcal{O}\left(\frac{g_{\parallel}(\varepsilon)}{\gamma_{0}}\right)$$

$$g_{\parallel}(\varepsilon) \approx g_{\perp}(\varepsilon) \approx \frac{9\gamma_{0}}{14 \ln \left[t^{*}/\ln |\tilde{\varepsilon}|/\varepsilon_{0}\right]}, \quad \boxed{\varepsilon_{0} \approx J \exp\left(-\pi v_{0}^{2}/\gamma_{0}\right)}$$

$$\tilde{\varepsilon} = \max(\varepsilon, \varepsilon_{0})$$

 \mathcal{E}_0 -is the energy at which the 1-loop approximation breaks down

All coupling constants are important. No chance for moving to chiral disorder!

Physical quantities with the "ultraviolet" logarithmic renormalization

Diffusion coefficient:

$$D(\epsilon) = \frac{v^2 \tau_{tr}(\epsilon)}{2}; \quad \frac{1}{\tau_{tr}(\epsilon)} = \frac{\pi \gamma_0 \nu(\epsilon)}{4}$$

Density of states:

$$\nu(\epsilon) = \frac{|\tilde{\varepsilon}|}{\pi \hbar^2 v^2(\epsilon)}$$

Conductivity:

$$\sigma(\epsilon) = 2e^2 \nu(\epsilon) D(\epsilon) = \frac{4e^2}{\pi^2 \hbar} \ln \left(\frac{|\tilde{\varepsilon}|}{\varepsilon_0} \right)$$

However, this is not the end of the story: localization effects!

Non-linear supermatrix σ -model for describing localization effects

General scheme: following "Supersymmetry in disorder and chaos", K.B. Efetov, Cambridge University Press, (1997)

$$\langle O \rangle = \int O(\hat{Q}) \exp\left(-F\left[\hat{Q}\right]\right) \mathcal{D}\hat{Q}$$

For any correlation function O

$$1 = \int \exp\left(-F\left[\hat{Q}\right]\right) \mathcal{D}\hat{Q}$$

Due to supersymmetry

For graphene: Q are 16x16 supermatrices (unity in the sublattice space), $Q^2 = 1$

Free energy functional F

$$F = \frac{\pi\hbar\nu(\epsilon)}{16}Str \int \left\{ D(\epsilon) \left(\nabla \hat{Q} \right)^2 + 2i\omega \hat{\Lambda} \hat{Q} \right.$$
$$- \frac{\pi\nu(\epsilon)g_{\parallel}(\epsilon)}{4} \left[\hat{\rho}_z, \hat{Q} \right]^2 - \frac{\pi\nu(\epsilon)g_{\perp}(\epsilon)}{4} \left(\left[\hat{\rho}_x, \hat{Q} \right]^2 + \left[\hat{\rho}_y, \hat{Q} \right]^2 \right) \right\} d\mathbf{r};$$
$$\hat{\rho}_j = \tau_j^{KK'} \otimes \mathbb{1}^{eh} \otimes \mathbb{1}^{AR} \otimes \mathbb{1}^g; \quad \hat{\Lambda} = \mathbb{1}^{KK'} \otimes \mathbb{1}^{eh} \otimes \tau_z^{AR} \otimes \mathbb{1}^g;$$

All the constants should be taken from the solution of the RG equations!

$$\begin{split} &D(\epsilon) = \frac{v^2(\epsilon)\tau_{tr}(\epsilon)}{2}; \quad \frac{1}{\tau_{tr}(\epsilon)} = \frac{\pi\gamma_0\nu(\epsilon)}{4} \\ &v\left(\varepsilon\right) = \left(\frac{\gamma_0}{\pi}\ln\frac{|\tilde{\varepsilon}|}{\varepsilon_0}\right)^{1/2}, \quad \gamma_0\left(\varepsilon\right) = \gamma_0 + \mathcal{O}\left(\frac{g_{\parallel}(\varepsilon)}{\gamma_0}\right) \\ &g_{\parallel}(\varepsilon) \approx g_{\perp}(\varepsilon) \approx \frac{9\gamma_0}{14\ln\left[t^*/\ln\left|\tilde{\varepsilon}\right|/\varepsilon_0\right]}, \qquad \tilde{\varepsilon} = \max\left(\epsilon, \#\varepsilon_0\right) \end{split}$$

1) If only diagonal disorder (γ_0) is present:

$$F = \frac{\pi v(\varepsilon)}{16} Str \int \left[D(\varepsilon) (\nabla Q)^2 + 2i\omega \Lambda Q \right] dr$$

$$\hat{Q} = \hat{C}\hat{Q}^T\hat{C}^T; \quad \hat{C} = i\tau_y^{KK'} \otimes \mathbb{1}^{AR} \otimes \left(\tau_-^{eh} \otimes \mathbb{1}^g - \tau_+^{eh} \otimes \tau_z^g\right)$$

The symmetry of Q corresponds to 2 replicas of the symplectic ensemble. Antilocalization!

The symplectic symmetry was first noticed by Suzuura and Ando (2002)

However, the terms with P_i break the symmetry and the ensemble becomes orthogonal!

At large distances:
$$\hat{Q} = \mathbb{1}_{KK'} \otimes \hat{Q}_o$$

-is 8x8 supermatrix of the orthogonal symmetry: localization!

Perturbation theory: first order gives the weak localization correction.

$$\Delta \sigma_{WL} = \frac{e^2}{2\pi^2 \hbar} \left[-\ln \frac{\tau_{\phi}}{\tau_{tr}} + 2\ln \frac{\tau_2}{\tau_{tr}} + \ln \frac{\tau_3}{\tau_{tr}} \right]$$

$$\frac{1}{\tau_2} = \frac{1}{\tau_\phi} + \frac{1}{\tau_\perp}; \ \frac{1}{\tau_3} = \frac{1}{\tau_\phi} + \frac{1}{2\tau_\parallel} + \frac{1}{\tau_\perp}; \qquad \tau_\phi \text{ -inelastic dephasing time}$$

Agrees with McCann et al (2006)

Ultraviolet logarithmic renormalization

$$\frac{\tau_{tr}}{\tau_1} \approx \frac{\tau_{tr}}{\tau_2} \approx \frac{36}{7 \ln[t^*/\ln|\tilde{\epsilon}/\varepsilon_0|]}$$

Qualitative picture (from Aleiner and Efetov (2006)).

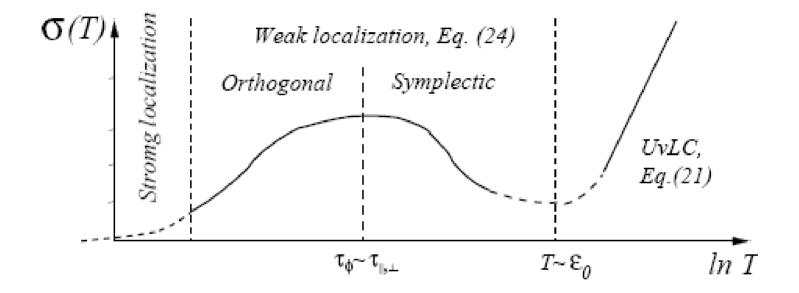


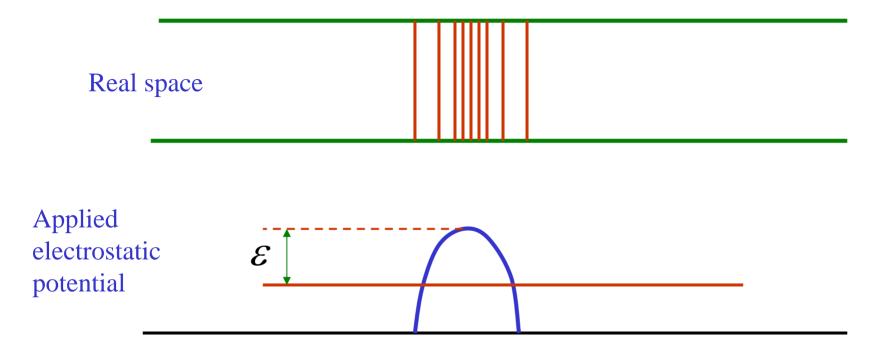
FIG. 2: Schematic dependence, $\sigma(T)$, for the undoped graphene and for $\tau_{\phi}^{-1} \propto T$.

How to make a quantum dot in graphene? Silvestrov and Efetov (2007)

Peculiarity of graphene with respect to conventional 2D electron gases: One barrier is sufficient!

Avoiding the reflectionless penetration (Klein paradox) through the barrier: making a strip with a barrier across it.

Transversal quantization removes the problem of the confinement!



Sufficient to make the quantum dot!

Quasiclassical treatment complemented by numerics.

Conductance as a function of the energy (gate voltage)

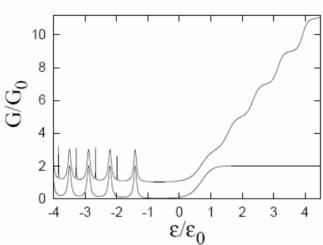


FIG. 1: Upper curve: Conductance of the graphene quantum dot as a function of the Fermi energy for the metallic armchair edges for $L=4\xi$, Eq. (4). The lower curve: the contribution to the conductance from the transmission channels with $p_y=\pm\pi\hbar/L$. All calculations are carried out for the zero temperature, T=0.

$$V = -\left(x/x_0\right)^2 U/2.$$

$$\xi = \left[\hbar c x_0^2 / U\right]^{1/3}$$
, $\varepsilon_0 = \hbar c / \xi$.

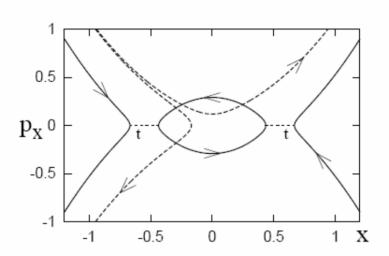


FIG. 2: Examples of trajectories described in the text drawn on the x, p_x plane (arbitrary units). Solid lines show the trajectories with $\varepsilon < 0$ either bouncing inside the barrier, or reflected by it from the left/right. Tunnelling events between the bounded and unbounded trajectories are shown schematically (t). Thick dashed lines show the trajectories with $\varepsilon > 0$ either transmitted for $|p_y| < \varepsilon/c$ (open channels) or reflected for $|p_y| > \varepsilon/c$ (closed channels).

Fabrication of quantum dots in graphene: essential detail for quantum computer! Experiments are being carried out!