Disordered Quantum Wires*

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*e=h=k_B=1

- Electrons in 3 dimensions
- Electrons in 1 dimension: what is different?
- Electrons in 1 dimension: experiments
- Quantum wire as a chain of quantum dots
- Long wires: Rare events
- Conclusions

1. Electrons in three dimensions

<u>Non-interacting electrons</u>: mass m, mean spacing a, Planck's constant \hbar

 \rightarrow energy scale $\hbar^2/ma^2 \sim E_F$

T << E_F : Pauli principle determines pressure $\rightarrow p \sim \hbar^2/ma^{d+2}$

Compressibility*: $\kappa = -a^{-2d} \partial \ln V / \partial p \sim 1/(E_F a^d)$

Specific heat : $C \sim T/E_F$



1. Electrons in three dimensions

Interacting spinless electrons:

charge e \rightarrow strength of interaction \sim (e²/a) / E_F \sim a/a_B!

but energy and momentum conservation reduce phase space ⇒ Landau's Fermi fluid of quasi-particles

with finite lifetime $\hbar/\tau \sim E_F(T/E_F)^2$

Disordered spinless electrons:

Metal insulator transition





2. Electrons in one dimension - what is different?



Electrons cannot avoid each other \rightarrow Landau's picture breaks down

 \Rightarrow density wave excitations : plasmons

(a) 1D clean wire: Luttinger liquid

 $G = dJ/dV = \sigma/L = (K) e^2/h$

(b)+ single impurity: K<1: impurity relevant,

 $ightarrow oldsymbol{G} \sim (extsf{max} (extsf{T}, extsf{eV}))^{2/ extsf{K-2}}$ Kane & Fisher '92, Furusaki & Nagaosa

K>1: impurity irrelevant

MoSe Nanowires

Venkataraman, PRL (2006)

J /T
$$^{\alpha+1}$$
 \sim max (V/T, V $^{\beta+1}$ /T $^{\alpha+1}$)





FIG. 1 (color online). (a) Structural model of a 7-chain MoSe nanowire along with the triangular Mo_3Se_3 unit cell. (b) and (c) AFM height images of MoSe nanowires between two Au electrodes. The wire heights are 7.2 nm and 12.0 nm, respectively. Scale bar = 500 nm.

short wires (L ~ 1 μ m):

"Temperature" Exponent (a) is close to "Voltage" Exponent (b) Agrees with the conventional "Luttinger-liquid" picture with

α=β=2/K-2

•<u>Multiwall carbon nanotubes</u>





conductance



FIG. 3. AFM images of junctions formed between two MWNTs: (a) end-bulk junction and (b) end-end junction. The arrows indicate the position of the junctions.

Variable Range Hopping conduction in polydiacetylene single crystals

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Polymer nanofibers

long wires: 10µm



FIG. 1 (color online). AFM image of a *R*-hel PA fiber; the inset shows the schematic of a two-probe device based on such a *R*-hel PA fiber on top of Pt electrodes.



"temperature" exponent **exceeds** "voltage" exponent!

J /T
$$^{\alpha+1}$$
 \sim max (V/T, V $^{\beta+1}$ /T $^{\alpha+1}$)

Disagrees with the conventional Luttinger-liquid picture

Observed Power-Law Exponents

L ~10 µm polymers	Sample	1	2	3	4	5	6
Aleshin et al.,	α	2.8	5.5	7.2	5.6	5.0	4.1
PRL (2004)	β	1.5	3.8	4.7	1.0	1.1	1.8

L ~ 100 µm InSb wires Zaitsev-Zotov et al., JPCM (2000)

Sample	1	2	3	4
α	2.3	3.4	4.5	4.6
β	1.3	3.4	2.8	2.0

• In long wires T-exponent exceeds V-exponent: $\alpha > \beta$

• Exponents are sample-dependent

4. Quantum wire as a chain of quantum dots



strong impurities, randomly (Poissonian) distributed

V=0 : <u>classical ground state</u> (K<<1)

integer number q_i of electrons between impurities i and i+1

$$q_i = [Q_i]_G$$
, $Q_i = a_i (2\pi k_F + \mu C)$ background charge



Transport from dot j to dot k by activation and <u>co-tunneling</u>



Transport from dot j to dot k by activation and <u>co-tunneling</u> $J_{jk} \sim e^{-E_{jk}/T} e^{-s|k-j|} \sinh((\xi_k - \xi_j)/T)$

Ohmic regime

typical energy mismatch $E_{jk} \sim \Delta / |k-j| \rightarrow |k-j| \sim (\Delta/Ts)^{1/2} \equiv x_{VRH}/a$

$$\Rightarrow$$
 J \sim e ^{(s Δ /T)^{1/2} F, Ts<< Δ VRH}

Non-ohmic regime

 $\begin{array}{ll} \text{typical } |\text{k-j}| aF \sim \Delta / |\text{k-j}| & \rightarrow & |\text{k-j}| \sim (\Delta/aF)^{1/2} \\ \\ \Rightarrow & \mathbf{J} \sim e^{s(\Delta/aF)^{1/2}}, & aF << \Delta & VRH \end{array}$

Large Voltage/temperature: power laws



Larkin and Lee '78

<u>Tunneling action</u> dominated by spreading of charge:

 $S_{tun} \sim \int d\tau \ E(\tau) \sim \int d\tau \ (C \ x(\tau))^{-1} \sim K^{-1} \ In \ (E_{initial}/E_{final})$

Large Voltage/temperature: power laws



Larkin and Lee '78

<u>Tunneling action</u> dominated by spreading of charge:

$$\begin{split} S_{tun} &\sim \int d\tau \; E(\tau) \sim \int d\tau \; (\mathcal{C} \; x(\tau))^{-1} \sim -K^{-1} \; ln \; (E_{final} / E_{initial}) & \frac{E_{initial} = k_F / \mathcal{C}}{E_{final} = max \; (\Delta, T, V)} \end{split}$$

 $s_{eff} \approx 2K^{-1} \ln \left[ak_{F}/max\left(1,T/\Delta,V/\Delta\right)\right] \approx s - 2K^{-1} \ln \left(max(T,V)\right) \text{ if } max(T,V) > \Delta$

 \rightarrow T > Δ : J_{k,k+1} \sim (max(T,V)/ Δ)^{2/K} Kane-Fisher 1992



 $\xi_{\text{loc}} \approx a/s$

Strong and weak pinning

T>T_{1,cr}: single impurity weak

 $T < T_2$: collective effects

Giamarchi, 2004: "Quantum Physics in One Dimension"



 $u < k_F$: SCHA $u \rightarrow u_{eff}$

4. Long wires: rare events

□ So far considered: typical quantum dots with $a_i \approx a$ □ Now: consider regions with many narrow dots with $a_i \ll a$



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Wire: non-overlapping breaks + low resistance connecting pieces

break resistance $R_0e^u \rightarrow R=L\int P_u(u) R_0e^u du$

 $P_u(u)$: prob. /unit length that break resistance at least e^u

 $\mathbf{I}_{jk} \sim e^{-s|j-k|} e^{-E_{jk}/T} \quad sinh ((\xi_j - \xi_k)/T) \qquad \qquad \xi_j = \mathbf{F} \mathbf{x}_j - \mu_j$

(a) Ohmic break of |j-k| dots

 $envelope: \quad \mathsf{E}_{\mathsf{i}}(\pm 1) \geq \epsilon_{\mathsf{i}}^{\pm} \ \rightarrow \ \mathsf{s}|\mathsf{j}\mathsf{-}\mathsf{k}| + \mathsf{max}(\epsilon_{\mathsf{j}}^{\pm}, \epsilon_{\mathsf{k}}^{\pm})/\mathsf{T} \geq \mathsf{u}$

simplification: rectangular break: s|j-k| \sim u, $\epsilon_{j}{}^{\pm}\!\sim$ uT



$$\begin{split} P_u(u) &\sim x_{VRH}^{-1} P_\epsilon(\epsilon)^{|j-k|} \sim x_{VRH}^{-1} P_\epsilon(uT)^{u/s} \begin{cases} & \ln P_u(u) \sim - u^2 T/\Delta \, s & \text{if } uT \ll \Delta \\ & \ln P_u(u) \approx - u/s \, \ln \left(2uT/\Delta\right) & \text{if } uT \gg \Delta \end{cases} \end{split}$$

 $\label{eq:lnP_u} \ln P_u(u) \sim - \ u^2 T / \Delta \ s \qquad \qquad \mbox{if } u T \mbox{--} \Delta \ s$

In P_u(u) \approx - u/s In (2uT/ Δ) if uT>> Δ



\rightarrow <u>infinite wire:</u>

Low T , s \ll 1 : $\rho \sim \exp (s\Delta/4T)$

Raikh Ruzin '89

High T, s \gg 1 : $\rho \sim \exp(s\Delta e^{(s-1)}/2T)$ new

finite wire:
$$P_u(u)(L/x_{VRH}) \approx 1 \rightarrow u_{max}$$

 $\label{eq:stimul} \text{sTln}(\text{L/x}_{\text{VRH}}) \ll \Delta : \qquad \rho \sim ~ \text{exp} ~ [~ \text{s} \Delta ~ \text{ln} ~ (\text{L/a}) ~ / ~ \text{T} ~]^{1/2} ~ ~ \text{VRH}$ Raikh Ruzin '89

 $sT \ln(L/x_{VRH}) \gg \Delta$:

$$\label{eq:rho} \begin{split} \rho &\sim ~ exp[s \, ln \, (L/x_{VRH}) \, / \, ln \, (Ts/\Delta)] \sim T^{\alpha} \\ \alpha &= s \, ln \, (L/x_{VRH}) \, / \, ln^2 \, (Ts/\Delta)] \end{split}$$

(b) <u>Non-Ohmic breaks of m dots</u> $V_b = \xi_i - \xi_k \gg T$ $T u_I + V_b/2$ $V_b/2$ $V_b/2$ - $V_b/2$ - $T u_I - V_b/2$ ϵ^- (main voltage drop across the breaks, but constant I = $I_0 e^{-u}$ current everywhere) $I_{ik} \sim e^{-s|j-k|} e^{-E_{jk}/T} \sinh((\xi_i - \xi_k)/T)$ $\rightarrow s|j-k| + max(\epsilon_i^{\pm}, \epsilon_k^{\pm})/T - V_b/2T \ge u$ <u>average electric field</u>: $V/L=F = \int dV_b V_b P_V(V_b) = x_{VRH}^{-1} 2Ts (Tu/\Delta)^{-u/s}$ $P_V(V_b) = x_{VRH}^{-1} [P_{\epsilon}(uT+V_b/2)]^{u/s} = x_{VRH}^{-1} (Tu/\Delta)^{-u/s} e^{-V_b/2Ts}$ V_⊾≪ u_IT u=-ln (I/I₀)= s ln (2Ts/Fx_{VRH})/ ln(Ts/ Δ)

 $I \sim V^{\beta}$, $\beta = s / \ln (Ts/\Delta)$, $\alpha/\beta = \ln(L/x_{VRH}) / \ln(Ts/\Delta) > 1$

Regime diagram



Conclusions:

- □ linear and non-linear conductivity
- field and temperature cross-over between single and many impurity tunneling
- Iow field and temperature: Mott-Shklovskii-VRH
- □ larger E,T: Kane-Fisher power law behavior
- Kane-Fisher "single-dominant-barrier" theory is not valid in long wires that contains many (> 100 ?) impurities
- true power-law exponents exceed the single-barrier ones by a "large" log-factor (perhaps, by 2 or 3 in practice)!
- resistance is controlled by "difficult spots" dense clusters of impurities
- global weak/strong pinning regime diagram