

Boltzmann equation in the presence of disorder

(47)

(Note that we consider scalar model! ~~H~~ H is not a matrix)

self-consistent

Born

$$\sum_{(1,2)} = \underbrace{\begin{array}{c} x_1 \xrightarrow{\text{---}} x_2 \\ | \quad \diagup \end{array}}_{\text{which already includes } \sum} = \langle v(1) \overset{1}{G}(1,2) v(2) \rangle_{\text{dis}}$$

↓
no convolutions
here!!!

$$\langle \dots \rangle_{\text{dis}} = \frac{1}{V^{N_i}} \int dR_1 \int dR_2 \dots \int dR_{N_i} \dots$$

R_1, R_2, \dots, R_{N_i} ← random positions of impurities

Impurity potential $V(r) = \sum_{i=1}^{N_i} \delta(r - R_i)$

$\delta(r)$ - impurity profile!

let's assume $\langle V \rangle_{\text{dis}} = 0$ (can always be achieved by subtracting a constant (for any given N_i))

Compute first

$$\langle V(r_1) V(r_2) \rangle_{\text{dis}} = \sum_{i,j} \langle \delta(r_1 - R_i) \delta(r_2 - R_j) \rangle_{\text{dis}} \quad (= 0 \text{ unless } i=j!)$$

since $\langle \sum \delta(r - R_i) \rangle = 0$

$$= \sum_{i=1}^{N_i} \langle \delta(r_1 - R_i) \delta(r_2 - R_i) \rangle_{\text{dis}} = \frac{N_i}{V} \int_V d^3 R \delta(r_1 - R) \delta(r_2 - R)$$

Volume

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$$\frac{N_i}{V} = n_{\text{imp}} - \text{impurity concentration:}$$

$$\Rightarrow \langle v(r_1) v(r_2) \rangle = n_{\text{imp}} \int d^3 R \, v(r_1 - R) \, v(r_2 - R) \quad (1)$$

use Fourier transform $v(r) = \int \frac{d^3 p}{(2\pi)^3} e^{ip \cdot r} v_p = \int \frac{d^3 p}{(2\pi)^3} e^{-ip \cdot r} v_p^*$

Since $v(r)$ - real potential

$$(1) \quad n_{\text{imp}} \underbrace{\int d^3 R}_{\int \frac{d^3 p_1}{(2\pi)^3} \int \frac{d^3 p_2}{(2\pi)^3}} e^{i p_2 \cdot r_2 - i p_1 \cdot r_1} v_{p_2} v_{p_1}^* \underbrace{e^{-i R \cdot (p_2 - p_1)}}_{}$$

$$\text{gives } (2\pi)^3 \delta(p_1 - p_2)$$

$$= n_{\text{imp}} \int \frac{d^3 p_1}{(2\pi)^3} e^{-ip_1 \cdot (r_1 - r_2)} |v_{p_1}|^2$$

$$\Sigma(1,2) = \underbrace{n_{\text{imp}} \int \frac{d^3 p_1}{(2\pi)^3} e^{-ip_1 \cdot (r_1 - r_2)} |v_{p_1}|^2}_{} G(1,2)$$

We need Wigner transform of this!

$$\sum(\varepsilon, p, t, R) = \int d(t_1 - t_2) e^{i\varepsilon(t_1 - t_2)} \underbrace{\int d^3(\gamma_1 - \gamma_2) e^{-iP(\gamma_1 - \gamma_2)}}_{\times} \times \\ \times n_{\text{imp}} \underbrace{\int \frac{d^3 p_1}{(2\pi)^3} e^{-ip_1(\gamma_1 - \gamma_2)}}_{|v_{p_1}|^2} G(t_1, t_2, p_2, R) e^{iP_2(\gamma_1 - \gamma_2)}$$

Integration over $(\gamma_1 - \gamma_2)$ gives $(2\pi)^3 \delta(p + p_1 - p_2)$

=

$$\Rightarrow p_1 = p_2 - p,$$

 \Rightarrow

$$\hat{\sum}(\varepsilon, p, t, R) = n_{\text{imp}} \int \frac{d^3 p^1}{(2\pi)^3} |v_{p-p^1}|^2 \hat{G}(\varepsilon, p^1, R, t)$$

all valid for matrices Γ_n

the zotated keldysh space!

$$\hat{\Sigma} = \begin{pmatrix} \Sigma^R & \Sigma^K \\ 0 & \Sigma^A \end{pmatrix}$$

$$\hat{G} = \begin{pmatrix} G^R & G^K \\ 0 & G^A \end{pmatrix}$$

Let us stick to ~~the~~ white-noise disorder for simplicity, i.e. $\nu_p = \nu_0 = \text{const}(\rho)$

$$\Rightarrow \hat{\Sigma} = \hat{\Sigma}(\varepsilon, t, R) = \underbrace{n_{\text{imp}} \nu_0^2}_{\begin{array}{c} \not{=} \\ \text{no } \rho \\ \text{dependence} \end{array}} \int \frac{d^3 p'}{(2\pi)^3} \hat{G}(\varepsilon, p'; R, t)$$

small parameter

We also ignore $\text{Re } \Sigma$ (spectrum shift)
 Then $\Sigma^R = -i\gamma$, $\Sigma^A = i\gamma$

This can be done more accurately by solving the Exercise on page 45

Remember:

$$\text{equation: } (\partial_t - (\nabla_R U) \nabla_p + (\nabla_p \varepsilon) \nabla_R) G^{R,A} = 0$$

this is solved by $G^{R,A} = \frac{1}{\varepsilon - \varepsilon_p - U(R) \pm i\gamma}$

However, when we compute $\hat{\Sigma}$ in collision integral $[\hat{\Sigma}, G]$ let us ignore the perturbation $U(R)$

i.e. disregard the driving force $\vec{F} = -\nabla_R U$ in the collision integral (standard approximation)

For Keldysh component we have:

$$(2_t - (\nabla_R V) \nabla_p + (\nabla_p \epsilon_p) \nabla_R) G^K = i(\Sigma^K (G^R - G^A) - G^K (\Sigma^R - \Sigma^A))$$

introduce distribution function by the integral

$$h_p = \oint d\epsilon \frac{G^K(\epsilon, t, R, p)}{-2\pi i}$$

also depends on
R and t



This correspond to the Ansatz

Note that

$$\oint d\epsilon \frac{G^R - G^A}{-2\pi i} = 1$$

also true
for any $V(R)$

$$G^K = h_p (G^R - G^A)$$

\uparrow
 $h_p(R, t) \leftarrow$ no explicit energy dependence!

Now, let us see what we have for $\Sigma^K, \Sigma^R, \Sigma^A$:

$$\Sigma^K = \cancel{\text{without}} n_{\text{imp}} v_0^2 \int \frac{d^3 p'}{(2\pi)^3} h_{p'} (G^R - G^A)_{(p', R, t) \epsilon}$$

$$\Sigma^R - \Sigma^A = n_{\text{imp}} v_0^2 \int \frac{d^3 p'}{(2\pi)^3} (G^R - G^A)_{(\epsilon, p', R, t)}$$

divide by $(-2\pi i)$ & integrate over ϵ

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$$(\partial_t - (\nabla_R \cdot) \nabla_p + (\nabla_p \cdot \epsilon_p) \nabla_R) h_p =$$

$$= i n_{imp} v_0^2 \left(\int d\epsilon \int \frac{d^3 p'}{(2\pi)^3} \left(h_{p'} (G_{p'}^R - G_{p'}^A) \frac{\epsilon_{p'}^R - \epsilon_{p'}^A}{-2\pi i} - (G_{p'}^R - G_{p'}^A) \frac{h_{p'} (G_{p'}^R - G_{p'}^A)}{-2\pi i} \right) \right)$$

$$= 2\pi n_{imp} v_0^2 \underbrace{\int d\epsilon \int \frac{d^3 p'}{(2\pi)^3} (h_{p'} - h_p) \frac{\epsilon_{p'}^R - \epsilon_{p'}^A}{-2\pi i} \frac{\epsilon_p^R - \epsilon_p^A}{-2\pi i}}$$

$$\text{Im } \sum \rightarrow 0 \Rightarrow \frac{\epsilon_p^R - \epsilon_p^A}{-2\pi i} \rightarrow \delta(\epsilon - \epsilon_p)$$

Consider the limit $\gamma \rightarrow 0$

In this limit we find

$$(\partial_t - (\nabla_R \cdot) \nabla_p + (\nabla_p \cdot \epsilon_p) \nabla_R) h_p = 2\pi n_{imp} v_0^2 \underbrace{\int \frac{d^3 p'}{(2\pi)^3} (h_{p'} - h_p) \delta(\epsilon_p - \epsilon_{p'})}_\text{Lorentz-Golden rule limit}$$

$$h_p = f - 2 \underbrace{f_p(R, t)}_\text{proper distribution function}$$

Some words on the interaction

Consider $H_{\text{int}} = g \Psi^\dagger \Psi \varphi$ \leftarrow bosonic field

Take a look at the self-energy  $= \Sigma$
due to interactions:

Define bosonic Green's function on the contours:

$$D(l, 2) = -i \langle T_{\text{cav}} \varphi(l) \varphi(2) \rangle \quad (\text{let it be real field})$$

Consider first the term $g \varphi(\gamma, t)$ as a general (quantum) potential for your electronic system: (see page 15)

Remember: $G_{\text{cav}}(1, 1') \varphi(1') G_{\text{cav}}(1', 2) \Rightarrow \underline{G}(1, 1') \tau_2 \varphi(1') \underline{G}(1', 2)$
in keldysh space
in rotated space:

$$\underline{\zeta}_2 \delta G L^+ = \underbrace{\zeta_2 \underline{G} L^+}_{\hat{G}} \underbrace{\zeta_2 \varphi \tau_2 L^+}_{\zeta Q L^+} \underbrace{\zeta_2 \underline{G} L^+}_{\hat{A}} = \hat{G} \hat{\varphi} \hat{G}$$

$$\hat{\varphi} = \begin{pmatrix} \varphi_{cl} & \varphi_q \\ \varphi_q & \varphi_{cl} \end{pmatrix}$$

So in the second order with respect to quantum potential you find:

$$\delta^{(2)} G = \hat{G}(1,2) \hat{\varphi}(2) \hat{G}(2,3) \hat{\varphi}(3) \hat{G}(3,1') = \delta^{(2)} G(1,1')$$

$$\hat{\varphi} = \begin{pmatrix} \varphi_{cl} & \varphi_q \\ \varphi_q & \varphi_{cl} \end{pmatrix}$$

$$\varphi_{cl} = \frac{1}{2} (\varphi(t_+) + \varphi(t_-))$$

t_+ = t (upper contour)

$$\varphi_q = \frac{1}{2} (\varphi(t_+) - \varphi(t_-))$$

t_- = t (lower contour)

To compute the diagram you need

to average $\langle \hat{\varphi}(2) \hat{\varphi}(3) \rangle$

\swarrow ^{upper} \swarrow ^{upper}

check that $\langle \varphi_{cl}(1) \varphi_{cl}(2) \rangle = \frac{1}{4} \langle \varphi_+(1) \varphi_+(2) + \varphi_-(1) \varphi_-(2) + \varphi_-(1) \varphi_+(2) + \varphi_+(1) \varphi_-(2) \rangle$

$$= \frac{i}{4} (D_{11} + D_{12} + D_{21} + D_{22}) = \frac{i}{2} D^K(1,2)$$

$$\langle \varphi_{cl}(1) \varphi_q(2) \rangle = \frac{i}{4} (D_{11} - D_{12} + D_{21} - D_{22}) = \frac{i}{2} D^R(1,2)$$

$$\langle \varphi_q(1) \varphi_{cl}(2) \rangle = \frac{i}{4} (D_{11} + D_{12} - D_{21} - D_{22}) = \frac{i}{2} D^A(1,2)$$

$$\langle \varphi_q(1) \varphi_q(2) \rangle = \frac{i}{4} (D_{11} - D_{12} - D_{21} + D_{22}) = 0$$

← hats
the rules
to compute
the self energy

Home work

Use the rules at the previous page

to compute

$$\hat{\Sigma}(1,2) = \text{Diagram} = \frac{i}{2} \sum_{i,j} z_i G(1,2) z_j D_{ij}^{(1,2)}$$

$z_0 = \begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix}, z_1 = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}$
 $D_{00} = D^K, D_{01} = D^R, D_{10} = D^A$
 $D_{11} = 0$

Go to Wigner coordinates for Σ

Derive corresponding cell/torus integral !!!

Consider first "bare" Coulomb interaction

Remember that bare Coulomb is instantaneous:

$$D_0(1,2) = -\frac{e^2}{4\pi r_{12}} \delta(t_1 - t_2)$$

Think of the RPA series:

in RPA finite D^K will emerge!

$$\omega = \omega_0 + \omega Q_m + \omega C_m +$$

and how to sum them up

$$D_0^R(q, \omega) = D_0^A(q, \omega) = -\frac{2\pi e^2}{q} \epsilon_{\text{ind}}$$

~~$D_0^R(q, \omega) = D_0^A(q, \omega) = -\frac{2\pi e^2}{q} \epsilon_{\text{ind}}$~~

$$D_0^K = 0 = \omega(\omega) (D^R(q, \omega) - D^A(q, \omega))$$