

## Solution to Problem Set 2

A system of fermions in the presence of impurities with concentration  $n_i = N/V$  is described by the Hamiltonian,

$$\mathcal{H} = \sum_k \epsilon_k a_k^\dagger a_k + \int dx U(x) \psi^\dagger(x) \psi(x) \quad (1)$$

where  $U(x) = \sum_i V(x - X_i)$  is the array of potentials due to impurities located at random positions  $X_i$ . The Fourier transform of the field operator is,

$$\psi(x) = \frac{1}{\sqrt{V}} \sum_k e^{ikx} a_k$$

and that of the impurity potential  $V(x)$  is,

$$V(q) = \int dx e^{-iqx} V(x)$$

The Hamiltonian can be rewritten as

$$\mathcal{H} = \sum_k \epsilon_k a_k^\dagger a_k + \frac{1}{V} \sum_q V(q) \rho_q \sum_p a_{p+q}^\dagger a_p \quad (2)$$

where  $\rho_q = \sum_i e^{-iqX_i}$ , is a function of the random variables  $X_i$ . We have to perform ensemble average over the random positions. This makes the Hamiltonian translationally invariant and therefore tractable. For a typical sample the ensemble average is justified only if the probability distribution over the ensembles is sharply peaked. This condition is satisfied in the limit of weak disorder. The simplest ensemble average (with uniform weight) is defined as,

$$\overline{F} = \prod_{i=1}^N \frac{1}{V} \int d^3 X_i F(X_1, \dots, X_N) \quad (3)$$

Then,

$$\overline{\rho_q} = \frac{N}{V} \int d^3 X e^{-iqX} = N \delta_{q,0} \quad (4)$$

This average will appear in the first order self energy diagram. For the second order term we have,

$$\overline{\rho_{q_1} \rho_{q_2}} = N^2 \delta_{q_1,0} \delta_{q_2,0} + N \delta_{q_1+q_2,0}$$

where we have approximated  $N(N-1)$  by  $N^2$ , assuming  $N$  to be large. For the calculation of self energy we need,

$$\overline{\rho_q \rho_{-q}} = N^2 \delta_{q,0} + N \quad (5)$$

The first term on the right side produce second order improper self energy diagram, which can be disregarded. The second term will be relevant for second order proper self energy diagram.

In the Keldysh formalism Green's function  $G$  and self energy  $\Sigma$  are matrices defined as,

$$G = \begin{bmatrix} G^t & G^< \\ G^> & G^{\tilde{t}} \end{bmatrix}, \quad \text{and} \quad \Sigma = \begin{bmatrix} \Sigma^t & \Sigma^< \\ \Sigma^> & \Sigma^{\tilde{t}} \end{bmatrix}$$

Dyson's equation in frequency-momentum space takes the form,

$$G = G_0 + G_0 \sigma_3 \Sigma \sigma_3 G \quad (6)$$

where  $\sigma_3$  is the Pauli matrix  $\begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}$ . Defining  $\tilde{G} = G \sigma_3$ , we can rewrite the above equation as

$$\tilde{G} = \tilde{G}_0 + \tilde{G}_0 \tilde{\Sigma} \tilde{G} \quad (7)$$

So, the Feynman rules to evaluate the diagrams are best written in terms of  $\tilde{G}$  and  $\tilde{\Sigma}$ . To understand what the Feynman rules are, let us look at the object  $\langle T \psi_\alpha \psi_\beta^\dagger(\sigma_3)_{\beta\mu} \rangle$ . In perturbation expansion we have,

$$\langle T \psi_\alpha \psi_\beta^\dagger(\sigma_3)_{\beta\mu} \rangle = \langle T \exp[-i \int_{\mathcal{C}} d\tau V(\tau) \psi^\dagger(\tau) \psi(\tau)] \psi_\alpha \psi_\beta^\dagger(\sigma_3)_{\beta\mu} \rangle_0 \quad (8)$$

where the subscript zero indicates average with respect to the free Hamiltonian, and  $\mathcal{C}$  indicates time integration over the Keldysh contour. The lower line of the Keldysh contour that runs from  $+\infty$  to  $-\infty$  can be reversed by introducing a negative sign. Then the expansion of the S-matrix looks like,

$$\begin{aligned} \exp[-i \int_{\mathcal{C}} d\tau V(\tau) \psi^\dagger(\tau) \psi(\tau)] = & 1 + (\sigma_3)_{\alpha\beta} \int_{-\infty}^{+\infty} d\tau (-i) V(\tau) \psi_\alpha^\dagger(\tau) \psi_\beta(\tau) \\ & + \frac{1}{2!} \left\{ (\sigma_3)_{\alpha\beta} \int_{-\infty}^{+\infty} d\tau (-i) V(\tau) \psi_\alpha^\dagger(\tau) \psi_\beta(\tau) \right\}^2 + \dots \end{aligned}$$

Using this expansion in eqn(8), and writing schematically we get,

$$\begin{aligned} \langle \psi_\alpha \psi_\beta^\dagger(\sigma_3)_{\beta\mu} \rangle = & \langle \psi_\alpha \psi_\beta^\dagger(\sigma_3)_{\beta\mu} \rangle_0 + \langle \psi_\alpha \psi_a^\dagger(\sigma_3)_{ab} \rangle_0 \int_{-\infty}^{+\infty} d\tau (-i) V(\tau) \langle \psi_a \psi_\beta^\dagger(\sigma_3)_{\beta\mu} \rangle_0 \\ & + \langle \psi_\alpha \psi_a^\dagger(\sigma_3)_{ab} \rangle_0 \int_{-\infty}^{+\infty} d\tau (-i) V(\tau) \langle \psi_b \psi_c^\dagger(\sigma_3)_{cd} \rangle_0 \int_{-\infty}^{+\infty} d\tau (-i) V(\tau) \langle \psi_d \psi_\beta^\dagger(\sigma_3)_{\beta\mu} \rangle_0 \\ & + \dots \end{aligned} \quad (9)$$

Now comparing eqn(7) and eqn(9) we arrive at the rules that each contraction gives a single line that denotes the free propagator  $i\tilde{G}_0$ . The scattering vertex is represented in momentum space by  $\frac{-iV(q)\rho_q}{V}$ , which has to be averaged according to equation (3). The self energy diagram represents  $-i\tilde{\Sigma}$ . Then, the first order self energy diagram, after the impurity average, yields,

$$\begin{aligned} -i\tilde{\Sigma}(k) &= -i \frac{N}{V} V(q=0) \\ \Sigma^t(k) &= n_i \int d^3x V(x) \\ \Sigma^{\bar{t}}(k) &= -n_i \int d^3x V(x) \end{aligned} \quad (10)$$

To first order there is no contribution to  $\Sigma^<$  and  $\Sigma^>$ . Then,  $\Sigma^A = \Sigma^t - \Sigma^> = \Sigma^t$ , and  $\Sigma^R = \Sigma^t - \Sigma^< = \Sigma^t$ . Also  $\Sigma^K = \Sigma^< + \Sigma^> = 0$ . Also to be noted is that the first order

contribution to self energy is constant, and it leads to a simple shift in the energy levels of the free theory. Therefore, this contribution is of no physical interest.

The second order proper self energy diagram, after impurity average, gives,

$$-i\tilde{\Sigma}(k) = \frac{N}{V^2} \sum_p (-iV(p-k))(-iV(k-p))(i\tilde{G}(p)) \quad (11)$$

To get a closed form expression let us assume a short range potential, namely,  $V(x) = V_0\delta(x)$ . Since,

$$\begin{aligned} G^t(k, \omega) &= \frac{1 - f(\epsilon_k)}{\omega - \epsilon_k + i\eta} + \frac{f(\epsilon_k)}{\omega - \epsilon_k - i\eta} \\ &= P \frac{1}{\omega - \epsilon_k} - i\pi\delta(\omega - \epsilon_k)(1 - 2f(\epsilon_k)) \end{aligned}$$

from equation (9) we get,

$$\Sigma^t(k, \omega) = \frac{n_i V_0^2}{V} \sum_p \left\{ P \frac{1}{\omega - \epsilon_p} - i\pi\delta(\omega - \epsilon_p)(1 - 2f(\epsilon_p)) \right\}$$

We define density of states as,

$$\rho(\epsilon) = \frac{1}{V} \sum_p \delta(\epsilon - \epsilon_p) \quad (12)$$

then, the above equation can be rewritten as,

$$\Sigma^t(k, \omega) = n_i V_0^2 \left\{ P \int_{-\infty}^{\infty} \frac{\rho(\epsilon)d\epsilon}{\omega - \epsilon} - i\pi\rho(\omega)(1 - 2f(\omega)) \right\} \quad (13)$$

Evaluation of the remaining self energies from equation (9) is on similar lines, and we get,

$$\begin{aligned} \Sigma^{\tilde{t}}(k, \omega) &= -n_i V_0^2 \left\{ P \int_{-\infty}^{\infty} \frac{\rho(\epsilon)d\epsilon}{\omega - \epsilon} + i\pi\rho(\omega)(1 - 2f(\omega)) \right\} \\ \Sigma^>(k, \omega) &= -i(1 - f(\omega))2\pi\rho(\omega)n_i V_0^2 \\ \Sigma^<(k, \omega) &= if(\omega)2\pi\rho(\omega)n_i V_0^2 \end{aligned} \quad (14)$$

And from the definition of  $\Sigma^R$  and  $\Sigma^A$  we get,

$$\begin{aligned} \Sigma^R &= n_i V_0^2 \left\{ P \int_{-\infty}^{\infty} \frac{\rho(\epsilon)d\epsilon}{\omega - \epsilon} - i\pi\rho(\omega) \right\} \\ \Sigma^A &= n_i V_0^2 \left\{ P \int_{-\infty}^{\infty} \frac{\rho(\epsilon)d\epsilon}{\omega - \epsilon} + i\pi\rho(\omega) \right\} \end{aligned} \quad (15)$$

Also

$$\Sigma^K = -i(1 - 2f(\omega))2\pi\rho(\omega)n_i V_0^2 \quad (16)$$