

Free energy of the electron gas:
 perturbation theory:

~~1st order~~ $F = -T \ln Z$

$$Z = \int D(\bar{\psi}, \psi) e^{-S[\bar{\psi}, \psi]}, \text{ where}$$

$$S[\bar{\psi}, \psi] = \sum_p \bar{\psi}_p \left(-i\omega_n + \frac{|\vec{p}|^2}{2m} - \mu \right) \psi_p - \frac{1}{2L^3} \sum_{p, p', q} \bar{\psi}_{p+q} \bar{\psi}_{p'-q} V(q) \psi_{p'} \psi_p$$

1st order contribution:

$$F^{(1)} = \frac{T^2}{2L^3} \sum_{p, p', q} \langle \bar{\psi}_{p+q} \bar{\psi}_{p'-q} V(q) \psi_{p'} \psi_p \rangle_0$$

Feynman rules:

* Interaction: 

* Contraction: $\langle \psi_p \bar{\psi}_p \rangle_0 = G_{0,p} = \frac{1}{-i\omega_n + \frac{|\vec{p}|^2}{2m} - \mu}$

* Sum over all 4-momenta at vertex is zero

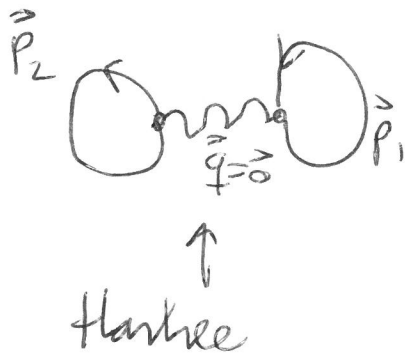
* Grassman var. anti commute: sign for each independent loop.

* Consider only connected diagrams (we're calculating $\ln Z$).

$$V(\vec{q}) = \frac{e^2}{q^2}, \quad V(z) = \frac{e}{z}, \quad \text{or:}$$

$$V(z) = \frac{e}{z} e^{-\lambda z} \sim \frac{e}{z} \frac{e^{-\lambda z}}{z^2}$$

There are two contributions at 1st order,



Hartree term: the Green's functions measure the electron density:

$$\begin{aligned} \text{Hartree term} &\propto \frac{1}{\beta} \sum_n \int \frac{d^3 p}{(2\pi)^3} \frac{1}{i\omega_n - \frac{\vec{p}^2}{2m} + \mu} \\ &= \frac{1}{\beta} \int \frac{d^3 p}{(2\pi)^3} n_F(\vec{p}) = \frac{N}{\beta}, \end{aligned}$$

↑
Metallic Fermi gas

which is the total charge. ~~The~~ The Hartree contribution cancels against a homogeneous background charge (not always the case!).

Hartree term: 'classical' density-density interaction.

[drop diagrams w/ $\vec{q} = \vec{0}$ interaction line]

The Fock contribution: (note: 1 fermion loop!)
 (spin: factor of 2)

$$- \frac{T^2}{L^3} \sum_{p, q} g_{0,p} g_{0,p+q} V(q)$$

$$= - \frac{T^2}{L^3} \sum_{p, p'} g_{0,p} g_{0,p'} V(p'-p)$$

($q = p' - p$)

$$= \frac{1}{L^3} \sum_{\vec{p}, \vec{p}'} n_F(\epsilon_{\vec{p}}) n_F(\epsilon_{\vec{p}'}) \frac{e^2}{|\vec{p} - \vec{p}'|^2}$$

$T \sum_{\omega_n} g_{\omega, p} = n_F(\epsilon_p)$

Evaluating this at zero temperature ($n_F \rightarrow$ step function)

$$F^{(1)}(T=0) = - \frac{1}{L^3} \sum_{\epsilon_p, \epsilon_{p'} < \mu} \frac{e^2}{|\vec{p} - \vec{p}'|^2}$$

$$= - \frac{e^2 L^3}{(2\pi)^4} \rho_F^4; \quad \text{Power of } \rho_F: \rho_F^{3+3-2}$$

Constant & evaluate integrals; (use Legendre polynomials)

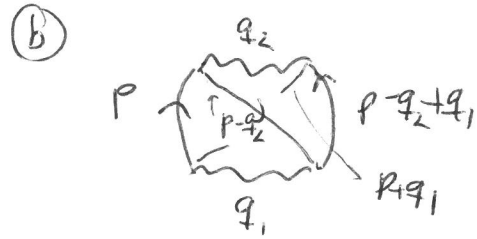
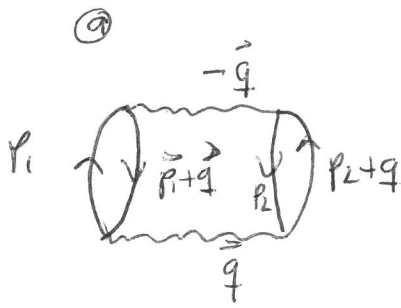
The contribution is negative, because the electrons avoid one another!

There is a problem, however! $F^{(0)} + F^{(1)}$ give rise to a divergence at ϵ_F , and an ρ that is

zero. Reason: long range Coulomb interaction.

Sol: screening or higher order PT,

2nd order terms (drop the ones w/ $\vec{q} = 0$ lines)



$$F^{2a} = - \frac{T^3}{L^6} \sum_{p_1, p_2, q} G_{\sigma, p_1} G_{\sigma, p_1+q} G_{\sigma, p_2} G_{\sigma, p_2+q} V(q)^2$$

$$F^{2b} = + \frac{T^3}{2L^6} \sum_{p, q_1, q_2} G_{\sigma, p} G_{\sigma, p+q_1} G_{\sigma, p+q_1-q_2} G_{\sigma, p-q_2} V(q_1) V(q_2)$$

b contrib: one particle hole pair

a contrib: two ind. particle hole pairs

How do these contribute: $\frac{1}{-i\omega_n + E(p) - \mu}$: suppressed away from $|\vec{p}| = p_F$

In (a) this means that $|\vec{q}|$ is small

(b) $|\vec{q}_1|, |\vec{q}_2|$ small,

In (a), there are two independent momentum integrals, in (b) only one. So, there is a much larger phase space for (a), which will dominate.

At higher order, we find that 'ring' diagrams dominate:



These diagrams correspond to:

$$F_{RPA}^{(n)} = \frac{T}{2n} \sum_q \left(\left(\frac{2T}{L^3} \right) V(q) \sum_p G_{Sp} G_{S+q} \right)^n$$

Combinatorics!

↳ $\frac{(n-1)!}{n!} \rightarrow$ # ways the int. can be arranged
 ↳ exponent

Important class of diagrams, which have large phase space!

We can actually sum over all of them:

$$F_{RPA} = \sum_{n=1}^{\infty} F_{RPA}^{(n)} = \frac{T}{2} \sum_q \ln (1 - V(q) \Pi_q),$$

where $\Pi_q \equiv \frac{2T}{L^3} \sum_p G_{Sp} G_{S+q}$ is the polarization operator, related to screening:

$$V_{eff} \equiv \frac{1}{V(q) - \Pi_q} = \frac{V(q)}{\epsilon(q)}, \text{ where } \epsilon(q) = 1 - V(q) \Pi_q$$


is a gen. dielectric function.

We see this from comparing $N^{(1)} = -\partial_\mu F^{(1)}$ and

$$N_{RPA} = -\partial_\mu F_{RPA}:$$

$$N^{(1)} = -\frac{2T^2}{L^3} \sum_{p,q} (G_{Sp})^2 G_{S+q} V(q)$$

$$N_{RPA} = \frac{T}{2} \sum_q \frac{V(q) \partial_\mu \Pi_q}{1 - V(q) \Pi_q} = \dots = -\frac{2T^2}{L^3} \sum_q V_{eff}(q) \sum_p G_{Sp} G_{S+q}$$

We can represent \tilde{G} as  :

$$\begin{aligned} \tilde{G} &= G + G \circ G + G \circ G \circ G + \dots \\ &= G + G \circ G \end{aligned}$$

Σ , symbolically

$$\begin{aligned} \tilde{G} - G \circ G &= G \\ \Rightarrow \tilde{G} &= \frac{G}{1 - G \circ G} = \frac{V(q)}{1 - V(q) \Pi(q)} \end{aligned}$$

We can do something similar for the self-energy
(ϕ^4 theory):

Write the greens function as:

$$\tilde{G} = G + G \circ \text{blob} + G \circ \text{blob} \circ \text{blob} + \dots$$

where  is 'one particle irreducible'!

$$\text{blob} = \text{blob} + G + \text{blob} + \dots \text{ etc.}$$

$$G = G_0 + G_0 \Sigma G_0 + G_0 \Sigma G_0 \Sigma G_0 + \dots$$

$$= G_0 + G_0 \Sigma G \quad \text{Dyson eq'n.}$$

$$\left[\tilde{G} = G + G \circ \text{blob} \right]$$

In momentum space: $G_p = G_{0,p} + G_{0,p} \Sigma_p G_p$, or

$$G_p = [1 - G_{0,p} \Sigma_p]^{-1} G_{0,p} = \frac{1}{G_{0,p}^{-1} - \Sigma_p} = \frac{1}{p^2 + m^2 - \Sigma_p}$$

Σ , to find greens function, we need to compute the self energy!