
Problem Set 2

1. Free electron gas in low dimensions (2+2+2+2+2 points)

Consider free electron gases confined in 1D and 2D boxes with “volume” V . The total number of electrons is N_e .

- Calculate the density of states (DOS) $D(\varepsilon)$, plot them and compare with the 3D result.
- Determine the Fermi energies ε_F by using the DOS obtained in (a). What are the corresponding Fermi wave vectors k_F ?
- Calculate the total energies of the Fermi gas at zero temperature. Compare your results with the 3D case ($E_{3D}/N_e = \frac{3}{5}\varepsilon_F$).
- Now consider the low temperature regime $T \ll T_F = \varepsilon_F/k_B$ and calculate the internal energies of the 1D and 2D free electron gases

$$U(T) = \int_{-\infty}^{\infty} d\varepsilon \varepsilon D(\varepsilon) f(\varepsilon, \mu, T),$$

where $f(\varepsilon, \mu, T) = 1/(e^{\beta(\varepsilon-\mu)} + 1)$ is the Fermi-Dirac distribution function.

Hint: You may use the Sommerfeld expansion to approximate the integral.

- Calculate the specific heat based on the results obtained in (d). How does the specific heat behave when $T \rightarrow 0$?

2. Density-density correlation function (3+3+2 points)

Consider the Fermi sphere of a *spinless* Fermi gas in 3D, described by the wave function

$$|\text{FS}\rangle = \prod_{|\mathbf{k}| < k_F} c_{\mathbf{k}}^{\dagger} |0\rangle.$$

- Calculate the Green's function

$$G(\mathbf{r}_1, \mathbf{r}_2) = \langle \psi^{\dagger}(\mathbf{r}_1) \psi(\mathbf{r}_2) \rangle,$$

where $\psi(\mathbf{r}) = \frac{1}{\sqrt{V}} \sum_{\mathbf{k}} c_{\mathbf{k}} e^{i\mathbf{k}\cdot\mathbf{r}}$ and $\langle \dots \rangle$ is the expectation value with respect to $|\text{FS}\rangle$. We note that $G(\mathbf{r}_1, \mathbf{r}_2)$ is often referred to as the one-particle reduced density matrix or correlation matrix.

- Calculate the density-density correlation function

$$C(\mathbf{r}_1, \mathbf{r}_2) = \langle n(\mathbf{r}_1) n(\mathbf{r}_2) \rangle - \langle n(\mathbf{r}_1) \rangle \langle n(\mathbf{r}_2) \rangle,$$

where $n(\mathbf{r}) = \psi^{\dagger}(\mathbf{r}) \psi(\mathbf{r})$ is the particle-density operator.

Hint: For calculating $\langle n(\mathbf{r}_1) n(\mathbf{r}_2) \rangle$, you may switch to momentum space or use the Wick's theorem $\langle \psi^{\dagger}(\mathbf{r}_1) \psi(\mathbf{r}_1) \psi^{\dagger}(\mathbf{r}_2) \psi(\mathbf{r}_2) \rangle = \langle \psi^{\dagger}(\mathbf{r}_1) \psi(\mathbf{r}_1) \rangle \langle \psi^{\dagger}(\mathbf{r}_2) \psi(\mathbf{r}_2) \rangle + \langle \psi^{\dagger}(\mathbf{r}_1) \psi(\mathbf{r}_2) \rangle \langle \psi(\mathbf{r}_1) \psi^{\dagger}(\mathbf{r}_2) \rangle$.

- Plot $C(\mathbf{r}_1, \mathbf{r}_2)$ as a function of the distance $|\mathbf{r}_1 - \mathbf{r}_2|$ and interpret the results (e.g. short/large distance behaviors and a characteristic oscillation known as “Friedel oscillation”).