5 BCS theory of superconductivity

Refs: [Mar] Section 27.3, [LP] E.M.Lifshitz et L.P.Pitaevskii, "Statistical Physics, Part 2" (vol.9 of "Landau et Lifshitz"), Sections 39,40.

As we have seen, phonons mediate an attraction between electrons. In this section, we will see how superconductivity emerges in an electron gas with attraction (the theory of Bardeen–Cooper–Schriffer).

5.1 Superconductivity as spontaneous symmetry breaking

Superconductivity is associated with developing nonzero anomalous averages

$$
\langle a_{\alpha} a_{\beta} \rangle \neq 0, \tag{5.1.1}
$$

where a_{α} and a_{β} are annihilation operators and α and β denote electron degrees of freedom (momentum/coordinate and spin). Such an average breaks the U(1) (electromagnetic) symmetry

$$
a \mapsto e^{i\alpha}a, \qquad a^+ \mapsto e^{-i\alpha}a^+.
$$
\n
$$
(5.1.2)
$$

The anomalous average [\(5.1.1\)](#page-0-0) may only be nonzero in a superposition of states with different particle numbers. Physically, the number of electrons in an isolated piece of a superconductor is fixed, in which case [\(5.1.1\)](#page-0-0) should be understood as a long-range order

$$
\lim_{|x-y|\to\infty} \langle (a_{\alpha}a_{\beta})_x (a_{\beta}^+ a_{\alpha}^+)_y \rangle \neq 0, \qquad (5.1.3)
$$

and the phase of an individual average $\langle a_{\alpha} a_{\beta} \rangle$ remains undetermined.

A good analogy to think of is the ferromagnetic transition: in a ferromagnet, the average magnetization is non-zero and points in a spontaneously chosen direction, even though in an isolated system, formally, the ground state is a superposition of states with all equivalent orientations of magnetization.

The phase of the average $\langle a_{\alpha} a_{\beta} \rangle$ is the spontaneously broken symmetry. It is not observable directly, but only in comparison with other such phases (the Josephson effect). A spatial modulation of this phase corresponds to the supercurrent (an electric current which propagates without dissipation).

Superconductors may be classified by the symmetry of indices in the anomalous average [\(5.1.1\)](#page-0-0). The most common symmetry (usually favored in superconductors with attraction due to phonons) is the *s-wave* superconductivity: the spin indices in $(5.1.1)$ form a singlet, and the pairing is isotropic in space.

5.2 Model Hamiltonian and mean-field approximation

We consider a model Hamiltonian of the form

$$
H = H_0 + H_{\text{int}} , \t\t(5.2.1)
$$

where

$$
H_0 = \sum_{k,\alpha} (\varepsilon_k - \mu) a_{k,\alpha}^+ a_{k,\alpha} \tag{5.2.2}
$$

is the free part, and

$$
H_{\rm int} = -\frac{1}{2V} \sum_{\tilde{k}_1, \tilde{k}_2, \tilde{k}_3} a_{\tilde{k}_1}^+ a_{\tilde{k}_2}^+ a_{\tilde{k}_3} a_{\tilde{k}_1 + \tilde{k}_2 - \tilde{k}_3} V_{\tilde{k}_1, \tilde{k}_2, \tilde{k}_3, \tilde{k}_1 + \tilde{k}_2 - \tilde{k}_3},\tag{5.2.3}
$$

where V is the system volume and $V_{\tilde{k}_1,\tilde{k}_2,\tilde{k}_3,\tilde{k}_4}$ are the interaction matrix elements (the spin indices are included in \tilde{k}_i for simplicity). In this chapter, we use the "sum" notation (the sum over k instead of integration over $d^3k/(2\pi)^3$, with the electronic states normalized as $\{a_k, a_{k'}^+\} = \delta_{kk'}$ (instead of the delta function of the continuous variable $k - k'$).

We will use the *mean-field approximation*: first, replace the products $a_k a_{k'}$ by their nonzero averages and then solve the *self-consistency* equation for those averages.

The structure of the non-zero anomalous averages depends on the interaction $V_{\tilde{k}_1,\tilde{k}_2,\tilde{k}_3,\tilde{k}_4}$. We assume that the superconductor is s-wave, with $\langle a_{k\uparrow}a_{-k\downarrow} \rangle \neq 0$. Correspondingly, we only consider the terms of this type in the interaction and neglect the k dependence of the interaction matrix elements (since, as we will see below, only k values around the Fermi surface are relevant). As a result, we simplify the interaction term to

$$
H_{\rm int} = -\frac{g_0}{\mathcal{V}} \sum_{k,k'} a_{k\uparrow}^+ a_{-k\downarrow}^+ a_{-k'} a_{k'\uparrow} = -\frac{g_0}{\mathcal{V}} \left(\sum_k a_{k\uparrow}^+ a_{-k\downarrow}^+ \right) \left(\sum_{k'} a_{-k'\downarrow} a_{k'\uparrow} \right) ,\qquad(5.2.4)
$$

where g_0 is some positive interaction constant.

We further define the complex numbers

$$
d_k = \langle a_{-k\downarrow} a_{k\uparrow} \rangle \qquad d_k^* = \langle a_{k\uparrow}^+ a_{-k\downarrow}^+ \rangle \,. \tag{5.2.5}
$$

These numbers will be later determined from the self-consistency conditions.

By replacing the products in the four-fermion operator by their averages, we get the quadratic Hamiltonian

$$
H_{\rm BCS} = \sum_{k} \left[(\varepsilon_k - \mu)(a_{k\uparrow}^+ a_{k\uparrow}^+ + a_{-k\downarrow}^+ a_{-k\downarrow}) + \Delta^* a_{-k\downarrow} a_{k\uparrow} + \Delta a_{k\uparrow}^+ a_{-k\downarrow}^+ \right],\tag{5.2.6}
$$

where

$$
\Delta = -\frac{g_0}{\mathcal{V}} \sum_k d_k \,. \tag{5.2.7}
$$

5.3 Bogoliubov quasiparticles and the BCS ground state

This quadratic Hamiltonian may be diagonalized by a rotation in the particle-hole space:

$$
\gamma_{k\uparrow}^+ = u_k a_{k\uparrow}^+ + v_k a_{-k\downarrow} \,. \tag{5.3.1}
$$

The coefficients u_k and v_k can be found, e.g., from the commutation relation

$$
[H_{\rm BCS}, \gamma_{k\uparrow}^+] = \tilde{\varepsilon}_k \gamma_{k\uparrow}^+ \,. \tag{5.3.2}
$$

We find the equation on the coefficients:

$$
\begin{pmatrix} \varepsilon_k - \mu & \Delta \\ \Delta^* & -(\varepsilon_k - \mu) \end{pmatrix} \begin{pmatrix} u_k \\ v_k \end{pmatrix} = \tilde{\varepsilon}_k \begin{pmatrix} u_k \\ v_k \end{pmatrix} . \tag{5.3.3}
$$

Figure 17: Left: The two sections across the Fermi surface contributing to the spectrum in the right panel. Each quasiparticle in the right panel corresponds to a linear combination of an electron with momentum k and a hole with momentum $-k$ and opposite spin. Right: The BCS spectrum [\(5.3.4\)](#page-2-0).

The eigenvalues give the spectrum:

$$
\tilde{\varepsilon}_k = \pm \sqrt{(\varepsilon_k - \mu)^2 + |\Delta|^2}.
$$
\n(5.3.4)

Thus $|\Delta|$ plays the role of the *qap* in the spectrum (see Fig. [17\)](#page-2-1).

The fermionic Fock space may be now represented in terms of the occupation numbers for (Bogoliubov) quasiparticles γ_{k}^+ κ_{\uparrow}^{+} . There are two ways to label quasiparticles:

- We can consider only $spin-up$ operators, as in $(5.3.2)$. In this case, we get two solutions for each k vector: one with positive, and one with negative energy.
- Alternatively, we can re-label the negative-energy solutions $(5.3.4)$ as annihilation operators $\gamma_{-k\downarrow}$. Then, for each k vector, we will have two quasiparticles γ_k^+ $\gamma_{k\uparrow}^+$ and $\gamma_{k\downarrow}^+$ $\stackrel{+}{k \downarrow},$ both with positive energies.

In any of these notations, the total number of quasiparticle states (the dimension of the Hilbert space) is the same as for original electrons: two single-particle states per k vector. We will use the second notation (with positive-energy quasiparticles).

It will also be convenient to normalize the coefficients so that

$$
|u_k|^2 + |v_k|^2 = 1.
$$
\n(5.3.5)

This would produce the canonical anticommutation relations for the quasiparticles:

$$
\{\gamma_{k\alpha}, \gamma_{k'\beta}^+\} = \delta_{kk'}\delta_{\alpha\beta} \,. \tag{5.3.6}
$$

The Hamiltonian [\(5.2.6\)](#page-1-1) can now be written in terms of quasiparticles as

$$
H_{\rm BCS} = \sum_{k} \tilde{\epsilon}_{k} \left(\gamma_{k\uparrow}^{+} \gamma_{k\uparrow} + \gamma_{k\downarrow}^{+} \gamma_{k\downarrow} \right) + E_{0} \,, \tag{5.3.7}
$$

Figure 18: The coherence factors [\(5.3.11\)](#page-3-0) as a function of energy.

The ground state of the superconductor $|GS\rangle$ may be found from the condition that it contains no quasiparticles:

$$
\gamma_{k\alpha}|\text{GS}\rangle = 0. \tag{5.3.8}
$$

Since sectors with different k vectors are decoupled in the Hamiltonian, this equation can be solved independently for each k vector:

$$
|GS\rangle_k = (u_k^* - v_k^* a_{k\uparrow}^+ a_{-k\downarrow}^+)|\star\rangle_k, \qquad (5.3.9)
$$

where $|\star\rangle$ is the state without electrons and the subscript k denotes that only states with a given k vector are considered. Combining all the k vectors together, we find the expression for the ground state of the superconductor:

$$
|\text{GS}\rangle = \prod_{k} (u_k^* - v_k^* a_{k\uparrow}^+ a_{-k\downarrow}^+) | \star \rangle. \tag{5.3.10}
$$

Note that this state is a superposition of states with different numbers of particles.

If we calculate u_k and v_k explicitly from diagonalizing the matrix [\(5.3.3\)](#page-1-2), we find

$$
u_k = e^{i\varphi} \sqrt{\frac{1}{2} \left(1 + \frac{\epsilon_k - \mu}{\tilde{\epsilon}_k} \right)},
$$

\n
$$
v_k = \sqrt{\frac{1}{2} \left(1 - \frac{\epsilon_k - \mu}{\tilde{\epsilon}_k} \right)},
$$
\n(5.3.11)

where φ is the phase of Δ . The absolute values of u_k and v_k are plotted in Fig. [18.](#page-3-1) We see that superconductivity changes the structure of the ground state only in the window of energies of the order Δ around the Fermi level (we usually have $\Delta \ll \mu$ in superconductors).

5.4 Self-consistency equations for the superconducting gap

The anomalous correlation functions d_k and the superconducting gap Δ are determined from the self-consistency conditions (5.2.5), where the averages are calculated in the quadratic system $(5.2.6)$ at a finite temperature T.

One of the possible ways to compute the anomalous average $\langle a_{-k\downarrow}a_{k\uparrow}\rangle$ is to re-express the a operators in terms of the quasiparticles γ and γ^+ and then use the equilibrium Fermi occupation numbers for the quasiparticles:

$$
\begin{cases}\n\gamma_{k\uparrow}^{+} = u_{k}a_{k\uparrow}^{+} + v_{k}a_{-k\downarrow} \\
\gamma_{-k\downarrow} = v_{k}^{*}a_{k\uparrow}^{+} - u_{k}^{*}a_{-k\downarrow}\n\end{cases}\n\Rightarrow\n\begin{cases}\na_{k\uparrow}^{+} = u_{k}^{*}\gamma_{k\uparrow}^{+} + v_{k}\gamma_{-k\downarrow} \\
a_{-k\downarrow} = v_{k}^{*}\gamma_{k\uparrow}^{+} - u_{k}\gamma_{-k\downarrow}\n\end{cases}\n(5.4.1)
$$

In terms of the quasiparticles γ_{k}^+ $\chi^+_{k\uparrow}$ and $\gamma_{-k\downarrow}$, the BCS Hamiltonian is diagonal, so we find

$$
\langle a_{-k\downarrow}a_{k\uparrow}\rangle_T = v_k^* u_k \langle \gamma_{k\uparrow}^+ \gamma_{k\uparrow} - \gamma_{-k\downarrow} \gamma_{-k\downarrow}^+ \rangle_T = v_k^* u_k \left[2n_F(\tilde{\varepsilon}_k) - 1 \right] = -v_k^* u_k \tanh\frac{\tilde{\varepsilon}_k}{2T}, \quad (5.4.2)
$$

where $\tilde{\varepsilon}_k$ is the quasiparticle energy given by Eq. (5.3.4).

Substituting this into Eq. $(5.2.7)$, we find the self-consistency equation for the gap

$$
\Delta = \frac{g_0}{\mathcal{V}} \sum_k v_k^* u_k \tanh \frac{\tilde{\varepsilon}_k}{2T}.
$$
\n(5.4.3)

Using Eq. $(5.3.11)$ for u_k and v_k , we find

$$
v_k^* u_k = \frac{\Delta}{2\tilde{\varepsilon}_k} \,. \tag{5.4.4}
$$

Note that this quantity is significant only in the vicinity of the Fermi surface (since far away form the Fermi surface either u_k or v_k tends to zero).

We remark that $\Delta = 0$ is always a formal solution to the equations $(5.4.3)$ – $(5.4.4)$. But one can show that at low temperatures this solution does not correspond to a minimum of a free energy, but to its maximum. In other words, at low temperatures the $\Delta = 0$ solution is unstable, and the physically relevant solution is a nontrivial one. To find this nontrivial solution, we divide the equation by Δ and replace the sum over k by integration over energies:

$$
\frac{1}{\mathcal{V}} \sum_{k} \quad \to \quad \nu_0 \int d\varepsilon \,, \tag{5.4.5}
$$

where ν_0 is the density of electronic states (for free electrons) per unit volume and per spin projection and ε is the free-electron energy. Substituting equation (5.3.4) for $\tilde{\varepsilon}_k$ and shifting the integration variable to $\varepsilon = \varepsilon_k - \mu$, we finally find the self-consistency equation in the closed form √

$$
1 = g_0 \nu_0 \int d\varepsilon \frac{\tanh \frac{\sqrt{\varepsilon^2 + |\Delta|^2}}{2T}}{2\sqrt{\varepsilon^2 + |\Delta|^2}}.
$$
 (5.4.6)

This equation, in principle allows to determine Δ as a function of temperature (see Fig. [19\)](#page-5-0).

5.5 Superconducting gap at zero temperature

A subtle point in this calculation is that the integral [\(5.4.6\)](#page-4-2) actually diverges logarithmically at large ε . Physically, this divergence is removed by introducing a cut-off at energies of the order of Debye energy ω_D (since the attraction mediated by phonons only extends to those energies).

Figure 19: A sketch of the gap dependence on the temperature.

At zero temperature, $tanh(...) \rightarrow 1$, and the equation [\(5.4.6\)](#page-4-2) reduces to

$$
1 = g_0 \nu_0 \int_0^{\infty} \frac{d\varepsilon}{\sqrt{\varepsilon^2 + \Delta_0^2}} = g_0 \nu_0 \left[\ln \frac{\omega_D}{\Delta_0} + \text{const} \right],
$$
 (5.5.1)

where const is a constant of order one. This gives the superconducting gap at zero temperature Δ_0 in the form

$$
\Delta_0 = \text{const } \omega_D \exp\left(-\frac{1}{g_0 \nu_0}\right). \tag{5.5.2}
$$

Note that the gap is exponentially small in g_0 .

5.6 Superconducting transition temperature

In a similar way we can find the superconducting transition temperature T_c , with the only difference that now we neglect Δ in the self-consistency equation [\(5.4.6\)](#page-4-2):

$$
1 = g_0 \nu_0 \int_0^{\infty} d\varepsilon \frac{\tanh \frac{\varepsilon}{2T_c}}{\varepsilon} = g_0 \nu_0 \left[\ln \frac{\omega_D}{T_c} + \text{const} \right],
$$
 (5.6.1)

with some const of order one (but different from that in the calculation of Δ_0 above!). In other words, T_c is of the same order of magnitude as Δ_0 .

Remarkably, one can determine the ratio T_c/Δ_0 without any ambiguity related to the cutoff. Namely, the difference of the integrals [\(5.5.1\)](#page-5-1) and [\(5.6.1\)](#page-5-2) is convergent and does not depend on the cut-off:

$$
0 = \int_0^\infty d\varepsilon \left[\frac{\tanh \frac{\varepsilon}{2T_c}}{\varepsilon} - \frac{1}{\sqrt{\varepsilon^2 + \Delta_0^2}} \right] = \int_0^\infty dx \left[\frac{\tanh(x/2)}{x} - \frac{1}{\sqrt{x^2 + (\Delta_0/T_c)^2}} \right].
$$
\n(5.6.2)

From this equation, one finds the *universal* value for the ratio T_c/Δ_0 :

$$
T_c \approx 0.57 \Delta_0 \,. \tag{5.6.3}
$$

[This value is easy to obtain by numerical methods. A more sofisticated analytic calculation gives

$$
T_c = \left(\frac{e^C}{\pi}\right) \Delta_0, \qquad (5.6.4)
$$

where $C = 0.577...$ is the Euler constant.

The relation [\(5.6.3\)](#page-5-3) is in a remarkably good agreement with experimental values on many conventional superconductors, despite the simplifications made in the BCS theory.

Problem Set 10

Problem 10.1

By diagonalizing the matrix in [\(5.3.3\)](#page-1-2), derive the spectrum [\(5.3.4\)](#page-2-0) and the eigenvectors $(5.3.11).$ $(5.3.11).$

Problem 10.2

(a) In superconductors, there is a characteristic length scale ξ called the *coherence* length. One of its possible definitions is the extent of the pair correlations. Consider the anomalous correlations in real space

$$
\Delta(x - y) = \langle a_{\downarrow}(x) a_{\uparrow}(y) \rangle
$$

It decays at a certain length scale ξ . Calculate this length in the BCS ground state.

Hint 1: In the BCS ground state, different wave vectors k are decoupled, so it is convenient to do a calculation at a given k vector, and then Fourier transform.

Hint 2: Only a vicinity of the Fermi surface contributes to this anomalous correlator, so you may linearize the electron spectrum near the Fermi surface.

Hint 3: You will find $\xi = v_F/\Delta$.

(b) For Aluminum, find in the literature the value of the gap Δ and estimate the superconducting coherence length ξ .