

Second Quantization

Appendix!

Describe occupation instead of describing the state of each particle

$$\Psi(0, 1, 0, 0, 1, 0, 1, 0, \dots)$$

Fermi exclusion principle 0, 1

Boxes
0, 1, 2, 3, ...

$$\phi_2, \phi_5, \phi_7 \quad (= \phi_0 e^{i\mathbf{k}\cdot\mathbf{r}})$$

not just product, Ψ has to be antisym for fermions

$$\Psi(0, 1, 0, 0, 1, 0, 0, \dots) = \frac{1}{\sqrt{N!}} \begin{vmatrix} \phi_2(a) & \phi_2(b) & \phi_2(c) \\ \phi_5(a) & \phi_5(b) & \phi_5(c) \\ \phi_7(a) & \phi_7(b) & \phi_7(c) \end{vmatrix}$$

10^{23} particle determinant!

$$C_i \Psi(\dots, \underset{i}{1}, \dots) = (-1)^{s(i)} \Psi(\dots, 0, \dots)$$

$$C_i \Psi(\dots, 0, \dots) = 0$$

$$C_i^+ \Psi(\dots, 0, \dots) = (i)^{s(i)}$$

$$C_i^+ \Psi(\dots, 1, \dots) = 0$$

$s(i)$
number of filled
rows below
(states)

Commutation relations

anticommutator

$$C_i C_j = -C_j C_i \Rightarrow \{C_i, C_j\} = 0$$

$$C_i^+ C_j = -C_j C_i^+$$

$$C_i^+ C_j^+ = -C_j^+ C_i^+$$

$$C_i C_i = C_i^+ C_i^+ = 0$$

Number operator

$$\hat{n}_i = C_i^+ C_i = \begin{cases} 1 & \text{if filled} \\ 0 & \text{if empty} \end{cases}$$

$$1 - \hat{n}_i = C_i C_i^+ = \begin{cases} 1 & \text{if empty} \\ 0 & \text{if filled} \end{cases}$$

$$\{C_i, C_j^+\} = \delta_{ij}$$

$$C_i C_j^+ + C_j^+ C_i = \begin{cases} 1 & \text{if } i=j \\ 0 & \text{if } i \neq j \end{cases}$$

Bosons

Commutation
instead of

anticommutation

$$C \psi_n = \sqrt{n} \psi_{n-1}$$

$$C^+ \psi_n = \sqrt{n+1} \psi_{n+1}$$

$$C^+ C = n$$

$$C C^+ = n+1$$

Fermi Groundst

$$\Psi = \prod_{k, k_f, \sigma} C_{k, \sigma}^+ |0\rangle$$

mixed occupation $\neq \Psi = (C_2^+ C_3^+ + C_4^+) |4\rangle$

Operators

Single particle

$$\hat{F}^{(1)} = \sum_{j,l} c_j^\dagger f_{jl} c_l$$

Example kinetic energy

$$\hat{E}_{kin} = \sum_{k\sigma} c_{k\sigma}^\dagger c_{k\sigma} E_k = \sum_{k\sigma} n_{k\sigma} E_k$$

Two Particle

$$\hat{F}^{(2)} = \frac{1}{2} \sum_{j,k,l} c_i^\dagger c_j^\dagger f_{ij,kl} c_l c_k$$

Coulomb interaction

describe scattering from state kl , to state ij

Fourier
Localized operators

$$\psi_\sigma(r) = \frac{1}{\sqrt{V}} \sum_k c_{k\sigma} e^{+ik\cdot r}$$

$$\psi_\sigma^\dagger(r) = \frac{1}{\sqrt{V}} \sum_k c_{k\sigma}^\dagger e^{-ik\cdot r}$$

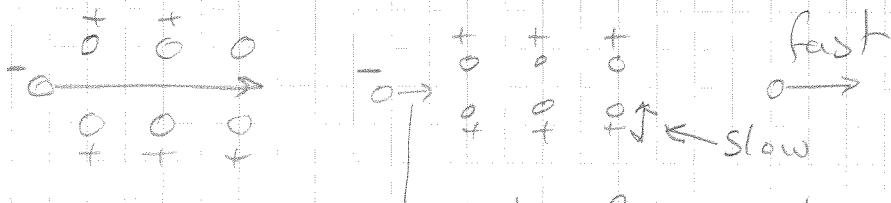
$$\{\psi_\sigma, \psi_\sigma^\dagger\} = \delta(\vec{r} - \vec{r}) \quad \text{Pauli principle}$$

$$F^{(1)} = \sum_j \int \psi_\sigma^\dagger f(r) \psi_\sigma d^3r$$

similar to the
expectation value

Simple minded picture

Normally electrons repel each other



attractive force due to
excess positive charge

This attractive force can be
more rigorously derived
describing the polarization of the
crystal due to lattice distortion
ie. electron phonon interaction

See sec 7.2 in the Book

This attractive force leads to
Cooper + pairing

Opposite k and σ

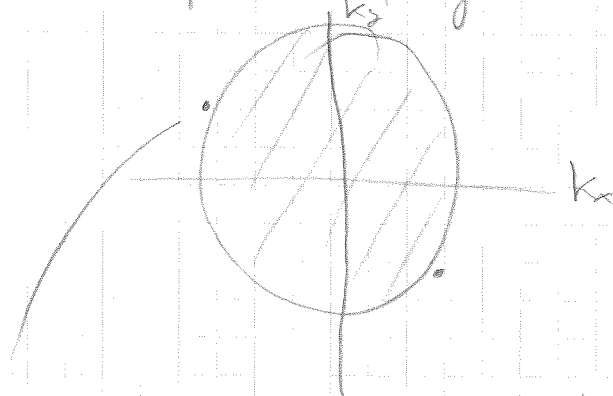
$$\vec{k}_1 = -\vec{k}_2 \quad \sigma_1 = -\sigma_2$$

$$H = H_0 + V(\vec{r}_1 - \vec{r}_2)$$

such that $V(\vec{r}_1 - \vec{r}_2) < 0$

The Cooper-pair model

Assume a filled Fermi sea
with all states for which $k < k_F$ are filled
if $k > k_F$ empty



Two electrons are added

Hamiltonian for the added pair

$$H = H_0 + V(\vec{r}_1 - \vec{r}_2)$$

non-interacting pair

interaction

Energies are measured with respect to E_F

SE

$$H\Psi = W\Psi$$

|
energy

$$(W - H_0)\Psi = V(\vec{r}_1 - \vec{r}_2)\Psi$$

non-interacting wave function

$$\phi_1 = e^{ik_1 \cdot \vec{r}_1} \quad \phi_2 = e^{ik_2 \cdot \vec{r}_2}$$

new coordinates

Center of mass

$$\bar{R} = \frac{1}{2}(\bar{r}_1 + \bar{r}_2)$$

$$\bar{K} = \frac{1}{2}(k_1 + k_2)$$

$$\bar{r}_1 = \bar{R} + \frac{r}{2}$$

$$\bar{r}_2 = \bar{R} - \frac{r}{2}$$

relative motion

$$r = \bar{r}_1 - \bar{r}_2$$

$$K = \frac{1}{2}(k_1 - k_2)$$

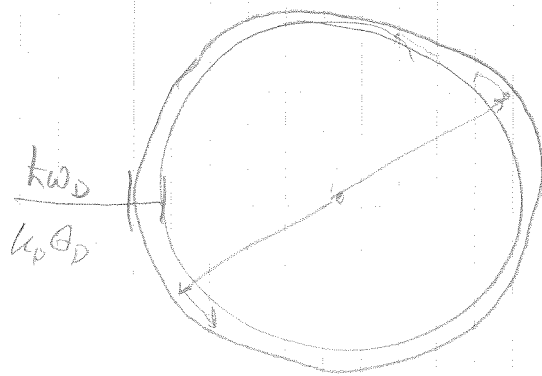
$$k_1 = \frac{1}{2}K + k$$

$$k_2 = \frac{1}{2}K - k$$

$$\Phi = \Phi_1 \cdot \Phi_2 = e^{ik_1 \bar{r}_1} \cdot e^{ik_2 \bar{r}_2} = e^{i(\bar{k}_1 \bar{r}_1 + \bar{k}_2 \bar{r}_2)}$$

$$= e^{i\left(\left(\frac{K}{2} + k\right)\left(\bar{R} + \frac{r}{2}\right) + \left(\frac{K}{2} - k\right)\left(\bar{R} - \frac{r}{2}\right)\right)}$$

$$= e^{i(\bar{K}\bar{R} + \bar{k}r)}$$



states with
opposite \bar{k}

$$\Rightarrow \bar{K} = 0$$

i.e

$$\Phi = e^{i\bar{k}(\bar{r}_1 - \bar{r}_2)} \quad \text{relative coordinates}$$

Stationary state is a superposition of different \bar{k} -states

$$\Psi = \sum_k a_k e^{ik(r_1 - r_2)}$$

$$(W - H_0) \Psi = V(\vec{r}_1 - \vec{r}_2) \Psi$$

$$(W - H_0) \sum_k a_k e^{ik(\vec{r}_1 - \vec{r}_2)} = V(\vec{r}_1 - \vec{r}_2) \sum_k a_k e^{ik(\vec{r}_1 - \vec{r}_2)}$$

Multiply by $e^{-ik(\vec{r}_1 - \vec{r}_2)}$ and \int over Volume Ω

Left hand side

$$\frac{1}{\Omega} \int_{\Omega} e^{-ik(\vec{r}_1 - \vec{r}_2)} (W - H_0) \sum_k a_k e^{ik(\vec{r}_1 - \vec{r}_2)} d(\vec{r}_1 - \vec{r}_2)$$

$$= \frac{1}{\Omega} W \sum_k a_k \int_{\Omega} e^{-ik(\vec{r}_1 - \vec{r}_2)} e^{ik(\vec{r}_1 - \vec{r}_2)} d(\vec{r}_1 - \vec{r}_2)$$

$$- \frac{1}{\Omega} \sum_k a_k \int_{\Omega} e^{-ik(\vec{r}_1 - \vec{r}_2)} H_0 e^{ik(\vec{r}_1 - \vec{r}_2)} d(\vec{r}_1 - \vec{r}_2)$$

$$= \frac{1}{\Omega} W a_k \Omega - \frac{1}{\Omega} \sum_k a_k E_k \int_{\Omega} e^{i(\vec{k}' - \vec{k})(\vec{r}_1 - \vec{r}_2)} d(\vec{r}_1 - \vec{r}_2)$$

$$W a_k - 2E_k a_k = (W - 2E_k) a_k$$

The right hand side

$$\frac{1}{\Omega} \int_{\Omega} e^{-ik(r_1-r_2)} \sum_k a_{k'} V(r_1-r_2) e^{+ik'(r_1-r_2)} d(r_1-r_2)$$

$$\left\{ V_{kk'} = \text{matrix element} = \frac{1}{\Omega} \int e^{-ik(r_1-r_2)} V e^{ik'(r_1-r_2)} d(r_1-r_2) \right\}$$

$$\sum_k a_k V_{kk}$$

$$(W - 2\epsilon_k) a_k = \sum_{k'} a_{k'} V_{kk'}$$

Assume V_{kk} independent of $k, k' = -V$
Calculate $W = \text{Energy of the Cooper pair}$

$$(W - 2\epsilon_k) a_k = -V \sum_{k'} a_{k'}$$

$$a_k = \frac{-V \sum_{k'} a_{k'}}{W - 2\epsilon_k}$$

$$\sum_{k'} a_{k'} = \sum_{k'} \left[\frac{-V \sum_{k''} a_{k''}}{W - 2\epsilon_{k'}} \right]$$

$$\downarrow = -V \sum_{k'} \frac{1}{W - 2\epsilon_{k'}}$$

convert sum to integral

$$I = -V \int_0^{\hbar\omega_D} \frac{N(\epsilon)}{W - 2\epsilon} \cdot d\epsilon$$

$$\hbar\omega_D \ll \epsilon_F \Rightarrow N(\epsilon) \approx N(0)$$

$$I = -VN(0) \int_0^{\hbar\omega_D} \frac{1}{W - 2\epsilon} d\epsilon$$

$$x = W - 2\epsilon$$

$$dx = -2d\epsilon$$

$$d\epsilon = -\frac{1}{2} dx$$

$$I = -VN(0) \left(-\frac{1}{2}\right) \int_0^{\hbar\omega_D} \frac{1}{x} dx$$

$$I = \frac{1}{2} VN(0) \left[\ln x \right]_{W-2\hbar\omega_D}^{W-2\cdot 0}$$

$$\frac{2}{VN(0)} = \ln \left(\frac{W - 2\hbar\omega_D}{W} \right)$$

$$e^{\frac{2}{VN(0)}} = 1 - \frac{2\hbar\omega_D}{W}$$

$$e^{\frac{2}{VN(0)}} - 1 = \frac{2\hbar\omega_D}{W}$$

$\gg 1$

$$W = -2\hbar\omega_D e^{-\frac{2}{VN(0)}}$$

Size of the Cooper pair

Limited number of \bar{k} states
is equivalent to

$$\Delta x \cdot \Delta p \approx \hbar$$

$$\Delta r \sim \frac{1}{\Delta k}$$

$$\Delta k = \frac{\partial k}{\partial E} \cdot \Delta E$$

$$E = \frac{\hbar^2 k^2}{2m}$$

$$\frac{\partial E}{\partial k} = \frac{\hbar^2 k}{m}$$

$$\Delta r = \frac{1}{\Delta k} = \frac{\partial E}{\partial k} \frac{1}{\Delta E} = \frac{\hbar^2 k}{m} \frac{1}{\Delta E} = \frac{\hbar v_F}{\Delta E} \sim$$

$$\approx \frac{\hbar v_F}{k_B T_c} \sim 1 \mu\text{m}$$