

# The BCS Model

Sara Changizi

This presentation closely follows parts of chapter 6 in  
Ring & Schuck “The nuclear many-body problem”.

# Outline

- Introduction to pairing
  - Essential experimental facts
- The BCS model
  - Pure pairing force

# Introduction to pairing

- In solid-state physics: A pair of electrons in metal with opposite spins and momentums close to Fermi surface interact with each other and make a pair. They have energy lower than Fermi surface which indicates that they are bound.
- BCS: Bardeen-Cooper-Schrieffer: microscopic model describes superconductivity (Nobel prize 1972)
- In nuclear physics: A pair of nucleons with total spin  $I_z=0$  ( $m_i, -m_i$ ). This is a short range (large spatial overlap) nucleon-nucleons interaction.

# Experimental observations

- Experimental observations that require short range interaction in the model :
  - The energy gap
  - The level density
    - Higher level density in low-lying excitation energies than found experimental values
  - Odd-even mass effect
  - Moment of inertia
    - Lowering of the moment of inertia compared to rigid body value
  - The low-lying  $2^+$  in even nuclei
    - Vibrate with low frequency :quadrupole oscillations

$$M_{(A \text{ odd})} > \frac{M_{A-1} + M_{A+1}}{2}$$

# The energy gap due to pairing

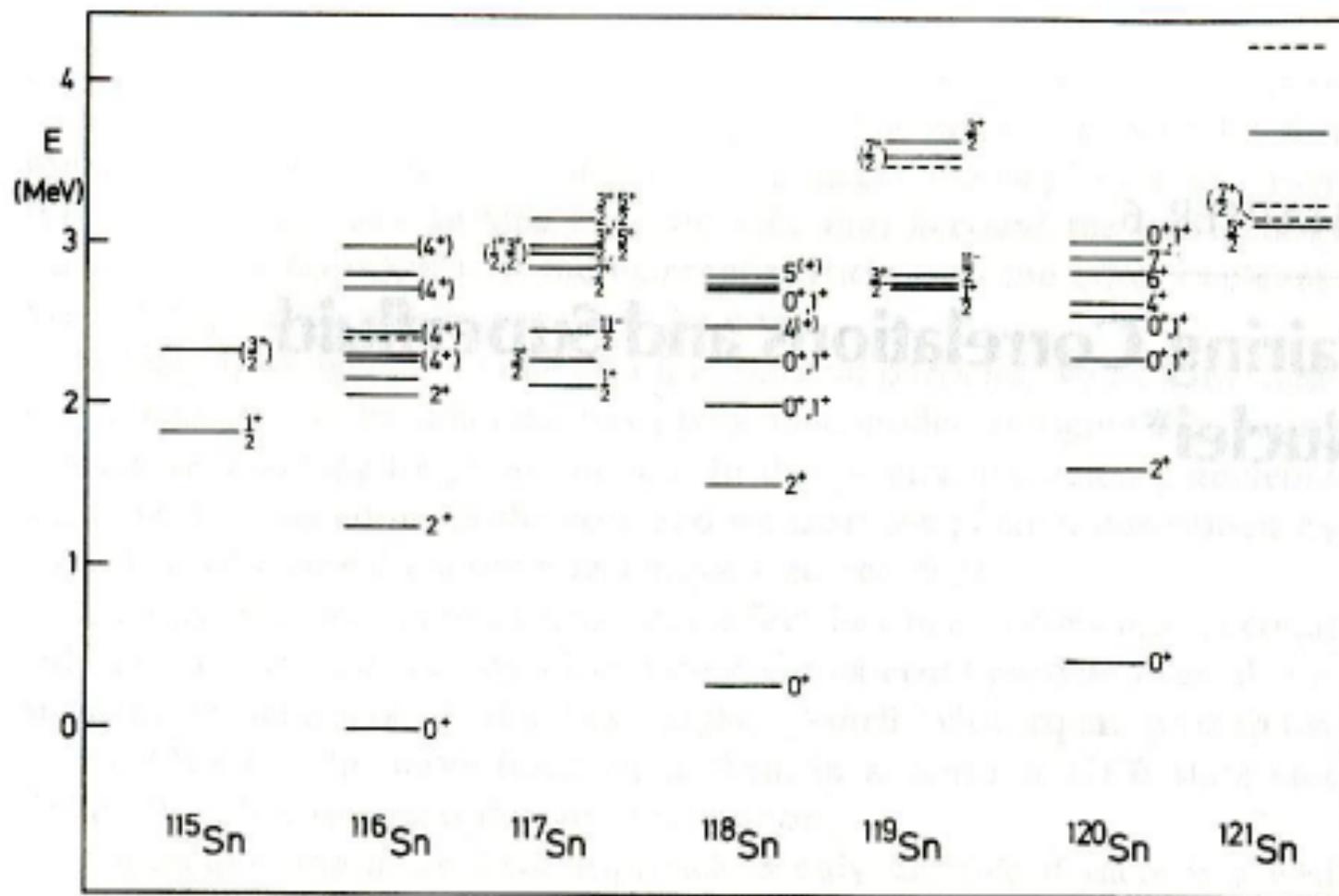


Figure 6.1. Excitation spectra of the  $^{50}\text{Sn}$  isotopes.

# Introduction to pairing

- To explain these phenomena we need to take into the account short range nucleon-nucleon interaction.
- The most effective pairing coupling is  $l=0$
- No exact solution, it is an approximation by variational principle

# Many-body system

A many-body system is described by following Hamiltonian in second-quantization form (particle-number representation):

$$H = \sum_{k_1 k_2} t_{k_1 k_2} a_{k_1}^+ a_{k_2} + \frac{1}{2} \sum_{k_1 k_2 k_3 k_4 <> 0} v_{k_1 k_2 k_3 k_4} a_{k_1}^+ a_{k_2}^+ a_{k_4} a_{k_3}$$
$$t_{k_1 k_2} \equiv \langle 1 : k_1 | t_1 | 1 : k_2 \rangle; \quad v_{k_1 k_2 k_3 k_4} \equiv \langle 1 : k_1 ; 2 : k_2 | v(1,2) | 1 : k_3 ; 2 : k_4 \rangle$$

First term: one body operator in second-quantization

Second term: two body operator in second-quantization

$V$ =matrix elements of the nucleon-nucleon interaction

$K_i$  are single-particle states and run over all available stats

Sums do not run on particles but on the infinite set of one-body states

# BCS model- BCS state

In BCS model we speculate that ground state should be built up from pair creation operators  $a_k^+ a_{-k}^+$

This is **the approximated solution:**

Trial wave function for even-even nuclei

$$|BCS\rangle = \prod_{k>0} (u_k + v_k a_k^+ a_{\bar{k}}^+) |-\rangle$$

$$|HF\rangle = \prod_{\alpha=1}^N a_{\alpha}^+ |-\rangle$$

K are the single particle levels

$V_k^2$  and  $U_k^2$  represent the probability that a certain pair state  $\{k, -k\}$  is or is not occupied.

$$|k\rangle = |nljm\rangle,$$

An example is a spherical basis

$$\text{Conjugate state: } |\bar{k}\rangle = |nlj - m\rangle \quad m > 0$$

# BCS model- BCS state

Example: For Hartree-Fock states we would have:

$v_k=1$  and  $u_k=0$  below Fermi level

$v_k=0$  and  $u_k=1$  above Fermi level

However in BCS model states over Fermi level can be occupied  
(energetically favored)

$u_k$  and  $v_k$  are variational parameters. We determine them in a way that the corresponding state has minimum energy.

Normalization of BCS state:

$$\langle BCS | BCS \rangle = \langle 0 | \prod u_k + v_k a_{\bar{k}} a_k \prod u_{k'} + v_{k'} a_{k'}^+ a_{\bar{k}'}^+ | 0 \rangle = \prod_{k>0} (u_k^2 + v_k^2)$$

We require:

$$u_k^2 + v_k^2 = 1$$

# BCS model

Great disadvantage: Particle number is not conserved in BCS! BCS state is the superposition of different number of pairs.

$$|BCS\rangle \propto |-\rangle + \sum_{k>0} \frac{v_k}{u_k} a_k^+ a_{-k}^+ |-\rangle + \frac{1}{2} \sum_{kk'} \frac{v_k v_{k'}}{u_k u_{k'}} a_k^+ a_{-k}^+ a_{k'}^+ a_{-k'}^+ |-\rangle + \dots$$

$$N = \langle BCS | \hat{N} | BCS \rangle = \langle BCS | a_k^+ a_k + a_{\bar{k}}^+ a_{\bar{k}} | BCS \rangle = 2 \sum_{k>0} v_k^2$$

This is fit to the interpretation of  $v_k$

Particle number uncertainty  $(\Delta N)^2 = \langle BCS | \hat{N}^2 | BCS \rangle - N^2 = 4 \sum_{k>0} u_k^2 v_k^2$

# BCS model

Hence, we need to restrict the variation by a supplementary condition. We define a parameter  $\lambda$  in the Hamiltonian to keep the expectation value of particle number to the desired particle number.

We add a term  $-\lambda N$  to Hamiltonian:

$$\hat{H}' = \hat{H} - \lambda \hat{N}$$

We call this Lagrange multiplier  $\lambda$  as Fermi energy or chemical energy, since it describes the energy variation in the system by changing the particle number.

# BCS model-Pure pairing force

Hamiltonian has a the form:

$$\sum_{k>0} \varepsilon_k a_k^+ a_k + \sum_{kk'>0} \langle k, \bar{k} | v | k', \bar{k}' \rangle a_k^+ a_{\bar{k}}^+ a_{\bar{k}'} a_{k'}$$

Single-particle part  $v_{kk\bar{k}\bar{k}'}$  Residual interaction acting only on pairs of nucleons

In this model we assume a constant matrix elements – G (pure pairing force):

$$H = \sum_{k>0} \varepsilon_k a_k^+ a_k - G \sum_{kk'>0} a_k^+ a_{\bar{k}}^+ a_{\bar{k}'} a_{k'}$$

# BCS model- Pure pairing force

Lets consider the Hamiltonian with the variational condition

$$H = \sum_{k>0} \varepsilon_k a_k^+ a_k - G \sum_{kk'>0} a_k^+ a_{k'}^+ a_{-k'}^- a_{-k}$$

$$H' = H - \lambda \hat{N}$$

$$\langle BCS | H - \lambda \hat{N} | BCS \rangle = \langle BCS | \sum_k (\varepsilon_k^0 - \lambda) a_k^+ a_k - G \sum_{kk'} a_k^+ a_{-k}^+ a_{-k'}^- a_{k'} | BCS \rangle$$

The expectation value:

$$\langle BCS | \hat{H}' | BCS \rangle =$$

$$\boxed{\Delta = G \sum_{k>0} u_k v_k}$$

$$2 \sum_{k>0} (\varepsilon_k^0 - \lambda) \underbrace{\langle BCS | a_k^+ a_k | BCS \rangle}_{v_k^2} - \sum_{k>0} G \underbrace{\langle BCS | a_k^+ a_{-k}^+ a_{-k} a_k | BCS \rangle}_{v_k^4} - G \underbrace{\left( \sum_{k>0} \underbrace{u_k v_k}_{\langle BCS | a_k^+ a_{-k}^+ a_{-k'}^- a_{k'} | BCS \rangle} \right)^2}_{\Delta}$$

# BCS model- Pure pairing force

$v_k$  determines the BCS wave function completely, we can express  $u_k$  in terms of  $v_k$  by the normalization condition

$$\delta\langle BCS|\hat{H}'|BCS\rangle = 0,$$

$$\left( \frac{\partial}{\partial v_k} + \frac{\partial u_k}{\partial v_k} \frac{\partial}{\partial u_k} \right) \langle BCS|H'|BCS\rangle = 0$$

$$u_k \frac{\partial u_k}{\partial v_k} + v_k \frac{\partial v_k}{\partial u_k} = 0 \Rightarrow \frac{\partial u_k}{\partial v_k} = -\frac{v_k}{u_k}$$

$$\left. \left( \frac{\partial}{\partial v_k} - \frac{v_k}{u_k} \frac{\partial}{\partial u_k} \right) \langle BCS|H'|BCS\rangle = 0 \right\}$$

$$\left( \frac{\partial}{\partial v_k} - \frac{v_k}{u_k} \frac{\partial}{\partial u_k} \right) \left( 2 \sum_{k>0} (\varepsilon_k^0 - \lambda) v_k^2 - \sum_{k>0} G v_k^4 - G \left( \sum_{k>0} u_k v_k \right)^2 \right) = 0$$

# BCS model

We get for **BCS equations** the following:

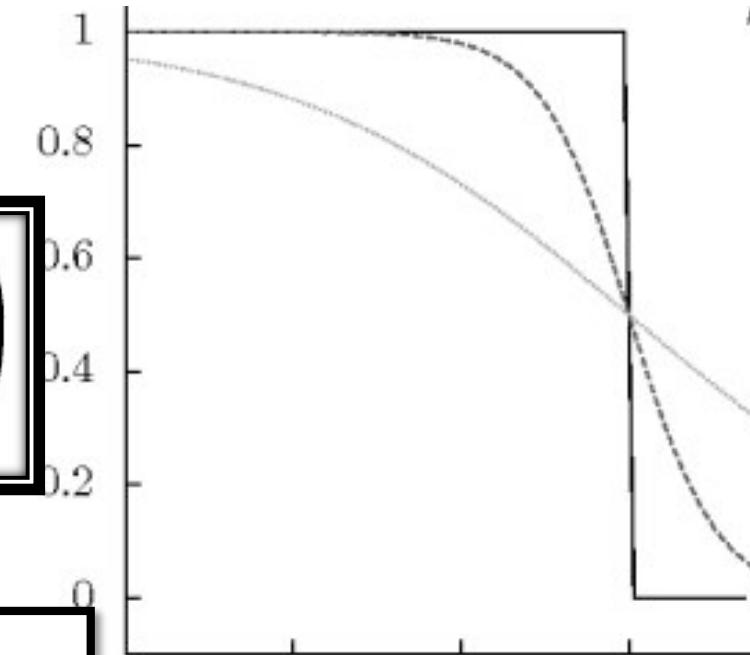
$$\varepsilon_k = \varepsilon_k^0 - \lambda - \underbrace{G \cdot v_k^2}_{\text{often neglected}}$$

$$2\varepsilon_k v_k u_k + \Delta(v_k^2 - u_k^2) = 0$$

The occupation probability in non interacting and

In the interacting case (assume we know  $\Delta$ ):

$$v_k^2 = \frac{1}{2} \left( 1 \pm \sqrt{1 - \frac{\Delta^2}{\varepsilon_k^2 + \Delta^2}} \right) = \frac{1}{2} \left( 1 \pm \frac{\varepsilon_k}{\sqrt{\varepsilon_k^2 + \Delta^2}} \right)$$



Insert this into definition of gap we can get this iterative equation.

Gap equation:

$$\Delta = \frac{G}{2} \sum_{k>0} \frac{\Delta}{\sqrt{\varepsilon_k^2 + \Delta^2}}$$

# Results of BCS model

- In the special case of a single  $j$ -shell: all  $\varepsilon_k$  are equals so all  $v_k^2$ 's are equals. From the particle-number condition we get

$$\langle BCS | \hat{N} | BCS \rangle = 2 \sum_{k>0} v_k^2 = N \Rightarrow v_k = \sqrt{\frac{N}{2\Omega}}, u_k = \sqrt{1 - \frac{N}{2\Omega}}$$

$$\Delta = G \cdot \sqrt{\frac{N}{2} \left( \Omega - \frac{N}{2} \right)}$$

- The gap has a parabolic dependence on the number of particle in the shell. It is zero for empty or filled shells. For  $N=\Omega$ :  $2\Delta = G \cdot \Omega$

$\Omega$  is the number of pairs

Thanks!