THERMODYNAMIC PROPERTIES OF HEAVY FERMION SUPERCONDUCTORS

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ABSTRACT

Certain types of interactions can lead to massive quasi-particles such that the resulting effective masses of the electrons may reach values between 100 and 1000 times the mass of the free electrons. The interactions between such heavy electrons are assumed to be singlet s-wave and triplet s-wave, and their contribution to s-wave superconductivity of heavy electrons has been studied. The effective Hamiltonian is diagonalized using the Bogoliubov-Valatin transformation, and the thermodynamic properties of such heavy fermion systems are calculated. The total energy is found to increase with temperature. There is a maximum in the specific heat at \( T_c \approx 4.8 \times 10^3 \) K suggesting a phase transition. The magnitude of specific heat is \( C_v = 4.8 \times 10^{33} \text{J/K} \) at \( T_c \). The entropy changes continuously through the transition temperature \( T_c \). Entropy of the system is found to decrease with temperature as is conventionally the case.

Keywords: Heavy Fermions, Super Fluid, Transition Temperature

INTRODUCTION

There are experimental observations in some cases that certain types of interactions can lead, at low temperatures, to the formation of a Fermi liquid-type state with massive quasi-particles (Andrei, 2004). Depending upon the detailed characteristics for the considered systems, the resulting effective masses of the electrons may reach values between 100 and 1000 times the mass of the free electrons. Such electrons are called heavy electrons. By definition, a common and independent Fermi liquid is stable and undergoes no phase transition. Experimental evidence for magnetic or superconducting transitions in heavy-electrons systems therefore imply that residual interactions driving these transitions need to be considered.

There is a common view that in the heavy-electron systems, the correlation effects or strong interactions are the dominating reasons behind the low-temperature properties of some materials. The heavy-electron state in most cases develops smoothly within a narrow temperature range upon cooling at low temperatures. In this sense, it may be regarded as a particular state of certain metals, caused by a delicate competition of interactions between ionic moments of atoms with incompletely-filled electronic shells (favorably f electrons) and itinerant electrons. Such a state may be unstable and undergo phase transitions, either to magnetic order or superconductivity. In fact, heavy-electron system involves many-body physics that deals with strongly interacting electrons in metals.

It is well known that for the standard BCS theory the obtained value for the size of the Cooper pair is of the order of \( 10^7 \text{Å}^2 \) whereas the experimentally extracted value for HTSC is of the order of \( 10^5 \text{Å}^2 \) (Riseman et al., 1995; Tifrea, 2003). Hence in the conventional superconductors (BCS type), the electrons in the Cooper pairs will be weakly bound because of the large size of the Cooper pairs, whereas the electrons in the HTSC will be strongly bound in the Cooper pairs because of the smaller size of the Cooper pair. Hence strong coupling between electrons in HTSC will lead to heavy fermions being present in HTSC. It is this fact that leads to the study of thermodynamic properties of heavy fermion superconductors.

Superconducting compounds which consist of one magnetic ion with 4f and 5f electrons, (generally Ce or U) and other constituents being s, p or electron metals are known as a family of heavy-fermion superconductors. A large number of heavy fermion materials superconduct exclusively under pressure.
The superconducting electron system is considered as being in some condensed phase and the scattered pairs of electrons are in s-state singlet pairing and s-state triplet pairing. The electrons are paired so as to minimize the ground state energy. With this assumption the Hamiltonian for the singlet and triplet pairing is written as,

$$H=\sum \varepsilon_k (c_k^*c_k + c_k^+c_k) - \sum V_{kk'} c_k^*c_{k'} + \sum U_{kk'} c_k^*c_k^*c_{k'}c_{k'}$$

(2)

Eq (2) will be expressed in terms of the new operators and then diagonalized to obtain the elements of the Hamiltonian. The new operators are related to the old fermion creation and annihilation operators such that,

$$\gamma_k = u_k c_k - v_k c_k^*$$

$$\gamma_k^* = u_k c_k^* + v_k c_k$$

(3)

Their conjugates are

$$\gamma_k^* = u_k c_k^* - v_k c_k$$

$$\gamma_k = u_k c_k^* + v_k c_k^*$$

(4)

In order to write the Hamiltonian in Eq (2) in terms of the new operators, the inverse transformation of Eq (3) and (4) are used. The inverse transformations of (3) and (4) are

$$c_k = u_k \gamma_k + v_k \gamma_k^*$$

$$c_k^* = u_k \gamma_k^* + v_k \gamma_k$$

$$c_k = u_k \gamma_k^* - v_k \gamma_k$$

$$c_k^* = u_k \gamma_k - v_k \gamma_k^*$$

(5)

The Hamiltonian expressed in terms of new operators will have three parts namely the Kinetic, Potential (Singlet) and Potential (Triplet) parts which are then diagonalized as explained in each step below.

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**Research Article**

**Kinetic Energy**

Eq (5) are used in Eq (2) to express the K.E of the Hamiltonian in terms of the new operators. Substituting the creation and annihilation operators using Eq (5) we get,

\[
H_{K} = \sum_{k} \varepsilon_{k} \left( u_{k}^{\dagger} i_{k} \gamma_{k} + u_{k} v_{k} \gamma_{k}^{\dagger} + u_{k} v_{k} \gamma_{k} - u_{k} v_{k} \gamma_{k}^{\dagger} + u_{k} v_{k} \gamma_{k}^{\dagger} + u_{k} v_{k} \gamma_{k} - u_{k} v_{k} \gamma_{k}^{\dagger} + u_{k} v_{k} \gamma_{k}^{\dagger} + v_{k} \gamma_{k} - v_{k} \gamma_{k}^{\dagger} + v_{k} \gamma_{k}^{\dagger} - v_{k} \gamma_{k}^{\dagger} \right)
\]  

(6)

Introducing particle number operators \( m_{k} \) that represent cases when particles are created we write,

\[
m_{k} = \gamma_{k}^{\dagger} \gamma_{k} \quad 1-m_{k} = \gamma_{k}^{\dagger} \gamma_{k}^{\dagger}
\]

(7)

Eq (8) are used in Eq (7) to get

\[
H_{KE} = \sum_{k} \varepsilon_{k} \left( u_{k}^{2} m_{k} + v_{k}^{2} (1-m_{k}) + v_{k}^{2} (1-m_{k}) + u_{k}^{2} m_{k} + u_{k} v_{k} \gamma_{k}^{\dagger} \gamma_{k}^{\dagger} + u_{k} v_{k} \gamma_{k} \gamma_{k} + u_{k} v_{k} \gamma_{k}^{\dagger} \gamma_{k} + u_{k} v_{k} \gamma_{k}^{\dagger} \gamma_{k}^{\dagger} + v_{k} \gamma_{k}^{\dagger} \gamma_{k}^{\dagger} - v_{k} \gamma_{k} \gamma_{k}^{\dagger} + v_{k} \gamma_{k}^{\dagger} \gamma_{k}^{\dagger} - v_{k} \gamma_{k} \gamma_{k}^{\dagger} - v_{k} \gamma_{k}^{\dagger} \gamma_{k}^{\dagger} \right)
\]

(8)

Using the relations,

\[
\gamma_{k}^{\dagger} \gamma_{k} = -\gamma_{k} \gamma_{k}^{\dagger} \quad \text{and} \quad \gamma_{k} \gamma_{k} = -\gamma_{k}^{\dagger} \gamma_{k}^{\dagger}
\]

(9)

We get,

\[
H_{KE} = \sum_{k} \varepsilon_{k} \left( 2v_{k}^{2} + (u_{k}^{2} - v_{k}^{2}) (m_{k} + m_{k}) + 2u_{k} v_{k} \left( \gamma_{k}^{\dagger} \gamma_{k}^{\dagger} + \gamma_{k} \gamma_{k} \right) \right)
\]

(10)

The first term of equation (11) is a constant, the terms with \( m_{k} \) and \( m_{k} \) generate particles and are called diagonal terms bearing \( \gamma_{k} \gamma_{k} \cdot \gamma_{k} \gamma_{k} \) operators, the third term is off-diagonal i.e. bearing \( \gamma_{k} \gamma_{k}^{\dagger} \) and \( \gamma_{k} \gamma_{k}^{\dagger} \) operators. They are without \( m_{k} \) and \( m_{k} \).

Eq (10) is the transformed form of the Kinetic Energy part.

**Potential Energy for Singlet Pairing**

Eq (5) are used to express the P.E of the singlet part of the Hamiltonian in terms of the new operators. From Eq(2)

\[
H_{P(E_{S})} = -\sum_{kk} V_{kk} c_{k}^{\dagger} c_{k}^{\dagger} c_{k} c_{k}
\]

(11)

Using Eq (5), Eq (11) and later transformed. The P.E Part thus becomes

\[
H_{P(E_{S})} = \sum_{kk} V_{kk} \left\{ \left( u_{k}^{2} \gamma_{k} \gamma_{k}^{\dagger} - u_{k} v_{k} \gamma_{k} \gamma_{k}^{\dagger} + u_{k} v_{k} \gamma_{k} \gamma_{k}^{\dagger} - u_{k} v_{k} \gamma_{k} \gamma_{k}^{\dagger} \right) \left( u_{k}^{2} \gamma_{k}^{\dagger} \gamma_{k}^{\dagger} + u_{k} v_{k} \gamma_{k} \gamma_{k} + u_{k} v_{k} \gamma_{k}^{\dagger} \gamma_{k}^{\dagger} + u_{k} v_{k} \gamma_{k} \gamma_{k}^{\dagger} - u_{k} v_{k} \gamma_{k}^{\dagger} \gamma_{k}^{\dagger} - v_{k} \gamma_{k} \gamma_{k}^{\dagger} + v_{k} \gamma_{k}^{\dagger} \gamma_{k}^{\dagger} - v_{k} \gamma_{k} \gamma_{k}^{\dagger} - v_{k} \gamma_{k}^{\dagger} \gamma_{k}^{\dagger} \right) \right\}
\]

(12)

Introducing particle number operators we write

\[
1-m_{k} = \gamma_{k} \gamma_{k}^{\dagger} \quad m_{k} = \gamma_{k} \gamma_{k}^{\dagger} \quad 1-m_{k} = \gamma_{k} \gamma_{k}^{\dagger} \quad 1-m_{k} = \gamma_{k} \gamma_{k}^{\dagger}
\]

(13)

Eq (13) are used to represent Eq (12) in terms of number operators. Thus

\[
H_{P(E_{S})} = \sum_{kk} V_{kk} \left\{ \left( u_{k}^{2} \gamma_{k} \gamma_{k}^{\dagger} - u_{k} v_{k} \gamma_{k} \gamma_{k}^{\dagger} + u_{k} v_{k} \gamma_{k} \gamma_{k}^{\dagger} - u_{k} v_{k} \gamma_{k} \gamma_{k}^{\dagger} \right) \left( u_{k}^{2} \gamma_{k}^{\dagger} \gamma_{k}^{\dagger} + u_{k} v_{k} \gamma_{k} \gamma_{k} + u_{k} v_{k} \gamma_{k}^{\dagger} \gamma_{k}^{\dagger} + u_{k} v_{k} \gamma_{k} \gamma_{k}^{\dagger} - u_{k} v_{k} \gamma_{k}^{\dagger} \gamma_{k}^{\dagger} - v_{k} \gamma_{k} \gamma_{k}^{\dagger} + v_{k} \gamma_{k}^{\dagger} \gamma_{k}^{\dagger} - v_{k} \gamma_{k} \gamma_{k}^{\dagger} - v_{k} \gamma_{k}^{\dagger} \gamma_{k}^{\dagger} \right) \right\} + 4OT
\]

(14)

(15)

where 4OT stands for "fourth order terms". The third term is an off-diagonal term which shall be considered later so as to leave a system of independent fermions.

**Potential Energy for Triplet Pairing**

From equation (2) we write

\[
H_{P(E_{T})} = \sum_{kk} U_{kk} c_{k}^{\dagger} c_{k}^{\dagger} c_{k} c_{k}
\]

(16)

Eq (5) are used to transform equation (16). The P.E (triplet) part thus becomes

\[
H_{P(E_{T})} = \sum_{kk} U_{kk} \left\{ \left( u_{k} \gamma_{k}^{\dagger} + v_{k} \gamma_{k} \right) \left( u_{k} \gamma_{k}^{\dagger} - v_{k} \gamma_{k} \right) \left( u_{k} \gamma_{k}^{\dagger} - v_{k} \gamma_{k} \right) \left( u_{k} \gamma_{k}^{\dagger} + v_{k} \gamma_{k} \right) \right\}
\]

(17)

or
The particle number operators in Eq(13) are then used in Eq (18) to get
\[ H_{ren} = \sum_{kk'} \left( u_k^+ u_{k'} v_{k'} - u_k v_{k'} u_{k'}^+ - u_k^+ v_{k'}^+ v_{k'} - u_{k'}^+ v_{k'} - u_k v_{k'} \right) \left( u_{k'}^+ v_{k'}^+ v_k - u_{k'} v_{k'} u_k^+ - u_{k'} v_k \right) \]
\[ + v_{k'}^+ u_k^+ m_{k'}^+ \gamma_k^+ + v_k^+ \gamma_k^+ m_{k'} u_{k'}^+ \left( 1-m_{k'} \right) \gamma_k^+ \gamma_{k'} + v_k^+ u_{k'}^+ \left( 1-m_{k'} \right) \gamma_{k'}^+ \gamma_k \]
\[ + u_k v_k^+ m_{k'} \gamma_k^+ - u_k^+ m_{k'}^+ \gamma_k - u_k^+ v_k v_{k'} \left( 1-m_{k'} \right) \gamma_{k'}^+ \gamma_k \]
\[ + u_{k'}^+ v_{k'} \left( m_{k'} - 1 + m_{k'} \right) \left( v_k^2 + u_k^2 \right) \left( \gamma_k^+ \gamma_{k'}^+ + \gamma_{k'} \gamma_k \right) \]
\[ + u_k v_k \left( m_k - 1 + m_k \right) \left( v_k^2 + u_k^2 \right) \left( \gamma_k^+ \gamma_{k'}^+ + \gamma_{k'} \gamma_k \right) \]  

(20)

Eq (20) is the transformed form of the Potential Energy of the triplet pairing. The last two terms are the off-diagonal terms.

Effective Hamiltonian

The effective Hamiltonian \( H = H_{KE} + H_{PES} + H_{PETF} \) is written by adding equations (10), (15) and (20) with the approximation that the 4OT can be neglected being negligible higher order terms i.e.
\[ H = H_{KE} + H_{PES} + H_{PETF} \]
\[ H = \sum_k \varepsilon_k \left( 2v_k^2 + \left( u_k^2 - v_k^2 \right) \left( m_k + m_k \right) \right) + 2u_k v_k \left( \gamma_k^+ \gamma_k \right) \]
\[ - \sum_k V_{kk'} \left( u_k v_k u_{k'}^+ v_{k'} + \left( 1 - m_{k'} - m_k \right) \left( 1 - m_{k'} - m_k \right) \right) \left( u_k^2 - v_k^2 \right) \left( \gamma_k^+ \gamma_{k'}^+ + \gamma_{k'} \gamma_k \right) \]
\[ + \sum_{kk'} V_{kk'} \left( u_k v_k u_{k'}^+ v_{k'} \left( m_{k'} + m_k - 1 \right) \left( m_{k'} + m_k - 1 \right) + u_k v_{k'} u_{k'}^+ v_k \left( 1 - m_{k'} - m_k \right) \left( u_k^2 - v_k^2 \right) \left( \gamma_k^+ \gamma_{k'}^+ + \gamma_{k'} \gamma_k \right) \right) \]
\[ + u_k v_k \left( m_k - 1 + m_k \right) \left( v_k^2 + u_k^2 \right) \left( \gamma_k^+ \gamma_{k'}^+ + \gamma_{k'} \gamma_k \right) \]  

(21)

Parametric Expressions for \( u_k \) and \( v_k \) for the Singlet State

A system of independent fermions is determined on the assumption that the off-diagonals terms of Eq (10), (15) and (20) vanish. At the lowest energy state of this system, both \( m_k \) and \( m_{k'} \) are zero. Hence to carry out the Bogoliubov-Valatin transformation for a superconductor in its ground state, \( m_k \) and \( m_{k'} \) are set to zero in Eq (10) and (15) and the 4OT neglected since the non-diagonal terms vanish. The modified form of combining equations (10) and (15) then gives
\[ \sum_k 2\varepsilon_k u_k v_k \left( \gamma_k^+ \gamma_{k'}^+ + \gamma_{k'} \gamma_k \right) - \sum_k V_{kk'} \left( u_k^2 - v_k^2 \right) \left( \gamma_k^+ \gamma_{k'}^+ + \gamma_{k'} \gamma_k \right) = 0 \]  

(22)

When the non-diagonal terms vanish Eq (22) becomes
\[ 2\varepsilon_k u_k v_k - \left( u_k^2 - v_k^2 \right) \sum_{k'} V_{kk'} u_k v_{k'} = 0 \]  

(23)

The condition \( u_k^2 + v_k^2 = 1 \) is used to express \( u_k \) and \( v_k \) in the form of a single variable \( \chi \) such that
\[ u_k = \frac{1}{2} \chi_k, \quad v_k = \frac{1}{2} + \chi_k \]  

(24)

and when used in Eq (23) gives
\[ 2\varepsilon_k \sqrt{\frac{1}{4} - \chi_k^2} + 2\chi_k \sum_{k'} V_{kk'} \sqrt{\frac{1}{4} - \chi_k^2} = 0 \]  

(25)

The quantity, \( \Delta_k \), is expressed as
\[ \Delta_k = \sum_{k'} V_{kk'} \sqrt{\frac{1}{4} - \chi_k^2} \]  

(26)

Hence on solving equation (25)
Quasi particle Spectrum for the Singlet State

Using Eq (27) in Eq (24) gives

\[ u_k^2 = \frac{1}{2} \left( 1 + \frac{\varepsilon_k}{\sqrt{\varepsilon_k^2 + \Delta_k^2}} \right) \]  
\[ v_k^2 = \frac{1}{2} \left( 1 - \frac{\varepsilon_k}{\sqrt{\varepsilon_k^2 + \Delta_k^2}} \right) \]  

The Hamiltonian consisting of diagonal terms is used to determine the quasi particle energy because the diagonal terms correspond to stationary states when the system is in equilibrium. The quasiparticles energy is found using equation (28) and equation (29).

The Energy \( E_{ks} \) necessary to create a quasi particle excitation will be obtained from the diagonal terms of the Hamiltonian \( H = \sum_k \varepsilon_k \left( 2v_k^2 + \left( u_k^2 - v_k^2 \right) \left( m_k + m_{k'} \right) \right) \sum_{kk'} V_{kk'} \left( u_{kk'} v_{k'} - v_{kk'} u_{k'} \left( 1 - m_{kk'} - m_{k'k} \right) \right) \) \( (30) \)

Now,

\[ E_{ks} = \frac{\partial \langle H_D \rangle}{\partial < m_k >} \]  \( (31) \)

\[ E_{ks} = \varepsilon_k \left( u_k^2 - v_k^2 \right) + 2u_k v_k \sum_{kk'} V_{kk'} u_{kk'} v_{k'} \left( 1 - m_{kk'} - m_{k'k} \right) \]  \( (32) \)

Approximating \( \Delta_k \approx \sum_{kk'} V_{kk'} u_{kk'} v_{k'} \left( 1 - m_{kk'} - m_{k'k} \right) \)

Equation (32) becomes

\[ E_{ks} = \varepsilon_k \left( u_k^2 - v_k^2 \right) + 2u_k v_k \Delta_k \]  \( (33) \)

Eq (28) and (29) are substituted in Eq (35) to obtain the quasi particle energy for the singlet pairing.

\[ E_{ks} = \sqrt{\varepsilon_k^2 + \Delta_k^2} \]  \( (34) \)

Equation (34) is the quasiparticles energy spectrum of the singlet state.

**Parametric expressions for \( u_k \) and \( v_k \) for the Triplet state**

To carry out the Bogoliubov-Valatin transformation for a superconductor in its ground state, the \( m_k, m_{k'} \) are set equal to zero in Eq (10) and (20) and the 4OT neglected since the non-diagonal terms vanish.

The modified form of combining equations (10) and (20) then gives

\[ \sum_k 2\varepsilon_k u_k v_k \left( \gamma_k^* \gamma_k^* + \gamma_{k'} \gamma_{k'} \right) + \sum_{kk'} U_{kk'} \left( u_k^2 v_{kk'} - v_k^2 u_{kk'} \right) \left( 1 - m_{kk'} - m_{k'k} \right) = 0 \]  \( (35) \)

or

\[ 2\varepsilon_k u_k v_k + u_k^2 \sum_{k'} U_{kk'} u_{k'} v_k = 0 \]  \( (36) \)

Eq (36) is solved using Eq (26) to get

\[ 2\varepsilon_k \left( \frac{1}{4} - X_k^2 \right)^{\frac{1}{2}} + \left( \frac{1}{2} - X_k \right) \sum_{k'} U_{kk'} \left( \frac{1}{4} - X_k^2 \right)^{\frac{1}{2}} = 0 \]  \( (37) \)
We now define a new quantity $\Delta$ as,

$$\Delta = \sum_{k} U_{kk} \left( \frac{1}{4} - X_{k}^2 \right)$$

(38)

Eq (37) then becomes

$$2\varepsilon_{k} \left( \frac{1}{2} - X_{k} \right)^{2} \left( \frac{1}{2} + X_{k} \right)^{2} + \left( \frac{1}{2} - X_{k} \right) \Delta = 0$$

(39)

or

$$X_{k} = \frac{1}{2} \left( \frac{\Delta^2 - 4\varepsilon_{k}^2}{4\varepsilon_{k}^2 + \Delta^2} \right)$$

(40)

The parametric expressions for $u_{k}$ and $v_{k}$ are obtained by substitution of Eq (40). Using Eq (24) and (40) we obtain

$$u_{k} = \pm \frac{2\varepsilon_{k}}{\sqrt{4\varepsilon_{k}^2 + \Delta^2}} \quad \text{and} \quad v_{k} = \pm \frac{\Delta}{\sqrt{4\varepsilon_{k}^2 + \Delta^2}}$$

(41)

Quasi particle Spectrum of the Triplet State

The Hamiltonian consisting of diagonal terms is used to determine the quasi particle energy because the diagonal terms correspond to stationary states when the system is in equilibrium. The quasiparticles energy is found using Eq (41). The Energy $E_{kT}$ necessary to create a quasi particle excitation will be obtained from the diagonal terms of the Hamiltonian

$$H_d = \sum_{k} 2\varepsilon_{k} v_{k}^2 + \sum_{k} \varepsilon_{k} (u_{k}^2 - v_{k}^2)(m_{k} + m_{k}) + \sum_{kk} U_{kk} u_{k} v_{k} u_{k} v_{k} ((m_{k} + m_{k} - 1)(m_{k} + m_{k} - 1))$$

(42)

Now

$$E_{kT} = \frac{\partial \langle H_d \rangle}{\partial <m_{k}>}$$

Approximating $\Delta \approx \sum_{k} U_{kk} u_{k} v_{k} ((m_{k} + m_{k} - 1))$ gives

$$E_{kT} = \varepsilon_{k} (u_{k}^2 - v_{k}^2) + 2u_{k} v_{k} \Delta$$

(43)

Eq (41) are substituted in Eq (43) to get

$$E_{kT} = \varepsilon_{k} \left( \frac{4\varepsilon_{k}^2}{4\varepsilon_{k}^2 + \Delta^2} \right) \Delta + 2 \left( \frac{2\varepsilon_{k}}{\sqrt{4\varepsilon_{k}^2 + \Delta^2}} \right) \Delta$$

$$E_{kT} = \frac{4\varepsilon_{k}^3 + 3\varepsilon_{k} \Delta^2}{4\varepsilon_{k}^2 + \Delta^2}$$

(44)

Eq (44) is the quasi particle energy for the triplet pairing.

Energy of the Assembly

The effective Hamiltonian given by Eq (2) is considered by summing the diagonal and non-diagonal parts of the various components of the Hamiltonian with the condition that the non-diagonal parts are set equal to zero or vanish whereas the 4OT are neglected i.e.

$$E_{k} = E_{\text{Singlet}} + E_{\text{Triplet}}$$

$$E_{k} = \sum_{k} \varepsilon_{k} \left( 2v_{k}^2 + (u_{k}^2 - v_{k}^2)(m_{k} + m_{k}) + 2u_{k} v_{k} (\gamma_{k} + \gamma_{-k}) \right)$$

$$- \sum_{kk} V_{kk} \left( u_{k} v_{k} u_{k} v_{k} (1-m_{k} - m_{k})(1-m_{k} - m_{k}) + u_{k} v_{k} (1-m_{k} - m_{k})(u_{k}^2 - v_{k}^2)(\gamma_{k} + \gamma_{-k}) \right)$$
The diagonal parts of Eq (45) correspond to quasi-particle states when the system is in equilibrium i.e.

\[ \text{Eq (45)} \]

The particle number operators \( m_k = m_{-k} = 0 \) (This means the quasi-particles represented by the operators \( \gamma \)'s are few in number or do not exist). Eq (46) then becomes

\[ \text{Eq (46)} \]

The singlet and triplet gaps are represented by

\[ \Delta_s = \sum_k V_{kk} u_k v_k \]  
\[ \Delta_u = \sum_k U_{kk} u_k v_k \]  

Eq (47) is then by substitution written as

\[ \text{Eq (47)} \]

To study the temperature dependence of the system the total energy \( E \) is found by multiplying energy of states \( E_k \) by the thermal activation factor \( e^{\frac{E_k}{k_B T}} \). Hence

\[ \text{Eq (51)} \]

Equation (50) is used in Eq (51) thus

\[ \text{Eq (52)} \]

Eq (52) is the total energy of the assembly with the triplet and singlet contributions. 

**Specific Heat Capacity**

The specific heat capacity is determined using the following relation

\[ \text{Eq (53)} \]

and hence

\[ \text{Eq (54)} \]

Eq (54) is the Specific heat Capacity equation for heavy fermion systems.

**Transition Temperature**

The transition temperature of the system is obtained by witing

\[ \text{Eq (55)} \]

To simplify equation (54) the larger terms are then represented by letting

\[ \rho = \frac{(2E_k \kappa_k^2 + E_k u_k v_k (\Delta_u - \Delta_s))}{K_B T^2} \]  
\[ \kappa_k = \beta. \]  

Then Eq (54) can be expressed as

\[ \text{Eq (57)} \]
Thus we get
\[
\left( -\frac{2\rho}{T^3} e^{\left(\frac{\beta}{T}\right)} + \frac{\beta\rho}{T^4} e^{\left(\frac{\beta}{T}\right)} \right)_{T=T_c} = 0
\]  \hspace{1cm} (58)

or
\[
T_c = \frac{\beta}{2} = \frac{E_K}{2K_B}
\]  \hspace{1cm} (59)

**Entropy**

The entropy of a system is determined from the relation
\[
S_f - S_i = \int_{T_i}^{T_f} \frac{dQ}{T} = \int_{V_i}^{V_f} mC dT
\]  \hspace{1cm} (60)

The equation for entropy from equation (60) then becomes
\[
S = m \left( \frac{2E_{K}E_{K}v_{K}^2 + E_{K}u_{K}E_{K}}{K_B} \right) \left( \frac{K_B e^{\frac{E_{K}}{K_B T}} + 3K_B^2 e^{\frac{E_{K}^2}{E_{K}^2}} + 6K_B^3 T e^{\frac{E_{K}^3}{E_{K}^3}} + 6K_B^4 T^2 e^{\frac{E_{K}^4}{E_{K}^4}}} {K_T e^{\frac{E_{K}}{K_T T}}} \right)
\]  \hspace{1cm} (61)

where \( m \approx 100m_e \).

**Essential Parameters**

The following values for different physical quantities have been used.

The numerical values of \( C^V \) are obtained using the condition \( u^2 + v^2 = 1 \)

\[
u_k = \frac{1}{\sqrt{2}} , \quad v_k = \frac{1}{\sqrt{2}}
\]

According to the experimental works of (Allan, et al., 2013) on heavy fermions

Single particle energy, \( \varepsilon_k = \hbar \omega = 300 \mu eV = 4.8 \times 10^{-23} J \)

The triplet and singlet energy gaps are given by the relation \( \Delta_v \approx 0.05 \Delta_u \) according to (Khanna, 2008).

The heavy fermion superconductivity energy window is equivalent to \(-600 \mu eV < E < 600 \mu eV\) according to (Allan, et al., 2013)

Energy gap for the triplet \( \Delta_u = 9.6 \times 10^{-23} J \) and \( E_{KS} = \sqrt{\varepsilon_k^2 + \Delta_k^2} \)

\( E_{KS} = 4.8239402 \times 10^{-23} J \), \( E_{K_T} = 9.6 \times 10^{-23} J \) and \( E_K = 1.44239 \times 10^{-23} J \)

Boltzmann constant \( K_B = 1.3806488 \times 10^{-23} JK^{-1} \)

**RESULTS AND DISCUSSION**

![Figure 1: Variation of Total Energy against Temperature](image)
Research Article

Eq (52) was used to compute the values of quasi-particle energy against changes in the temperature. A graph depicting the variation of quasi-particle energy versus temperature is shown in figure 1.

Eq (54) is used to compute the values of specific heat against temperature changes.

![Graph of Specific Heat Against Temperature](image)

Figure 2: The graph of specific heat against temperature.

The shape of the specific heat curve is similar to the specific heat curve of UBe$_{13}$ (Ott H.R., 1992). The curve shows a turning point at $T_C=5.2K$.

Equation (61) gives the variation of Entropy with temperature.

![Graph of Entropy Against Temperature](image)

Figure 3: Variation of Entropy against Temperature

Discussion

The value of E decreases below $T_C$ (5.2K) and becomes zero at T=0 K and this is consistent with the nature of the super-fluid state. The total energy of the system increases with increase in temperature of the system.

In figure 1, it is observed that at T=0, E=0 which implies that the states below the superconducting gap are filled and the states above the superconducting gap are empty. Now it is much harder for impurities to scatter the pair of electrons because they need a finite amount of energy. If there is not enough energy to get to an unoccupied state, then the electron does not change its state, there is no dissipation, and hence there is no electrical resistance. To get across the superconducting gap requires a minimum amount of energy of about $1.2 \times 10^{-23}$. The gap doesn’t appear suddenly at $T_C$ (5.2K); it opens up gradually as T decreases below $T_C$, just like the order parameter. When the temperature increases, the energy is observed to increase. This implies thermal fluctuations will populate the f-level by conduction electrons, thereby increasing the f-electron number. This will involve the breaking of Cooper pairs and hence the
suppression of the superconducting gap. At high temperatures the f-electrons are localized on their atomic sites and do not contribute to conduction.

Superconductivity is a true reversible thermodynamic state and in moving from the normal to the superconducting state, a material undergoes a thermodynamic phase transition. In order for this to happen, the overall free energy must be lower in the superconducting state than in the normal state as seen in figure 1 and this energy difference, which depends on the temperature, is known as the condensation energy. At high temperatures heavy fermion compounds behave like normal metals and the electrons can be described as Fermi gas, in which the electrons are assumed to be non-interacting fermions. The specific heat capacity of heavy fermion superconductors jumps at $T_C$, but the scale remains set by the large effective mass of the quasiparticles (Coleman, 2007). If the superconducting state did not have heavy fermion properties, the specific heat below the jump would be much smaller (Coleman, 2007).

In figure 2, the graph of specific heat assumes nearly perfect Gaussian shape. A peak in specific heat occurs at about 5.2K. At this point the system is unstable and there is a second order phase transition from a normal metal into a superconducting state that is much like the super fluid transition. At $T_C$, there is a phase transition, the fermions pair up and the specific heat depicting a second order transition meaning there is no latent heat.

The weak interaction of fermions is mediated by the lattice and approaches zero at absolute zero. Super fluidity in fermions is exhibited on the basis of interaction and is a phenomenon that is realized at any temperature below a finite transition temperature. At room temperatures, the $f$-electrons of the magnetic ions behave as localized spins; the conduction electrons are the $s$, $p$ or $d$ electrons and have quite ordinary effective masses. As the temperature is lowered, the $f$-electrons begin to couple to the conduction electrons, resulting in very large effective masses for the hybridized carriers (Andrei, 2004). At low temperatures the electronic specific heat shows a continuous increase until superconductivity sets in at $T_C=5.2K$.

Furthermore, the temperature dependence of the heat capacity below $T_C$ is not exponential. Instead, it follows a power law, as seen from figure 2 indicating that the energy gap at the Fermi surface has nodes in certain directions. Thus, the energy gap is highly anisotropic (Andrei, 2004).

However, contrary to conventional materials, some of the $f$-electrons become itinerant at low temperatures. Crudely speaking, it is these itinerant $f$-electrons that lead to the uncommon behaviors of heavy fermion materials at low temperatures (Hua, 2008). The large magnitude of the specific heat jump ($4.8\times10^{23} J/K$) that occurs at the superconducting transition temperature, $T_C$, shows that the electrons that take part in the formation of the heavy–fermion state form the Cooper pairs.

In figure 3 shown above, entropy increases continuously with increase in temperature. This shows excitation of the particles in the system. The entropy approaches zero at $T\rightarrow0$, as it should. The gradient $dS/dT$ goes to zero sharply as $T=0K$. The rate of change $dS/dT$ (gradient of the graph) consistently decreases at temperatures below $T_C=5.2K$ and becomes zero from $T=1.5K$ i.e. system becomes more orderly. When molecules are cooled their total energy decreases resulting into less and less vigorous molecules.

The entropy of the system at $T_C$ is $3.5\times10^{-21} J/K$. The graph confirms this observation by depicting that entropy decreases with decrease in temperature. At extremely low temperatures as the entropy of the material drops, an electron passing by lattices in the crystal will cause an electron- electron interaction making efforts to form Cooper pairs and at a particular temperature i.e. $T_C$ the material changes electric state. This is when the material becomes superconducting. But if a field is applied, the electrons would have some component counteracting the electron-electron interaction and decreases entropy. At $T=0$, $S=0$, it seems reasonable to regard this entropy as being carried by a set of localized f spins which emerge above temperatures. To ensure that $S = 0$ at $T = 0$, localized spins order magnetically due to spin-spin interaction. Similarly, a system with the fermion condensate must experience some sort of a low-temperature phase transition eliminating the excessive entropy $S$. Particle disorder decreases with decrease in the total energy of the system.
Conclusion

Figure 3 suggests that particles settle and interact less as the system gives out energy. This is in agreement with conventional knowledge and concurs with (Khanna et al., 2010). It is also in good agreement with (Ayodo, 2008), (Ayodo et al., 2010) on low temperature statistical thermodynamics of binary bose-fermi system. Entropy is a measure of molecular disorder (Ayodo, 2008) when the system cools, the internal energy of the particles decreases resulting in less and less particle motion. The graph in figure 3, confirms this observation by predicting that entropy decreases with decrease in temperature even for a system of heavy fermions

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