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Entropy and spin susceptibility of s-wave type-II superconductors near $H_{c2}$

Takahumi Kita
Division of Physics, Hokkaido University, Sapporo 060-8610, Japan
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A theoretical study is performed on the entropy $S_s$ and the spin susceptibility $\chi_s$ near the upper critical field $H_{c2}$ for s-wave type-II superconductors with arbitrary impurity concentrations. The changes of these quantities through $H_{c2}$ may be expressed as $[S_s(T,B)−S_s(T,0)]/[S_s(T)−S_s(T,0)]=1−a_s(1−B/H_{c2})=(B/H_{c2})^{\alpha_s}$, for example, where $B$ is the average flux density and $S_s$ denotes entropy in the normal state. It is found that the slopes $\alpha_s$ and $\chi_s$ at $T=0$ are identical, connected directly with the zero-energy density of states, and vary from 1.72 in the dirty limit to 0.5−0.6 in the clean limit. This mean-free-path dependence of $\alpha_s$ and $\chi_s$ at $T=0$ is quantitatively the same as that of the slope $\alpha_s(T=0)$ for the flux-flow resistivity studied previously. The result suggests that $S_s(B)$ and $\chi_s(B)$ near $T=0$ are convex downward (upward) in the dirty (clean) limit, deviating substantially from the linear behavior $\propto B/H_{c2}$. The specific-heat jump at $H_{c2}$ also shows fairly large mean-free-path dependence.

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I. INTRODUCTION

This paper considers the changes of the entropy $S_s$ and the spin susceptibility $\chi_s$ through $H_{c2}$ for clean s-wave type-II superconductors. These quantities were calculated by Maki$^{1,2}$ in the dirty limit for superconducting alloys nearly 40 years ago. However, detailed studies on clean systems are still missing even for s-wave superconductors. Writing these quantities as

$$S_s(T,B)−S_s(T,0)/S_s(T)−S_s(T,0)=1−\alpha_s(1−B/H_{c2})=(B/H_{c2})^{\alpha_s},$$

(1a)

$$\chi_s(T,B)−\chi_s(T,0)/\chi_s(T)−\chi_s(T,0)=1−\alpha_x(1−B/H_{c2})=(B/H_{c2})^{\alpha_x},$$

(1b)

the slopes $\alpha_s$ and $\alpha_x$ will be obtained quantitatively for arbitrary impurity concentrations. The results near $H_{c2}$ will also be useful for getting an insight into the behaviors over $0\leq B \leq H_{c2}$. Indeed, $\alpha>1$ ($\alpha<1$) indicates overall field dependence which is convex downward (upward), as seen from Eq. (1).

It seems to have been widely accepted that various physical quantities of clean s-wave type-II superconductors follow the linear field dependence with $\alpha=1$ at low temperatures. A theoretical basis for it is the density of states for a single vortex calculated by Caroli, de Gennes, and Matricon.3,4 However, few quantitative calculations have been carried out so far on the explicit field dependence. Recently, Ichioka et al.$^5$ performed a numerical study on the density of states of clean two-dimensional s-wave superconductors with $\kappa>1$ at $T=0.5T_c$. They found the exponent $\alpha=0.67$ for the overall field dependence of the zero-energy density of states. Also, experiments on the $T$-linear specific-heat coefficient $\gamma_s(B)$ for clean $V_x$Si$^6$, NbSe$_2$7–10 and CeRu$_2$$^{11}$ showed marked upward deviations from the linear behavior $\gamma_s(B/H_{c2})$. Even early experiments on $\gamma_s(B)$ for clean V and Nb indicate similar deviations,$^{12,13}$ although not recognized explicitly in those days due to the absence of a theory on clean systems. These results indicate that the field dependence with $\alpha<1$ is a general feature of clean s-wave superconductors, as suggested by Ramirez.$^6$

Following the preceding works on the Maki parameters$^{14}$ and the flux-flow resistivity,$^{15}$ which will be referred to as I and II, respectively, I here present a detailed study on $S_s$ and $\chi_s$ near $H_{c2}$ at all temperatures. I thereby hope to clarify the $\kappa$ and mean-free-path ($l_u$) dependence of $\alpha_s$ and $\alpha_x$. Calculations are performed for both two- and three-dimensional isotropic systems to see the dependence of $\alpha_s$ and $\alpha_x$ on detailed Fermi-surface structures. I also calculate the specific-heat jump at $H_{c2}$ for various values of $\kappa$ and $l_u$. To my knowledge, this kind of a systematic study has not been performed even for clean s-wave superconductors.

Unlike the convention, I adopt the average flux density $B$ in the bulk as an independent variable instead of the external field $H$. An advantage for it is that the irrelevant region $H < H_{c1}$ is automatically removed from the discussion on the field dependence. This distinction between $B$ and $H$ becomes important for low-$\kappa$ materials where $H < H_{c1}$ occupies a substantial part of $0 < H < H_{c2}$. Any experiment on the $B$ dependence should be accompanied by a careful measurement on the magnetization, especially for low-$\kappa$ materials such as Nb and V.

Section II provides the formulation. Sec. III presents numerical results, and Sec. IV summarizes the paper. The main analytic results are tabulated in Table I for an easy reference. I put $k_B=1$ throughout.

II. FORMULATION

A. Entropy and Pauli paramagnetism

As before,$^{14,15}$ I consider the s-wave pairing with an isotropic Fermi surface and s-wave impurity scattering in an external magnetic field $H \parallel \mathbf{z}$. The formulation proceeds in exactly the same way for both the three-dimensional system and the two-dimensional system placed in the $xy$ plane perpendicular to $H$. The vector potential in the bulk can be written as

$$A(r)=B\hat{y}+\hat{A}(r),$$

(2)
where \( B \) is the average flux density produced jointly by the external current and the supercurrent inside the sample, and \( \mathbf{A} \) expresses the spatially varying part of the magnetic field satisfying \( \int \nabla \times \mathbf{A} \, dr = 0 \).

I first write down the expressions for the entropy and the magnetization in the presence of Pauli paramagnetism. This effect can be included in the Eilenberger equations\(^{22}\) for the quasiclassical Green’s functions \( f, f^\dagger \), and \( g \) by the replacement\(^{23-25}\)

\[
\epsilon_n \rightarrow \epsilon_n^\prime = \epsilon_n - i \mu_B \mathbf{z} \cdot (\nabla \times \mathbf{A}),
\]

where \( \epsilon_n = (2n + 1) \pi T \) is the Matsubara energy and \( \mu_B \) is the Bohr magneton. Choosing \( B \) as an independent variable and measuring the energy from the normal state at the same temperature \( T \) in zero field, the corresponding Eilenberger’s free-energy functional\(^22\) is given by

\[
F = \int dr \left\{ \frac{1}{8\pi} \left[ (\nabla \times \mathbf{A})^2 + \frac{\chi_0}{2} (\nabla \times \mathbf{A})^2 + N(0) |\Delta(\mathbf{r})|^2 \right] \ln \frac{T}{T_c}ight. + \left. \pi TN(0) \sum_{n=-\infty}^{\infty} \left( \frac{|\Delta(\mathbf{r})|^2}{|\epsilon_n|^2} - \langle I(\epsilon_n, \mathbf{k}_F, \mathbf{r}) \rangle \right) \right\},
\]

Here \( \chi_0 = 2 \mu_B^2 N(0) \) is the normal-state spin susceptibility with \( N(0) \) the density of states per one spin and per unit volume at the Fermi level, \( \Delta \) is the pair potential, \( \mathbf{k}_F \) is the Fermi wave vector, and \( \langle \cdots \rangle \) denotes the Fermi-surface average with \( 1 \) = \( 1 \). The quantity \( I \) is defined by\(^{14}\)

\[
I = \Delta^f + \Delta f^\dagger + 2 \epsilon_n^\prime [g - \text{sgn}(\epsilon_n)] + \hbar f(f^\dagger) + (f^\dagger)f^\dagger + \frac{g(g^\prime - 1)}{2 \tau} - \frac{\hbar f^\dagger \mathbf{v}_F \cdot \nabla f - f \mathbf{v}_F \cdot \nabla^\dagger f^\dagger}{2 [g + \text{sgn}(\epsilon_n)]},
\]

where \( \tau \) is the relaxation time in the second-Born approximation, \( \mathbf{v}_F \) is the Fermi velocity, and \( \nabla \) denotes

\[
\nabla = \nabla - \frac{2e}{\hbar c} \mathbf{A}.
\]

The quasiclassical Green’s functions \( f \) and \( g \) are connected by \( g = (1 - ff^\dagger)^{1/2} \text{sgn} \epsilon_n \) with \( f(\epsilon_n, \mathbf{k}_F, \mathbf{r}) = f^\dagger(-\epsilon_n, \mathbf{k}_F, \mathbf{r}). \) The functional derivatives of Eq. (4) with respect to \( f^\dagger, \Delta^\dagger, \) and \( \mathbf{A} \) lead to the Eilenberger equation for \( f \), the self-consistency equation for \( \Delta(\mathbf{r}) \), and the Maxwell equation for \( \mathbf{A} \), respectively.

The expression of the entropy \( S_s \) is obtained from Eq. (4) by the thermodynamic relation. \( S_s = S_n - \partial F/\partial T \). Considering the stationarity with respect to \( f, \Delta, \mathbf{A} \), we only have to differentiate \( F \) with respect to the explicit temperature dependence. We thereby obtain

\[
S_s = S_n = \frac{N(0)}{T} \int dr \left[ |\Delta(\mathbf{r})|^2 - \pi T \sum_{n=-\infty}^{\infty} \langle I(\epsilon_n, \mathbf{k}_F, \mathbf{r}) \rangle \right] - 2 \pi T \sum_{n=-\infty}^{\infty} \epsilon_n \langle g - \text{sgn}(\epsilon_n) \rangle,
\]

where \( S_n = 2 \pi^2 N(0) VT/3 \) with \( V \) the volume of the system. In contrast, the expression of the external field \( H \) may be derived by applying the Dorla-Guberman-Rainer scaling to Eq. (4).\(^{26}\) The details are given in Appendix A of I, and we obtain

\[
H = -4 \pi M_{ap} + B + \frac{1}{BV} \int dr (\nabla \times \mathbf{A})^2 + \frac{\pi^2 TN(0)}{BV} \sum_{n=-\infty}^{\infty} \int dr \left\{ \frac{\hbar^2 f^\dagger \mathbf{v}_F \cdot \nabla f - f \mathbf{v}_F \cdot \nabla^\dagger f^\dagger}{g + \text{sgn} \epsilon_n} \right\},
\]

where \( M_{ap} = \chi_{\alpha \beta} B \) denotes the normal-state magnetization due to Pauli paramagnetism. We thus arrive at the expression of the magnetization from Pauli paramagnetism as

\[
M = M_{ap} - \frac{2 \pi^2 TN(0) \mu_B}{BV} \sum_{n=-\infty}^{\infty} \int dr \langle g \rangle \mathbf{z} \cdot (\nabla \times \mathbf{A}),
\]

When Pauli paramagnetism is negligible compared with the orbital diamagnetism by supercurrent, we can take the limit \( \mu_B \rightarrow 0 \) in Eqs. (7) and (9) and retain only the leading-order terms. This procedure yields \( \epsilon_n^\prime \rightarrow \epsilon_n \) for Eq. (7). On the other hand, Eq. (9) is transformed by noting Eq. (3) into
\[ M_{\text{sp}} = M_0 \left[ 1 - \frac{\pi T}{V} \sum_{n=-\infty}^{\infty} \int d\mathbf{r} \frac{\partial (g)}{\partial e_n} \left( \mathbf{V} \times \mathbf{A} \right) \right]^2. \] (10)

If the zero-field expression \( g = e \sigma / \sqrt{\epsilon_0^2 + |\Delta|^2} \) is substituted into Eq. (10) with \( \mathbf{V} \times \mathbf{A} = \mathbf{B} \mathbf{z} \), the terms in the square bracket reduce to the Yosida function.\(^{27}\)

**B. Expressions near \( H_{c2} \)**

I now consider the cases where Pauli paramagnetism is small and provide explicit expressions to Eqs. (7) and (10) near \( H_{c2} \). From now on I adopt the units used previously\(^{14,15}\), where the energy, the length, and the magnetic field are measured by the zero-temperature energy gap \( \Delta(0) \) at \( H = 0 \), the coherence length \( \xi_0 = \hbar v_F / \Delta(0) \) with \( v_F \) the Fermi velocity, and \( B_0 = \phi_0 / 2 \pi \xi_0^2 \) with \( \phi_0 = h c / 2 e \) the flux quantum, respectively, with \( \hbar = 1 \).

First, \( f, g, \) and \( \tilde{A} \) are expanded up to the second order in \( \Delta(r) \) as

\[
\begin{align*}
f & = f^{(1)}, \\
g & = (1 - \frac{1}{2} f^{(1)} f^{(1)}) \text{sgn}(e_n), \\
\tilde{A} & = \tilde{A}^{(2)}.
\end{align*}
\] (11)

Substituting them into Eqs. (7) and (10) and using the Eilenberger equations for \( f^{(1)} \) and \( f^{(1)*} \) to remove terms with \( v_F \cdot \hat{\mathbf{a}} \), we obtain

\[
\begin{align*}
S_n & = 1 - \frac{3}{2 \pi^2 T^2} \sum_n \int d\mathbf{r} \left[ \left| \Delta(r) \right|^2 - \frac{\pi T}{2} \sum_n \left( f^{(1)*} \Delta + f^{(1)} \Delta^* \right) \right] \\
& \quad + \pi T \sum_n \left| e_n (f^{(1)} f^{(1)}) \right|,
\end{align*}
\] (12a)

\[
\begin{align*}
\frac{M_{\text{sp}}}{M_0} & = 1 + \pi T \sum_n \int d\mathbf{r} \left\{ \frac{\partial f^{(1)*}}{\partial e_n} f^{(1)} + f^{(1)} \frac{\partial f^{(1)*}}{\partial e_n} \right\} \text{sgn}(e_n).
\end{align*}
\] (12b)

Further, \( \Delta(r) \) and \( f^{(1)} \) near \( H_{c2} \) can be expanded in the basis functions \( \psi_{Nq}(r) \) of the vortex lattice as\(^{14}\)

\[
\Delta(r) = \sqrt{\Delta_0} \psi_{0q}(r),
\] (13a)

\[
f^{(1)}(e_n, \mathbf{k}_F, \mathbf{r}) = \sqrt{\Delta_0} \sum_{N=0}^{\infty} f^{(1)}_N(e_n, \mathbf{r}) e^{iNq \cdot \mathbf{r}} \psi_{Nq}(\mathbf{r}),
\] (13b)

where \( (\theta, \varphi) \) are the polar angles of \( \mathbf{v}_F \) with \( \sin \theta = 1 \) in two dimensions, \( N \) denotes the Landau level, and \( q \) is an arbitrary chosen magnetic Bloch vector characterizing the broken translational symmetry of the flux-line lattice and specifying the core locations.\(^{21}\) The coefficients \( \Delta_0 \) and \( f^{(1)}_N \) are both real for the relevant hexagonal lattice. Substituting these expressions into Eqs. (12a) and (12b) and using the orthonormality of \( \psi_{Nq}(r) \) and \( e^{iNq \cdot r} \), we obtain

\[
\begin{align*}
\frac{S_n}{S_0} & = 1 - \frac{3 \Delta_0^2}{2 \pi^2 T^2} \sum_{n=-\infty}^{\infty} \left( f^{(1)}_0 \right)^2 \\
& \quad + \pi T \sum_{n=-\infty}^{\infty} \left| e_n \right| \sum_{N} \left( -1 \right)^N \left( f^{(1)}_N f^{(1)}_N \right) \text{sgn}(e_n).
\end{align*}
\] (14a)

\[
\begin{align*}
\frac{M_{\text{sp}}}{M_0} & = 1 + \pi T \sum_{n=-\infty}^{\infty} \left| e_n \right| \sum_{N} \left( -1 \right)^N \left( \frac{\partial f^{(1)}_N}{\partial e_n} \right) \text{sgn}(e_n).
\end{align*}
\] (14b)

Except \( \Delta_0^2 H_{c2} - B \), all the quantities in Eqs. (14a) and (14b) are to be evaluated at \( H_{c2} \).

It is possible to give an alternative convenient expression to Eq. (14a). To this end, we make use of the equation for \( H_{c2} \) obtained as Eq. (33) of I:

\[
\ln \frac{T_c}{T} + \pi T \sum_{n=-\infty}^{\infty} \left( f^{(1)}_0 \right)^2 - \frac{1}{\left| e_n \right|^2} = 0.
\] (15)

Differentiating Eq. (15) with respect to \( T \) yields

\[
-1 + \pi T \sum_{n} \left( f^{(1)}_0 \right) + \left( \frac{\partial f^{(1)}_0}{\partial e_n} \right) \left( \frac{\partial \left( f^{(1)}_0 \right)}{\partial e_n} \right) \frac{dH_{c2}}{dT} = 0.
\] (16)

The quantity \( \frac{\partial f^{(1)}_0}{\partial H_{c2}} \) has been calculated as Eqs. (31) and (32) of I:

\[
\frac{\partial f^{(1)}_0}{\partial H_{c2}} = \sum_{N} \left( -1 \right)^{N+1} \frac{N+1}{8 H_{c2}} \left( f^{(1)}_N f^{(1)}_N \right) \sin \theta.
\] (17)

A similar procedure leads to the analytic expressions for \( \frac{\partial f^{(1)}_0}{\partial e_n} \) and \( \frac{\partial f^{(1)}_N}{\partial e_n} \) in Eq. (14b) as

\[
\begin{align*}
\frac{\partial f^{(1)}_0}{\partial e_n} & = -\sum_{N} \left( -1 \right)^N \left( f^{(1)}_N f^{(1)}_N \right) \text{sgn}(e_n), \\
\frac{\partial f^{(1)}_N}{\partial e_n} & = -\sum_{N} K_N \left( f^{(1)}_N f^{(1)}_N \right) \frac{\partial f^{(1)}_0}{\partial e_n} \text{sgn}(e_n) \frac{\partial f^{(1)}_0}{\partial e_n}.
\end{align*}
\] (18a)

(18b)

where \( K_N \) is defined by Eq. (25) of I. Using Eqs. (16) and (18a) in Eq. (14a), we obtain

\[
\begin{align*}
\frac{S_n}{S_0} & = 1 - \frac{3 \Delta_0}{2 \pi^2 T} \sum_{n=-\infty}^{\infty} \left( f^{(1)}_0 \right) \frac{dH_{c2}}{dT} \\
& \quad + \pi T \sum_{n=-\infty}^{\infty} \left| e_n \right| \sum_{N} \left( -1 \right)^N \left( \frac{\partial f^{(1)}_N}{\partial e_n} \right) \text{sgn}(e_n),
\end{align*}
\] (19)

with

\[
\begin{align*}
\frac{dH_{c2}}{dT} & = \left( \frac{\partial f^{(1)}_0}{\partial e_n} \right) \frac{dH_{c2}}{dT} \\
& \quad + \pi T \sum_{n=-\infty}^{\infty} \left| e_n \right| \sum_{N} \left( -1 \right)^N \left( \frac{\partial f^{(1)}_N}{\partial e_n} \right) \text{sgn}(e_n).
\end{align*}
\] (20)

Equation (20) also enables us to calculate the specific-heat jump at \( H_{c2} \). Indeed, it is given in conventional units as\(^{1}\)
\[ \Delta C = \frac{T}{4\pi} \left( \frac{dH_c}{dT} \right)^2 \frac{1}{(2\kappa^2-1)|\beta_A|}. \]  

(21)

where \( \kappa \) is the Maki parameter\(^{1,2}\) and \( \beta_A = 1.16 \).

Equations (14) and (21) with Eqs. (17), (18), and (20) are the main analytic results of the paper. The quantities \( \Delta_0 \), \( \tilde{f}^{(1)}_N \), and \( \kappa \) have been obtained in I. The explicit expression for \( \tilde{f}^{(1)}_N \) is given by

\[ \tilde{f}^{(1)}_N = \frac{\tilde{K}^0_N \text{sgn}(e_n)}{1 - \langle \tilde{K}^0_N \rangle \text{sgn}(e_n)/2\tau}, \]  

(22)

where \( \tilde{K}^N \) may be calculated efficiently by the procedure in Sec. IIIF of I, with a change of definition of \( \tilde{e}_n \) as

\[ \tilde{e}_n = \left( |e_n| + \frac{1}{2\tau} \right) \text{sgn}(e_n). \]  

(23)

C. Analytic results at \( T = 0 \)

Now it will be shown that Eqs. (14a) and (14b) reduce to an identical expression at \( T = 0 \) for arbitrary impurity concentrations, which has the physical meaning of the zero-energy density of states.

Let us start from Eq. (14a) where \( e_n > 0 \) and \( e_n < 0 \) yield the same contribution. Using this fact and Eq. (18a), it is transformed into

\[ \frac{S_s}{S_n} = 1 - \frac{3\Delta_0^2}{2\pi^2T} \left[ 1 - \frac{2\pi T}{0} \sum_{n=0}^\infty \left( \tilde{f}^{(1)}_0 + \epsilon_n \frac{\partial \tilde{f}^{(1)}_0}{\partial \epsilon_n} \right) \right]. \]  

(24)

The summation over \( n \) for \( T \to 0 \) may be performed by using the Euler-Maclaurin formula and the asymptotic property \( \tilde{f}^{(1)}_0(e_n) \to e_n^{-1} \) (\( e_n \to \infty \)).\(^{14}\) For example,

\[ 2\pi T \sum_{n=0}^\infty \langle \tilde{f}^{(1)}_0(e_n) \rangle \approx \int_0^\infty \langle \tilde{f}^{(1)}_0(e) \rangle \, de + \pi T \langle \tilde{f}^{(1)}_0 \rangle (\pi T) \]

\[ - \frac{(\pi T)^2}{3} \langle \tilde{f}^{(1)}_0 \rangle (\pi T) \]

\[ + \int_0^\infty \langle \tilde{f}^{(1)}_0(e) \rangle \, de + \frac{(\pi T)^2}{6} \langle \tilde{f}^{(1)}_0 \rangle (0). \]

(25)

We thereby obtain

\[ \frac{S_s}{S_n} \to 1 + \frac{\Delta_0^2}{2} \langle \tilde{f}^{(1)}_0 \rangle (0). \]  

(26a)

Equation (14b) may be transformed similarly as

\[ \frac{M_{sp}}{M_{ap}} = 1 - \pi T \Delta_0^2 \sum_{n=0}^\infty \frac{\partial^2 \langle \tilde{f}^{(1)}_0(e_n) \rangle}{\partial \epsilon_n^2} \to 1 + \frac{\Delta_0^2}{2} \langle \tilde{f}^{(1)}_0 \rangle (0). \]  

(26b)

Thus, \( S_s/S_n = M_{sp}/M_{ap} \), or equivalently, \( \alpha_s = \alpha \) at \( T = 0 \) for arbitrary impurity concentrations.

Equations (26a) and (26b) have a simple physical meaning. Indeed, noting Eqs. (11), (13b), and (18a), we find an alternative expression at \( T = 0 \):

\[ \frac{S_s}{S_n} = \frac{M_{sp}}{M_{ap}} = \frac{1}{V} \int \langle g(e_n, 0, \kappa_r) \rangle \, dr, \]  

(27)

which is nothing but the normalized density of states at \( e = 0 \). Thus, the entropy and the spin susceptibility at \( T = 0 \) are both determined by the zero-energy density of states.

The coefficient of \( \Delta_0 \approx H_c - B \) have been obtained in I. Also, \( \tilde{f}^{(1)}_0(0) \) in Eqs. (26a) and (26b) may be calculated efficiently from Eq. (22) by using the analytic expression\(^{14}\)

\[ \langle \tilde{K}^0_N \rangle (\tilde{e}_n) = \sqrt{\frac{2}{\pi}} \int_0^\infty \tilde{e}_n \tilde{e}_n^2 + x^2 \beta^2 e^{-x^2/\beta^2} \, dx, \]  

(28)

with \( \beta = \sqrt{H_c \sin \theta / 2} \). Hence, Eqs. (26a) and (26b) at \( T = 0 \) can be evaluated easily.

D. Analytic results in the dirty limit

I here summarize analytic results in the dirty limit \( \tau \to 0 \). First, the key quantities \( \tilde{R}^0_N \) are calculated by choosing \( N_{on} = 1 \) in the procedure in Sec. IIIF of I. The results are given by

\[ \tilde{R}^0_N = \frac{\tilde{e}_n}{\tilde{e}_n^2 + \beta^2}, \quad \tilde{R}^1_N = \frac{\beta}{\tilde{e}_n^2 + \beta^2}. \]  

(29)

Since \( \beta^2 \) is of the order of \( 1/\tau \), as shown below, \( \langle \tilde{R}^0_N \rangle \) may be approximated as \( \langle \tilde{R}^0_N \rangle \approx \langle \tilde{R}^0_N \rangle \approx \langle \tilde{e}_n \rangle / \langle \tilde{e}_n^2 + \beta^2 \rangle \). Using this \( \langle \tilde{R}^0_N \rangle \) in Eq. (22) and retaining only the leading-order contributions, we obtain

\[ \tilde{f}^{(1)}_0 = \frac{1}{|\epsilon_n| + 2\tau \beta^2}, \quad \tilde{f}^{(1)}_0 = \frac{2\tau \beta \text{sgn}(\epsilon_n)}{|\epsilon_n| + 2\tau \beta^2}. \]  

(30)

Notice that \( \tilde{f}^{(1)}_0 \) is smaller than \( \tilde{f}^{(1)}_0 \). Substitution of Eq. (30) into Eq. (15) leads to the equation for \( H_c \) obtained by Maki\(^{28}\) and de Gennes\(^{29}\):

\[ \ln(T_c/T) + \psi(1/2) - \psi(x) = 0, \]  

(31)

where \( \psi \) is the digamma function, and \( x \) is defined by

\[ x = \frac{1}{2} + \frac{\tau \beta^2}{\pi T \epsilon} = \frac{1}{2} + \frac{\tau H_c}{4\pi T d}, \]  

(32)

with \( d = 2.3 \) the dimension of the system. As shown by Maki\(^{28}\), Eq. (31) can be solved near \( T = 0 \) by using the asymptotic expression of \( \psi(x) \) as

\[ H_c \approx \frac{d}{\tau} \left[ 1 - \frac{2}{3} (\pi T)^2 \right]. \]  

(33)

Thus \( \beta^2 \approx H_c \sim \tau^{-1} \), as assumed at the beginning. Differentiating Eq. (31) with respect to \( T \), we obtain
Finally, $\kappa_2$ and $[\Delta_0(B)]^2$ are calculated from Eqs. (34b) and (36) of I as

$$\kappa_2 = \frac{d}{\sqrt{2}} \frac{\langle \hat{\psi}^{(2)}(x) \rangle}{\sqrt{2} \tau H_{c2} \psi'(x)} \left. \kappa_0 \right| \rightarrow \frac{\kappa_0}{\sqrt{2}}$$

(35)

$$\Delta_0^2 = \frac{(H_{c2} - B)^2 \kappa_0^2 (2 \kappa_2^2 - 1) \beta_A + 1}{(2 \kappa_2^2 - 1) \beta_A + 1} - \frac{4 \pi T d \tau H_{c2} \psi'(x)}{\kappa_0^2}$$

(36)

where $\kappa_0$ is defined by $\kappa_0 = \phi_0 / 2 \pi \xi_0^2 H_{c2}(0)$ with $H_{c2}(0)$ the thermodynamic critical field at $T=0$. Equation (35) agrees with the result by Caroli, Cyrot, and de Gennes.30

Now, let us substitute Eq. (30) into Eqs. (14b) and (19) and use Eq. (33). We thereby obtain

$$S_n = S_n + \frac{d H_{c2}}{d T} + \frac{3}{8 \pi^2 T^2} \psi'(x)_{\tau=0} - 2 \Delta_0^2$$

(37a)

$$M_{d} = M_{d} + \frac{1}{8 \pi^2 T^2} \psi'(x)_{\tau=0} - 2 \Delta_0^2$$

(37b)

Thus, $M_d / M_{d0}$ and $S_n / S_n$ are the same at $T=0$, in agreement with Eq. (27); they both determine the zero-energy density of states. Equation (37b) is the result first obtained by Maki.2 Also, the expression $1 - 2 \Delta_0^2$ for the normalized zero-energy density of states at $T=0$ agrees with the result for the local density of states by de Gennes.42

Equation (36) tells us that $\Delta_0^2 = (1 - B/H_{c2}) \beta_A$ as $T \rightarrow 0$ for $\kappa_2 \gg 1$. We hence find from Eqs. (1), (37a), and (37b) that the initial slopes at $T=0$ for $\kappa_2 \gg 1$ are given by

$$\alpha_s = \alpha_c = 2 \beta_A = 1.72.$$  

(38)

The results suggest the overall field dependence of $S_n$ and $\chi_s$ at $T=0$ which is convex downward. Notice that the flux-flow resistivity $\rho_f$ at $T=0$ also has the same initial slope $\alpha_f = 1.27$ in the dirty limit.15,31,32 These results strongly suggest that the density of states at $e=0$ is mainly relevant to the physical properties of the vortex state at $T=0$.

E. The case with $p$-wave impurity scattering

If the $p$-wave impurity scattering is relevant, the following additional terms appear on the right-hand side of Eq. (5):

$$d \left( \hat{k} \cdot \hat{k}' \right) \frac{1}{4 \tau_1} \left( \hat{k}' \cdot \hat{k}' \right) + d \left( \hat{k} \cdot \hat{k}' \right) \frac{1}{2 \tau_1}$$

(39)

where $\left( \hat{k}' \cdot \hat{k} \right) = \hat{k} \cdot \hat{k} \left( e_n, k_{\Psi}, r \right)$, for example, $\tau_1$ is the $p$-wave relaxation time, and $\hat{k}$ is the unit vector along $k_F$. However, Eqs. (7), (9), and (10) remain unchanged once $I$ is modified as above.

The corresponding calculations near $H_{c2}$ may be performed as described in Appendix A of I. It thereby follows that Eqs. (14) and (21) are also valid together with Eqs. (17), (18a), and (20), where $\tilde{f}^{(1)}_N$ is now given by

$$\tilde{f}^{(1)}_N = \frac{1}{D} \left[ 1 - \frac{d}{4 \tau_1} \left( \tilde{K}_1 \sin ^2 \theta' \right) \left( \tilde{K}_n \right) \tilde{K}_n \right]$$

(40)

with

$$D = \left[ 1 - \frac{1}{2 \tau} \left( \tilde{K}_0 \right) \left( \tilde{K}_0 \right) \tilde{K}_0 \sin ^2 \theta' \right]$$

(41)

In addition, Eq. (18b) is to be replaced by

$$\frac{\partial \tilde{f}^{(1)}_N}{\partial \epsilon_n} = - \sum_N \tilde{K}_N \frac{\partial \tilde{f}^{(1)}_N}{\partial \epsilon_n}$$

(42)

where

$$\frac{\partial \tilde{f}^{(1)}_N \sin \theta'}{\partial \epsilon_n} = - \sum_N \left( -1 \right) \tilde{K}_N \frac{\partial \tilde{f}^{(1)}_N \sin \theta'}{\partial \epsilon_n}$$

(43)

with

$$\tilde{f}^{(1)}_N = \frac{1}{D} \left[ 1 - \frac{1}{2 \tau} \left( \tilde{K}_0 \right) \left( \tilde{K}_0 \right) \tilde{K}_n \sin \theta' \right]$$

(44)

Finally, the analytic results in the dirty limit are the same as those given in Sec. II D with a replacement of $\tau$ by the transport lifetime $\tau_u$ defined through

$$\frac{1}{\tau_u} = \frac{1}{\tau} - \frac{1}{\tau_1}$$

(45)

F. Numerical procedures

I have adopted the same parameters as I and II to express different impurity concentrations:

$$\xi_{r} / l_{w} = 1 / 2 \pi T c r_{u} \ L_{w} / \tau$$

(46)

Numerical calculations of Eqs. (14) and (21) with Eqs. (17), (18a), and (20) have been performed for each set of parameters by restricting every summation over the Matsubara frequencies to those satisfying $|\epsilon_n| \leq \epsilon_c$. Choosing $\epsilon_c = 200$ is sufficient to obtain an accuracy of $\sim 0.1\%$ for Eqs. (14b) and (21), whereas $\epsilon_c = 20\,000$ (4000) is required for Eq. (14a) in the dirty (clean) limit. Summations over Landau levels have been truncated at $N = N_{cut}$ where I put $N_{cut} = 1$ in the calculation of $\tilde{K}_N$; see Sec. II F of I for the details. Enough
curves start from the same value $a_{100, 200, 1500, and 4000}$ for $j$ 

a is small, as expected from $N$. The linear behavior

In particular, the curves of $a$ have been calculated in three dimensions for $\nu$ 

~purity concentrations parametrized by Eq. $T_{1.72}$ in the dirty limit to around 0.6 for $j$. We thus realize that the zero-energy density of states is mainly determined by the mean free path, and may not depend much on the Fermi-surface structures nor the details of the impurity scattering.

For comparison, Fig. 3 presents the slope $\alpha_p$ for the flux-flow resistivity $\rho_f$ calculated previously$^{15}$ for the same parameters as in Fig. 1. At finite temperatures in the dirty limit, $\alpha_p$ is much larger than $\alpha_s$ and $\alpha_x$, indicating a steeper decrease of $\rho_f$ just below $H_{c2}$. However, the difference is seen to diminish as the temperature is reduced, and it has been checked numerically that $\alpha_s = \alpha_x = \alpha_p$ holds at $T = 0$ for arbitrary impurity concentrations. This fact suggests that $\rho_f$ at $T = 0$ is also determined by the zero-energy density of states.

Next, we examine the dependence of the slopes on the Ginzburg-Landau parameter $\kappa_{GL}$. Figure 4 shows the same curves as in Fig. 1 near the type-I–type-II boundary of $\kappa_{GL} = 1$. Each curve is shifted upwards from the corresponding one in Fig. 1 for $\kappa_{GL} = 50$, but the quantitative difference is

III. RESULTS

Figure 1 shows temperature dependence of $\alpha_s$ and $\alpha_x$ defined by Eqs. (1a) and (1b), respectively, for different impurity concentrations parametrized by Eq. (46). They have been calculated in three dimensions for $l_u/l = 1.0$ and $\kappa_{GL} = 50$, where $\kappa_{GL}$ is the Ginzburg-Landau parameter.$^{33}$ All the curves start from the same value $a_{0.3} = a_{0.862}$ at $T = T_c$ and develop differences among different impurity concentrations at lower temperatures. The equality $a_{0.3} = a_{0.862}$ holds at $T = 0$, as shown by Eq. (27), and the value decreases from 1.72 in the dirty limit to around 0.6 for $\xi_E/l_u = 0.1$. According to Eq. (27), this variation in the slope at $T = 0$ can be attributed to the mean-free-path dependence of the zero-energy density of states $N_c(0,B)$. In particular, $N_c(0,B)$ in the dirty (clean) limit decreases more rapidly (mildly) than the linear behavior $N(0)B/H_{c2}$ near $H_{c2}$. From this result, we expect the overall field dependence of the entropy and the spin susceptibility at $T = 0$ which is convex downward (upward) in the dirty (clean) limit, as seen from Eq. (1).

The difference between $a_{0.3}$ and $a_{0.862}$ at finite temperatures is small, as expected from $a_{0.3} = a_{0.862}$ holding at $T = 0$ and $T_c$. In particular, the curves of $\alpha_x$ and $\alpha_x$ in the dirty limit depend neither on the dimensions nor $l_u/l$. However, the dependenc...
FIG. 4. Slope $\alpha_s$ as a function of $T/T_c$ for different impurity concentrations with $d=3$, $l_u/l=1$, and $\kappa_{GL}=1$.

rather small. This is also the case for $\alpha_x$. Thus, the slopes $\alpha_s$ and $\alpha_x$ defined in terms of $B$ do not have large $\kappa_{GL}$ dependence.

Finally, Fig. 5 plots the specific-heat jump $\Delta C$ over $T$ at $H_{c2}$ as a function of $T/T_c$ for different impurity concentrations with $d=3$, $l_u/l=1$, and $\kappa_{GL}=50$. It is normalized by the corresponding quantity at $T=T_c$ and $H=0$, i.e., $\Delta C(T_c)/T_c=1.43$ in the weak-coupling model. The curves change gradually from almost $T$-linear overall temperature dependence in the dirty limit to $T^2$ dependence in the clean limit, and approach zero as $\propto T^2$ at lowest temperatures. Although the ratio near $T_c$ is strongly dependent on $\kappa_{GL}$ at $0.5-0.6$ in the clean limit. This change is due completely to the mean-free-path dependence of the zero-energy density of states. The fact also suggests variation of the overall field dependence at $T=0$ from convex downward in the dirty limit to upward in the clean limit. (v) The slopes have only small dependence on the dimensions and the details of the impurity scattering. (vi) The slope $\alpha_B$ for the flux-flow resistivity $\rho_f$, which has been calculated previously, also shows a complete numerical agreement at $T=0$ with $\alpha_s$ and $\alpha_x$. This fact indicates that the zero-energy density of states is also responsible for $\rho_f$ at $T=0$.

The $T$-linear specific-heat coefficient $\gamma_s(B)$ observed in clean materials presents curves with $\gamma<1$, which is in a qualitative agreement with the present calculation. On the other hand, $\gamma_s(B)$ for dirty samples follows the well-accepted linear field dependence $\propto B/H_{c2}$ and apparently in contradiction with the present result in the dirty limit. However, it should be noted that a careful experiment on $\rho_f$ shows field dependence near $T=0$ which is convex downward, and experimentally obtained $\alpha_B$ agrees quantitatively with the dirty-limit theory. Detailed experiments on the mean-free-path dependence of $\gamma_s(B)$ and $\rho_f(B)$ are desired to remove these discrepancies.

IV. SUMMARY

The entropy and the spin susceptibility near $H_{c2}$ have been calculated for $s$-wave type-II superconductors with arbitrary impurity concentrations. The results have been expressed conveniently with respect to the initial slopes $\alpha_s$ and $\alpha_x$ defined by Eq. (1). The main conclusions are summarized as follows: (i) $\alpha_s=\alpha_x$ holds both at $T=0$ and $T=T_c$. (ii) $\alpha_s=\alpha_x=0.862$ at $T=T_c$ for all impurity concentrations. (iii) At $T=0$, the slope $\alpha_s$ decreases from 1.72 in the dirty limit to $0.5-0.6$ in the clean limit. This change is due completely to the mean-free-path dependence of the zero-energy density of states. The fact also suggests variation of the overall field dependence at $T=0$ from convex downward in the dirty limit to upward in the clean limit. (iv) The slopes have only small dependence on the dimensions and the details of the impurity scattering. (v) The slope $\alpha_B$ for the flux-flow resistivity $\rho_f$, which has been calculated previously, also shows a complete numerical agreement at $T=0$ with $\alpha_s$ and $\alpha_x$. This fact indicates that the zero-energy density of states is also responsible for $\rho_f$ at $T=0$.

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