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Specific heat of REBaSrCu₃O₇ compounds (R = Nd, Sm, Dy, Er)

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Abstract

Specific heat measurements have been performed on polycrystalline REBaSrCu₃O₇ (R = Nd, Sm, Dy, Er) in the temperature range from 0.3 to 3 K. Antiferromagnetic order is indicated by the λ -type peak in the specific heat at 1.2 and 0.6 K for Dy and Er compounds, respectively, whereas only a broad anomaly around 0.6 K is observed for Nd and Sm compounds. The specific heat data for Nd and Sm compounds are well fit with a Schottky-type anomaly with an energy splitting of 1.4 K, indicating the existence of an internal field of 2.42 T at the RE site. The magnetic singlet ground state of ErBaSrCu₃O₇ and DyBaSrCu₃O₇ is produced by exchange interactions of -0.05 and -0.035 T f.u./ μ_B , respectively.

Keywords: Crystal field; High- T_C ; REBaSrCu₃O₇; Specific heat

High- T_C superconductors REBa₂Cu₃O₇ (RE-123O₇, RE = most rare earths) crystallize in the orthorhombic structure. The c -axis (and the cell volume) increases continuously with increasing radius of the RE³⁺ ion. The superconducting temperature T_C (~ 90 K) exhibits a very small change for different RE ions. Compounds REBaSrCu₃O₇ are known to belong to the RE-123O₇ family with T_C between 54 and 86 K. However, with Sr substitution, compounds with ionic radius $r_{RE} < r_{Dy}$ preserve an orthorhombic structure, while those with $r_{RE} > r_{Dy}$ have a tetragonal structure [1]. The c -axis (and the cell volume) was found to decrease from RE = La to Dy, and then increase again. The maximum T_C is obtained for RE = Gd and Dy, i.e. close to the tetragonal/orthorhombic phase transition. The Dy compound can be stabilized in both the orthorhombic and tetragonal structure with the same oxygen content, depending on the heat treatment [2].

In this paper, we present specific heat measurements performed on polycrystalline REBaSrCu₃O₇ (RE = Nd, Sm, Dy, Er) in the temperature range from 0.3 to 3 K.

REBaSrCu₃O_{7- δ} (R = Nd, Sm, Dy, Er; $\delta = 0.06$ – 0.07) samples were prepared by a standard ceramic technique. Details of sample preparation, crystal structure and oxygen content determination, and studies of the superconducting properties have been published elsewhere [1,2].

Specific heat measurements in the temperature range 0.3–3 K were performed in a ³He cryostat. Measurements were carried out on five samples with Nd, Sm (tetragonal), Dy (both tetragonal and orthorhombic) and Er (orthorhombic). For comparison, data for the (same) RE-123O₇ compounds [3,4] are also reported.

The specific heat $C(T)$ of NdBaSrCu₃O₇ is shown in Fig. 1. The broad anomaly around 0.6 K becomes more visible in the Sr-substituted compound, whereas the small peak indicating the antiferromagnetic order of the Nd³⁺

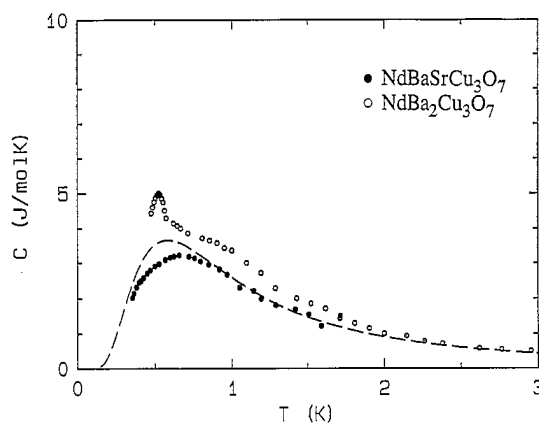


Fig. 1. Specific heat plotted as C vs. T for NdBaSrCu₃O₇ (●). The data for Nd-123O₇ (○) are taken from Ref. [4]. (---) The Schottky anomaly with an energy splitting ΔE of 1.4 K.

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ions in the Nd-123O₇ compound at $T_N = 0.5$ K [4] disappears. The entropy amounts to 4.8 J/kmol at 3 K for Nd-123O₇, whereas it reaches only 4 J/kmol for the Sr-substituted compound. The $C(T)$ data for SmBaSrCu₃O₇ are shown in Fig. 2. The sharp peak at $T_N = 0.6$ K with an entropy of almost $R \ln 2$ disappears with Sr substitution and only a broad anomaly is observed for this compound around 0.6 K. This indicates that the substitution of Ba by Sr causes not only a change from the orthorhombic to the tetragonal structure, but also affects the magnetic properties of the Nd and Sm compounds. Our calculations show that the $C(T)$ data for the two Sr-substituted compounds with Nd and Sm exhibiting a broad anomaly around 0.6 K follow a Schottky-type curve for a two-level system with an energy splitting ΔE of 1.4 K. Such a two-level system is expected to originate from the Kramers doublet ground state as found for other Nd and Sm systems [5]. The splitting indicates the presence of an internal field of 2.42 T at the RE site.

The specific heat $C(T)$ of Dy and ErBaSrCu₃O₇ compounds is shown in Fig. 3. The formation of antiferromagnetic order is indicated by the λ -type peak at 1.2 and 0.6 K, respectively. Sr substitution results in an enhancement of T_N (from 0.9 to 1.2 K) for the Dy compound. The $C(T)$ data are almost the same for Dy compounds with tetragonal and orthorhombic structure except for a slight downward shift of T_N of 0.1 K. The entropy reaches a value of $R \ln 2$ in these compounds. Moreover, superconducting properties are found to be very similar for both phases. The Curie behaviour of the paramagnetic susceptibility, with the effective paramagnetic moment close to the Dy³⁺ free-ion value, was not found to be influenced by the different structures [2]. The Dy magnetic moments lie along the c -axis, and thus the change of anisotropy in the a - b plane due to the tetragonal–orthorhombic phase transition would not affect the magnetic properties. The Er

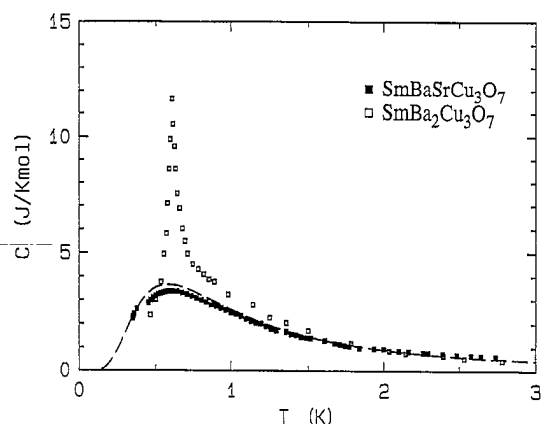


Fig. 2. Specific heat plotted as C vs. T for SmBaSrCu₃O₇ (■). The data for Sm-123O₇ (□) are taken from Ref. [4]. (---) The Schottky anomaly with an energy splitting ΔE of 1.4 K.

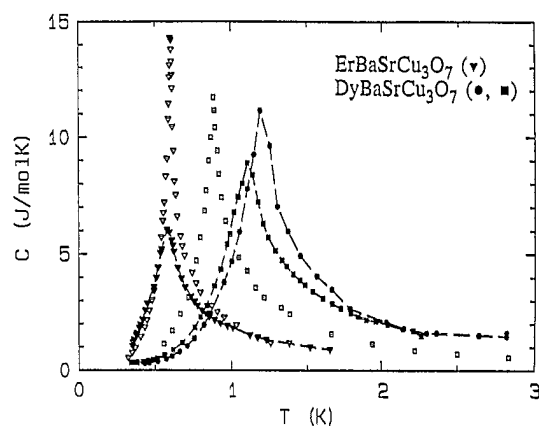


Fig. 3. Specific heat plotted as C vs. T for tetragonal (■) and orthorhombic (●) DyBaSrCu₃O₇, and for orthorhombic ErBaSrCu₃O₇ (▼). The data for Dy- (□) and Er-123O₇ (▽) are taken from Refs. [4,3].

compound forms the orthorhombic structure. Sr substitution causes only a broadening of the λ -type peak at T_N , which is mostly related to the change of the anisotropy ratio in the a - b plane. Moreover, the entropy only reaches 70% of $R \ln 2$ in this case. As the Néel temperature is the same at 0.6 K, we expect that the magnetic singlet ground state, produced by weak exchange interactions of -0.05 T f.u./ μ_B as for the Er-123O₇ compound [3], would also be appropriate for ErBaSrCu₃O₇. The exchange interaction was estimated to be -0.027 and -0.035 T f.u./ μ_B for Dy-123O₇ and DyBaSrCu₃O₇, respectively.

In conclusion, the substitution of Ba by Sr in RE-123O₇ superconductors has a significant influence on the structural and magnetic properties for light RE (Nd, Sm), whereas only a small effect is observed for heavy RE (Dy, Er). The specific heat data for Nd and Sm compounds indicate the presence of an internal field of 2.42 T at the RE site. All considered RE ions are Kramers systems and lifting of the Kramers degeneracy may be the reason for the large low-temperature specific heat [5].

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