

Research Article

Amorphization Effect for Kondo Semiconductor $\text{CeRu}_2\text{Al}_{10}$

Yusuke Amakai,^{1,2} Yasuhiro Shiojiri,¹ Kei Ishihara,¹ Hiroto Hitotsukabuto,¹
Shigeyuki Murayama,¹ Naoki Momono,^{1,2} and Hideaki Takano¹

¹Graduate School of Engineering, Muroran Institute of Technology, 27-1 Mizumoto-cho, Muroran, Hokkaido 050-8585, Japan

²Research Center for Environmentally Friendly Materials Engineering, Muroran Institute of Technology,
27-1 Mizumoto-cho, Muroran, Hokkaido 050-8585, Japan

Correspondence should be addressed to Yusuke Amakai; a-rain@mmm.muroran-it.ac.jp

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We measured the magnetic susceptibility χ , electrical resistivity ρ , and specific heat C_p of a sputtered amorphous (*a*-) $\text{CeRu}_2\text{Al}_{10}$ alloy. χ value for *a*- $\text{CeRu}_2\text{Al}_{10}$ alloy follows a Curie-Weiss paramagnetic behavior in the high-temperature region, and magnetic transition was not observed down to 2 K. The effective paramagnetic moment p_{eff} is $1.19 \mu_B/\text{Ce-atom}$. The resistivity shows a typical disordered alloy behavior, that is, small temperature dependence for the whole temperature range. We observed an enhancement of ρ and C_p/T in the low-temperature region of $T < 10$ K. The enhancement in ρ is suppressed by applying a magnetic field. It is suggested that this behavior is caused by the Kondo effect.

1. Introduction

The ternary rare-earth compound, $\text{CeRu}_2\text{Al}_{10}$, exhibits an unusual antiferromagnetic phase transition at $T_0 \approx 27$ K [1, 2]. The resistivity for $\text{CeRu}_2\text{Al}_{10}$ exhibits a semiconducting behavior in the paramagnetic phase. However, this behavior is suppressed by substituting La in Ce-site of $\text{CeRu}_2\text{Al}_{10}$. In addition, T_0 decreases rapidly with an increase in La concentration and disappears when about half of Ce is substituted by La. The resistivity for Ce dilute region of La-substituted $\text{CeRu}_2\text{Al}_{10}$ exhibits a metallic behavior at high temperatures, and the resistivity exhibits a minimum in the low-temperature region. Tanida et al. proposed that the long-range magnetic order was suppressed randomly in Ce dilute region for La substitution of $\text{CeRu}_2\text{Al}_{10}$. The authors pointed out that the resistivity minimum can be explained by a typical impurity Kondo effect, where Ce exists as a magnetic impurity [3, 4]. Moreover, the semiconducting band gap of $\text{CeRu}_2\text{Al}_{10}$ is broken by La substitution.

However, several studies on the Kondo effect of structural-disordered Ce-alloys such as bulk metallic glasses and amorphous alloys, where Ce-atom is arranged randomly, have been recently performed. For example, in Ce-Al bulk metallic glasses, the tunable competition between the Kondo

effect and the Ruderman-Kittel-Kasuya-Yoshida (RKKY) interaction with the variation in Ce-concentration and the magnetic field [5] is suggested. We studied the low-temperature properties of several binary amorphous Ce-alloys. Amorphous (*a*-)Ce-Mn and *a*-Ce-Ru alloys in Ce high-concentration region exhibit a large electronic specific heat coefficient γ (>200 mJ/molK²) and T^2 law with a large coefficient A ($>0.02 \mu\Omega\text{cm}/\text{K}^2$) in the low-temperature resistivity [6–10]. From these results, we show that an itinerant heavy-fermion state occurs at low temperatures as a Fermi-liquid ground state in the structure-disordered system after the formation of a dense Kondo state.

In this study, to investigate the influence of the structural-disordered effect on the electrical resistivity and magnetic properties for $\text{CeRu}_2\text{Al}_{10}$, we prepared *a*- $\text{CeRu}_2\text{Al}_{10}$ alloy and measured its magnetic susceptibility, resistivity, and specific heat. In addition, we prepared *a*- $\text{LaRu}_2\text{Al}_{10}$ alloy that does not have 4*f*-electrons with a rare-earth element for comparison.

2. Experimental

Bulk ingots of $\text{CeRu}_2\text{Al}_{10}$ and $\text{LaRu}_2\text{Al}_{10}$ were prepared by using the arc-melting method with a stoichiometric

composition of Ce 99.9%, La 99.9% (Nippon Yttrium Co., Ltd.), Ru 99.95% (Rare Metallic Co., Ltd.), and Al 99.99% (Mitsui Chemicals Co., Ltd.), in Ar atmosphere. The amorphous alloy was prepared using a dc high-rate sputtering method with arc-melt ingots on a water-cooled Cu substrate (30 mm ϕ). The sample thickness was ~ 200 μm . The structure of the obtained samples was confirmed using X-ray diffraction measurements. Measurements were performed with an as-sputtered film. The chemical compositions of the present amorphous alloys were determined using scanning electron microscope energy dispersive X-ray spectroscopy (SEM-EDS) to be $\text{Ce}_{10}\text{Ru}_{15}\text{Al}_{75}$ and $\text{La}_8\text{Ru}_{17}\text{Al}_{75}$ (suffixes represent at%). Thus, we will use the notations $a\text{-CeRu}_2\text{Al}_{10}$ and $a\text{-LaRu}_2\text{Al}_{10}$ for the samples, hereafter. The magnetic susceptibility was measured using a commercial SQUID magnetometer (Quantum Design MPMS) from 2 to 300 K. The electrical resistivity was measured by using a typical four-terminal method (Quantum Design PPMS) from 2 to 300 K. The resistivity measurements of $a\text{-CeRu}_2\text{Al}_{10}$ alloy were performed in a magnetic field (0, 20 kOe, 40 kOe, and 60 kOe) and the temperature range was 2–60 K. The specific heat was measured by PPMS from 2 to 300 K.

3. Results and Discussion

Figure 1 shows the X-ray diffraction patterns for $a\text{-CeRu}_2\text{Al}_{10}$ and $a\text{-LaRu}_2\text{Al}_{10}$ alloys. The diffraction patterns for both alloys exhibit two broad peaks at the center at about 25° and 42° , and definite Bragg peaks are not observed. Therefore, the samples are identified as amorphous materials.

Figure 2 shows the temperature dependence of the magnetic susceptibility χ (left axis) and the inverse susceptibility $1/(\chi - \chi_0)$ (right axis) for $a\text{-CeRu}_2\text{Al}_{10}$ and $a\text{-LaRu}_2\text{Al}_{10}$ alloys at $H = 10$ kOe. We calculated χ using the composition ratio of $\text{Ce}_{10}\text{Ru}_{15}\text{Al}_{75}$ obtained by SEM-EDS as 1 mol. χ value for $a\text{-CeRu}_2\text{Al}_{10}$ alloy increases monotonically with decreasing temperature, and a magnetic transition is not observed in the measurement temperature region. However, χ value for $a\text{-LaRu}_2\text{Al}_{10}$ alloy is almost independent of the temperature and is very small value ($< 10^{-6}$ emu/mol). $1/(\chi - \chi_0)$ value for $a\text{-CeRu}_2\text{Al}_{10}$ alloy exhibits a linear behavior in the high-temperature region of $T > 30$ K following the Curie-Weiss law,

$$\frac{1}{(\chi - \chi_0)} = \frac{1}{C} (T - \theta), \quad (1)$$

where χ_0 is a constant for the independence of temperature, C is the Curie constant, and θ is the Weiss temperature. χ_0 value for $a\text{-CeRu}_2\text{Al}_{10}$ alloy obtained from (1) was $\sim 1.0 \times 10^{-6}$ emu/mol. The value of θ is -20 K. C is obtained as

$$C = \frac{N_{\text{Ce}} \mu_{\text{eff}}^2}{3k_B}, \quad (2)$$

where N_{Ce} is the number of Ce atoms, μ_{eff} is the paramagnetic effective magnetic moment, and k_B is the Boltzmann constant. Here, μ_{eff} is calculated from Ce-concentration (10 at%) of composition ratio $\text{Ce}_{10}\text{Ru}_{15}\text{Al}_{75}$ obtained from SEM-EDS.

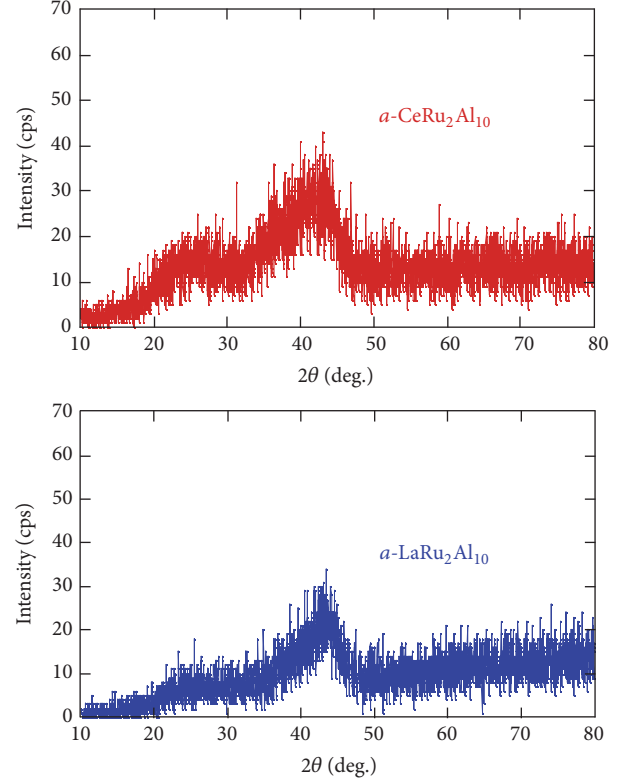


FIGURE 1: X-ray diffraction pattern of $a\text{-CeRu}_2\text{Al}_{10}$ and $a\text{-LaRu}_2\text{Al}_{10}$ alloys.

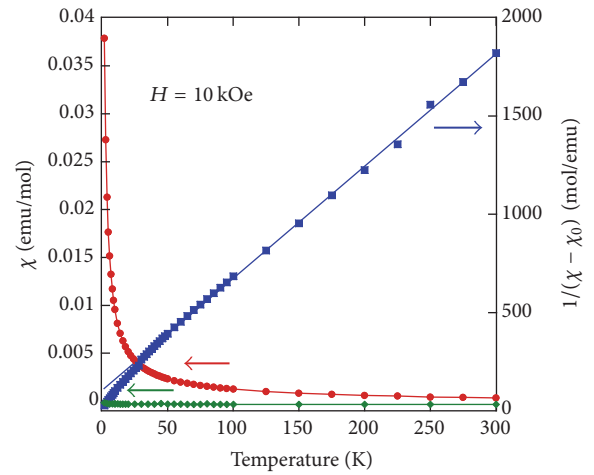


FIGURE 2: Temperature dependence of the magnetic susceptibility χ and inverse susceptibility $1/(\chi - \chi_0)$ for $a\text{-CeRu}_2\text{Al}_{10}$ alloy. The green points and line are the temperature dependence of the magnetic susceptibility χ for $a\text{-LaRu}_2\text{Al}_{10}$ alloy.

The estimated μ_{eff} obtained using (2) is $1.19 \mu_B/\text{Ce-atom}$. Since the effective magnetic moment expected for trivalent Ce ($J = 5/2$) is $2.54 \mu_B$, the obtained value of $1.19 \mu_B/\text{Ce}$ is about half this value. Therefore, it is expected that about half of Ce of $a\text{-CeRu}_2\text{Al}_{10}$ alloy exists as nonmagnetic tetravalent Ce and the remaining half exists as magnetic trivalent Ce in the alloy.

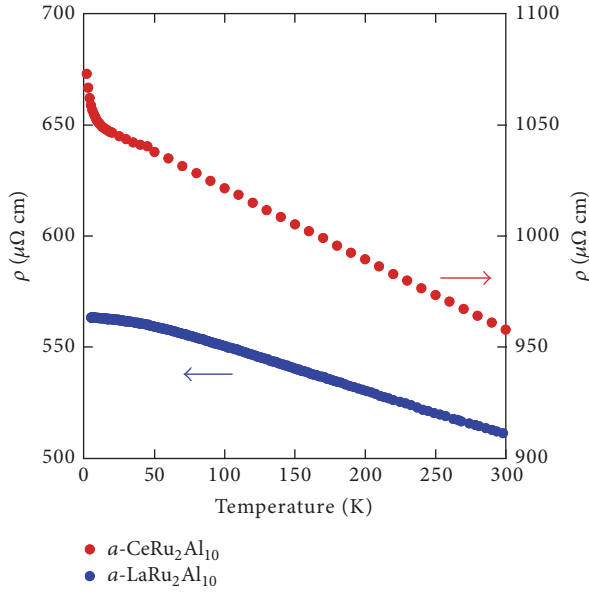


FIGURE 3: Temperature dependence of the resistivity ρ for a -CeRu₂Al₁₀ and a -LaRu₂Al₁₀ alloys.

Figure 3 shows the temperature dependence of the resistivity ρ for a -CeRu₂Al₁₀ and a -LaRu₂Al₁₀ alloys. The value of ρ at 300 K for a -CeRu₂Al₁₀ alloy is about 2 times greater than that of a -LaRu₂Al₁₀ alloy. ρ value for both alloys exhibits small temperature dependence less than 10% in the whole temperature region, and it increases with decreasing temperature. Such temperature dependence is one of the characteristics of disordered alloys. In contrast to ρ value of a -LaRu₂Al₁₀ alloy, which increases monotonically with decreasing temperature, ρ value of a -CeRu₂Al₁₀ alloy exhibits an increase in the low-temperature region. Generally, in the case of disordered alloys, the resistivity is larger than that for the crystalline counterparts owing to the random arrangement of atoms. However, the absolute value of the low-temperature ρ in a -CeRu₂Al₁₀ alloy is on about the same order as that for the polycrystalline CeRu₂Al₁₀. In addition, although ρ value of a -CeRu₂Al₁₀ alloy increases with decreasing temperature, a small temperature dependence similar to that for the typical disordered alloy was observed. Therefore, it is considered that the semiconductor band for crystalline CeRu₂Al₁₀ exhibited a metallic behavior as a result of amorphization.

Figure 4 shows the low-temperature specific heat C_p over T versus T^2 plots for a -CeRu₂Al₁₀ and a -LaRu₂Al₁₀ alloys. The inset shows the temperature dependence of C_p for the alloys. C_p values of both alloys are almost in agreement in the whole temperature region, as shown in the inset. The low-temperature specific heat of a usual metal can be expressed by

$$C_p = \gamma T + \beta T^3, \quad (3)$$

where γT is the electronic specific heat term and βT^3 is the phonon specific heat term. This relation can also be applied to amorphous alloys [11]. As shown in Figure 4, C_p/T of a -LaRu₂Al₁₀ alloy follows a linear relation as a function

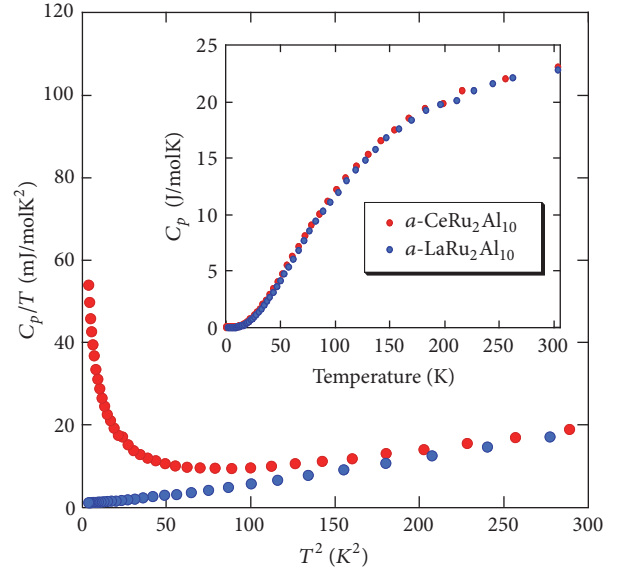


FIGURE 4: C_p/T versus T^2 plot for a -CeRu₂Al₁₀ and a -LaRu₂Al₁₀ alloys. The inset shows the temperature dependence of the specific heat C_p for a -CeRu₂Al₁₀ and a -LaRu₂Al₁₀ alloys.

of T^2 below 50 K². However, C_p/T of a -CeRu₂Al₁₀ alloy increases rapidly with decreasing T^2 below 100 K². The value of C_p/T at the lowest temperature is ~ 54 mJ/molK². The temperature region where the enhancement in C_p/T occurs agrees with the temperature region of the upturn for ρ . Such enhancement in C_p/T and ρ at low-temperature does not occur for a -LaRu₂Al₁₀. Therefore, it is considered to be due to the magnetic contribution of the 4f-electron of Ce. In order to clarify the magnetic contribution of the 4f-electron of Ce for a -CeRu₂Al₁₀, we measured the magnetoresistance of a -CeRu₂Al₁₀.

Figure 5 shows the magnetic field dependence of the transverse magnetoresistance $\Delta\rho/\rho$ at 2 K, 10 K, and 60 K for a -CeRu₂Al₁₀ alloy. We calculated $\Delta\rho/\rho$ as follows:

$$\frac{\Delta\rho}{\rho} = \frac{\rho(H) - \rho(0)}{\rho(0)}, \quad (4)$$

where $\rho(H)$ is the resistivity in the magnetic field and $\rho(0)$ is the resistivity at zero field. $\Delta\rho/\rho$ increases negative with increasing magnetic field strength. $\Delta\rho/\rho$ at 2 K exhibits a large negative enhancement compared with that at other temperatures. Figure 6 shows the temperature dependence of $\Delta\rho/\rho$ at 20 kOe, 40 kOe, and 60 kOe for a -CeRu₂Al₁₀ alloy. The $\Delta\rho/\rho$ in the high-temperature region ($T > 10$ K) is almost independent of the temperature for all the magnetic fields. However, $\Delta\rho/\rho$ in the low-temperature region increases rapidly with decreasing temperature at $T < 10$ K for all the fields. Thus, the enhancement of ρ at low temperatures is suppressed by the applied magnetic field.

The present a -CeRu₂Al₁₀ alloy seems reasonable that the magnetic order of the crystalline counterpart have disappeared because they lose a long-range crystallographic order by amorphization. However, the local structure of

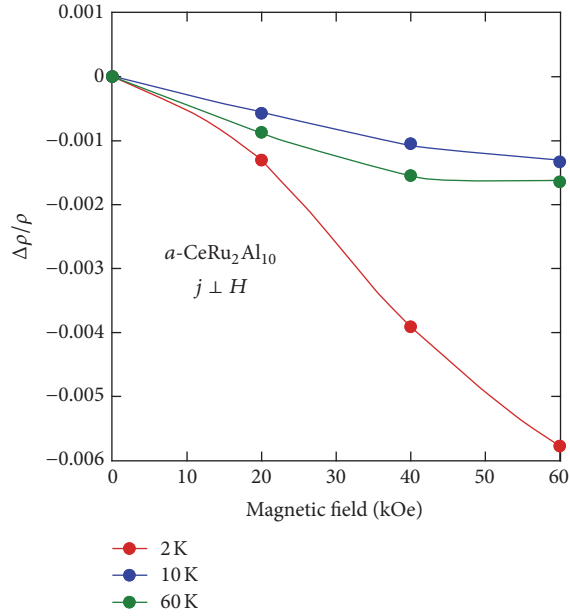


FIGURE 5: Field dependence of the magnetoresistance $\Delta\rho/\rho$ at various temperatures for $a\text{-CeRu}_2\text{Al}_{10}$ alloy.

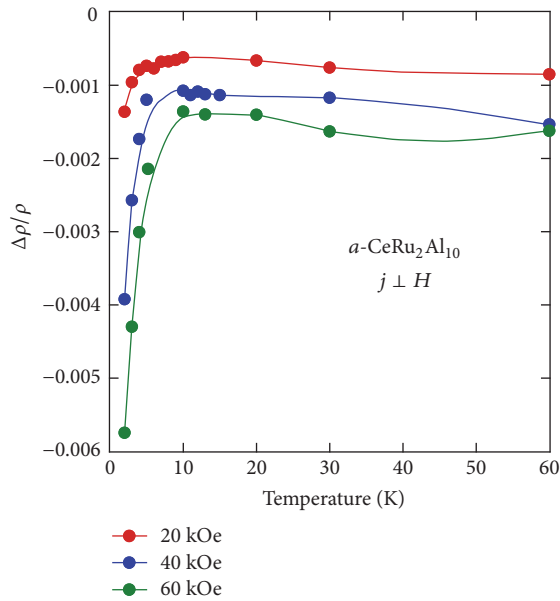


FIGURE 6: Temperature dependence of the magnetoresistance $\Delta\rho/\rho$ under various magnetic fields for $a\text{-CeRu}_2\text{Al}_{10}$ alloy.

the amorphous alloy is generally considered to be similar to that of the crystalline counterpart. Even in $a\text{-CeRu}_2\text{Al}_{10}$ alloy, a local structure similar to that of the crystalline $\text{CeRu}_2\text{Al}_{10}$ is expected to be realized. In addition, based on μ_{eff} , we have shown that half of Ce in $a\text{-CeRu}_2\text{Al}_{10}$ alloy is in the trivalent state. Ce^{3+} ions in $a\text{-CeRu}_2\text{Al}_{10}$ alloy are distributed randomly in the amorphous sample. In this case, it is considered that the random distribution of Ce^{3+} in the disordered structure realizes the impurity Kondo effect at low temperatures. Thus, the enhancement of

ρ and C_p/T is observed in the low-temperature region for $a\text{-CeRu}_2\text{Al}_{10}$ alloy. Since the Kondo scattering is suppressed by the applying a magnetic field, $\Delta\rho/\rho$ increases negatively. Therefore, the negative increase of $\Delta\rho/\rho$ for $a\text{-CeRu}_2\text{Al}_{10}$ alloy at low temperatures and high magnetic fields indicates that the scattering center of the resistivity is due to Kondo scattering.

4. Conclusion

We prepared $a\text{-CeRu}_2\text{Al}_{10}$ and $a\text{-LaRu}_2\text{Al}_{10}$ alloys and measured their susceptibility, resistivity, and specific heat. χ value for $a\text{-CeRu}_2\text{Al}_{10}$ alloy follows the Curie-Weiss law in the high-temperature region of $T > 30$ K with no magnetic transition. The effective paramagnetic moment and the Weiss temperature are $1.19 \mu_B/\text{Ce-atom}$ and -20 K, respectively. ρ value for $a\text{-CeRu}_2\text{Al}_{10}$ and $a\text{-LaRu}_2\text{Al}_{10}$ alloys shows a small temperature dependence. However, an enhancement of ρ was observed in the low-temperature region for $a\text{-CeRu}_2\text{Al}_{10}$ alloy. C_p/T value of $a\text{-CeRu}_2\text{Al}_{10}$ alloy increases rapidly with decreasing T^2 below 100 K^2 . The magnetoresistance $\Delta\rho/\rho$ for $a\text{-CeRu}_2\text{Al}_{10}$ alloy increases rapidly with decreasing temperature for $T < 10$ K. Therefore, it is considered that $a\text{-CeRu}_2\text{Al}_{10}$ alloy is formed the impurity Kondo state in the low-temperature region.

Conflicts of Interest

The authors declare that there are no conflicts of interest regarding the publication of this paper.

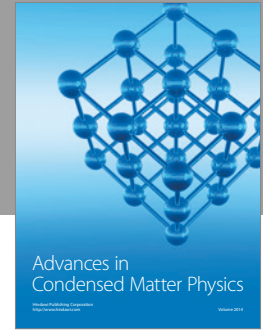
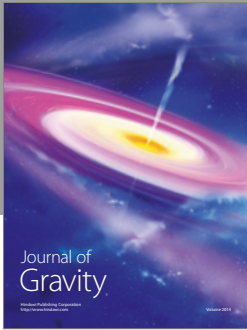
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