

Anomalous Nernst effect of epitaxial Fe₄N films grown on SrTiO₃(001) substrates
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Recently, the development of thermoelectric power generation (TEG) devices utilizing the anomalous Nernst effect (ANE) is attracting increasing interest. However, a current critical technical issue for practical application to ANE-based TEG is a much small value of the thermoelectric power of ANE (S_{ANE}) compared with that of the conventional Seebeck effect (SE)-based TEG. We focus on Fe₄N as a new material with a large S_{ANE} . Fe₄N is composed of earth abundant elements and relatively large S_{ANE} of 2.2 $\mu\text{V/K}$ has been experimentally reported in a Fe₄N film on a MgO(001) substrate [1]. In addition, large transverse Peltier coefficient (α_{xy}) of 2.4 A/(m·K) is theoretically predicted [2]. In this study, ANE of epitaxial Fe₄N films grown on SrTiO₃(STO)(001) substrates were measured in order to experimentally evaluate its α_{xy} .

The epitaxial Fe₄N film with a thickness of 27 nm was grown at 400 °C on STO(001) substrates by molecular beam epitaxy (MBE). Subsequently, a 2-nm-thick Al capping layer was formed on the Fe₄N layer in a sputtering chamber connected to the MBE chamber. The sample was microfabricated into a Hall bar shape using photolithography and ion milling for the measurements of ANE, SE, and the anomalous Hall effect (AHE). External magnetic fields were applied in the direction perpendicular to the film surface and temperature gradient (∇T) was applied to the in-plane Fe₄N[110] direction.

The S_{ANE} of 2.5 $\mu\text{V/K}$ was obtained and this value is larger than that of the Fe₄N film on the MgO(001) substrate [1]. The lattice mismatches of MgO(001)/Fe₄N and STO(001)/Fe₄N are -9.9 and -2.8%, respectively, and the Fe₄N films on the STO(001) substrates have better crystal orientation [3]. That is considered to be the reason why a larger S_{ANE} was obtained in STO(001)/Fe₄N sample. By using the results of ANE, SE, and AHE measurements, α_{xy} of the sample was estimated to be 1.5 A/(m·K). This value is smaller than the theoretically predicted 2.4 A/(m·K) [2]. The difference between the experimental and the theoretically predicted values might be due to the imperfect chemical ordering of N atoms in the Fe₄N film. The anisotropy of S_{ANE} in Fe₄N films depending on the crystallographic ∇T direction will be presented.

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