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HS-04. Effect of Cr doping on Mössbauer studies of double perovskite Sr_2FeMoO_6 . H. Ryu¹, S. Kim¹, C. Kim¹ and B. Lee²1. Physics, Kookmin University, Seoul, South Korea; 2. Physics, Hankuk Univ. of Foreign Studies, Young In, Kyungki, South Korea

We investigated the crystallographic and magnetic properties of double perovskite $Sr_2Fe_{0.9}Cr_{0.1}MoO_6$ and compared these with those of Sr_2FeMoO_6 [1]. We prepared single phases of the polycrystalline Sr₂Fe_{0.9}Cr_{0.1}MoO₆ powder using the solid-state reaction method, and confirmed the chemical composition of Sr, Fe, Cr, and Mo of the sample as stoichiometric using Rutherford backscattering spectrometry analysis. We found the structure of the $Sr_2Fe_{0.9}Cr_{0.1}MoO_6$ to be tetragonal, with lattice constants $a_0=5.5789$ Å, c_0 =7.9129 Å. The magnetoresistance magnitude ($\Delta \rho/\rho_0$) was 14.9 % and 1.4 % at 77 K and 300 K, respectively, under the applied field with 1 T. The Curie temperature of the Cr doped sample ($T_{\rm C} = 415$ K) was slightly smaller than that of the pure sample ($T_{\rm C}$ = 425 K). Mössbauer spectra of the Sr₂Fe_{0.9}Cr_{0.1}MoO₆ have been taken at various temperatures ranging from 15 to 415 K. As the temperature increased towards $T_{\rm C}$, the Mössbauer spectra showed line broadening and 1, 6 and 3, 4 line width differences because of anisotropic hyperfine field fluctuation. The Mössbauer spectra indicated that an anisotropic field fluctuation of $+H(P_{+}=0.85)$ was greater than $-H(P_{-}=0.85)$ 0.15). We calculated the temperature dependence of anisotropy energy from its relaxation rate. The field fluctuation frequency factor and the anisotropy energy were calculated as $8.2 \Gamma/h$ and 143.52 erg/cm^3 , respectively, using the relatively accurate data for T = 260 K that is associated with the large line broadening. We interpreted of effect of $Cr(t^3_{2g})$ doping as a decrease in the anisotropy energy.

[1] Sung Baek Kim, Key Taeck Park, and Chul Sung Kim, J. Appl. Phys. 89, 7681 (2001).