

## Site occupancy and anisotropy distribution of Al substituted Ba-ferrite with high coercivity

Dong Hyeok Choi, Sung Yong An, Sang Won Lee, In-Bo Shim, and Chul Sung Kim\*

Department of Physics, Kookmin University, Seoul 136-702, Korea

\*Corresponding author: e-mail: cskim@phys.kookmin.ac.kr, Phone: +82 2 910 4752, Fax: +82 2 910 4728

M-type barium hexagonal ferrite has been intensively investigated as a material for permanent magnets, high-density recording media, and microwave device. Recently, most of the research has emphasized the modification of magnetic properties by the substitution of  $\text{Fe}^{3+}$ , such as  $\text{Co}^{2+}$ - $\text{Ti}^{4+}$ ,  $\text{Co}^{2+}$ - $\text{Sn}^{4+}$ ,  $\text{Cr}^{3+}$ ,  $\text{Mn}^{3+}$  and  $\text{Al}^{3+}$ . Especially, when the substituted  $\text{Al}^{3+}$ , changes of structural and magnetic properties is very enlarged. In this study, the site occupancy and anisotropy distribution of Al substituted  $\text{BaFe}_{12-x}\text{Al}_x\text{O}_{19}$  ( $0.0 \leq x \leq 4.0$ ) have been studied with Mössbauer spectroscopy, x-ray diffraction, and vibrating sample magnetometry. Nanocrystalline  $\text{BaFe}_{12-x}\text{Al}_x\text{O}_{19}$  ( $0.0 \leq x \leq 4.0$ ) powders were fabricated by the sol-gel method. The result of XRD measurement shows that the lattice  $a$  and  $c$  parameters are decreased with increasing  $x$  from  $a = 5.901 \text{ \AA}$  and  $c = 23.243 \text{ \AA}$  for  $x = 0.0$ , to  $a = 5.818 \text{ \AA}$  and  $c = 22.754 \text{ \AA}$  for  $x = 4.0$ . With increasing  $x$ , the saturation magnetization,  $M_s$  decreased linearly but the coercivity,  $H_c$  greatly increased up to  $x = 2.0$ , and then slightly decreased over  $x = 2.0$ . Mössbauer spectra obviously changes with Al doping with a decrease in intensity from the  $4f_1+2a$  and  $12k$  sites. As can be seen, the linewidths broaden, especially for the  $4f_1+2a$  sites, and the relative area for the  $12k$  site gets larger as  $x$  increases. It can be seen that  $\text{Al}^{3+}$  ions have a strong preference for the  $4f_1+2a$  sites. However, when  $x > 2.0$  the  $4f_1+2a$  sites are only slightly involved in the substitution.

### References

- [1] C. S. Kim, S. W. Lee, S. Y. An, and I. B. Shim, Phys. Stat. Sol. (a), **189**, 903 (2002).
- [2] S. Diaz, J. L. Sanchez, F. Leccabue, B. E. Watts, G. Bocelli, and G. Albanese, J. Phys. IV France **7**, C1-331 (1997).

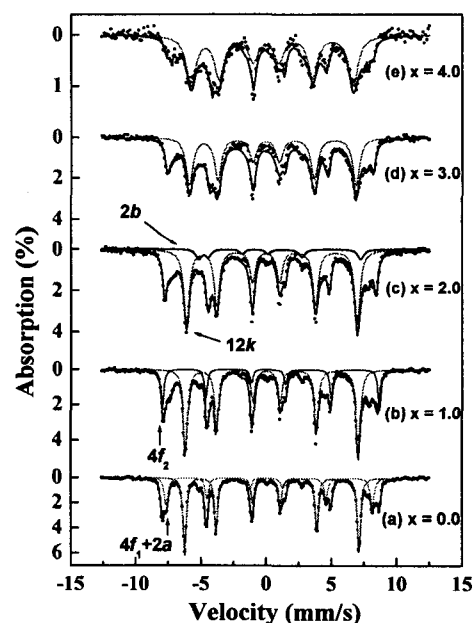


Fig. 1. Mössbauer spectra of  $\text{BaFe}_{12-x}\text{Al}_x\text{O}_{19}$  for  $x=0.0$  to  $x=4.0$  at room temperature. Solid circles are experimental data. Solid lines are fits according to models described in the text.