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## CARRIER DOPING DEPENDENCE OF $T_c$ IN DOUBLE PEROVSKITES $A_2FeMoO_6$ (A = Sr AND Ba)

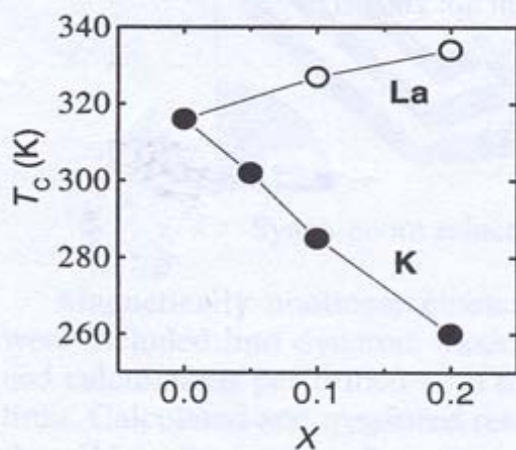
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We have studied effects of the carrier doping on the magnetic transition of double perovskites  $A_2FeMoO_6$  (A = Sr and Ba). The substitution of  $La^{3+}$  (or  $K^+$ ) in  $A_2FeMoO_6$  changes the average ionic radius ( $\langle r_A \rangle$ ) without distorting the crystal symmetry and the valence state of Fe/Mo ions via carrier doping.

Polycrystalline  $A_{2-x}B_xFeMoO_6$  (B = La and K;  $x = 0, 0.1$  and  $0.2$ ) samples were prepared by standard solid-state reaction. X-ray diffraction patterns reveal that samples are single-phase with cubic  $Fm3m$  symmetry for A = Ba and tetragonal  $I4/mmm$  symmetry for A = Sr. Lattice parameter of  $Ba_{2-x}La_xFeMoO_6$  decreases from  $8.076 \text{ \AA}$  for  $x = 0$  to  $8.045 \text{ \AA}$  for  $x = 0.2$  upon La substitution. This considerable reduction of lattice parameter arises from the substitution of the smaller  $La^{3+}$  ions into the  $Ba^{2+}$  ionic sites. However, the lattice parameter of  $Ba_{2-x}K_xFeMoO_6$  decreases slightly to  $8.074 \text{ \AA}$  for  $x = 0.2$ , which is due to the similar ionic radius between  $K^{1+}$  and  $Ba^{2+}$ . In  $Sr_{2-x}K_xFeMoO_6$  the lattice parameters increase monotonically with increasing  $x$  due to the substitution of bigger  $K^+$  ions.

A plot of the  $T_c$  vs carrier doping concentration for  $Ba_{2-x}B_xFeMoO_6$  (B = La and K) is shown in figure. Open symbols represent data for  $Ba_{2-x}La_xFeMoO_6$  and filled symbols correspond to data for  $Ba_{2-x}K_xFeMoO_6$ . The magnetic transition temperature  $T_c$  is defined as the temperature of the inflection point of  $M(T)$  curve. In  $Ba_{2-x}La_xFeMoO_6$  the  $T_c$  increases at a rate of  $1.8 \text{ K/\%}$  with increasing La doping. The  $T_c$  is  $316 \text{ K}$  for  $x = 0$  and  $334 \text{ K}$  for  $x = 0.2$ . However, the  $T_c$  of  $Ba_{2-x}K_xFeMoO_6$  decreases at a rate of  $5.6 \text{ K/\%}$  with K doping. The  $T_c$  for  $Ba_{1.8}K_{0.2}FeMoO_6$  ( $x = 0.2$ ) is  $260 \text{ K}$ .



Similar to  $Ba_{2-x}K_xFeMoO_6$ , the  $T_c$  for  $Sr_{2-x}K_xFeMoO_6$  decreases from  $379 \text{ K}$  for  $x = 0$  to  $368 \text{ K}$  for  $x = 0.1$ . The partial substitution of  $La^{3+}$  ( $K^+$ ) for  $Ba^{2+}$  in  $Ba_2FeMoO_6$  changes the  $\langle r_A \rangle$  within the cubic symmetry and the valence state of Fe/Mo ions via electron (hole) doping. The lattice parameters of  $Ba_{2-x}La_xFeMoO_6$  and  $Ba_{2-x}K_xFeMoO_6$  are found to vary linearly with the  $\langle r_A \rangle$ . As the  $\langle r_A \rangle$  decreases, the electronic bandwidth becomes broader, and thus the  $T_c$  increases [1]. In  $Ba_{2-x}K_xFeMoO_6$  system, the lattice parameter decreases with K doping; this should lead to the raise of  $T_c$ . However,  $T_c$  decreases slightly with increasing K concentration. This indicates that the carrier doping can promote some modification of  $T_c$  in these oxides.

- [1] RITTER, C. – IBARRA, M. R. – MORELLON, L. – BLASCO, J. – GARCIA, J. – De TERESA, J. M.: J. Phys.: Condens. Matter, **12**, (2000), 8295.

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