

Comparative Study of Impurity Scattering on Al- and C-doped MgB₂ Single Crystals

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We investigated the temperature dependence of the upper critical fields $H_{c2}(T)$ of Al- and C-substituted MgB₂ single crystals along the two main crystallographic directions. Despite similar suppression of T_c with the Al and C concentrations, the effects of the two substitutions on $H_{c2}(T)$ are found to be different. $H_{c2}(T)$ of the Al- and the C-doped MgB₂ can be well explained within the dirty-limit two-gap theory though $H_{c2}(0)$ of two substitutions show an opposite behavior, indicating very weak interband scattering. For the C-doping case, the ratio of in-plane electron diffusivities D_σ/D_π is much smaller than that for the Al-doping case indicating that C-doping enhances impurity scattering mainly in the σ bands as anticipated. The ratio of H_{c2} along the two directions, $\gamma_H(T)$ ($= H_{c2}^{ab} / H_{c2}^c$) systematically decreases with both Al and C dopings, however, $\gamma_H(T)$ for the C-doped shows a stronger temperature dependence than that for the Al-doped MgB₂ crystals.

Keywords : MgB₂ single crystals, Al doping, C doping, impurity scattering