# INFLUENCE OF CHEMICAL SUBSTITUTIONS ON ANISOTROPIC UPPER CRITICAL FIELD IN $MgB_2$ : IMPACT OF FERMI SURFACE CHANGES

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Specific band structure of MgB<sub>2</sub>, with two bands  $\pi$  and  $\sigma$  involved in superconductivity, leads to high critical temperature,  $T_c$ , of 39 K and temperature and field dependent anisotropy of superconducting parameters. Chemical substitutions lead to modification of band structure and therefore influence all superconducting parameters, especially  $T_c$ , the upper critical field,  $H_{c2}$ , and its anisotropy,  $\gamma_H$ . Magnetic investigations of  $Mg_{1-x}Al_xB_2$  crystals show the slight increase of  $H_{c2||c}$  for the samples with small x, significant reduction of  $\gamma_H$  at lower temperatures for Al substituted samples as compared to this of unsubstituted crystals. In  $Mg(B_{0.94}C_{0.06})_2$  single crystals  $H_{c2||c}(0) \cong 85$  kOe is more than twice as large as that one of  $\cong$  31 kOe in unsubstituted MgB<sub>2</sub>. Anisotropy of  $H_{c2}$  decreases to about 4 at low temperatures, the value considerably lower than that in MgB<sub>2</sub>, and its temperature dependence is much less pronounced. The corresponding  $H_{c2||ab}(0) \approx 330-350$  kOe is likely close to the maximum enhancement of  $H_{c2}$  due to chemical substitutions. The enhancement of  $H_{c2}$  can be explained as a disorder effect only if the main result of disorder is to make the  $\pi$  bands more dirty while not affecting the  $\sigma$  bands as much. However, in addition to disorder and weakened electron-phonon coupling, the impact of the Fermi level shifting into a region with lower  $\sigma$  Fermi surface velocities has to be taken into account in the analysis of  $H_{c2}$  data as well.

← 13.4 cm —

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 $9.7~\mathrm{cm}$