

Structure/Properties Relationships in doped MgB₂ Single Crystals

Götz Schuck^a, M. Wörle^b, N.D. Zhigadlo^a, K. Rogacki^a, J. Karpinski^a,

^a*Solid State Physics Laboratory ETH 8093 Zürich, Switzerland.*

^b*Laboratory of Inorganic Chemistry ETH 8093 Zürich Switzerland.* E-mail: schuck@solid.phys.ethz.ch

MgB₂ is a two electronic-bands, two energy gaps superconductor with a high T_c of 39K and unusual properties such as temperature and field dependent anisotropy.

Superconducting hexagonal single crystals of pure and Al, C, Mn and Fe doped MgB₂ phase have been grown at a pressure of 30 kbar using cubic anvil technique [1-2] to study the intrinsic properties of MgB₂. The superconducting transition of doped MgB₂ single crystals can be tuned in a wide temperature range between 10 and 39 K by adjustment of the nominal composition. Al [1], Mn and Fe are substituted on the Mg position of MgB₂ and C [2] on the B position. Introduction of disorder by substitution is partly observed.

In order to elucidate structure/properties relationships we have carried out x-ray single crystal measurements on doped MgB₂ single crystals (the composition varies between 1 and 15 % doping material). Additional temperature dependent measurements on pure MgB₂ and Mn doped MgB₂ has been carried out.

- [1] Karpinski J., Zhigadlo N. D., Schuck G., Kazakov S. M., Batlogg B., Rogacki K., Puzniak R., Jun J., Müller E., Wagli P., Gonnelli R., Daghero D., Ummarino G. A., Stepanov V. A., *Phys. Rev. B, submitted*, cond mat/0411449.
[2] Kazakov S. M., Puzniak R., Rogacki K., Mironov A. V., Zhigadlo N. D., Jun J., Soltmann Ch., Batlogg B., Karpinski J., *Phys. Rev. B*, 2005, **71**, 024533.

Keywords: high-T_c superconductivity, structure-physical properties relationships, x-ray crystal structure analysis