

Superconductivity of β -Type Salts under Uniaxial Compression

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Among organic superconductors, β -type salts attract attention for their high transition temperature (T_C). Due to weak dimerization of molecules compared to the κ -type salt, study on the β -type salts is intriguing in order to clarify whether the dimerization is a prerequisite for the high- T_C in organic superconductors. Uniaxial compression is a powerful technique to get insight into the mechanism of superconductivity since it controls the transfer integrals directly. Using epoxy encapsulation method, we have measured the T_C for β -(BDA-TTP)₂X [X=SbF₆, AsF₆] under in-plane uniaxial compression parallel and perpendicular to the molecular stack [1]. For X=SbF₆, T_C increases and takes a maximum at 3 kbar under the parallel compression. On the other hand, T_C decreases monotonically under the perpendicular compression. This behavior is understood as a competition between spin fluctuation and spin frustration effects on the dimerized triangular-lattice Hubbard model like κ -type salts. However, for the X=AsF₆, a T_C minimum is observed as a function of uniaxial pressure, which contradicts to the dimerized model [1].

In this presentation, we report the uniaxial compression effect on β -(BEDT-TTF)₂I₃. By embedding the crystal into epoxy resin, 8 K superconducting state is stabilized. By applying uniaxial compression, further T_C increase taking a maximum at 3-4 kbar is observed for both directions parallel and perpendicular to the molecular stack. In order to understand the different behavior against uniaxial compression, we calculate the T_C within the fluctuation-exchange approximation on dimerized single-band or non-dimerized two-band Hubbard models, by assuming displacements of molecules under uniaxial compression for each salt. For β -(BDA-TTP)₂SbF₆, both models are consistent with the experiment, but for β -(BDA-TTP)₂AsF₆, the non-dimerized model explains the T_C minimum. On the other hand, β -(BEDT-TTF)₂I₃ is rather understood by the dimerized model. These results are understood in terms of the difference in the strength of the dimerization among the three salts, that is, β -(BDA-TTP)₂AsF₆ < β -(BDA-TTP)₂SbF₆ < β -(BEDT-TTF)₂I₃.

[1] H. Ito *et al.*, Phys. Rev. B 78 (2008) 172506.