

Synthesis, Structure and Physical Properties of the REIr₃ and RE₂Ir₇ Binary Iridates

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The family of compounds characterized by the general formula RE_{2m+n}T_{4m+5n}, where RE is a rare earth element and T is a transition metal, is of significance due to its structural complexity and wide range of observed physical properties. The m and n parameters determine the number of MgCu₂-type and CaCu₅-type structural blocks, respectively, thereby providing a systematic framework for understanding structural variations and their influence on the properties. The REIr₃ system (m = 1 and n = 1) is known to form in different crystal structure types, depending on the choice of the RE element and the post-synthesis annealing temperature.

In this lecture we will present synthesis, structural and physical studies of the rhombohedral (PuNi₃ type) CeIr₃, PrIr₃, NdIr₃ and ThIr₃. The key element in this family, which affects the physical properties, is the rare earth metal. While LaIr₃, CeIr₃ and ThIr₃ are superconductors with T_c = 3.5 K [1], 2.5 K [2] and 4.4 K [3], respectively. In contrast, the remaining two are classified as ferromagnets, with the Curie temperature T_{Curie} = 7.5 K (PrIr₃) [4] and 10.6 K (NdIr₃) [5]. Interestingly, at around 70 K, the both magnetic compounds reveal a structural phase transition. The rhombohedral R-3m crystal structure undergoes a transformation to a monoclinic C2/m structure upon cooling, and further cooling reveals temperature-independent behaviour of the unit cell volume.

We also report on the synthesis and study of the m = 1, n = 2 variant of this family, i.e. La₂Ir₇ and Ce₂Ir₇, which also form in the rhombohedral crystal structure (Gd₂Co₇-type). Superconductivity was revealed through electrical resistivity and heat capacity measurements, with transition temperatures of T_c = 1.4 K and 1.6 K for La₂Ir₇ and Ce₂Ir₇, respectively. Thermodynamic and transport measurements demonstrate that both compounds are type-II superconductors, exhibiting a fully opened superconducting gap and weak electron-phonon coupling. The experimental results will be analyzed alongside the theoretical calculations.

[1] N. Haldolaarachchige, et al., J. Phys.: Condens. Matter **29**, 475602 (2017).

[2] K. Górnicka, et al., Supercond. Sci. Technol. **32**, 025008 (2019).

[3] K. Górnicka, et al., Phys. Rev. B **100**, 214514 (2019).

[4] K. Górnicka, et al., Phys. Rev. B **110**, 064431 (2024).

[5] K. Górnicka, et al., Phys. Rev. B **99**, 14430 (2019).