

Contribution ID: 348 Type: Oral Presentation

Optical spectroscopic investigations of antiferromagnetic semiconducting BaMn2P2

Friday 11 July 2025 10:50 (20 minutes)

Materials with ThCr₂Si₂-type crystal structure (Space Group: I4/<i>mmm</i>) have sparked scientific interest for several decades now owing to their novel properties and exotic ground states. The ThCr₂Si₂-type compounds have physical properties that are highly tunable, making this family of compounds ideal for investigating the structure-property relationship. Unconventional high <i>T</i>_c superconductivity was observed in iron-based arsenides, Ba_{1-<i>x</i>}K_{<i>x</i>}F [1] and BaFe < sub > 2 - < i > x < / i > < / sub > Co < sub > < i > x < / i > < / sub > As < sub > 2 < / sub > [2]. The quest for higher < i > T < / i > < sub > c < / sub > Co < sub > (a > x < / i > < / sub > Co < sub > (a > x < / i > < / sub > Co < sub > (a > x < / i > < / sub > Co < sub > (a > x < / i > < / sub > Co < sub > (a > x < / i > < / sub > Co < sub > (a > x < / i > < / sub > Co < sub > (a > x < / i > < / sub > Co < sub > (a > x < / i > < / sub > Co < sub > (a > x < / i > < / sub > Co < sub > (a > x < / i > < / sub > Co < sub > (a > x < / i > < / sub > Co < sub > (a > x < / i > < / sub > Co < sub > (a > x < / i > < / sub > Co < sub > (a > x < / i > < / sub > (a > x < / i > < / sub > (a > x < / i > < / sub > (a > x < / i > < / sub > (a > x < / i > x < / i > < / sub > (a > x < / i > < / sub > (a > x < / i > < / sub > (a > x < / i > < / sub > (a > x < / i > < / sub > (a > x < / i > < / sub > (a > x < / i > < / sub > (a > x < / i > < < / sub > (a > x < / i > < < / sub > (a > x < / i > < < / sub > (a > x < / i > < < / sub > (a > x < / i > < < / sub > (a > x < / i > x < / i > < < / sub > (a > x < / i > < < / sub > (a > x < / i > x < / i > < < / sub > (a > x < / i > < < > x < / i > < < > x < / i > < < > x < / i > < < > x < / i > < < > x < / i > < < > x < / i > < < > x < / i > < < > x < / i > < < > x < / i > < < > x < / i > < < > x < / i > < < > x < / i > < < > x < / i > < < > x < / i > < < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > x < > xsuperconductivity led research into other Ba<i>T</i>₂<i>Pn</i>₂ compounds, where <i>T</i> = transition metal and <i>Pn</i> = P, As, Sb, Bi, revealing a variety of physical properties. BaMn₂P₂ is one such compound that was recently investigated by us [3]. Electrical resistivity and heat capacity measurements on single crystals of BaMn₂P₂ revealed an insulating ground state with a small band gap [3]. Anisotropic magnetic susceptibility measurements confirmed that BaMn₂P₂, like its As-, Sb- and Bi- counterparts, has collinear Néel type antiferromagnetism below <i>T</i>_N = 795(15) K, which is the highest value for the family of 122-pnictide compounds thus far. Moreover, the magnetic susceptibility increases above <i>T</i>_N, like in the As-, Sb-, and Bi-based compounds, suggesting that antiferromagnetic correlations persist above the magnetic ordering temperature. In this contribution, we probe the properties of this interesting compound using Raman spectroscopy, investigating its structure around <i>T</i>_N. We also attempt to probe the effect of tweaking the ground states by changing charge-carrier concentrations.

References

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Session Classification: Physics of Condensed Matter and Materials

Track Classification: Track A - Physics of Condensed Matter and Materials