## 第 534 回物性セミナー・

## 創発的物性物理研究拠点セミナー

## Electronic structure of EuTGe3 (T: transition metal) studied by x-ray spectroscopies

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\*This seminar talk will be given in English.

Quantum critical phase transitions have been one of the central issues in strongly correlated 4f-electron systems. At absolute zero temperature, tuning non-thermal control parameters, such as pressure and magnetic field, tilts the balance between the competing ground states associated with conflicting Kondo effects and Ruderman-Kittel-Kasuya-Yoshida interactions [1]. The competition between Kondo and RKKY interactions in Ce- and Yb-compounds are often discussed using the Doniach phase diagram [2], whereas Eu compounds are believed to have a first-order phase transition with a Eu valence transition and not to have a quantum critical point (QCP). Most of the reported Eu-compounds favor divalent electronic states ( $Eu^{2+}$ , J=7/2) with an antiferromagnetic ground state. By applying pressure or a small element substitution, the system undergoes a first-order phase transition by changing into a non-magnetic ground state with a trivalent (or valence fluctuating) state ( $Eu^{3+}$ , J=0). Very recently, an unconventional behavior has been discovered in antiferromagnetic EuRhSi<sub>3</sub> and  $Eu_2Ni_3Ge_5$  [3] and indicates a potential QCP.

EuTGe<sub>3</sub> (T= Co, Ni, Rh, Ir) and EuRhSi<sub>3</sub> are isostructural, possessing a non-centrosymmetric BaNiSn<sub>3</sub>-type crystal structure. The Eu ions expect to be in a divalent electronic state and antiferromagnetic ordering appears at ~10-15 K [4]. We have performed hard x-ray photoelectron spectroscopy [5] and partial fluorescence yield x-ray absorption spectroscopy on EuTGe<sub>3</sub>. Details of the changes in the electronic structure as a function of transition metal substitution and pressure will be discussed in the talk.

## References

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- [4] O. Bednarchuk, et al., J. Alloys Comp. 622, 432-439 (2015).
- [5] Y. Utsumi, et al., Phys. Rev. B 97, 115155 (2018).

5 研究科共同セミナーの認定科目です

担当:木村 昭夫(理学研究科)



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