## Electronic properties across the first-order phase transition in $Fe_{1.05}$ Te

**S.** Rößler<sup>*a*</sup>, Dona Cherian<sup>*b*</sup>, Harikrishnan S. Nair<sup>*b*</sup>, Handady L. Bhat<sup>*b*</sup>, Suja Elizabeth<sup>*b*</sup>, Frank Steglich<sup>*a*</sup>, and Steffen Wirth<sup>*a*</sup>

 $^a{\rm Max}$ Planck Institute for Chemical Physics of Solids, Nöthnizer Straße 40, 01187, Dresden, Germany  $^b{\rm Department}$  of Physics, Indian Institute of Science, Bangalore 560012, India

We present here resistivity, magnetization, specific heat, scanning tunneling microscopy, and spectroscopy (STM/S) studies on Fe<sub>1.05</sub>Te single crystals grown by a horizontal Bridgman method. In this compound, the superconductivity appears upon Se substitution and the physical properties are found to be extremely sensitive to non-stoichiometry and disorder [1]. In our crystals, a first-order phase transition is observed around 57 K in the resistivity, magnetization and the specific heat measurements. This transition is associated with a structural change from the tetragonal P4/nmm to the monoclinic P  $2_1$ /m space group. At this temperature, the compound becomes antiferromagnetic and the temperature dependence of the resistivity changes from log (-T) to T<sup>2</sup>. This observation suggests that the material becomes a Fermiliquid metal at low temperatures. Metallic behavior is also confirmed in the I-V characteristics of the STM measurements taken on an atomically resolved surface.

[1] S. Rößler et al., Phys. Rev. B, 82 144523(2010).