## Preliminary resistivity results on U<sub>2</sub>Ni<sub>2</sub>Sn single crystals

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 $U_2Ni_2Sn$  is a member of a large family of intermetallic compounds with the tetragonal  $Mo_2FeB_2$  crystal structure, which has been studied intensively over the past years [1-3]. It orders antiferromagnetically at 25 K with a propagation vector q=(0,0,1/2). Magnetization, magneto-acoustic, and neutron-diffraction experiments on a single crystal provide evidence that the uranium moments align parallel to the c-axis with the anisotropy energy of  $\approx 170$  K, indicating that  $U_2Ni_2Sn$  can be classified as an Ising system. This behavior is rather exceptional, majority of the isostructural uranium ternaries have U moments confined into the basal plane.

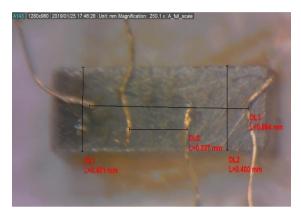
Last results [1] are actually at variance with previous studies on polycrystals, which indicated different magnetic structure, and which were incompatible with the 5f-5f two-ion anisotropy model dominant in most of U band systems. High-field magnetization studies [1] exhibit a weak linear response for fields along the basal plane up to the highest field applied (60 T), while the c-axis magnetization curve exhibits three metamagnetic transitions at approximately 30, 39 and 50 T.

Few single crystals of  $U_2Ni_2Sn$  were grown by the Czochralski method from a stoichiometric mixture of the pure elements (99.9% U, 99.99% Ni and 99.9999% Sn) in a triarc furnace with a water-cooled copper crucible under protective argon atmosphere. A tungsten rod was used as a seed. The pulling speed was 10 mm/h. We cut two differ-

ent single crystal pieces for resistivity measurements with orientation of [001] and [110], respectively. The small size of crystals was chosen to be compatible with high-pressure study envisaged.

The four-point technique, as shown in Fig. 1, was used to measure the resistivity. 25  $\mu m$  Au wire was used as leads for the measurement.

Although the primary goal is the high-pressure study (as  $U_2Ni_2Sn$  is an itinerant antiferromagnet, we expect fast suppression of magnetic order with pressure), already the ambient pressure data (see below) yield an interesting insight.



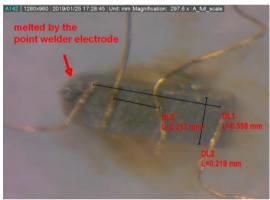


Figure 1. Single crystals of  $U_2N_{12}Sn$  mounted for the resistivity measurement in the [001] (up) and [110] (down) direction.

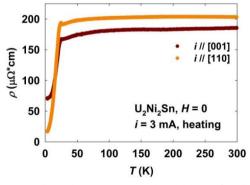


Figure 2a. The temperature dependence of resistivity of  $U_2Ni_2Sn$  single crystal with i // [110] and i // [110] (left).

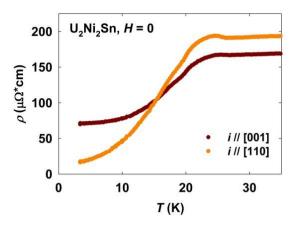


Figure 2b. The temperature dependence of resistivity of  $U_2Ni_2Sn$  single crystal with i // [110] and i // [110] (left) The low-temperature detail.

The first finding is that single crystal resistivity has better RRR value, reaching 10 for i // [110] and 3 for the [001] direction. The polycrystal value is about 3 [2]. None of the two directions exhibits the negative resistivity slope,  $d\Box/dT < 0$ , found for the polycrystal, and  $\Delta$  (T) shows merely a saturation. The magnetic propagation vector implies an influence of magnetic superzone gapping for [001], which can explain much higher residual resistivity in this direction. There is AF coupling between the U moments within the same basal-plane sheet, but it all happens within one unit cell, accommodating 4 U atoms.

Further differences are revealed by a numerical analysis. Fitting for i // [110] indicates the quadratic dependence, compatible with the high uniaxial anisotropy, not permitting a population of magnons with energy lower than the anisotropy gap. On the other hand, i // [001] exhibits an additional gap excitation term,  $\rho = \rho_0 + AT^2 + C * exp\left(-\frac{\Delta}{T}\right)$  with  $\Delta = 40\text{-}60$  K, i.e. much smaller than the anisotropy gap 170K. Such exchange gap has to be identified as a spin gap, suggesting an easy spin flipping along c.

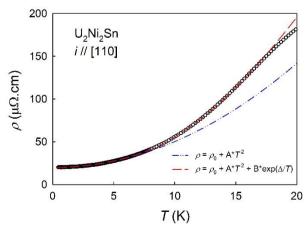
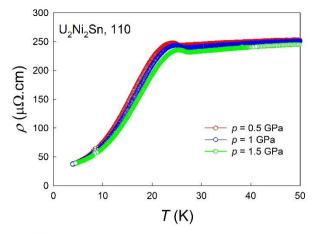


Figure 3. The temperature dependence of resistivity of  $U_2N_1^2S_1$  single crystal with i // [110] and i // [110] fitting.

The temperature dependence on pressure and field of resistivity of  $U_2Ni_2Sn$  single crystal with i // [110] and i // [110], are shown in figers 4 and 5. on boths transition temperatures goes down on higher fields or pressure.



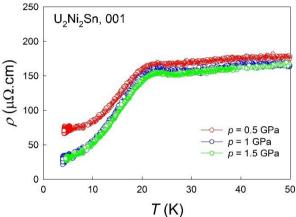


Figure 4. The temperature and pressure dependence of resistivity of  $U_2Ni_2Sn$  single crystal with i // [110] and i // [110]

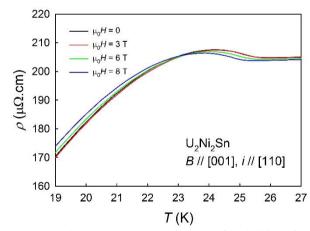


Figure 5. The temperature dependence of resistivity of  $U_2N_{12}Sn$  single crystal with i // [110] and B // [001].

## References

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