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## Phase Transition studies of $\text{MnSb}_{2-x}\text{Ta}_x\text{Sb}_2\text{O}_6$

Recently, the trigonal modification (sg. P321) of  $\text{MnSb}_2\text{O}_6$  has drawn significant attention as it could be an unusual type of multiferroic behaviour and weakly polar material [1]. The magnetic susceptibility of  $\text{MnSb}_2\text{O}_6$  shows a short range ordering below 200 K and long range ordering is observed below the Neel temperature ( $T_N = 12.5$  K) resulting in an incommensurately ordered three-dimensional Heisenberg antiferromagnet [2].  $\text{MnTa}_2\text{O}_6$  adopts the orthorhombic  $\text{MgNb}_2\text{O}_6$  structure type (sg. Pbcn) and it shows a monoclinic magnetic structure:  $P2_1/c$  symmetry at 4.2 K ( $T_N = 4.4$  K) [3]. In this solid solution, Ta could occupy the MnO interlayers and it will induce the decrease of the magnetic inter layer coupling. According to the Lab X-ray diffraction studies, a new tetragonal modification can be observed between  $x = 0.2$  and 1.8. The powder can be refined as mixtures of the trigonal  $\text{MnSb}_2\text{O}_6$  structure and a tetragonal tri-rutile modification, which is known for other  $\text{MSb}_2\text{O}_6$  compounds, for the refinement from  $x = 0.6$  to  $x = 1.2$ . The tri-rutile modification could be refined as the sole phase between  $x = 1.4$  to 1.6. The orthorhombic  $\text{MnTa}_2\text{O}_6$  modification could only be observed from  $x = 1.7 - 2.0$ . Interestingly, the tri-rutile modification has previously been described as a meta-stable modification for  $\text{MnTa}_2\text{O}_6$  [4].

### References

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**Primary author(s) :** KANG, Hyung-Been (The University of Auckland)

**Co-author(s) :** SUZUKI, Furitsu (The University of Auckland); Dr SOEHNEL, Tilo (The University of Auckland)

**Presenter(s) :** KANG, Hyung-Been (The University of Auckland)