

## Mn<sub>3</sub>MnB<sup>'</sup><sub>2</sub>O<sub>9</sub> high pressure A-site manganites with 1:2 B-site ordered triple perovskite structure

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ABO<sub>3</sub> oxides with perovskite related structures have largely been studied due to their chemical and physical versatility. Different types of cation order (1:1, 1:2 and 1:3) can be induced to enhance their physical properties. [1-3] The introduction of Mn into the A sites of the perovskite structure using high-pressure and high-temperature (HPHT) synthesis, provides a rich playground of potential spin, charge and orbital orderings. This throws attention to the growing family of functional A-site manganites, with non-ordered simple (Pv), 1:1 double (DPv, A or B) and doubly ordered (or double double DDPv, A and B) and 1:3 quadruple (QPv, A) perovskite structures. These compounds, with increasing chemical and structural complexity, show a wide range of promising physical properties as spintronics (MnVO<sub>3</sub> Pv), colossal magnetoresistance (Mn<sub>2</sub>FeReO<sub>6</sub> DPv), crystal field induced magnetic anisotropy (MnRMnSbO<sub>6</sub> DDPv) and magneto orbital coupling (MnMn<sub>3</sub>Mn<sub>4</sub>O<sub>12</sub> QPv). [4-7]

A new family of 1:2 B-site ordered A-site manganites Mn<sub>3</sub>MnB<sup>'</sup><sub>2</sub>O<sub>9</sub> (B' = Nb, Ta) with triple perovskite (TPv) structure (Fig. 1a) has recently been reported [8,9] to show a complex monoclinic distortion of the parent cubic Pv with Cc space group and  $a = \sqrt{2}\sqrt{3}a_c$ ,  $b = \sqrt{2}a_c$  and  $c = 2\sqrt{3}a_c$  cell parameters Their complex magnetic behaviour includes three subsequent transitions where Mn<sup>2+</sup> spins order into a collinear antiferromagnetic (AFM) structure, modulate in a complex spin density wave (SDW) with incommensurate [ $k_x$  0  $k_z$ ] propagation vector continuously evolving with temperature and lock at low temperatures with  $k_L = [\frac{1}{3}$  0 -  $\frac{1}{6}]$  (Fig.1b). A high temperature phase transition for B' = Ta, allows to recover the room pressure multiferroic Mn<sub>4</sub>B<sup>'</sup><sub>2</sub>O<sub>9</sub>. A 25% band gap reduction between both phases is due to their different structural connectivity.

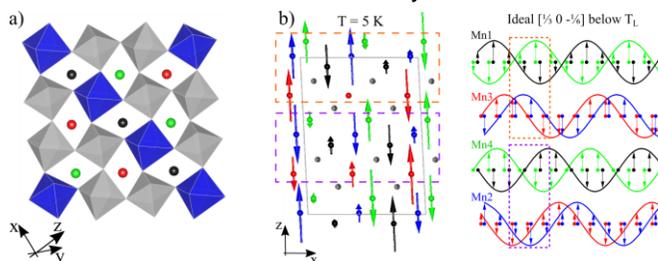


Fig.1. Crystal (a) and low temperature magnetic (b) structures of HP-Mn<sub>3</sub>MnNb<sub>2</sub>O<sub>9</sub> TPv.

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