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CORRECTION

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Correction: Defect chemistry and enhancement of thermoelectric performance in Ag-doped $Sn_{1+\delta-x}Ag_xTe$

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Correction for 'Defect chemistry and enhancement of thermoelectric performance in Ag-doped $Sn_{1+\delta-x}Ag_xTe$' by Min Ho Lee *et al.*, *J. Mater. Chem. A*, 2017, **5**, 2235–2242.

The authors regret an error in Fig. 4 of the original manuscript. The correct version of Fig. 4 is as below.

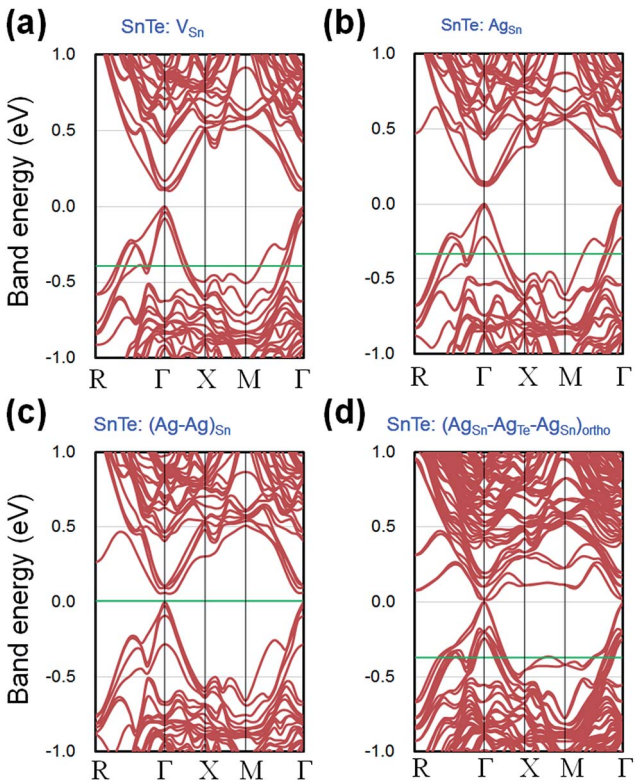

Fig. 4 Electronic band structures of SnTe supercells with defects: Sn-vacancy (V_{Sn}) (a), Ag-substitution at the Sn-site (Ag_{Sn}) (b), Ag–dimer complex defect $(Ag–Ag)_{Sn}$ (c), and Ag complex defect $(Ag_{Sn}–Ag_{Te}–Ag_{Sn})_{orthogonal}$ (d). Valence band maximum energy is set to zero and the Fermi level is denoted by a green line.

The Royal Society of Chemistry apologises for these errors and any consequent inconvenience to authors and readers.

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