Correction: Significant average ZT enhancement in Cu₃SbSe₄-based thermoelectric material via softening p–d hybridization

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The authors regret the following errors in the labelling of Cu and Ag orbitals in the published article:

In the Results and discussion section, the text “The calculated partial density of states (PDOS) for pristine Cu₃SbSe₄ (Fig. S3†) shows the valence band is dominated by strong hybridization between Cu 4d orbitals and Se 4p orbitals, whereas the conduction band edge is mainly formed of Sb 5s orbitals and p orbitals of neighboring Se. As Ag 5d orbitals are located at much lower energies (Fig. 4e)...” should instead read as follows: “The calculated partial density of states (PDOS) for pristine Cu₃SbSe₄ (Fig. S3†) shows the valence band is dominated by strong hybridization between Cu 3d orbitals and Se 4p orbitals, whereas the conduction band edge is mainly formed of Sb 5s orbitals and p orbitals of neighboring Se. As Ag 4d orbitals are located at much lower energies (Fig. 4e)...”

In Fig. 4e, the labels ‘Cu 4d’ and ‘Ag 5d’ should instead read ‘Cu 3d’ and ‘Ag 4d’, respectively. A corrected version of Fig. 4 is provided below.

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In the originally published version of Fig. S3, the label ‘Cu 4d’ should instead have read ‘Cu 3d’. This error has now been corrected in the ESI which is available online.

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