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CORRECTION

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Cite this: *Phys. Chem. Chem. Phys.*, 2021, **23**, 15386

Correction: A first-principles study of the stability, electronic structure, and optical properties of halide double perovskite $Rb_2Sn_{1-x}Te_xI_6$ for solar cell applications

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DOI: 10.1039/d1cp90137k

rsc.li/pccp

Correction for 'A first-principles study of the stability, electronic structure, and optical properties of halide double perovskite $Rb_2Sn_{1-x}Te_xI_6$ for solar cell applications' by Muhammad Faizan *et al.*, *Phys. Chem. Chem. Phys.*, 2021, **23**, 4646–4657, DOI: 10.1039/D0CP05827K.

The published version of this manuscript included errors in the text in the last paragraph of page 4653. The correct sentences are given below:

For $Rb_2Sn_{0.75}Te_{0.25}I_6$, the strongest absorption ($\sim 6.9 \times 10^5$ cm^{-1}) occurs at approximately 2.93 eV. Similarly, for $Rb_2Sn_{0.50}Te_{0.50}I_6$, the maximum absorption (5.82×10^5 cm^{-1}) occurs at ~ 3.1 eV.

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

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