

Erratum: Density Functional Theory and Experimental Determination of Band Gaps and Lattice Parameters in Kesterite $\text{Cu}_2\text{ZnSn}(\text{S}_x\text{Se}_{1-x})_4$

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The authors regret that one of their co-authors, Dr Oliver S. Hutter, was omitted from the original publication *J. Phys. Chem. Lett.* 2020, 11, 10463–10468 doi.org/10.1021/acs.jpcllett.0c03205 due to an oversight. Dr Hutter was responsible for performing a large part of the experiments on bulk crystalline materials. The authors would therefore like to take this opportunity to include Dr Hutter on the author list and to extend their apologies to him for the earlier oversight.

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