



Correction to Influence of Rb/Cs Cation-Exchange on Inorganic Sn Halide Perovskites: From Chemical Structure to Physical Properties

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Supporting Information

We are issuing a correction for eq 3. The correct version should appear as shown below:

$$\Delta S = -k_B[x \ln x + (1 - x) \ln(1 - x)]$$

Additionally, we are issuing a correction for Table S1 in the Supporting Information. The correct values for the lattice parameters (a_0 , b_0 , and c_0) of the cubic α -CsSnI₃ phase should appear as 6.28 Å.

ASSOCIATED CONTENT

Supporting Information

The Supporting Information is available free of charge on the ACS Publications website at DOI: [10.1021/acs.chemmater.7b01616](https://doi.org/10.1021/acs.chemmater.7b01616).

Detailed bulk and surface properties of the polymorphs (CsSnI₃ and RbSnI₃) and their solid solutions and their thermodynamic stabilities, surface band structures, and band level alignment under the influence of SOC effect and various xc functionals (PDF)