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Electronic Supplementary Information

High thermoelectric performance of In-doped Cu₂SnSe₃ prepared by fast

combustion synthesis

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Fig.S1 XRD pattern of the monoclinic Cu₂SnSe₃ prepared by combustion synthesis in air.



Fig. S2 Photographs of the $Cu_2Sn_{0.9}In_{0.1}Se_3$ sample prepared by high-pressure combustion synthesis in an Ar atmosphere with a pressure of 2 MPa.



Fig.S3 XRD patterns of the $Cu_2Sn_{1-x}In_xSe_3$ samples after spark plasma sintering: (a) an overview; (b) an enlarged view to clearly show the small peaks. No significant difference was observed in the XRD patterns of the samples just synthesized and after spark plasma sintering.



Fig.S4 Back-scattered electron images and EDS results of the $Cu_2Sn_{1-x}In_xSe_3$ samples: (a) and (b) prepared by high-pressure combustion synthesis, x=0.05, with SnSe as the secondary phase; (c) and (d) prepared by high-pressure combustion synthesis, x=0.20, with a Cu-rich secondary phase; (e) and (f) prepared by high-pressure combustion synthesis followed with spark plasma sintering, x=0.1, with much more grain boundaries compared with the just synthesized samples before spark plasma sintering.



Fig.S5 Lattice thermal conductivities of $Cu_2Sn_{1-x}In_xSe_3$ (x=0.05, 0.10) samples: (a) prepared by onestep HPCS, (b) prepared by HPCS-SPS.

The lattice thermal conductivities are calculated according to the Wiedemann-Franz Law ($\kappa = \kappa_e + \kappa_L = L_0 \sigma T + \kappa_L$), where κ_e is the thermal conductivity contributed by carriers, L_0 is the Lorentz number, σ is the electrical conductivity, and T is the absolute temperature. In our calculation, the Lorentz number of $L_0=2.0\times10^{-8}$ W Ω K⁻² is used according to the report by Shi et al. (X. Shi, L. Xi, J. Fan, W. Zhang and L. Chen, Chemistry of Materials, 2010, 22, 6029-6031.)