

# Decoupled electron-phonon transport in Ag<sub>2</sub>Se thermoelectric materials through constructing TiO<sub>2</sub>/MoS<sub>2</sub> co-decorated cell-membrane-mimic grain boundaries

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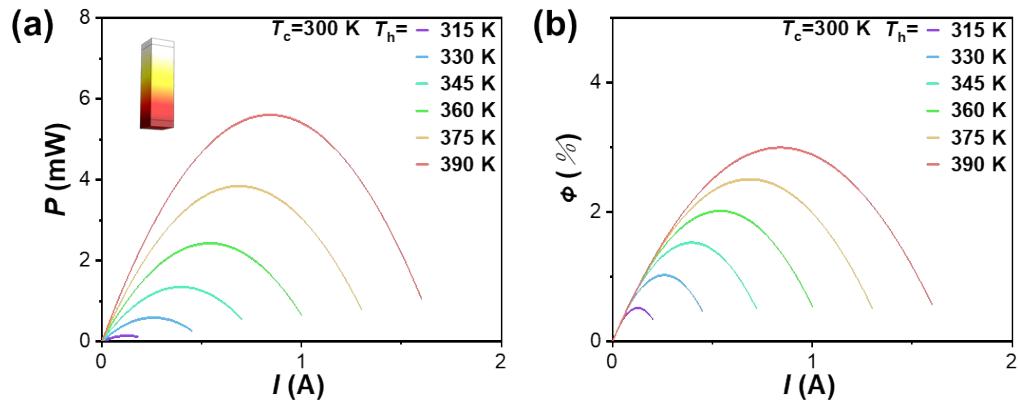
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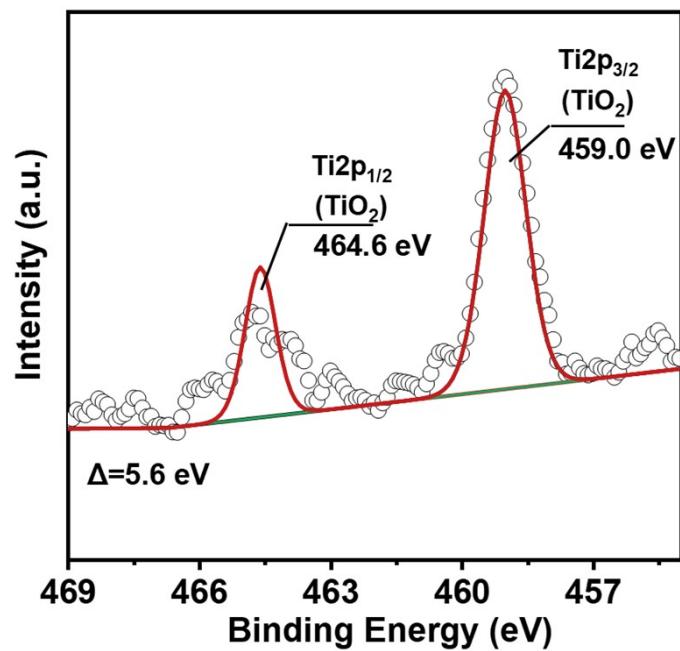
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**Fig. S1.**



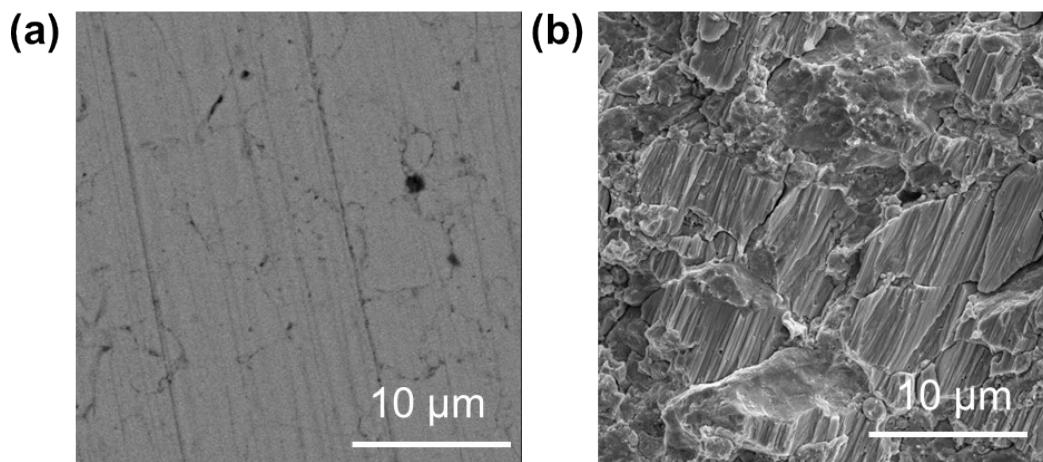
**Fig. S1.** (a-b) The results of the finite-element analysis of the  $\text{Ag}_2\text{Se}$  based module detail the output power ( $P$ ) and conversion efficiency ( $\Phi$ ) of the  $\text{Ag}_2\text{Se}$  based module as a function of current ( $I$ ) with a fixed cold side temperature ( $T_c$ ) and varied hot side temperature ( $T_h$ ).

**Fig. S2.**



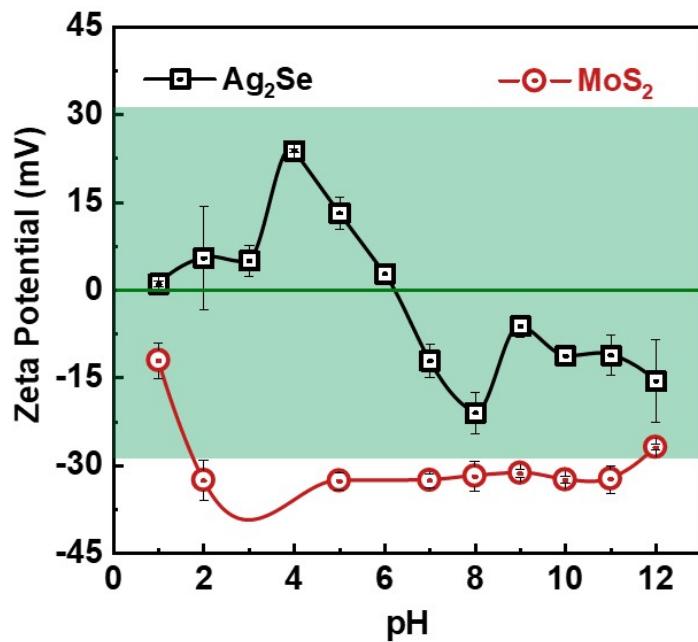
**Fig. S2.** High resolution XPS spectra of the AS-T-0.5 MoS<sub>2</sub> powder sample, inset is the Ti 2p spectrum.

**Fig. S3.**



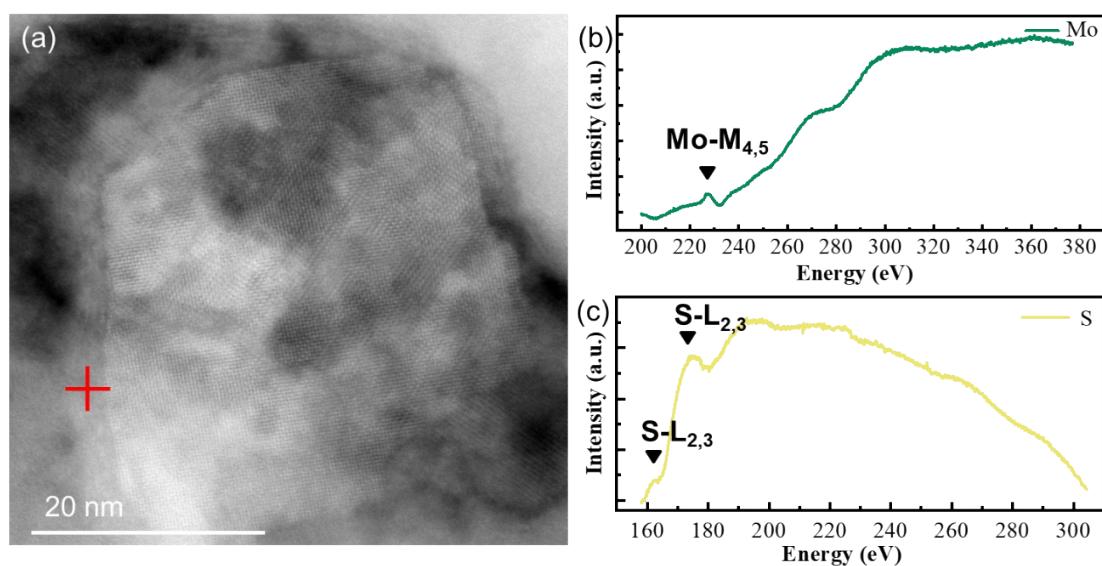
**Fig. S3.** (a) As-synthesized  $\text{Ag}_2\text{Se}$  to the as-decorated bulk samples. (b) The cross-section SEM of  $\text{Ag}_2\text{Se}$  bulk samples.

**Fig. S4.**



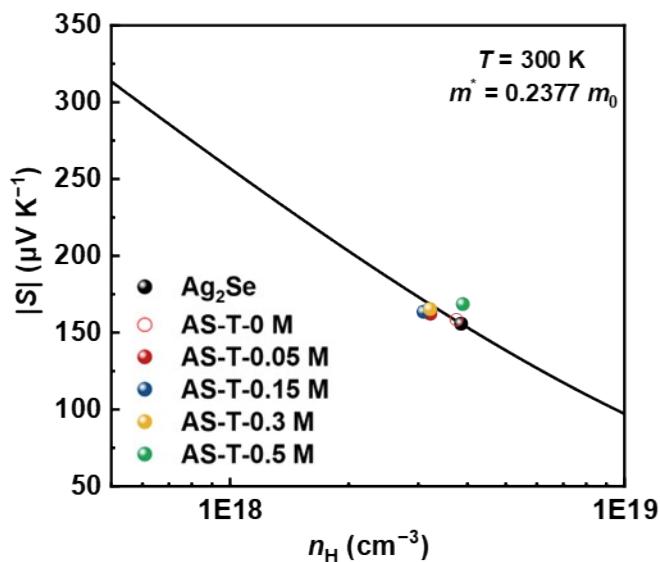
**Fig. S4.** Zeta potential of  $\text{MoS}_2$  and  $\text{Ag}_2\text{Se}$  dispersion at different pH.

**Fig. S5.**



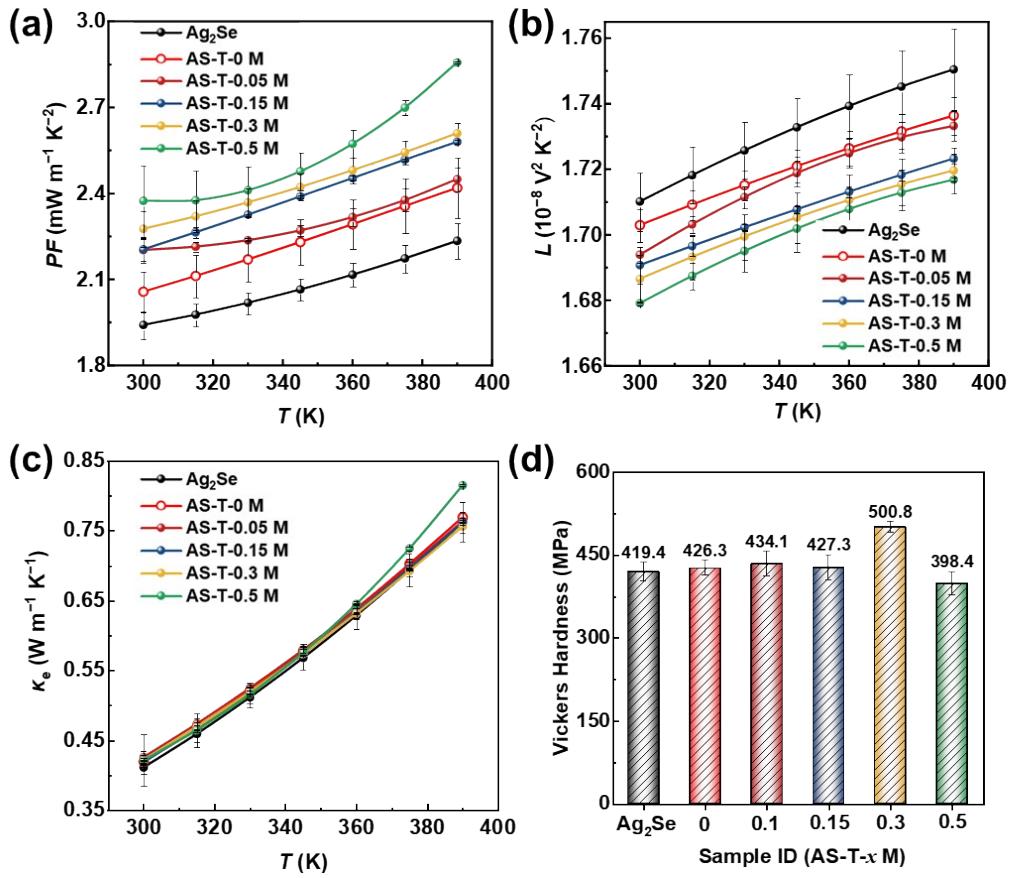
**Fig. S5.** (a) STEM-HAADF image of the TiO<sub>2</sub>/MoS<sub>2</sub> co-doped Ag<sub>2</sub>Se nanocomposites (b-c) EELS spectra of the element of Mo and S collected from a MoS<sub>2</sub> area (red cross).

**Fig. S6.**



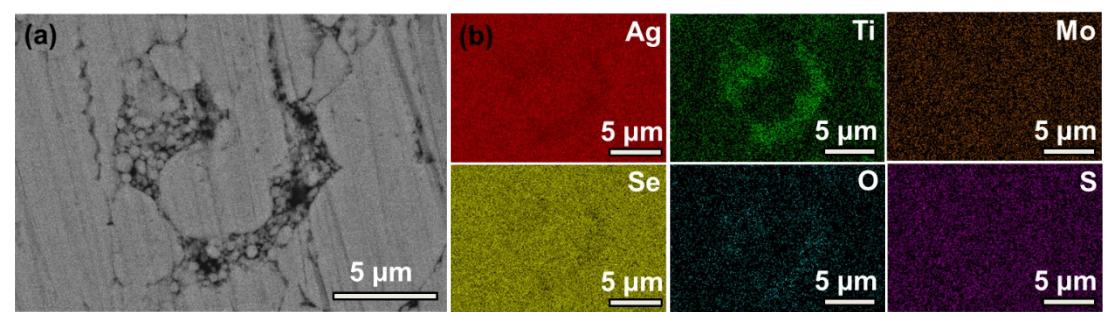
**Fig. S6.** The variation of  $S$  with  $n_H$  carrier concentration for of AS-T- $x$  M (where  $x = 0, 0.05, 0.15, 0.3$ , and  $0.5$  in wt%) compared to the Pisarenko line, which is caused by the change in the increased effective density of states mass ( $m^*$ ).

**Fig. S7.**



**Fig. S7.** Temperature-dependent (a) power factor ( $PF$ ), (b) lorentz number ( $L$ ), (c) electrons thermal conductivity ( $\kappa_e$ ) and (d) the Vickers hardness ( $H_v$ ) of  $\text{AS-T}-x \text{ M}$  (where  $x = 0, 0.05, 0.15, 0.3$ , and  $0.5$  in wt%).

**Fig. S8.**



**Fig. S8.** (a) Low magnification SEM image of AS-T-0.5M sample. (b) Corresponding elemental distribution of Se, Ag, O, Ti, S, and Mo.

**Table.S1**

PH	1	2	5	7	8	9	10	11	12
Zeta	-12.033	-	-	-	-	-	-	-	-26.9
Potential (mV)		32.566	32.666	32.566	31.833	31.333	32.466	32.333	
Error range	-3.0666	-	-	-	2.5333	-	-	-	0.5
		3.4333	1.5333	1.1333		0.7666	0.5666	2.3666	

**Table.S2**

Sample	Actual density	Relative density (%)
Ag <sub>2</sub> Se	8.137	98.15
AS-T-0.05 M	8.020	96.74
AS-T-0.15 M	8.026	96.82
AS-T-0.3 M	8.007	96.59
AS-T-0.5 M	8.110	97.83