## N-type Cu<sub>0.008</sub>Bi<sub>2</sub>Te<sub>2.7-x</sub>Se<sub>0.3</sub> 열전 합금에서 Te 공공 형성에 의한 열전 성능 연구

## AEML 홍석원 (신소재), 이민형(신소재), 안지우(신소재), 한유진(신소재), 전지훈(화공) 지도교수: 김상일



Advanced Energy Materials Lab.

에너지 나노소재 연구실



electrical conductivity  $\sigma$  [S/cm] Seebeck coefficient, S [ $\mu V/K$ ] Thermal Conductivity [ $W/m \cdot K$ ] Absolute Temperature, T [K]

 $zT = \frac{\sigma S^2}{\kappa} T$ 

 $P.F. = \sigma S^2$ 

Power Factor, P. F.  $[mW/mK^2]$ 





Fig. 5. X-ray diffraction patterns of (a)  $Cu_{0.008}Bi_2Te_{2.7-x}Se_{0.3}$  and (b) Calculated lattice parameters a and c of the  $Cu_{0.008}Bi_2Te_{2.7-x}Se_{0.3}$  with x=0, 0.005, 0.01 and 0.02.



Fig. 6. (a) The measured electrical conductivity  $\sigma$ , (b) Seebeck coefficient S and (c) power factor  $\sigma$ S<sup>2</sup> of the samples as a function of temperature for the Cu<sub>0.008</sub>Bi<sub>2</sub>Te<sub>2.7-x</sub>Se<sub>0.3</sub> with x=0, 0.005, 0.01, 0.02.

(d) (e) Estimated carrier concentrations and mobilities from the Hall measurement. (f) Pisalenko plot.



Fig. 7. (a) The electron ( $n_e$ ) and hole ( $n_h$ ) concentration, (b) weighted mobility(U) of CB and VB and (c) weighted mobility ratio (A=U<sub>CB</sub>/U<sub>VB</sub>) calculated from SPB at 300K.

## Table 1. Band parameters obtained using the two-band model

	Cu	$I_{0.008}Bi_2Te_{2.7-x}Se_{0.3}$			
Band parameters		x=0	x=0.005	x=0.01	x=0.02
Conduction band (CB)	$\operatorname{CB} E_{\operatorname{def}}(\operatorname{eV})$	18.4	18.0	17.7	17.5
	CB $m^*$ (in $m_0$ )	1.06	1.07	1.08	1.09
	$U_{\rm CB}~({\rm cm}^2/{\rm Vs})$	291	299	306	313
	Electron conc. $n_{\rm e} (10^{19} {\rm cm}^{-3})$	2.98	3.28	3.24	4.01
Valence band (VB)	$\operatorname{VB} E_{\operatorname{def}}(\operatorname{eV})$	29.5	29.0	29.1	27.1
	VB $m^*$ (in $m_0$ )	1	1	1	1
	$U_{\rm VB}~({\rm cm}^2/{\rm Vs})$	119	124	123	142
	Hole conc. $n_{\rm h} (10^{16} {\rm cm}^{-3})$	4.74	4.25	4.41	3.30
A (U <sub>CB</sub> /U <sub>VB</sub> )		2.44	2.42	2.50	2.21

Table 1. Band parameters obtained using the two-band model.

 $E_{def} =$  deformation potential

 $m^*$  = density-of-states effective mass (m<sub>0</sub> = electron mass)

 $\mu_0$  = nondegenerate mobility

U = weighted mobility

A = weighted mobility ratio



Fig. 8. (a) The total thermal conductivity  $\kappa_{tot}$ , (b) the electronic thermal conductivity  $\kappa_{elec}$ , (c) the bipolar thermal conductivity  $\kappa_{bp}$ , and (d) the lattice thermal conductivity  $\kappa_{latt}$  of the  $Cu_{0.008}Bi_2Te_{2.7-x}Se_{0.3}$  with x=0, 0.005, 0.01 and 0.02.

Fig. 5. (a) Temperature dependence of the dimensionless figure of merit zT (b) zT520K,  $zT_{average}$  and  $zT_{max}$  of the  $Cu_{0.008}Bi_2Te_{2.7-x}Se_{0.3}$  with x=0, 0.005, 0.01 and 0.02.

## **Thank You**