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Single and Double-Doping Effects on the Thermoelectric Properties of Two Zintl Compounds: $Eu_{11}Bi_{8.07(2)}Sn_{1.93}$ and $Eu_{10.74(2)}K_{0.26}Bi_{9.14(2)}Sn_{0.86}$

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Table S1 Detailed structural Information including atomic coordinates for the three structure models

model	$Eu_{11}Bi_{10}$	$Eu_{11}Bi_8Sn_2$	$Eu_{10}KBi_8Sn_2$
space group	*I4/mmm* (No.139)	*I4/mmm* (No.139)	*I4/mmm* (No.139)
unit cell dimensions (Å)	*a* =12.571 *c* =18.173	*a* =12.538 *c* =18.125	*a* =12.750 *c* =18.433
volume ($Å^3$)	2871.9	2849.3	2996.5

Atomic coordinates

atom	*x*	*y*	*z*
		$Eu_{11}Bi_{10}$	
Eu1	0	0.2515	0.3110
Eu2	0	0.3334	0.1015
Eu3	0.3259	0.3259	0
Eu4	0	0	0.1655
Bi1	0.2055	0.2055	0.1730
Bi2	0.1484	0.5	0
Bi3	0.1199	0.1199	0
Bi4	0	0	0.3775
Bi5	0	0.5	0.25
		$Eu_{11}Bi_8Sn_2$	
Eu1	0	0.2515	0.3110
Eu2	0	0.3334	0.1015
Eu3	0.3259	0.3259	0
Eu4	0	0	0.1655
Bi1	0.2055	0.2055	0.1730
Bi2	0.1484	0.5	0
Sn1	0.1199	0.1199	0
Bi4	0	0	0.3775
Bi5	0	0.5	0.25
		$Eu_{10}KBi_8Sn_2$	
Eu1	0	0.2515	0.3110
Eu2	0	0.3334	0.1015
Eu3	0.3259	0.3259	0
K4	0	0	0.1655
Bi1	0.2055	0.2055	0.1730
Bi2	0.1484	0.5	0
Sn1	0.1199	0.1199	0
Bi4	0	0	0.3775
Bi5	0	0.5	0.25

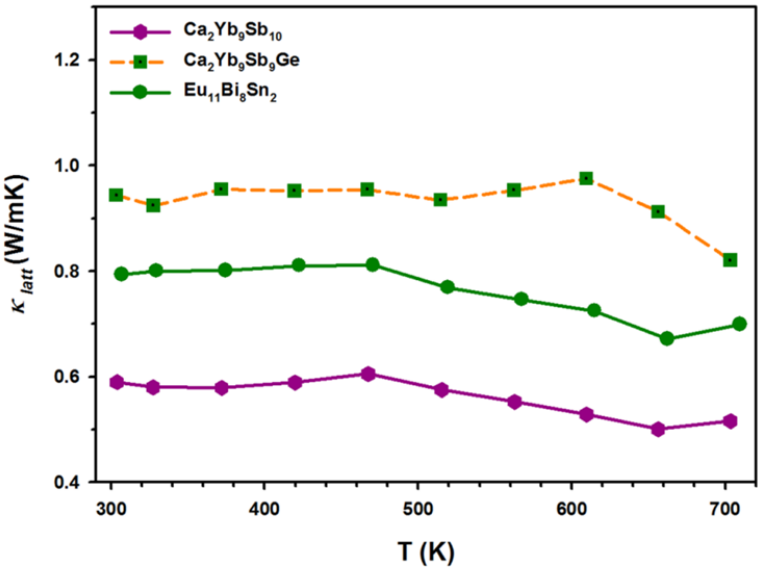

Fig. S1. Temperature-dependent lattice thermal conductivities κ_{latt} of the three compounds measured over the temperature range of 300-700 K.