

Supplemental Material for
“Functionality-directed Screening of Pb-free
Hybrid Organic-inorganic Perovskites with
Desired Intrinsic Photovoltaic Functionalities”

Dongwen Yang,^{†,||} Jian Lv,^{†,||} Xingang Zhao,[†] Qiaoling Xu,[†] Yuhao Fu,[†] Yiqiang
Zhan,[‡] Alex Zunger,^{*,¶} and Lijun Zhang^{*,†,§}

[†]Department of Materials Science and Engineering and Key Laboratory of Automobile
Materials of MOE, Jilin University, Changchun 130012, China

[‡]State Key Laboratory of ASIC and System, Department of Microelectronics, SIST, Fudan
University, Shanghai 200433, China

[¶]University of Colorado, 13067d..30672 11)0.13067(n)0.30621(d)349.928(K)R.14(3 2)1.2171(S)0.30





Figure 6 Band structures of the selected $AM^{IV}X_3^{VII}$ perovskites having indirect band gaps. The valence band maximum and conduction band minimum are marked by red circles. The actual band gaps E_g and direct band gaps E_g^d .

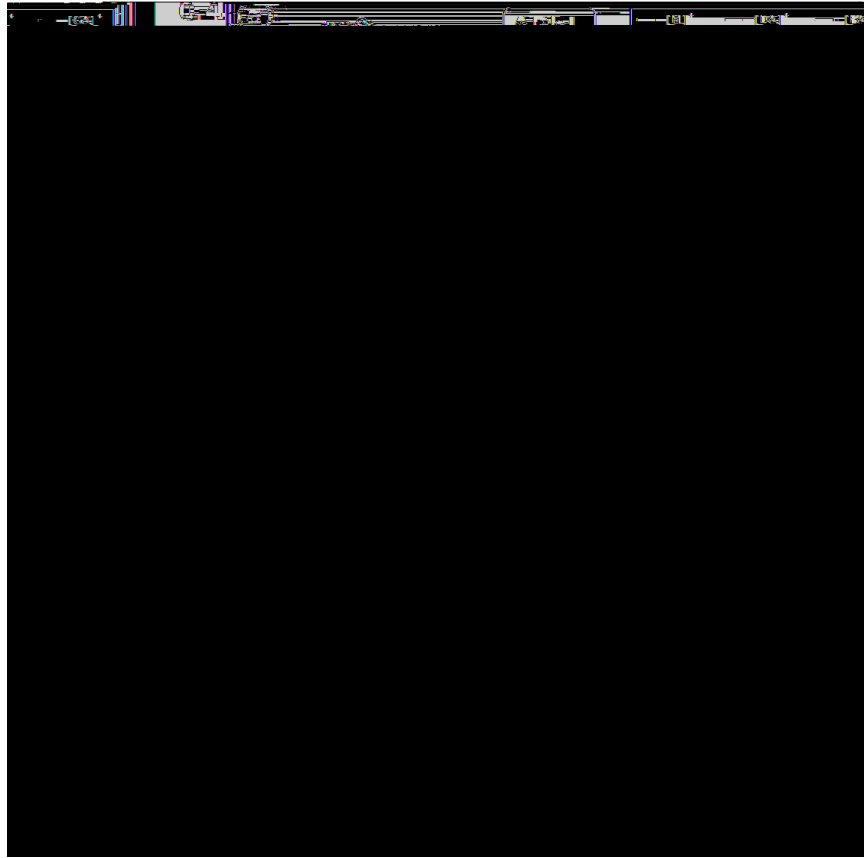


Figure 1. Crystal orbital overlap populations (COOP) of the $AM^{IV}X_3^{VII}$ perovskites. For comparison, the CBM of each material is set to energy zero.



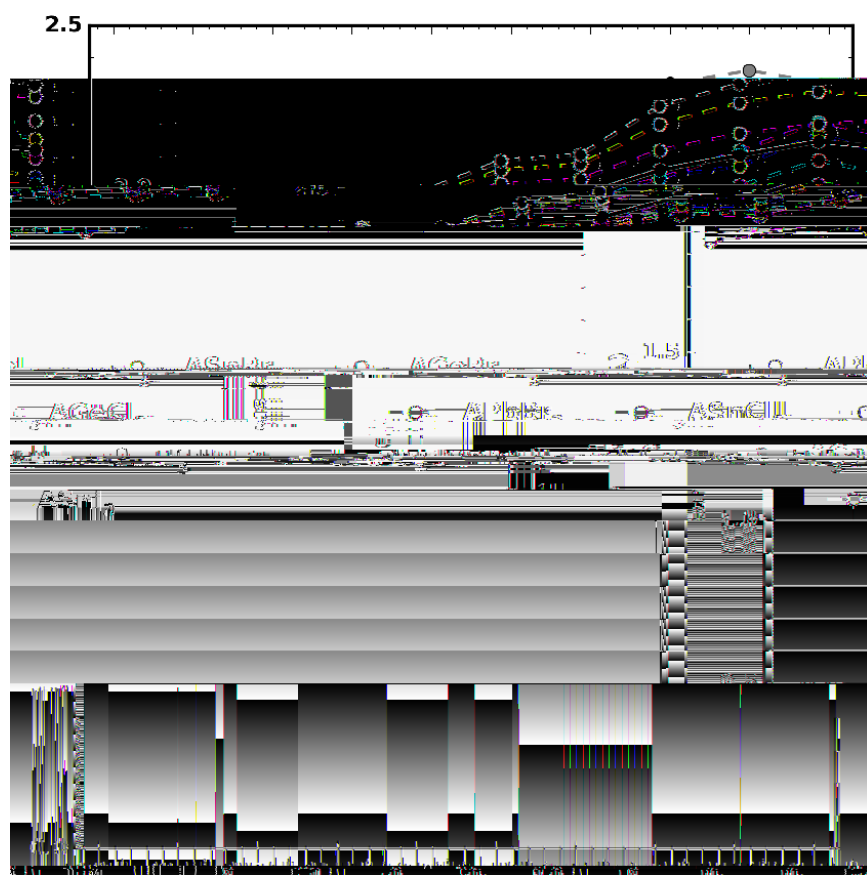


Figure 10 Evaluation of the steric sizes of organic cations within the idealized solid sphere model (see the Experimental section for details)

Table 1: Calculated decomposition enthalpies H of the candidate $AM^{IV}X_3^{VII}$

Table 1: Calculated direct band gaps E_g^d of the candidate $AM^{IV}X_3^{VII}$ perovskites. The green shading indicates the compounds passing the current DM $E_g^d < E_{DM}$, as we use the DM in Table 1. The light blue shading indicates the compounds passing only the current DM.

E_g^d (eV)	PbI ₃	PbBr ₃	PbC ₃	nI ₃	nBr ₃	nC ₃	GeI ₃	GeBr ₃	GeC ₃
[M ⁺]	/	/	/	/	/	/	/	/	/

Table 1: Calculated exciton binding energies E_B of the candidate $AM^{IV}X_3^{VII}$ perovskites. The green shading indicates the compounds passing the current DM $E_B < 1$ eV, as well as the DMs in Tables 1, 2, and 4. The light blue shading indicates the compounds passing on by the current DM.

E_B (eV)	PbI ₃	PbBr ₃	PbC ₃	nI ₃	nBr ₃	nC ₃	GeI ₃	GeBr ₃	GeC ₃
M ⁺	0.4	1.0	0.4	0.9	0.9	1.4	0.9	0.9	1.1
Cs ⁺	0.9	0.0	0.4	0.0	0.0	0.0	1.4	0.4	0.4
HA ⁺	0.9	0.9	0.9	0.9	0.4	1.9	0.0	0.0	1.4
DA ⁺	0.9	1.4	0.4	0.0	0.9	0.9	0.4	1.4	0.9
MA ⁺	0.4	0.9	0.0	0.0	0.9	0.9	0.0	0.9	0.4
FM ⁺	0.0	0.9	0.9	0.9	0.0	0.4	0.9	0.0	0.4
FA ⁺	0.0	0.9	1.9	0.9	1.9	0.4	0.4	0.9	0.0
EA ⁺	1.9	1.9	0.0	0.0	1.1	0.0	1.4	0.0	0.0
GA ⁺	0.0	0.0	0.4	0.0	0.0	0.9	0.0	0.9	0.0
DEA ⁺	0.9	1.1	0.9	0.4	1.9	0.0	0.9	0.4	0.0

Table 1. Calculated various DMs for the $AM^{IV}X_3^{VII}$ perovskites containing pseudo halogen anions, $AM^{IV}BF_4$ and $AM^{IV}CN_3$ with $A = Cs^+, MA^+, FA^+$ and $M^{IV} = Pb^{2+}, Sn^{2+}$

Materials	H	E_g^d	ϵ_e^*	ϵ_h^*	E_{Be}	α_{ex}
Compounds based on BF_4^-						
CsPb BF_4	1	1	1 ϵ_0	1 4		1 4
Cs $SnBF_4$			1 ϵ_0	1 4		1 4
MAPb BF_4	1	3 4	ϵ_0 3		1	
MA $SnBF_4$		3 1	4 ϵ_0	3	3	ϵ_0
Compounds based on CN^-						
CsPb CN_3	3 1					