

**Figure S1.** (a) XRD patterns of  $(PMA)_2PbI_4$  and  $(pFPMA)_2PbI_4$  films. The inset is the enlarged XRD patterns with 2 $\theta$  between 5.8° and 6.8°. (b) UV-vis absorption spectra of the  $(PMA)_2PbI_4$  and  $(pFPMA)_2PbI_4$  films. The inset is the band gap of films.



**Figure S2.** Calculated electronic band structure for (PMA)<sub>2</sub>PbI<sub>4</sub>. (b) Calculated electronic band structure for (pFPMA)<sub>2</sub>PbI<sub>4</sub>.



Figure S3. (a) Band gaps of  $(PMA)_2PbI_4$  and  $(pFPMA)_2PbI_4$  calculated using PBE and PBE + vdW. (b) Effective masses of electrons and holes in  $(PMA)_2PbI_4$  and  $(pFPMA)_2PbI_4$ .