

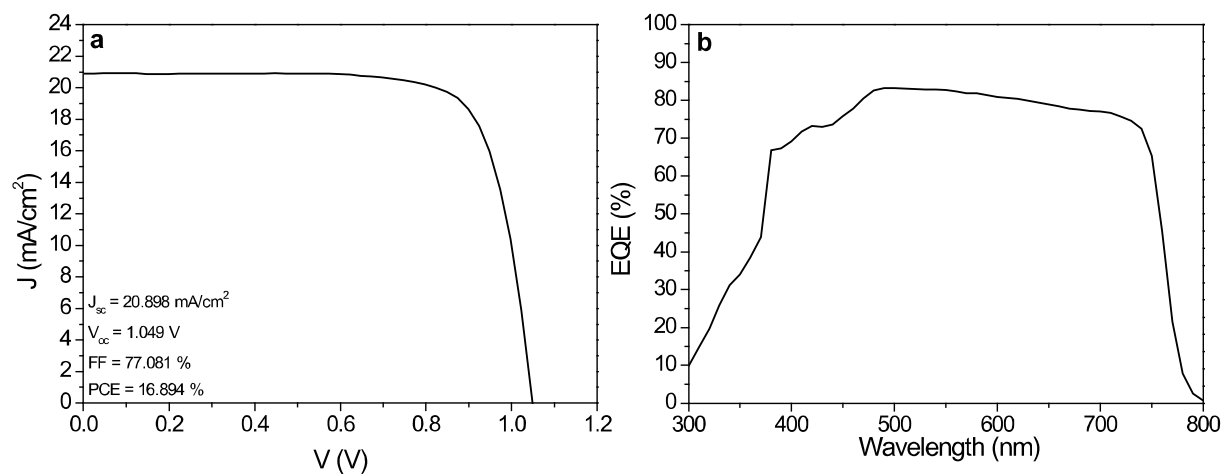
## Supporting Information

### Understanding Effects of Precursor Solution Aging in Triple Cation Lead Perovskite

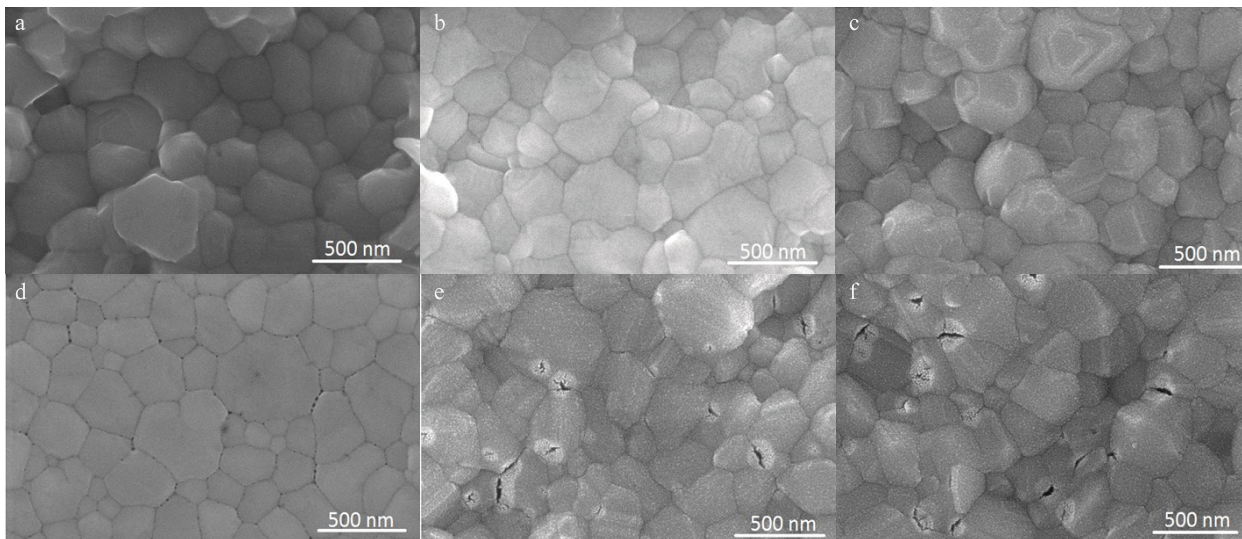
Passarut Boonmongkolras<sup>a</sup>, Daehan Kim<sup>a</sup>, Esra M. Alhabshi<sup>b</sup>, Issam Gereige<sup>b</sup>, Byungha Shin<sup>a\*</sup>

<sup>a</sup> Department of Material Science and Engineering, Korea Advanced Institute of Technology, Daejeon 34141, Korea

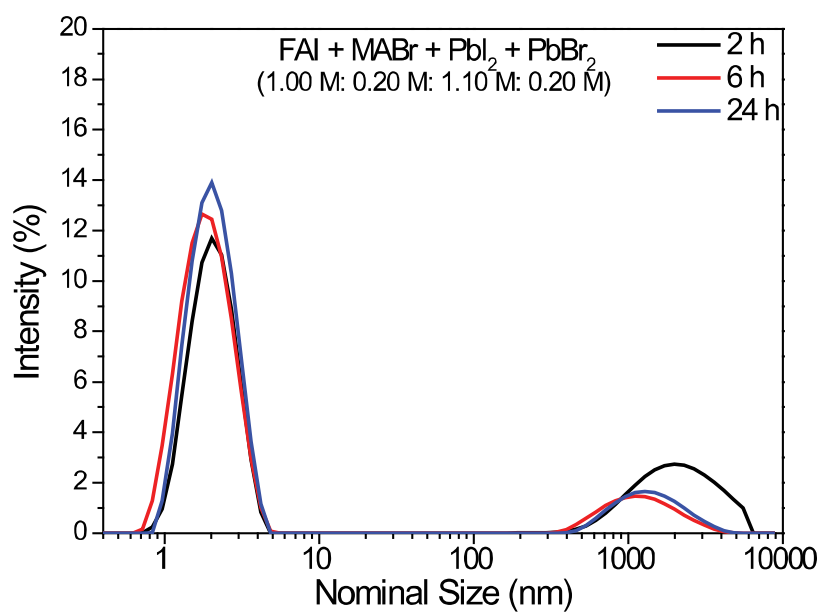
<sup>b</sup> Saudi Aramco Research & Development Center, Dhahran 31311, Kingdom of Saudi Arabia



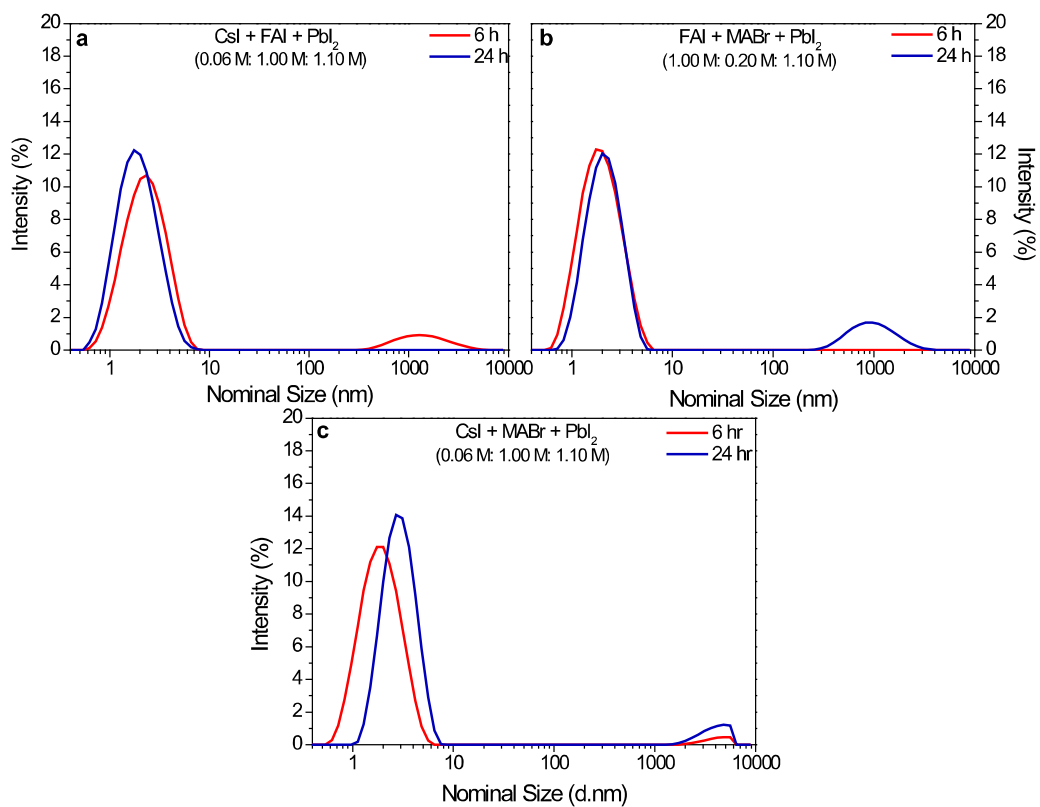
**Figure S1:** Device performance (a) JV curve and (b) EQE of the best performing cell.



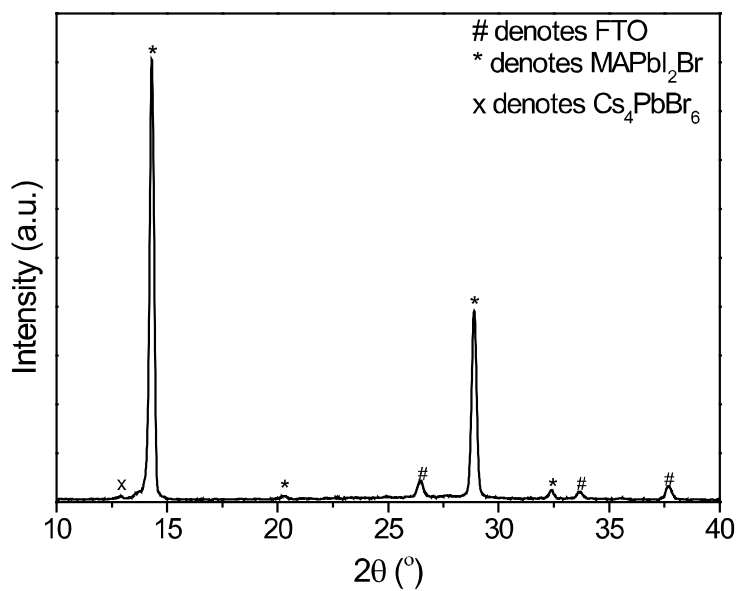
**Figure S2:** SEM images of the films fabricated from different triple cation lead halide perovskite solution where (a) – (f) are 2, 6, 24, 72, 168, and 720 hr, respectively.



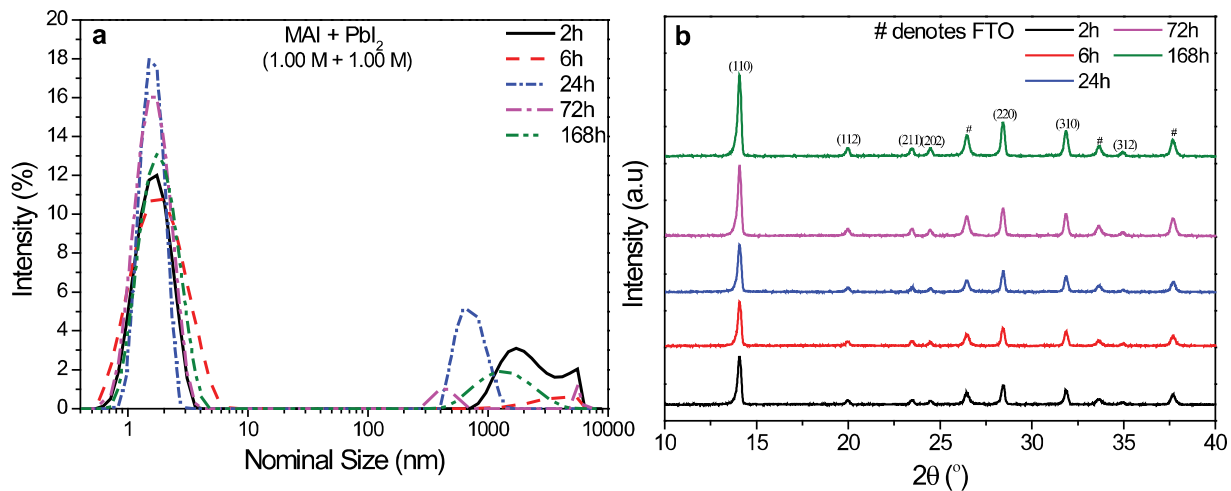
**Figure S3:** DLS data of the precursor solution containing FAI + MABr + PbI<sub>2</sub> + PbBr<sub>2</sub> with respect to its aging time.



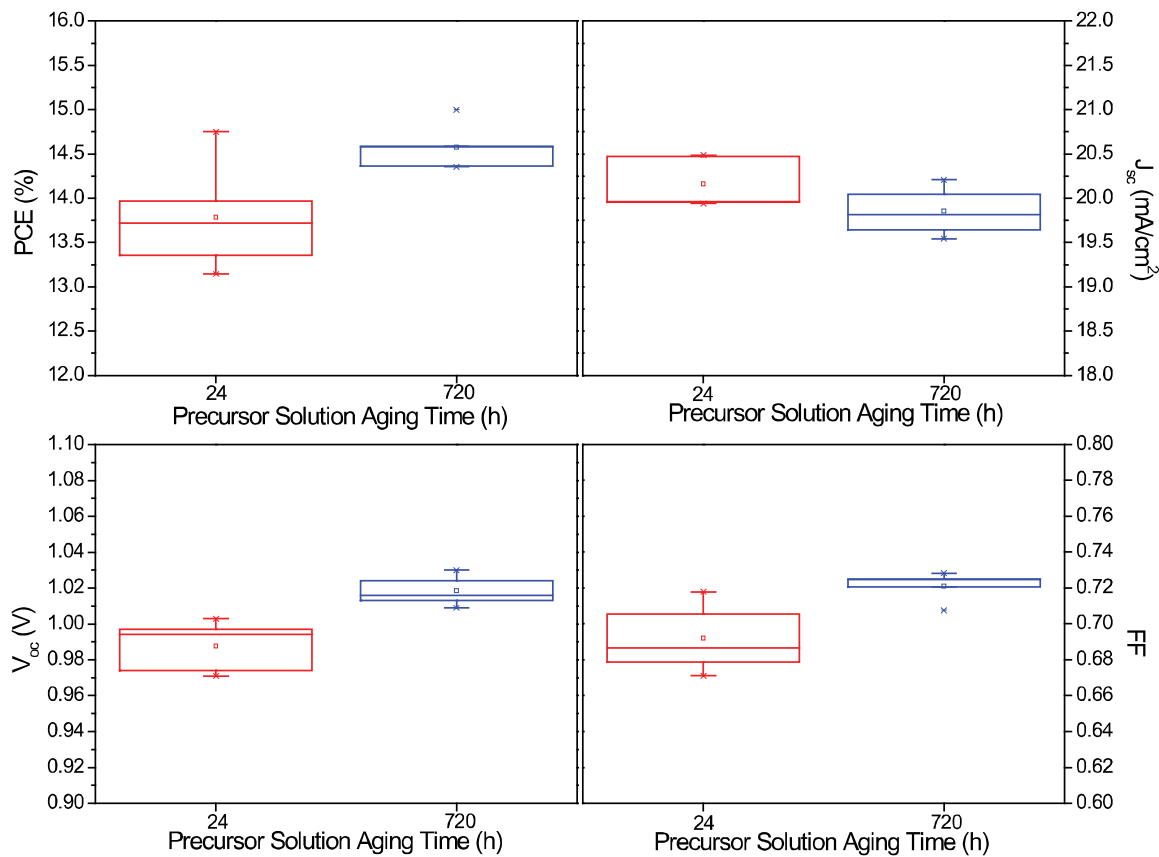
**Figure S4:** DLS data of the precursor solution containing (a) CsI + FAI + PbI<sub>2</sub>, (b) FAI + MABr + PbI<sub>2</sub>, and (c) CsI + MABr + PbI<sub>2</sub> with respect to its aging time.



**Figure S5:** XRD pattern of MAPbI<sub>2</sub>Br perovskite film with addition of CsI. The amount of CsI added was equal to the amount added in triple cation precursor solution.



**Figure S6:** (a) DLS and (b) XRD data of MAPbI<sub>3</sub> with respect to its perovskite precursor solution aging time.



**Figure S7:** Device distribution of MAPbI<sub>3</sub> perovskite with different precursor solution aging time.

Input Precursors	Cation: Lead: Halide Molar Ratio (A: B: X Molar Ratio)	6 h	24 h	72 h or more
FAI + PbI <sub>2</sub> + MABr + PbBr <sub>2</sub> + CsI (Triple-cation Lead Halide)	1.26: 1.30: 3.86	X	O	O
FAI + PbI <sub>2</sub> + MABr + PbBr <sub>2</sub>	1.20: 1.30: 3.80	O	O	O
FAI + PbBr <sub>2</sub>	1.00: 1.10: 3.20	O	O	O
MABr + PbBr <sub>2</sub>	1.00: 1.10: 3.20	X	O	O
FAI + PbI <sub>2</sub>	1.00: 1.10: 3.20	O	O	O
MABr + PbI <sub>2</sub>	1.00: 1.10: 3.20	X	O	O
CsI + FAI + PbI <sub>2</sub>	1.06: 1.10: 3.26	O	X	O
MABr + FAI + PbI <sub>2</sub>	1.20: 1.10: 3.40	X	O	O

**Table 1:** Summary of DLS experiments with different combinations of the precursors. The label “O” indicates the existence of the large aggregates within the precursor solution and “X” indicates no large aggregates.