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Exploring stability of formamidinium lead trihalide for solar cell application

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**Abstract:** Formamidinium lead triiodide (FAPbI<sub>3</sub>) is a promising photoactive perovskite for low-cost and efficient solar cells. This article reports on an experimental investigation on the stability of FAPbI<sub>3</sub> by comparison with that of widely-used methylamidinium lead triiodide (MAPbI<sub>3</sub>). A hydration of the FAPbI<sub>3</sub> with moisture could be the dominant mechanism for its degradation in air, rather than a common thermal decomposition in the MAPbI<sub>3</sub>. This can be mainly contributed to a relatively strong bond formation between formamidinium ions (FA<sup>+</sup>) and I<sup>-</sup>. Consequently, the stability of FAPbI<sub>3</sub> based devices can be greatly enhanced by removal moisture in the surrounding. This conclusion renders FAPbI<sub>3</sub> extremely attractive for stable perovskite solar cells with fine encapsulation.

Keywords: Perovskite; Solar cell; Stability; Cation

#### 1. Introduction

A broad organic-inorganic series of hybrid metal iodide perovskites with the general formulation ABX<sub>3</sub> have been reported for solar cell application, where A is the organic cation such as methylammonium  $(CH_3NH_3^+)$  or formamidinium  $(HC(NH_2)_2^+)$  at the corner of cubic unit cell, B being a divalent metal  $(Pb^{2+}, Sn^{2+})$  at the body center, and X being an halogen anion  $(Cl^-, Br^-, \Gamma)$  locating at the face center. Following the pioneer work of methylamidinium lead triiodide  $(MAPbI_3)$  perovskite

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