

How Important is the Organic Part of Lead Halide Perovskite Photovoltaic Cells? Efficient CsPbBr₃ Cells

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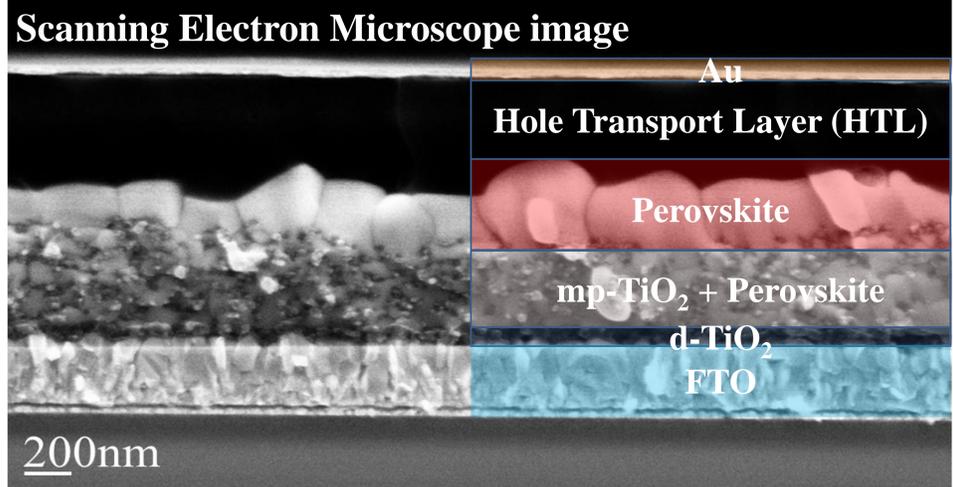
Background

To date, there are no reports on high band gap (> 2 eV) all-inorganic perovskite solar cells. CH₃NH₃PbBr₃-based devices have shown promising device parameters with open-circuit voltages surpassing 1.5 V [1], but is the organic cation really necessary to obtain high quality devices? CsPbBr₃, a high bandgap semiconductor with perovskite structure at standard temperature and pressure, has been estimated to have an electron mobility of ~1000 cm²/(V s) and comparable electron/hole lifetimes of 2.5 μs [2], emphasizing that it has potential as a promising absorber in solar cells. Here we discuss if perovskites with an inorganic A cation can form light absorbers with PV properties, comparable to those with an organic (e.g., alkyl ammonium) one, in particular in terms of the high open circuit voltages that are an important feature of these cells [3].

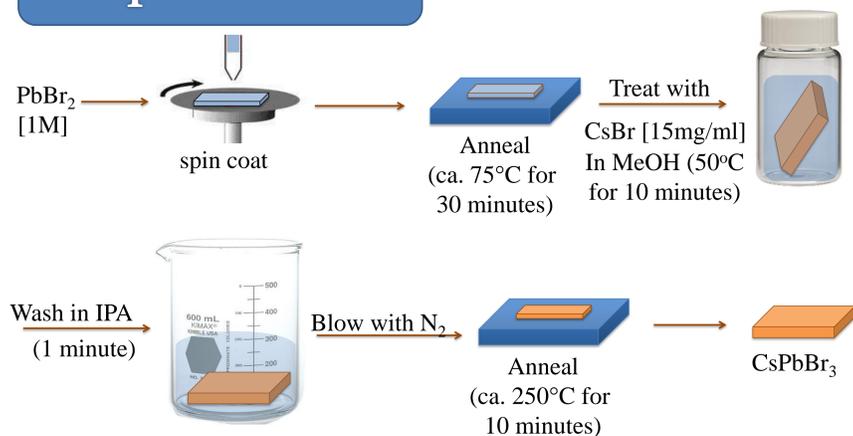
- [1] E. Edri et al., *J. Phys. Chem. Lett.* (2014), 5, 429-433.
 [2] C.C. Stoumpos et al., *Cryst. Growth Des.* (2013), 13, 2722-2727.
 [3] M. Kulbak, D. Cahen and G. Hodes, *J. Phys. Chem. Lett.* (2015), 6, 2452-2456.

Device Structure

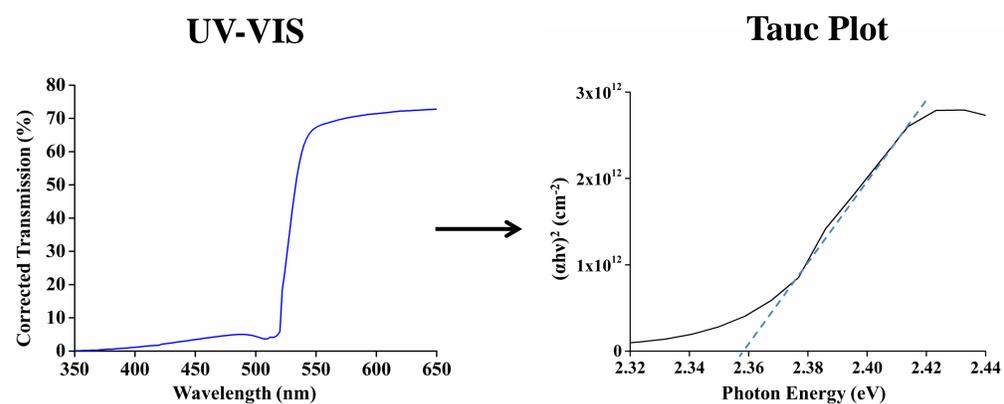
FTO/d-TiO₂/mp-TiO₂/CsPbBr₃/HTM/Au



Experimental

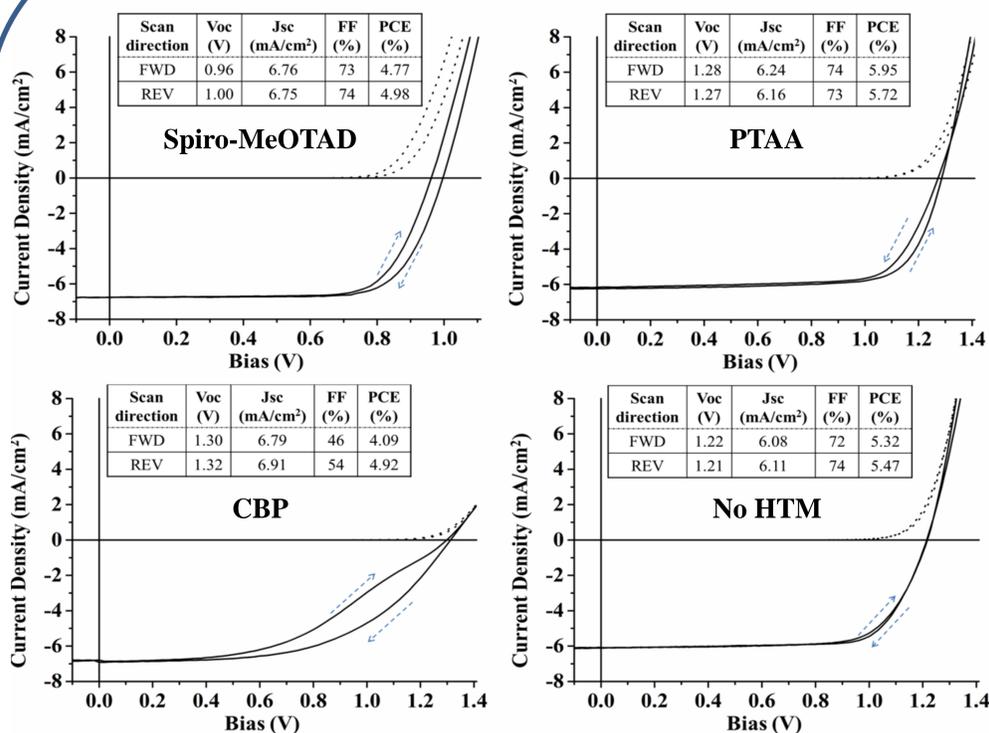


Optical properties of CsPbBr₃



- Films of CsPbBr₃ on FTO/dTiO₂/mp-TiO₂ show a drop in transmission at ca. 520 nm, correlating to a calculated direct bandgap of 2.36 eV.

Device I-Vs with different HTLs



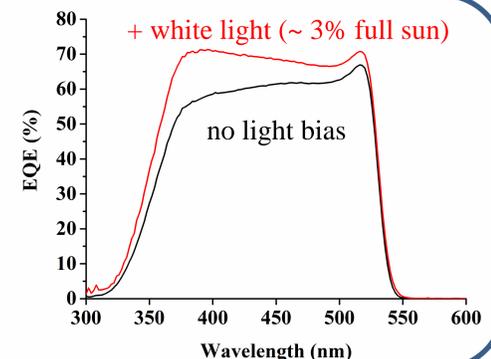
HTM	Spiro	PTAA	CBP
HOMO [eV]	5.0	5.2	5.7

- The deeper lies the HTL HOMO level, the higher is the V_{OC}.
- Devices without HTL layer work just as well as cells with it.
- No significant hysteresis is seen except with CBP.
- 1.3 V V_{OC} values are still below the top, 1.55 V, MAPbBr₃ one.

EQE

EQE of CsPbBr₃/PTAA device:

- White light increases the EQE close to 70 %.
- The EQE value yields an equivalent J_{SC} that agrees with the directly measured result (IV measurements).



Conclusions

We show CsPbBr₃ can have a PV performance that does not fall below that of CH₃NH₃Br₃, incl. encouraging V_{OC} values, comparable to our early results with the hybrid perovskite.

Acknowledgements

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