

Potential Hole Transport Materials for Perovskite Solar Cell (PSC)

Mrinal K. Das,^a Sinchan Maji,^b Pritam Dey,^c Tanmoy Maiti,^{c*} Debaprasad Mandal,^{b*} and Alakesh Bisai^{a*}^cDepartment of Material Science and Engineering, IIT Kanpur; ^bDepartment of Chemistry, IIT Ropar; ^aDepartment of Chemistry, IISER Bhopal*E-Mail: alakesh@iiserb.ac.in

Truxene(**1**) is a fluorescent polyaromatic structure that exhibits a poor solubility in most of the common organic solvents in its unsubstituted form.¹ To address this issue, introduction of solubilizing groups have been reported.^{1b} In fact, Gratzel and Nazeeruddin *et. al.* revealed that molecularly engineered *aza*-truxene(**2**) exhibit surface interaction with perovskite material resulting efficient hole injection from the valence band of perovskite into the HOMO of Hole Transport Material (HTM).² On the basis of the CV measurements, UV-vis, and PL spectra, the HOMO and LUMO energy levels of **2** was found to be in good alignment with the energy levels of perovskite material. Remarkable power conversion efficiency of 17.7% was realized using **2** as HTM and compositive perovskite as absorber, which is on par with that of spiro-OMeTAD (17.1%).³ Recently, Jen *et. al.* have reported a strong dipolar D-A (*Donor-Acceptor*) chromophore (**4**) with good hole-transporting ability as an efficient *dopant-free* HTM to realize high performance conventional PSCs, which shows power conversion efficiency of 17%.⁴

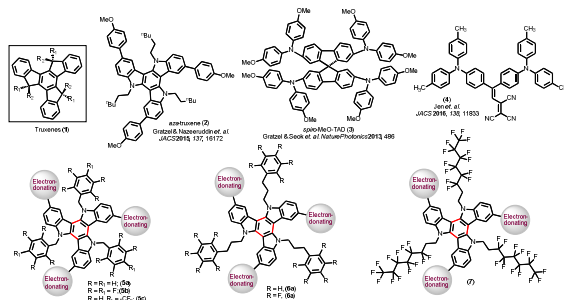


Figure. Potential Hole Transport Materials (HTMs) for Perovskite Solar Cell (PSC).

We hypothesized that *aza*-truxenes anchored with electron-deficient fragments, such as **5-7**, would not only improve interfacial properties and efficiency but also protect from humidity and UV induced degradation of PSCs. Our synthetic endeavours towards the development of two dimensional fluorinated *aza*-truxene materials and their spectroscopic behaviors will be discussed.⁵

References:

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