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S.no	Material	Structure	E _g (ev)	PCE (%)	Reference Number	year
1	MASnI₂Br	Glass/FTO/c-TiO2/m- TiO2/MASnIBr/Spiro-OMeTAD/Au	1.56	5.48	1	2014
2	CsGel ₃	Glass/FTO/c-TiO2/m-TiO2/CsGel/Spiro- OMeTAD/Au	1.64	0.11	2	2015
3	CsPb _{0.9} Sn _{0.1} IBr ₂	Glass/c-TiO2/m-TiO2/CsPbSnIBr/C	1.79	11.3	3	2017
4	CsSnIBr	Glass/FTO/c-TiO2/mp- TiO2/CsSnIBr/PTAA/Au	1.79	3.04	4	2017
5	CsPbl ₂ Br	Glass/FTO/NiOx/CsPbIBr/ZnO/C60/Ag	1.92	13.3	5	2018
6	CsPbIBr ₂	Glass/FTO/c-TiO2/CsPbIBr/Au	2.05	4.7	6	2016
7	MASnBr ₃	Glass/FTO/c-TiO2/m- TiO2/MASnBr/Spiro-OMeTAD/Au	2.15	4.56	1	2014
8	CsPbBr ₃	Glass/FTO/TiO2/CsPbBr/PTAA-TPB- LiTFSI/Au	2.36	6.24	7	2016

 Table S 1: High band gap value of organic and inorganic based perovskite materials for multi junction solar cell application.

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Fig.S1 Solution pictures for (a) Cesium chloride with solvent DMF (b) Tin chloride with solvent DMF (c) Both Cesium and Tin chloride with solvent DMF (d) Cesium chloride with solvent Water (e) Tin chloride with solvent Water (f) Both Cesium and Tin chloride with solvent Water (g) Tin chloride with solvent Water (0.1M HCl) & (h) Both Cesium and Tin chloride with solvent Water (0.1M HCl)



Fig. S2 Reitveld refinement using FullProf Software for Cs₂SnCl₆



Fig. S3 XRD Pattern obtained for (MA_{0.5}Cs_{0.5})₂SnCl6 at 14.79°



Fig. S4 PL spectrum for (a) Cs_2SnCl_6 (b) MA_2SnCl_6 and (c) $(MA_{0.5}Cs_{0.5})_2SnCl_6$





Fig.S5 SAED Pictures for (a) Cs_2SnCl_6 (b) MA_2SnCl_6and (c) $(MA_{0.5}Cs_{0.5})_2SnCl_6$



Fig.S6 EDAX analysis for synthesized Cs_2SnCl_6



Fig.S7 EDAX analysis for synthesized MA₂SnCl₆



Fig.S8 EDAX analysis for synthesized (MA_{0.5}Cs_{0.5})₂SnCl_{6.}