Supporting Information

Diammonium Spacer-Induced Stable Zigzag Type 2D Dion-Jacobson Lead/Tin-Based Perovskite Solar Cells

Mi-Hee Jung*, †

Department of Nanotechnology and Advanced Materials Engineering, Sejong University, 209, Neungdong-ro, Gwangjin-gu, Seoul 05006, Republic of Korea; Fax: +82-2-3408-4342; Tel: +82-2-6935-2597

*To whom correspondence should be addressed. E-mail: mhjung@sejong.ac.kr



Figure S1. Crystal structure of $C_{18}H_0I_{16}N_6Sn_4$



Figure S2. The enlarged SEM images of the (a) 10% and (b) 15% of $3API_2$ added $FA(Pb_{0.5}Sn_{0.5})I_3$ perovskite films.

		10	15	20
0.095	0.097	0.098	0.094	0.099
0.123	0.121	0.121	0.120	0.124
	0.095	0.095 0.097 0.123 0.121	0.0950.0970.0980.1230.1210.121	0.0950.0970.0980.0940.1230.1210.1210.120

Table S1. Full width half maximum (FWHM) values calculated from the (110) diffraction peak in Figure



Figure S3. J-V curves of the 2D DJ (3AP) $FA_{n-1}(Pb_{0.5}Sn_{0.5})_nI_{3n+1}$ (n = 1, 2, 3, 4) based devices

Table S2. The photovoltaic parameters of the 2D DJ $(3AP)FA_{n-1}(Pb_{0.5}Sn_{0.5})_nI_{3n+1}$ (n = 1, 2, 3, 4) based devices

Perovskites	Jsc [mA/cm ²]	Voc [V]	FF [%]	PCE [%]
$(3AP)(Pb_{0.5}Sn_{0.5})I_4$	15.47	0.25	35.04	1.38
$(3AP)(FA)(Pb_{0.5}Sn_{0.5})_2I_7$	17.42	0.29	38.27	1.92
$(3AP)(FA)_2(Pb_{0.5}Sn_{0.5})_3I_1$	20.10	0.39	49.39	3.86
$(3AP)(FA)_3(Pb_{0.5}Sn_{0.5})_4I_1$	21.37	0.40	51.81	4.42