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Aqueous Electrolytes Based Dye-Sensitized Solar Cells Using Organic Sensitizers

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Figure S1. J-V curve and IPCE spectra of **JK-2.**

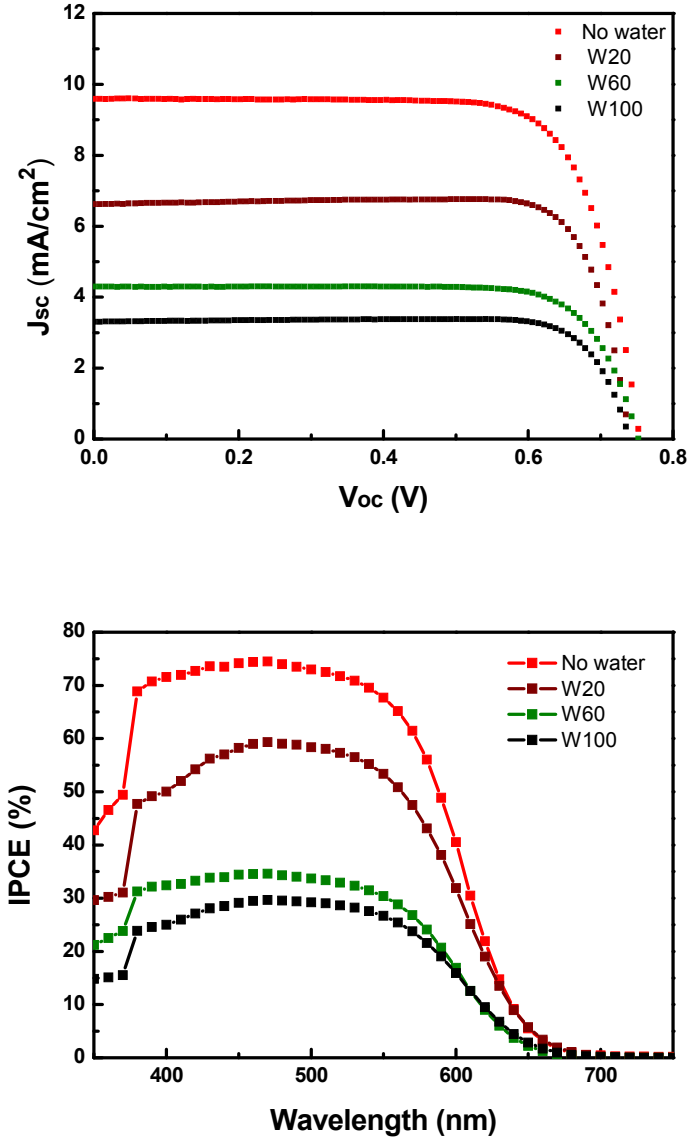

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Table S1. Calculated excitation energy characteristics of the **JK-259** and **JK-262.**

Table S1. Calculated excitation energy characteristics of the **JK-259** and **JK-262**[a]

Compounds	energy (eV)	F[b]	Composition (%)[c]
JK-259	1.8894	0.4167	71 (HOMO → LUMO)
	2.6419	1.4480	63 (HOMO-1 → LUMO)
			31 (HOMO → LUMO+1)
	2.7705	0.1608	31 (LUMO → HOMO-1)
			63 (HOMO → LUMO+1)
JK-262	2.0512	0.5538	71 (HOMO → LUMO)
	2.9659	0.8528	68 (HOMO-1 → LUMO)
			16 (LUMO+1 → HOMO)
	3.3641	0.2336	23 (LUMO → HOMO-2)
			13 (HOMO-1 → LUMO)
			65 (HOMO → LUMO+1)

[a] The characteristics were calculated by the time dependent-density functional theory (TD-DFT) using the B3LYP functional/6-31G* basis set.
[b] The oscillator strength (*f*) of a transition is a measure of its intensity and is related to the molar absorption coefficient.
[c] The composition means contribution of each transition for excitation energies.