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Corrigendum

Highly efficient new indoline dye having strong electron-withdrawing group for zinc oxide dye-sensitized solar cell [Tetrahedron 67 (34) (2011) 6289–6293]

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The authors have identified some errors in the published version of their paper. The corrected text is provided below. Page 6289, left hand column, second paragraph:

The $E_{\rm ox}$ level should be more positive than ca. 0.2 V versus Fc/Fc^+ to show high IPCE, corresponding to HOMO level more stable than -4.9 eV by the density functional theory (DFT) calculations. Page 6291:

Table 1 Physical properties of indoline dyes

Dye	$\lambda_{\max} (\varepsilon)^a / nm$	λ_{max} on ZnO/nm	$F_{\rm max}^{\ a}/{\rm nm}$	$E_{\rm ox}^{\ \ b}/V$	$E_{\text{ox}} - E_{0-0}^{\text{c}} / \text{V}$	HOMO ^d /eV	LUMO ^d /eV
D205	395 (38,100), 554 (74,700)	540	641	+0.35	-1.73	-5.06	-2.36
DN317	373 (28,200), 521 (61,900)	505	608	+0.35	-1.85	-4.99	-2.15
DN319	400 (36,700), 566 (68,000)	542	662	+0.37	-1.66	-5.18	-2.53

 $^{^{\}rm a}$ Measured on 1.0×10 $^{\rm -5}$ mol dm $^{\rm -3}$ of substrate in chloroform at 25 $^{\circ}$ C.

Page 6291, left hand column, first complete paragraph:

The HOMO level of DN319 was calculated to be -5.18 eV, being sufficiently stable to show high IPCE.

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b Versus Fc/Fc⁺ in DMF.

^c Calculated on the basis of $E_{\rm ox}$ and $\lambda_{\rm int}$.

^d Calculated by the B3LYP/6-31G(d,p)//B3LYP/3-21G level.

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