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CORRECTION

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Correction: Twisted organic TADF triads based on a diindolocarbazole donor for efficient photoisomerization of stilbene and photo-arylation of heteroarenes

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Correction for 'Twisted organic TADF triads based on a diindolocarbazole donor for efficient photoisomerization of stilbene and photo-arylation of heteroarenes' by Sushil Sharma *et al.*, *Org. Chem. Front.*, 2023, **10**, 6087–6095, https://doi.org/10.1039/D3QO01542D.

The authors regret that the singlet and triplet energy values were incorrectly reported in Table 1. The corrected Table 1 appears below.

Table 1 Absorption wavelength (λ_{abs}), emission wavelength (λ_{em}), Stokes shift, molar extinction coefficient (ε) and singlet energy (S_1), triplet energy (T_1) and singlet–triplet energy gap (ΔE_{ST}) of **DI-PF** and **DI-PI**

Compounds	λ_{ex} (nm)	λ_{em} (nm)	Stokes shift (nm)	ε (M^{-1} cm^{-1})	S_1 (eV)	T_1 (eV)	ΔE_{ST} (eV)
DI-PF	279, 373, 394	590	196	48 268	2.39	2.38	0.01
DI-PI	328, 363, 393	519	126	80 783	3.14	2.77	0.37

The Royal Society of Chemistry apologises for these errors and any consequent inconvenience to authors and readers.

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