ORGANIC CHEMISTRY FRONTIERS

CORRECTION

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Cite this: *Org. Chem. Front.*, 2025, **12**, 3942

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

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DOI: 10.1039/d5qo90048d

rsc.li/frontiers-organic

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 ($\lambda_{abs}$), emission wavelength ($\lambda_{em}$), Stokes shift, molar extinction coefficient ($\varepsilon$) and singlet energy ($S_1$), triplet energy ($T_1$) and singlet–triplet energy gap ($\Delta E_{ST}$) of **DI-PF** and **DI-PI**

| Compounds | $\lambda_{ex}$ (nm) | $\lambda_{em}$ (nm) | Stokes shift (nm) | $\varepsilon$ ($M^{-1}$ $cm^{-1}$) | $S_1$ (eV) | $T_1$ (eV) | $\Delta E_{ST}$ (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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