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Correction: Solvent dependent triplet state delocalization in a co-facial porphyrin heterodimer

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Correction for 'Solvent dependent triplet state delocalization in a co-facial porphyrin heterodimer' by Susanna Ciuti et al., *Phys. Chem. Chem. Phys.*, 2022, <https://doi.org/10.1039/D2CP04291F>.

The authors would like to correct an error in Fig. 4 of the published article, which did not contain the simulated spectra. The correct figure is shown below:

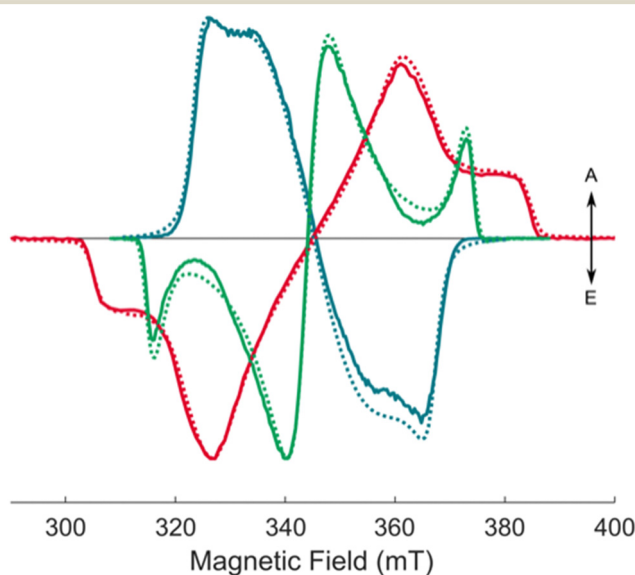


Fig. 4 Experimental (solid) and calculated (dashed) X-band transient EPR spectra of the AlPor–O–PPor–PF₆ dimer (teal) and the two reference monomers AlPor–OH (red), and PPor–OMe–PF₆ (green) in 3 : 1 MeTHF : CH₂Cl₂ at 80 K and ~1 μs after the laser flash (λ_{ex} = 575 nm).

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

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