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CORRECTION

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# Correction: An exchange coupled *meso–meso* linked vanadyl porphyrin dimer for quantum information processing

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Correction for 'An exchange coupled *meso–meso* linked vanadyl porphyrin dimer for quantum information processing' by Davide Ranieri *et al.*, *Chem. Sci.*, 2022, https://doi.org/10.1039/d2sc04969d.

During the editorial production of the finished article, the graphics associated with Fig. 3(a) and (b) were inadvertently switched. The intended version of Fig. 3 is shown below, and replaces that of the original article:

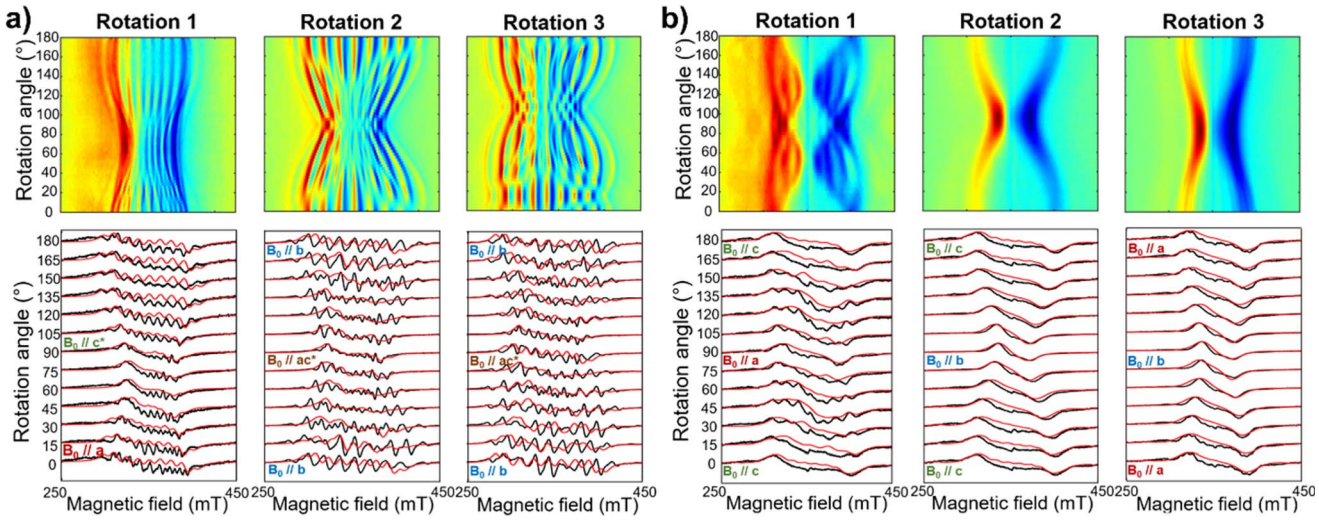


Fig. 3 Room temperature angular-dependent CW EPR X-band spectra of (a) *m*-[VO(TrPP)]$_2$ and (b) *o*-[VO(TrPP)]$_2$ for crystal rotations around three orthogonal axes. For both panels, the upper row shows the 2D experimental EPR contour plots for the three rotations, acquired with a 3° step; the lower row shows representative EPR spectra (black lines) for the three rotations – from 0° to 180° every 15° – together with the best spectral simulations (red lines) obtained by using $|J| = 0.01$ (0.005) cm$^{-1}$ and $|J| = 0.05$ (0.01) cm$^{-1}$ for (a) and (b), respectively. Experimental frequency: 9.40 GHz for (a), 9.87 GHz for (b).

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

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