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CORRECTION

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Correction: ^{57}Fe Mössbauer spectroscopy and high-pressure structural analysis for the mechanism of pressure-induced unique magnetic behaviour in (cation)[$Fe^{II}Fe^{III}$(dto)$_3$] (cation = Ph_4P and $^nPrPh_3P$; dto = 1,2-dithiooxalato)

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Correction for '^{57}Fe Mössbauer spectroscopy and high-pressure structural analysis for the mechanism of pressure-induced unique magnetic behaviour in (cation)[$Fe^{II}Fe^{III}$(dto)$_3$] (cation = Ph_4P and $^nPrPh_3P$; dto = 1,2-dithiooxalato)' by Ryosuke Taniai *et al.*, *Dalton Trans.*, 2023, **52**, 8368–8375.

The authors regret that several errors were published in the original article. The correct details are given below.

In the first paragraph of "Synthetic procedures" on page 8369, "$C_{30}H_{20}Fe_2O_6PS_6$" should read "$C_{27}H_{22}Fe_2O_6PS_6$".

In the third paragraph of the section "Crystal structures" on page 8370, the sentence "the nearest C6···S2^b and C7···O2^c distances were 3.62(3) and 3.28(3) Å, respectively", should read "the nearest C6···S2^b and C4···O1^a distances were 3.62(3) and 3.26(3) Å, respectively".

In addition, the authors regret that an incorrect version of Fig. 2 was included in the original article. The correct version of Fig. 2 is presented here.

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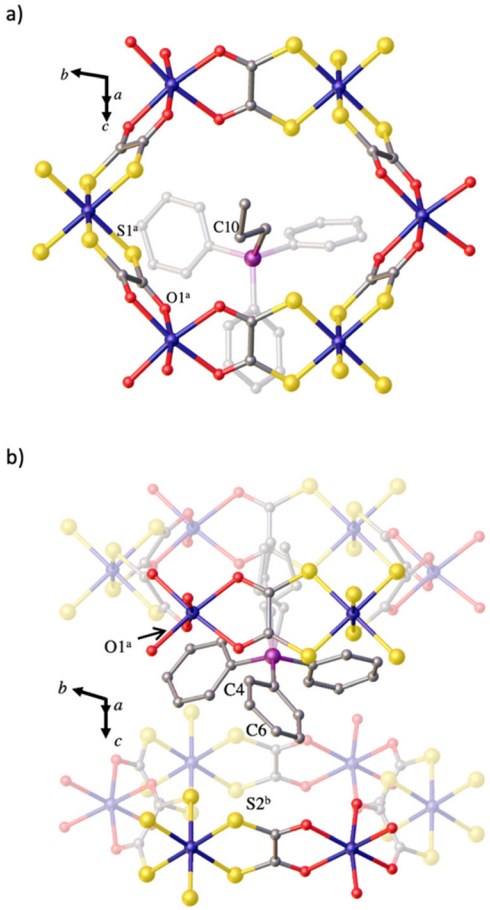

Fig. 2 (a) Coordination environment of O_6 and S_6 atoms for the two Fe ions in **3**. Colour codes: gray, red, yellow, purple and navy represent C, O, S, P, and Fe atoms, respectively. A 2-D honeycomb layer in the *ab* plane is shown. (b) Layered structure along the *c* axis. One side of two disordered configurations is shown in the $^{n}PrPh_3P$ position. Symmetry codes: (a) 1 + *x*, 1 + *y*, +*z*; (b) 1 + *y*, 1−*x* + *y*, 1/2 + *z*.

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