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Correction: 2D → 2D parallel interpenetration of (4,4) sheets constructed from a ditopic bis(4,2':6',4"-terpyridine)

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Correction for '2D → 2D parallel interpenetration of (4,4) sheets constructed from a ditopic bis(4,2':6',4"-terpyridine)' by Edwin C. Constable et al., *CrystEngComm*, 2014, **16**, 3494–3497.

The crystallographic data for the compound $\{[\text{Zn}_2\text{Cl}_4(1)] \cdot 4\text{H}_2\text{O}\}_n$ reported in footnote 28 was incorrectly given as that of the free ligand: $\text{C}_{52}\text{H}_{56}\text{N}_6\text{O}_2$, $M = 797.03$, yellow block, monoclinic, space group $P2_1/c$, $a = 13.3440(8)$, $b = 11.0423(7)$, $c = 14.6607(9)$ Å, $\beta = 90.571(3)^\circ$, $U = 2160.1(2)$ Å³, $Z = 2$, $D_c = 1.225$ Mg m⁻³, $\mu(\text{Cu-K}\alpha) = 0.589$ mm⁻¹, $T = 123$ K. Total 31 623 reflections, 3923 unique, $R_{\text{int}} = 0.0353$. Refinement of 3425 reflections (272 parameters) with $I > 2\sigma(I)$ converged at final $R_1 = 0.0393$ (R_1 all data = 0.0447), $wR_2 = 0.1046$ (wR_2 all data = 0.1096), gof = 1.034. CCDC 986027.

The correct data for $\{[\text{Zn}_2\text{Cl}_4(1)] \cdot 4\text{H}_2\text{O}\}_n$ are as follows:

$\text{C}_{52}\text{H}_{56}\text{Cl}_4\text{N}_6\text{O}_2\text{Zn}_2 \cdot 4\text{H}_2\text{O}$, $M = 1141.64$, yellow block, monoclinic, space group $C2/c$, $a = 20.6102(11)$, $b = 11.5999(6)$, $c = 23.8198(12)$ Å, $\beta = 90.978(3)^\circ$, $U = 5693.9(5)$ Å³, $Z = 4$, $D_c = 1.322$ Mg m⁻³, $\mu(\text{Cu-K}\alpha) = 3.169$ mm⁻¹, $T = 123$ K. Total 39 090 reflections, 5115 unique, $R_{\text{int}} = 0.0345$. Refinement of 4828 reflections (382 parameters) with $I > 2\sigma(I)$ converged at final $R_1 = 0.0591$ (R_1 all data = 0.0616), $wR_2 = 0.1643$ (wR_2 all data = 0.1659), gof = 1.192. CCDC 986028.

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

