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CORRECTION

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Correction: [3+3] Imine and β-ketoenamine tethered fluorescent covalent-organic frameworks for CO_2 uptake and nitroaromatic sensing

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Correction for '[3+3] Imine and β-ketoenamine tethered fluorescent covalent-organic frameworks for CO_2 uptake and nitroaromatic sensing' by D. Kaleeswaran *et al.*, *J. Mater. Chem. C*, 2015, **3**, 7159–7171.

In Table 1 (page no. 7166), the unit for average pore diameter should be Å instead of nm. A suitably corrected table is shown below.

Table 1 N_2 (at 77 K), H_2 (at 77 K) and CO_2 (at 273 K) sorption data for COFs TAPB-TFPB, TAPB-TFP, *i*PrTAPB-TFPB and *i*PrTAPB-TFP at 1.0 bar

COF(s)	SA_{BET} ($m^2\ g^{-1}$)	SA_{Lang} ($m^2\ g^{-1}$)	Average pore diameter (Å)	H_2 uptake (wt%)	CO_2 uptake (mg g^{-1})	CO_2 uptake (wt%)
TAPB-TFPB	229.4	539.5	40	0.68	40.1	4.0
*i*PrTAPB-TFPB	390.6	1191.0	50	0.43	31.2	3.1
TAPB-TFP	567.0	989.0	26	1.08	180.0	18.0
*i*PrTAPB-TFP	756.0	1515.0	34	1.15	105.2	10.5

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

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