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## Correction: [3+3] Imine and $\beta$ -ketoenamine tethered fluorescent covalent-organic frameworks for CO<sub>2</sub> uptake and nitroaromatic sensing

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Correction for '[3+3] Imine and  $\beta$ -ketoenamine tethered fluorescent covalent-organic frameworks for CO<sub>2</sub> 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<sub>2</sub> (at 77 K), H<sub>2</sub> (at 77 K) and CO<sub>2</sub> (at 273 K) sorption data for COFs TAPB-TFPB, TAPB-TFP, *i*PrTAPB-TFPB and *i*PrTAPB-TFP at 1.0 bar

COF(s)	S <sub>A</sub> BET (m <sup>2</sup> g <sup>-1</sup> )	S <sub>A</sub> Lang (m <sup>2</sup> g <sup>-1</sup> )	Average pore diameter (Å)	H <sub>2</sub> uptake (wt%)	CO <sub>2</sub> uptake (mg g <sup>-1</sup> )	CO <sub>2</sub> uptake (wt%)
TAPB-TFPB	229.4	539.5	40	0.68	40.1	4.0
<i>i</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>i</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.

