MSDE



Check for updates

Cite this: Mol. Syst. Des. Eng., 2022, 7, 1162

Correction: Engineering metal-organic frameworks for adsorption-based gas separations: from process to atomic scale

Marco Taddei^{*a} and Camille Petit^{*b}

DOI: 10.1039/d2me90023h

rsc.li/molecular-engineering

Correction for 'Engineering metal–organic frameworks for adsorption-based gas separations: from process to atomic scale' by Marco Taddei and Camille Petit, *Mol. Syst. Des. Eng.*, 2021, **6**, 841–875, DOI: https://doi.org/10.1039/D1ME00085C.

The authors regret an error made in a sentence found in section 3.3. The sentence found in the original manuscript reads "A high heat capacity limits the temperature swing during sorption. Therefore, in a TSA process, for the same energy input used to drive desorption, the working capacity increases, which helps the CAPEX". The corrected sentence reads as follows: "A high heat capacity limits the temperature swing during sorption. Therefore, in a TSA process, for the same energy input used to drive desorption, the working capacity decreases, which increases the CAPEX".

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

^a Department of Chemistry and Industrial Chemistry, University of Pisa, Via Giuseppe Moruzzi, 13, Pisa, Italy. E-mail: marco.taddei@unipi.it

^b Barrer Centre, Department of Chemical Engineering, Imperial College London, South Kensington Campus, London SW7 2AZ, UK. E-mail: camille.petit@imperial.ac.uk



View Article Online View Journal | View Issue