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## Correction: Thermally induced top-down nanostructuring for the synthesis of a core/shell-structured CoO/CoS<sub>x</sub> electrocatalyst

Min Soo Kim,<sup>a</sup> Muhammad Awais Abbas,<sup>b</sup> Raju Thota<sup>b</sup> and Jin Ho Bang<sup>\*abc</sup>Correction for 'Thermally induced top-down nanostructuring for the synthesis of a core/shell-structured CoO/CoS<sub>x</sub> electrocatalyst' by Min Soo Kim *et al.*, *J. Mater. Chem. A*, 2019, 7, 26557–26565.

The authors regret a mistake in Fig. 5, where Co<sub>3</sub>O<sub>4</sub> and CoO curves in Fig. 5(b) were mistakenly interchanged. The corrected figure is presented below:

There were also errors in the sentence "The calculated TOFs of Co<sub>3</sub>O<sub>4</sub>, CoO, CoO/CoS<sub>x</sub>, and IrO<sub>2</sub> at an overpotential of 400 mV were 0.81, 2.79, 3.46, and 5.03 s<sup>-1</sup>, respectively." on page 26563. The corrected sentence is copied below:

"The calculated TOFs of Co<sub>3</sub>O<sub>4</sub>, CoO, CoO/CoS<sub>x</sub>, and IrO<sub>2</sub> at an overpotential of 400 mV were 0.081, 0.279, 0.346, and 0.503 s<sup>-1</sup>, respectively."

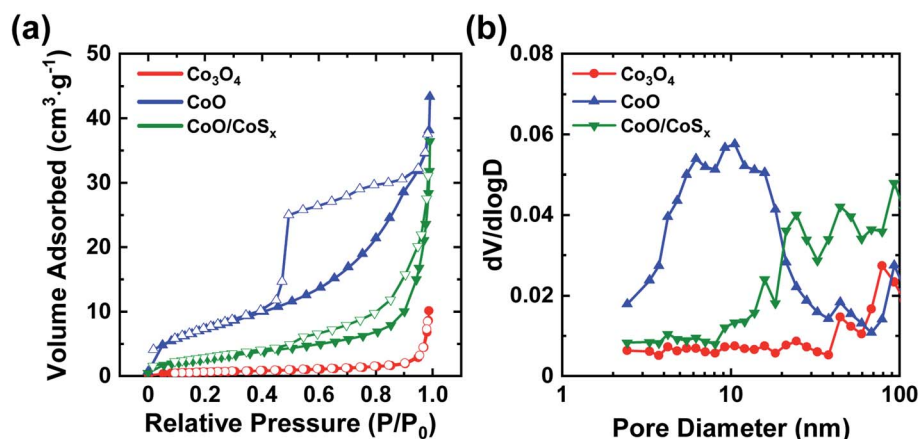


Fig. 5 (a) N<sub>2</sub> physisorption isotherms and (b) pore size distributions of Co<sub>3</sub>O<sub>4</sub>, CoO, and CoO/CoS<sub>x</sub> particles.

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

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