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## CORRECTION

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## Correction: An *in situ* growth route towards antiperovskite Ni<sub>3</sub>InN nanoparticles embedded within amorphous silicon nitride

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Correction for 'An *in situ* growth route towards anti-perovskite Ni<sub>3</sub>InN nanoparticles embedded within amorphous silicon nitride' by Shotaro Tada *et al., J. Mater. Chem. A*, 2024, 12, 3689–3699, https://doi.org/10.1039/D3TA06212K.

The authors apologise for an error in Fig. 6. The graphs for Fig. 6b and c were incorrectly swapped so that they appeared in the wrong position. The corrected figure is shown here.

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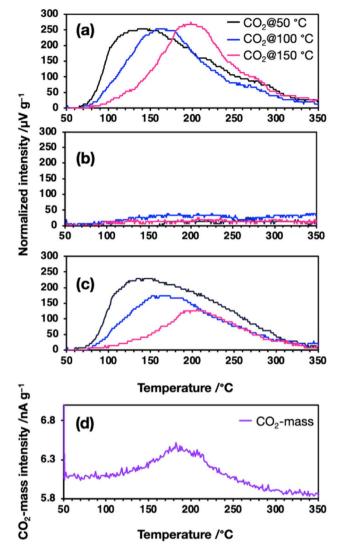


Fig. 6  $CO_2$ -TPD spectra under flowing He recorded for Ni<sub>3</sub>InN/a-SiN nanocomposites: normalized TCD curves of (a) the Ni<sub>0.05</sub>In<sub>0.1</sub>-DRZ600 sample, (b) bulk Ni<sub>3</sub>InN obtained by ammonolysis of an oxide precursor and (c) microporous amorphous SiN derived from In<sub>0.1</sub>-DRZ600 synthesized by pyrolysis under NH<sub>3</sub> at 600 °C and (d) CO<sub>2</sub>-mass spectra of the Ni<sub>0.05</sub>In<sub>0.1</sub>-DRZ600 sample recorded after CO<sub>2</sub> treatment at 150 °C.

In addition, the authors regret a mistake in the text of the manuscript. In the left column on page 3696 the section that begins "In contrast, the  $CO_2$ -TPD spectra of the bulk  $Ni_3InN$  sample (Fig. 6b)..." and ends "...which decreases in intensity with increasing  $T_{CO_2}$ , indicating  $CO_2$  physisorption behavior." should be as shown below:

"In contrast, the CO<sub>2</sub>-TPD spectra of the bulk Ni<sub>3</sub>InN sample (Fig. 6b) exhibits no pronounced peak under the same measurement conditions. Interestingly, the CO<sub>2</sub> desorption peak intensity of the Ni<sub>0.05</sub>In<sub>0.1</sub>-DRZ600 sample increases consistently with increasing  $T_{CO_2}$ , suggesting the CO<sub>2</sub> chemisorption behavior (Fig. 6a). In contrast, the In<sub>0.1</sub>-DRZ600 sample exhibits a broad CO<sub>2</sub> desorption curve which decreases in intensity with increasing  $T_{CO_2}$ , indicating CO<sub>2</sub> physisorption behavior (Fig. 6c)".

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