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CORRECTION





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# Correction: X-ray absorption near edge structure simulation of $LiNi_{0.5}Co_{0.2}Mn_{0.3}O_2$ *via* first-principles calculation

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Correction for 'X-ray absorption near edge structure simulation of $LiNi_{0.5}Co_{0.2}Mn_{0.3}O_2$ *via* first-principles calculation' by Toshiharu Ohnuma *et al.*, *RSC Adv.*, 2019, **9**, 35655–35661.

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

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