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Correction: X-ray absorption near edge structure simulation of $\text{LiNi}_{0.5}\text{Co}_{0.2}\text{Mn}_{0.3}\text{O}_2$ via first-principles calculation

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Correction for 'X-ray absorption near edge structure simulation of $\text{LiNi}_{0.5}\text{Co}_{0.2}\text{Mn}_{0.3}\text{O}_2$ 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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