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CORRECTION

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Correction: Reactive simulations of the activation barrier to dissolution of amorphous silica in water

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Correction for 'Reactive simulations of the activation barrier to dissolution of amorphous silica in water' by Michael Kagan *et al.*, *Phys. Chem. Chem. Phys.*, 2014, **16**, 9294–9301.

On page 9300, in the last sentence of the first paragraph of the Conclusions, the first data range is incorrectly shown as 18–24 kcal mol^{-1}, instead of 14–24 kcal mol^{-1}. The sentence should therefore read as follows:

This result is within the lower end of the experimental data, which varies from 14–24 kcal mol^{-1}, while various *ab initio* calculations using small cluster models obtain values that vary from 18–39 kcal mol^{-1}.

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

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