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CORRECTION

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# Correction: Accelerating materials discovery using integrated deep machine learning approaches

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Correction for 'Accelerating materials discovery using integrated deep machine learning approaches' by Weiyi Xia *et al.*, *J. Mater. Chem. A*, 2023, **11**, 25973–25982, https://doi.org/10.1039/d3ta03771a.

The authors apologise for an error in Fig. 5c and d. The figure previously mistakenly portrayed the electronic band structure and electronic density of states, rather than the described phonon dispersion and density of states. The corrected figure is shown below.

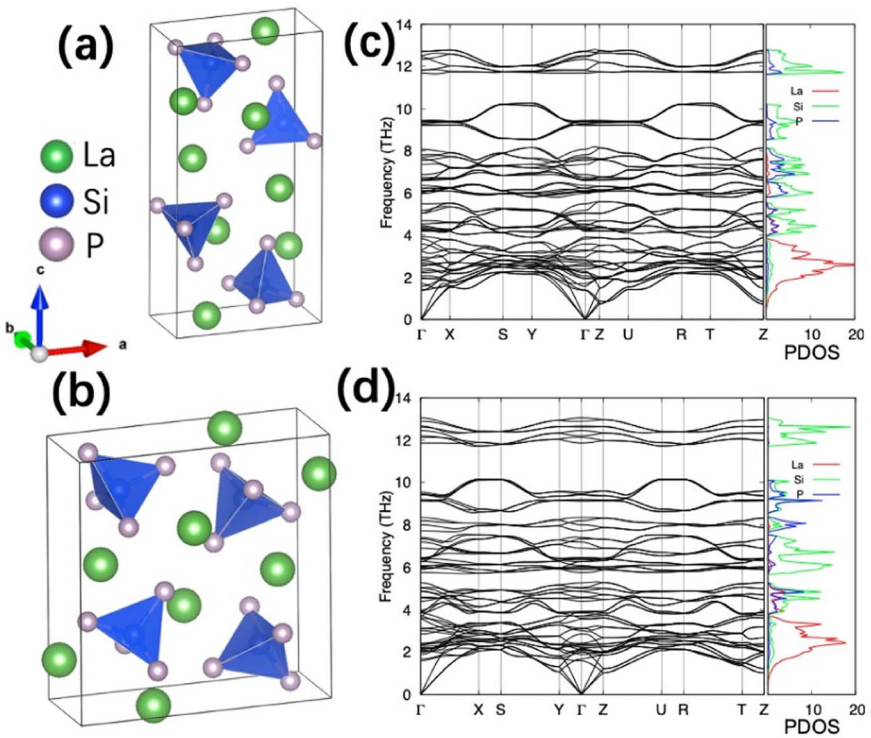


Fig. 5 (a and b) The structures of the predicted two $La_2SiP_3$ phases. The formation energies above the convex hull are 1 meV per atom and 33 meV per atom, respectively. (c and d) The phonon dispersion and density of states of the two predicted $La_2SiP_3$ phases.

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

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