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Correction: Accelerating materials discovery using integrated deep machine learning approaches

Weiye Xia,^a Ling Tang,^b Huaijun Sun,^c Chao Zhang,^d Kai-Ming Ho,^e
Gayatri Viswanathan,^{af} Kirill Kovnir^{af} and Cai-Zhuang Wang^{*ae}

Correction for 'Accelerating materials discovery using integrated deep machine learning approaches' by Weiye 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.

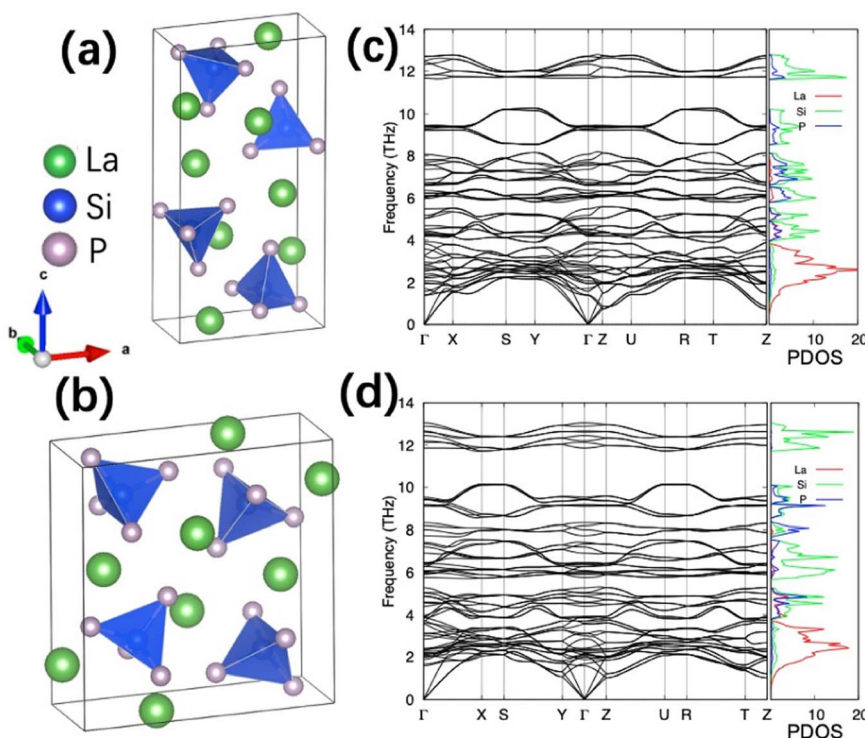


Fig. 5 (a and b) The structures of the predicted two La₂SiP₃ phases. The formation energies above the convex hull are 1 meV per atom and 33 meV per atom, respectively. (c and d) The phonon dispersion and density of states of the two predicted La₂SiP₃ phases.

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

^aAmes National Laboratory, U.S. Department of Energy, Ames, IA 50011, USA

^bDepartment of Applied Physics, College of Science, Zhejiang University of Technology, Hangzhou, 310023, China

^cJiyang College of Zhejiang Agriculture, Forestry University, Zhuji 311800, China

^dDepartment of Physics, Yantai University, Yantai 264005, China

^eDepartment of Physics and Astronomy, Iowa State University, Ames, IA 50011, USA

^fDepartment of Chemistry, Iowa State University, Ames, IA 50011, USA

