Journal of Materials Chemistry A

ROYAL SOCIETY OF CHEMISTRY

CORRECTION

Cite this: *J. Mater. Chem. A*, 2024, **12**, 16929

Correction: Graph theory and graph neural network assisted high-throughput crystal structure prediction and screening for energy conversion and storage

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DOI: 10.1039/d4ta90115k

rsc.li/materials-a

Correction for 'Graph theory and graph neural network assisted high-throughput crystal structure prediction and screening for energy conversion and storage' by Joshua Ojih *et al.*, *J. Mater. Chem. A*, 2024, **12**, 8502–8515, https://doi.org/10.1039/D3TA06190F.

On page 8511 in section 3.2 of the published article, the text "The single values of bonding and antibonding for each structure are obtained by performing integration over COHP curves for each atomic pair as evaluated by the LOBSTER package" should have included citations to six references. The missing references are listed below as ref. 1–6.

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

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