



Cite this: *New J. Chem.*, 2023, 47, 17608

DOI: 10.1039/d3nj90137h

rsc.li/njc

Correction: The theoretical prediction of the structural characteristics and SO₂ adsorption-sensing properties of pristine HfS₂ and TM-doped HfS₂ monolayers (TM = Ni, Pd, or Pt)

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Correction for 'The theoretical prediction of the structural characteristics and SO₂ adsorption-sensing properties of pristine HfS₂ and TM-doped HfS₂ monolayers (TM = Ni, Pd, or Pt)' by Tuan V. Vu et al., *New J. Chem.*, 2023, <https://doi.org/10.1039/d3nj01053h>.

The authors regret that one of the affiliations (affiliation b) was incorrectly shown in the original manuscript. The corrected list of affiliations is as shown here.

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

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