Correction: A high-entropy B₄(HfMo₂TaTi)C and SiC ceramic composite

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Correction for ‘A high-entropy B₄(HfMo₂TaTi)C and SiC ceramic composite’ by Hanzhu Zhang et al., Dalton Trans., 2019, DOI: 10.1039/c8dt04555k.

The authors regret that there is an error in writing the crystal structure in the article. The authors would like to address as follows:

The presented XRD and TEM results revealed a hexagonal crystal structure. The following analyses including the identification of the lattice parameters and the DFT calculation were based on a hexagonal lattice. Therefore, the HCP (hexagonal close-packed) structure mentioned in the article should be regarded as a hexagonal structure. The HCP term used in the introduction, where the article from Joshua Gild et al. was cited, should also be regarded as hexagonal AlB₂ structure.

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