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## Correction: Theoretical study of large-scale graphene on the Cu(111) surface using machine learning potential

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Correction for 'Theoretical study of large-scale graphene on the Cu(111) surface using machine learning potential' by Jingli Han *et al.*, *Phys. Chem. Chem. Phys.*, 2025, **27**, 17717–17729, <https://doi.org/10.1039/D5CP02042E>.

The authors regret that in the originating article, the affiliations of Yongpeng Yang and Jingli Han were incorrect. Their correct affiliation is Henan Institute of Advanced Technology, Zhengzhou University, Zhengzhou 450052, PR China, and School of Material and Chemical Engineering, Zhengzhou University of Light Industry, Zhengzhou, 450001, China, respectively.

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

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