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CORRECTION

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# Correction: Enhanced 5f-δ bonding in $[U(C_7H_7)_2]^-$: C K-edge XAS, magnetism, and *ab initio* calculations

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Correction for 'Enhanced 5f-δ bonding in $[U(C_7H_7)_2]^-$: C K-edge XAS, magnetism, and *ab initio* calculations' by Yusen Qiao *et al.*, *Chem. Commun.*, 2021, **57**, 9562–9565, DOI: 10.1039/D1CC03414F.

The authors regret that two references to the multi-configurational pair-density functional theory methods used in the wavefunction calculations did not appear in the main article. The references were originally included in the Methods section of the Electronic Supplementary Information as ref. 17*a* and *b*, but should also have appeared in the second paragraph on page 9564. The references are listed below as ref. 1 and 2.

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

## References

1 G. Li Manni, R. K. Carlson, S. Luo, D. Ma, J. Olsen, D. G. Truhlar and L. Gagliardi, *J. Chem. Theory Comput.*, 2014, **10**, 3669–3680.
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