

Correction to “Polarizable Density Embedding: A New QM/QM/MM-based Computational Strategy”

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We inadvertently reported the experimental vacuum excitation energy of the permanganate ion as 2.19 eV where it should be 2.09 eV.¹ The relevant sentence in the first column on page 5350 should therefore read:

“The calculated vacuum excitation energy ($E_{\text{exc}}^{\text{vac}}$) is 1.956 eV, in good agreement with the experimental value of 2.09 eV.⁸³”

■ REFERENCES

(1) Houmøller, J.; Kaufman, S. H.; Stöckel, K.; Tribedi, L. C.; Brøndsted Nielsen, S.; Weber, J. M. On the Photoabsorption by Permanganate Ions in Vacuo and the Role of a Single Water Molecule. New Experimental Benchmarks for Electronic Structure Theory. *ChemPhysChem* **2013**, *14*, 1133–1137.