

CORRECTION

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Correction: Conceptual design of tetraazaporphyrin- and subtetraazaporphyrin-based functional nanocarbon materials: electronic structures, topologies, optical properties, and methane storage capacities

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Correction for 'Conceptual design of tetraazaporphyrin- and subtetraazaporphyrin-based functional nanocarbon materials: electronic structures, topologies, optical properties, and methane storage capacities' by Rodion V. Belosludov *et al.*, *Phys. Chem. Chem. Phys.*, 2016, **18**, 13503–13518.

In the published version of this paper, a small number of TDDFT-predicted excited states were used for nanocages 3 and 4 in Fig. 11. The correct figure with a larger number of TDDFT-predicted excited states for nanocages 3–4 is provided below. The results and conclusions described in the paper are unchanged.

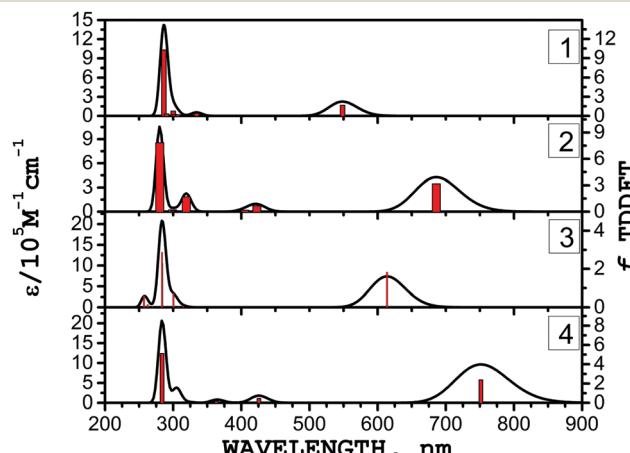


Fig. 11 TDDFT-predicted UV-vis spectra for nanocages 1–4.

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

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