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Correction: Mechanistic insights into Ag⁺ induced size-growth from [Au₆(DPPP)₄]²⁺ to [Au₇(DPPP)₄]²⁺ clusters

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Correction for 'Mechanistic insights into Ag⁺ induced size-growth from [Au₆(DPPP)₄]²⁺ to [Au₇(DPPP)₄]²⁺ clusters' by Ying Lv *et al.*, *Nanoscale Adv.*, 2022, 4, 3737–3744, <https://doi.org/10.1039/D2NA00301E>.

The authors regret that Fig. 2 and Fig. 3 in the published paper are identical. The correct Fig. 2 is given here.

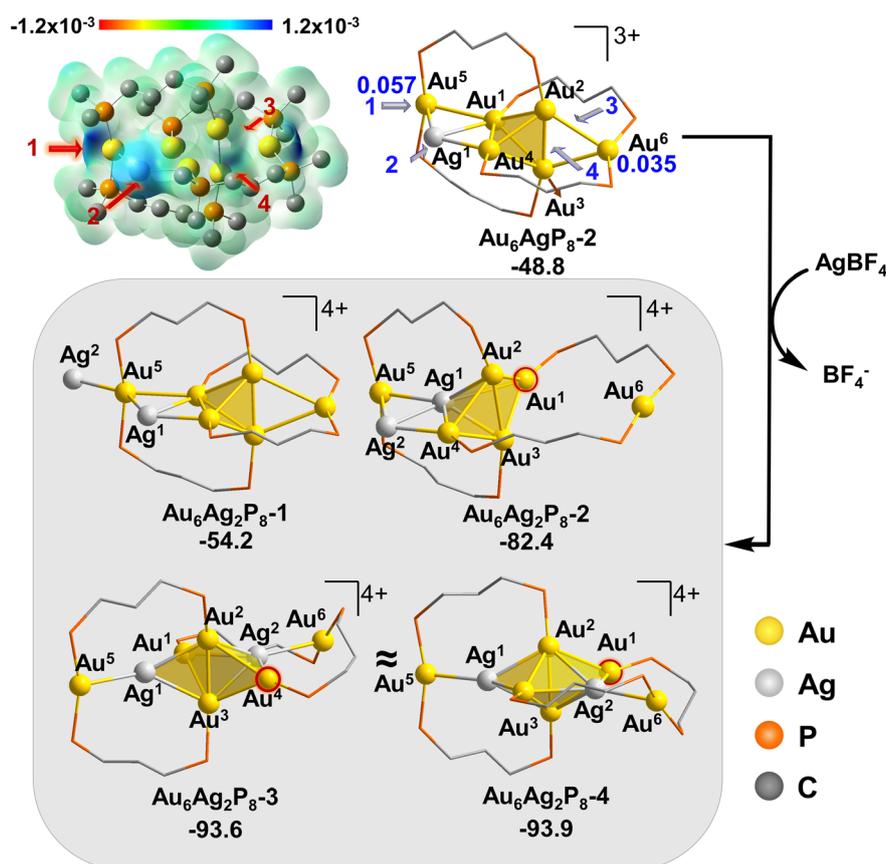


Fig. 2 The isosurface of the f^- for $\text{Au}_6\text{AgP}_8-2$, using the width of Gaussian function of 0.01 au and the energy (in kcal mol^{-1}) and structural changes for the doping of second Ag^+ into $\text{Au}_6\text{AgP}_8-2$. The Hirshfeld charge of $\text{Au}^{5/6}$ in starting structure is given in blue and bold.

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

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