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## Correction: Mechanistic insights into $\text{Ag}^+$ induced size-growth from $[\text{Au}_6(\text{DPPP})_4]^{2+}$ to $[\text{Au}_7(\text{DPPP})_4]^{2+}$ clusters

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Correction for 'Mechanistic insights into  $\text{Ag}^+$  induced size-growth from  $[\text{Au}_6(\text{DPPP})_4]^{2+}$  to  $[\text{Au}_7(\text{DPPP})_4]^{2+}$  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.

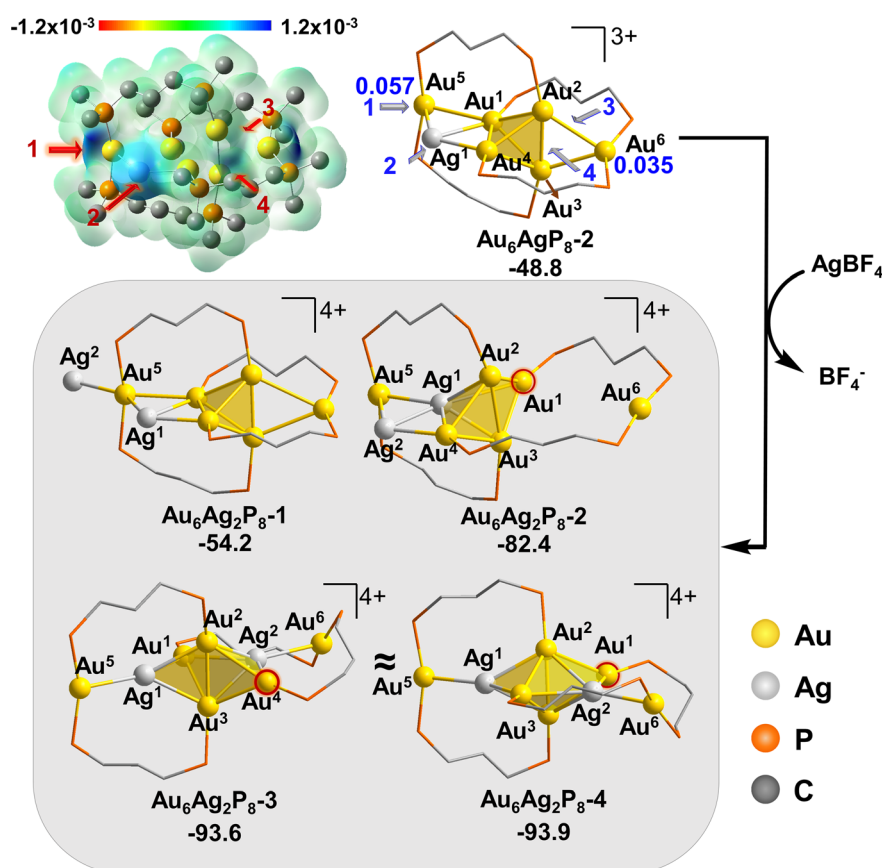


Fig. 2 The isosurface of the  $f^-$  for  $\text{Au}_6\text{AgP}_8-2$ , using the width of Gaussian function of 0.01 au and the energy (in  $\text{kcal mol}^{-1}$ ) and structural changes for the doping of second  $\text{Ag}^+$  into  $\text{Au}_6\text{AgP}_8-2$ . The Hirshfeld charge of  $\text{Au}^{5/6}$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