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CORRECTION

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# Correction: Mechanistic insights into $Ag^+$ induced size-growth from $[Au_6(DPPP)_4]^{2+}$ to $[Au_7(DPPP)_4]^{2+}$ clusters

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Correction for 'Mechanistic insights into $Ag^+$ induced size-growth from $[Au_6(DPPP)_4]^{2+}$ to $[Au_7(DPPP)_4]^{2+}$ 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.

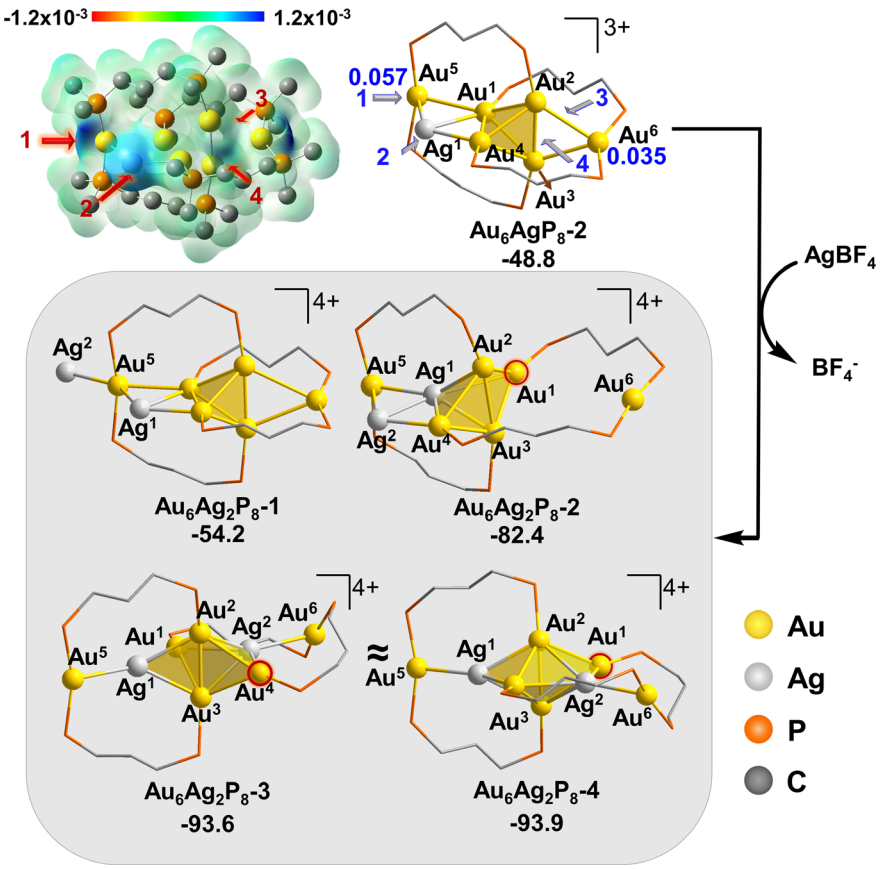


Fig. 2 The isosurface of the $f^-$ for $Au_6AgP_8$-2, using the width of Gaussian function of 0.01 au and the energy (in kcal $mol^{-1}$) and structural changes for the doping of second $Ag^+$ into $Au_6AgP_8$-2. The Hirshfeld charge of $Au^{5/6}$ in starting structure is given in blue and bold.

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

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