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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.

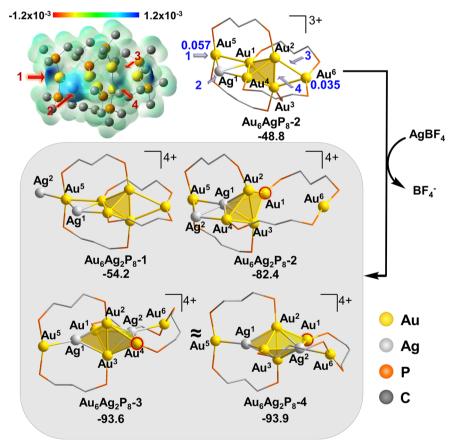


Fig. 2 The isosurface of the  $f^-$  for  $Au_6AgP_8-2$ , using the width of Gaussian function of 0.01 au and the energy (in kcal mol<sup>-1</sup>) and structural changes for the doping of second  $Ag^+$  into  $Au_6AgP_8-2$ . The Hirshfeld charge of  $Au^{5/6}$  in starting structure is given in blue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

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

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