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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

Ying Lv,[a] Xiaohang Wu,[a] Shuping He[a] and Haizhu Yu*[ab]

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.

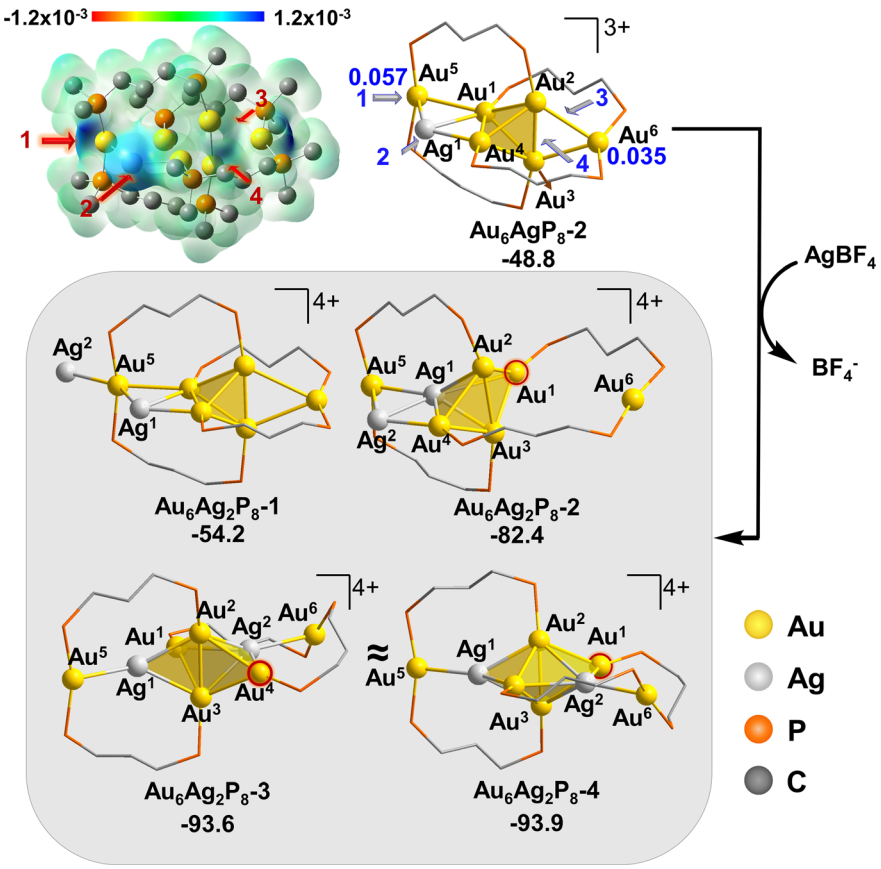

Fig. 2 The isosurface of the f^- for Au_6AgP_8-2, using the width of Gaussian function of 0.01 au and the energy (in kcal mol^{-1}) and structural changes for the doping of second Ag^+ into Au_6AgP_8-2. The Hirshfeld charge of $Au^{5/6}$ in starting structure is given in blue and bold.

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

[a]*Department of Chemistry, Centre for Atomic Engineering of Advanced Materials, Anhui Province Key Laboratory of Chemistry for Inorganic/Organic Hybrid Functionalized Materials, Key Laboratory of Structure and Functional Regulation of Hybrid Materials of Ministry of Education, Anhui University, Hefei 230601, Anhui, P. R. China. E-mail: yuhaizhu@ahu.edu.cn*

[b]*Institute of Energy, Hefei Comprehensive National Science Center, Hefei, Anhui, 230031, P. R. China*