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CORRECTION

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Correction: Sterically controlled 5-*exo*-dig cyclization enables synthesis of non-benzenoid polycyclic aromatic hydrocarbons with intriguing (anti)aromaticity and diradical properties

Liangliang Chen,*[a] Zhichun Shangguan,[b] Tianyu Shi,[c] Liyuan Qin,[c] Yiyun Zeng,[a] Qingqiu Zhu,[a] Jin Chen,[a] Junhong Liang,[a] Wentao Miao,[a] Yurong He,[a] Xiaosong Qiu,[a] Xunchang Wang,[a] Deqing Zhang[c] and Renqiang Yang*[a]

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Correction for 'Sterically controlled 5-*exo*-dig cyclization enables synthesis of non-benzenoid polycyclic aromatic hydrocarbons with intriguing (anti)aromaticity and diradical properties' by Liangliang Chen *et al.*, *Chem. Sci.*, 2026, https://doi.org/10.1039/d6sc00121a.

The authors regret that due to an error in figure assembly part of Fig. 2b was obscured in the final version of the manuscript. The corrected Fig. 2 is shown below. This does not alter the interpretations or conclusions of the paper.

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

[a]*Key Laboratory of Flexible Optoelectronic Materials and Technology (Ministry of Education), School of Optoelectronic Materials & Technology, Jianghan University, Wuhan, 430056, China. E-mail: chenliangliang@jhun.edu.cn; yangrq@jhun.edu.cn*

[b]*College of Chemistry and Materials Engineering, Wenzhou University, Key Lab of Biohealth Materials and Chemistry of Wenzhou, Wenzhou, 325027, China*

[c]*Beijing National Laboratory for Molecular Sciences, Institute of Chemistry, Chinese Academy of Sciences, Beijing, 100190, China*

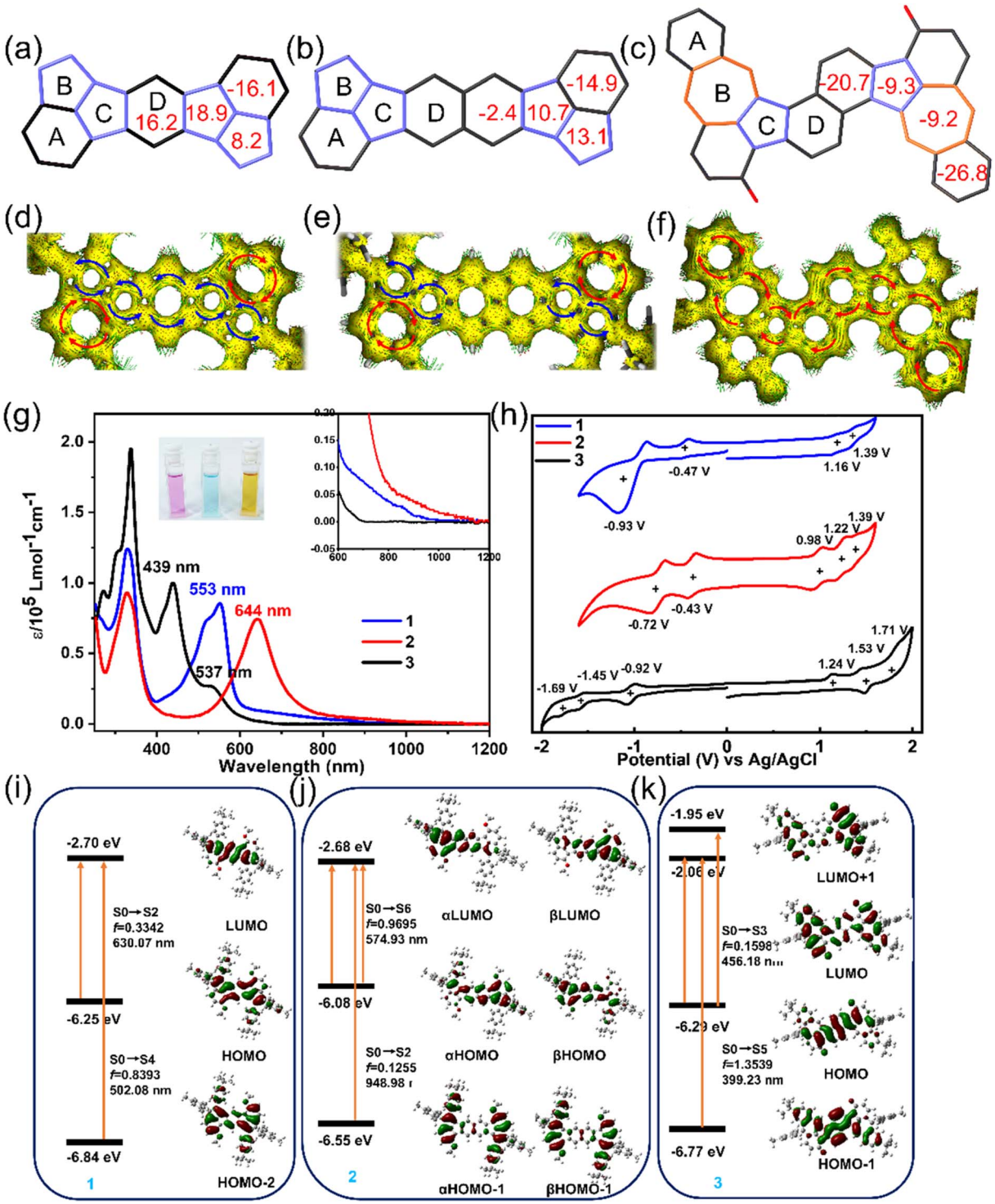

Fig. 2 The calculated NICS(1)$_{zz}$ values at 1 Å of *Z* axis of (a) **1**, (b) **2** and (c) **3**. The calculated ACID plots of (d) **1**, (e) **2** and (f) **3**. The clockwise ring current represents aromaticity and the counter-clockwise ring current represents antiaromaticity. (g) The UV-vis absorption spectra of **1**, **2** and **3** (10^{-5} mol L^{-1} in DCM solution). (h) Cyclic voltammograms (CV) curves of **1**, **2** and **3**. The calculated absorption transitions of (i) **1**, (j) **2** and (k) **3** based on TD-DFT calculations. (i) and (k) were based on M06-2X/6-311G(d,p) level of theory with closed-shell single state. (j) was based on UM06-2X/6-311G(d,p) level of theory with open-shell singlet state.