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Addition/Correction

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Correction to Polyradicals of Polycyclic Aromatic Hydrocarbons as Finite Size Models of Graphene: Highly Open-Shell Nature, Symmetry Breaking, and Enhanced-Edge Electron Density

Amarjeet Yadav and P. C. Mishra*

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Typographical errors need correction as follows. There is no change in the discussion or any other part of the paper.

(1) In Table 1, for molecule 4R4, multiplicity quartet, DFT Functional B3LYP, the total energy should be −1839.522 00 au instead of −1151.522 00 au. For the same system and multiplicity, with DFT functional M06-2X, the total energy should be −1838.808 34 au instead of −1839.808 34 au.

(2) In Table 2, for radical 1R1−5H (1−5), multiplicity quartet, DFT Functional B3LYP, the total energy should be −228.858 71 au instead of −229.858 71 au.

(3) In Table 3, for radical 2R2−2H (1−2), multiplicity quintet, DFT Functional M06-2X, the total energy should be −614.071 43 au instead of −614.151 58 au.

For radical 2R2−7H (1−7), multiplicity quartet, DFT Functional M06-2X, the total energy should be −610.802 99 au instead of −612.802 99 au.

(4) In Table 4: System 2R2, in the column of singlet: Total energy is negative (−609.01175 au).

(5) In Section 3.9, Spin Density Distribution, last sentence: Figure 4a,b should read Figure 5a,b.

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