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## Correction: Tetraaryl pyrenes: photophysical properties, computational studies, crystal structures, and application in OLEDs

Tarek H. El-Assaad,<sup>a</sup> Manuel Auer,<sup>b</sup> Raul Castañeda,<sup>c</sup> Kassem M. Hallal,<sup>a</sup> Fadi M. Jradi,<sup>a</sup> Lorenzo Mosca,<sup>d</sup> Rony S. Khnayzer,<sup>e</sup> Digambara Patra,<sup>a</sup> Tatiana V. Timofeeva,<sup>c</sup> Jean-Luc Brédas,<sup>f</sup> Emil J. W. List-Kratochvil,<sup>bg</sup> Brigitte Wex<sup>\*e</sup> and Bilal R. Kaafarani<sup>\*a</sup>

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Correction for 'Tetraaryl pyrenes: photophysical properties, computational studies, crystal structures, and application in OLEDs' by Tarek H. El-Assaad *et al.*, *J. Mater. Chem. C*, 2016, DOI: 10.1039/c5tc02849c.

The *x* coordinate for compound 3 in Table 7 is missing and only the *y* coordinate is given. The correct coordinates are 0.148, 0.243 and the correct version of this table is as follows:

**Table 7** Electroluminescent characteristics of the investigated compounds in a single-layer geometry

| Compound | $V_{on}^a$ [V] | $L_{max}$ [cd m <sup>-2</sup> ] | $H^b$ [cd A <sup>-1</sup> ] | CIE1931 [x, y] |
|----------|----------------|---------------------------------|-----------------------------|----------------|
| 2        | 2.8            | 13 542                          | 2.0000                      | 0.163, 0.200   |
| 3        | 2.9            | 6902                            | 2.6000                      | 0.148, 0.243   |
| 4        | 2.9            | 85                              | 0.0050                      | 0.148, 0.244   |
| 7        | 8.6            | 7                               | 0.0039                      | 0.153, 0.124   |

<sup>a</sup> Voltage at a luminance of 1 cd m<sup>-2</sup>. <sup>b</sup> Value of the maximum efficiency.

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

<sup>a</sup> Department of Chemistry, American University of Beirut, Beirut 1107-2020, Lebanon. E-mail: bilal.kaafarani@aub.edu.lb

<sup>b</sup> NanoTecCenter Weiz Forschungsgesellschaft mbH, Franz-Pichler-Straße 32, A-8160 Weiz, Austria

<sup>c</sup> Department of Biology & Chemistry, New Mexico Highlands University, Las Vegas, NM 87701, USA

<sup>d</sup> Department of Chemistry and Center for Photochemical Sciences, Bowling Green State University, Bowling Green, Ohio, 43403, USA

<sup>e</sup> Department of Natural Sciences, Lebanese American University, Byblos, Lebanon. E-mail: brigitte.wex@lau.edu.lb

<sup>f</sup> Solar & Photovoltaics Engineering Research Center, Physical Science and Engineering Division, King Abdullah University of Science & Technology, Thuwal 23955-6900, Kingdom of Saudi Arabia

<sup>g</sup> Institute of Solid State Physics, Graz University of Technology, A-8010 Graz, Austria

