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

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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 [ <i>x</i> , <i>y</i> ] |
|----------|----------------|---------------------------------|-----------------------------|---------------------------------|
| 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.

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