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4, 3059**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>  
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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.

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