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CORRECTION

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Correction: A computational study of tri-*s*-triazine-based molecules as ambipolar host materials for phosphorescent blue emitters: effective geometric and electronic tuning

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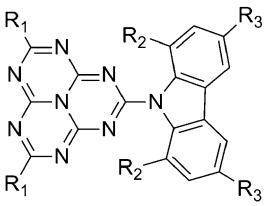
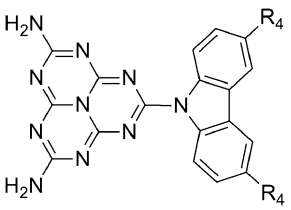
Correction for 'A computational study of tri-*s*-triazine-based molecules as ambipolar host materials for phosphorescent blue emitters: effective geometric and electronic tuning' by Jing Li *et al.*, *J. Mater. Chem. C*, 2015, **3**, 4859.

There were errors in the structures depicted in Schemes 1 and 2. These have been revised and the correct versions of Schemes 1 and 2 are:



Scheme 1 The structure of designed tri-*s*-triazine derivatives (R = F, Cl, Br, CN, NO_2, COOH, NH_2, OH, OPh, SPh, methyl, ethyl, propyl, i-propyl, butyl, i-butyl and *t*-butyl).



R_4			
I1	F,	I6	i-propyl,
I2	NH_2,	I7	butyl,
I3	CH_3,	I8	i-butyl,
I4	ethyl,	I9	t-butyl,
I5	propyl,		

Scheme 2 Sketch map of the structures of series I.

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

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