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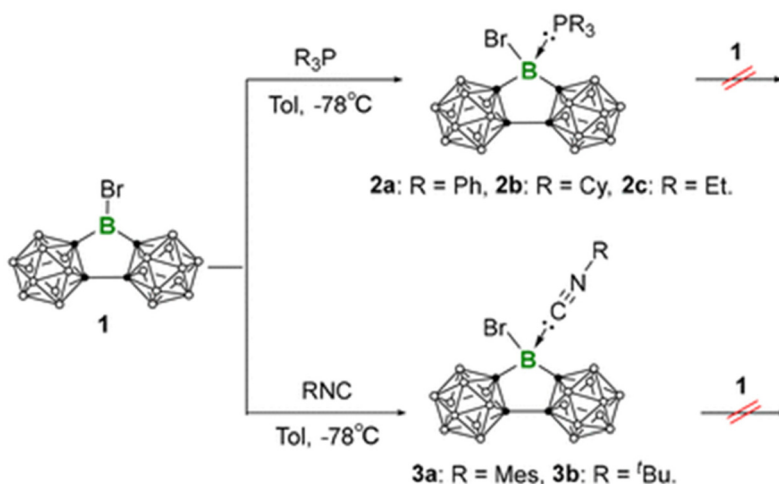
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Correction: Reactivity study of Lewis superacidic carborane-based analogue of 9-bromo-9-borafluorene towards Lewis bases

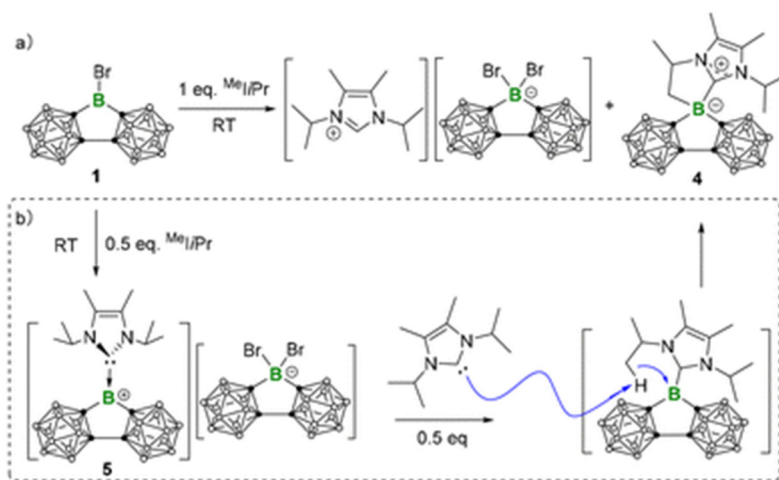
Libo Xiang,^{a,b} Alexander Matler,^{a,b} Leibo Tan^{a,b} and Qing Ye^{*a,b}Correction for 'Reactivity study of Lewis superacidic carborane-based analogue of 9-bromo-9-borafluorene towards Lewis bases' by Libo Xiang *et al.*, *Dalton Trans.*, 2024, **53**, 11655–11658, <https://doi.org/10.1039/D4DT01615G>.

The authors regret that there was a mismatch between the images and captions for Schemes 1 and 2 and Fig. 2 and 3 in the original article. The correct figures are shown here.



Scheme 1 Synthesis of 2–3.

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Scheme 2 (a) Synthesis of 4; (b) synthesis of 4 via 5.

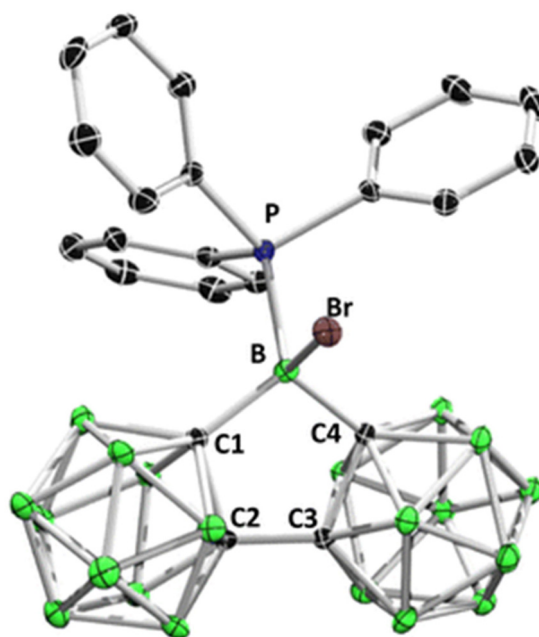


Fig. 2 Single crystal structure of 2a. Hydrogen atoms were omitted for clarity. Thermal ellipsoids are drawn at the 50% probability level. Selected bond lengths [Å]: B1–P 2.032(2), B1–C1 1.671(3), B1–C4 1.667(3), C1–C2 1.667(3), C3–C4 1.670(3), B–Br 2.036(2).



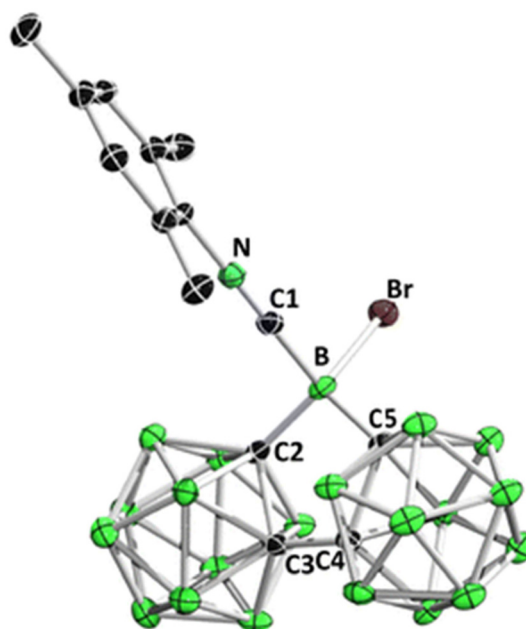


Fig. 3 Single crystal structure of **3a**. Hydrogen atoms were omitted for clarity. Thermal ellipsoids are drawn at the 50% probability level. Selected bond lengths [Å]: B–C1 1.605(3), B–C2 1.648(3), B–C5 1.643(3), C2–C3 1.656(3), C4–C5 1.653(3), C–N1 1.143(3), B–Br 2.009(2).

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

