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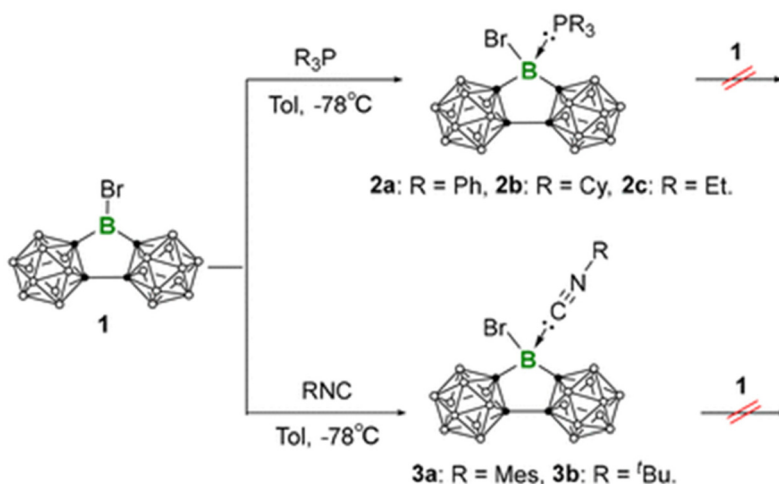
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## Correction: Reactivity study of Lewis superacidic carborane-based analogue of 9-bromo-9-borafluorene towards Lewis bases

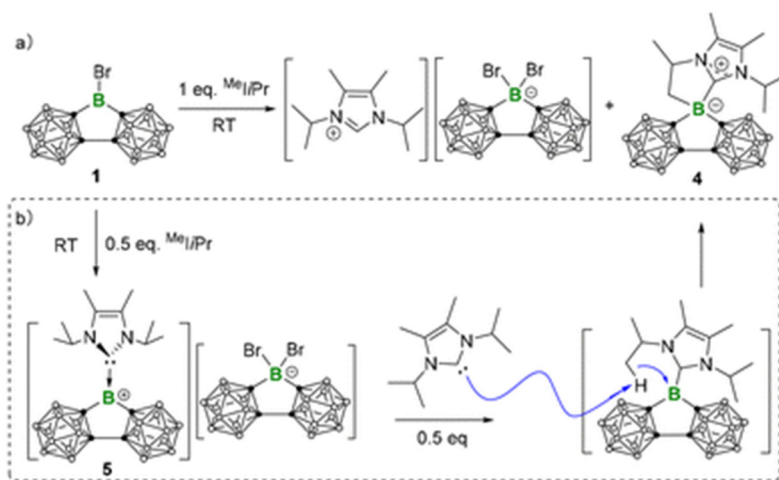
Libo Xiang,<sup>a,b</sup> Alexander Matler,<sup>a,b</sup> Leibo Tan<sup>a,b</sup> and Qing Ye<sup>\*a,b</sup>Correction for 'Reactivity study of Lewis superacidic carborane-based analogue of 9-bromo-9-borafluorene towards Lewis bases' by Libo Xiang *et al.*, *Dalton Trans.*, 2024, **53**, 11655–11658,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

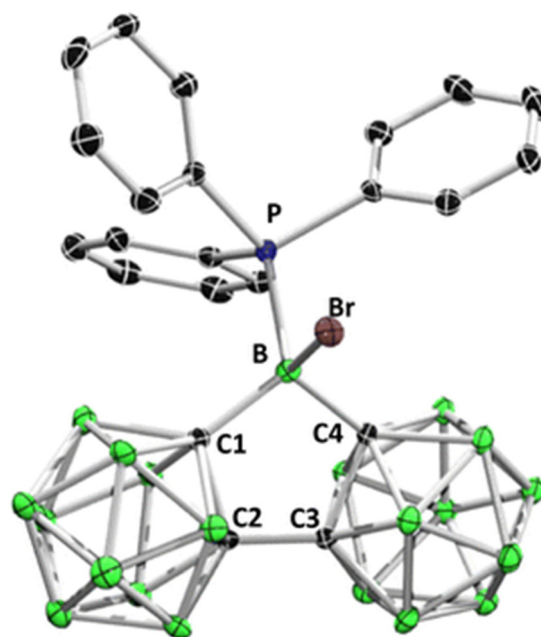


Scheme 1 Synthesis of 2–3.

<sup>a</sup>Institute for Inorganic Chemistry, Julius-Maximilians-Universität Würzburg, Am Hubland, 97074 Würzburg, Germany. E-mail: qing.ye@uni-wuerzburg.de<sup>b</sup>Institute for Sustainable Chemistry & Catalysis with Boron, Julius-Maximilians-Universität Würzburg, Am Hubland, 97074 Würzburg, Germany

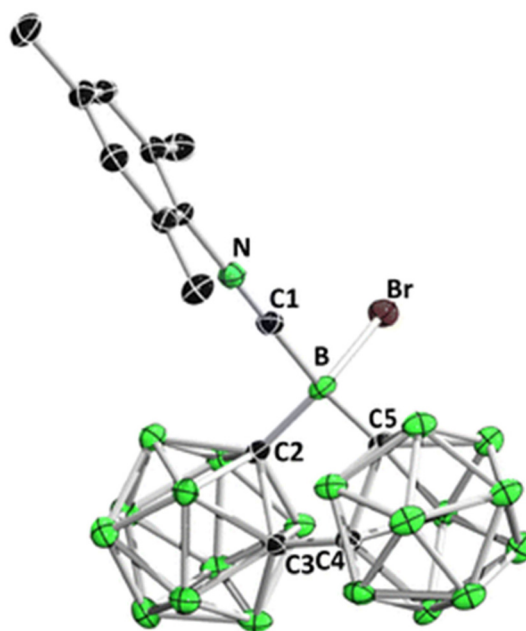


**Scheme 2** (a) Synthesis of 4; (b) synthesis of 4 via 5.



**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]: 2, 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.

