

# Correction to Catching the First Oligomerization Event in the Catalytic Formation of Polyaminoboranes: H<sub>3</sub>B·NMeHBH<sub>2</sub>·NMeH<sub>2</sub> Bound to Iridium

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# Supporting Information

Page 11077 and caption of Figure 3. Reanalysis of the single crystal X-ray diffraction data corresponding to compound 2, H<sub>3</sub>B·NMeHBH<sub>2</sub>·NMeH<sub>2</sub>, has indicated an incorrect assignment of C1 and B1, which are in fact inverted. As a consequence, the discussion on page 11077 and the caption of Figure 3 are incorrect. The molecule is found in fact in a gauche conformation, as for its Ir-bound analogue, and the N1–B1 bond length measures at 1.590 Å, which is in close agreement to the value of 1.576 Å for the analogous bond in the metal-bound species. The corrected value for N1–B1 is also in close agreement with the B–N bond distance in the related amine—borane, MeNH<sub>2</sub>·BH<sub>3</sub>, of 1.594 Å.<sup>1</sup>

The previously reported value for N1–B1 in 2 of 1.487 Å corresponds in fact to the N1–C1 bond distance and is in close agreement with the value of 1.483 Å attributed to the N2–C2 bond within the same molecule. Correcting this misassignment also reduces the R factor of the structure from to 0.071 to 0.035, as expected. The authors apologize for this mistake.

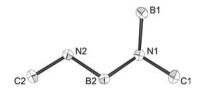


Figure 1. Corrected molecular structure of  $H_3B \cdot NMeHBH_2 \cdot NMeH_2$ .

# ASSOCIATED CONTENT

### S Supporting Information

Full experimental data for the syntheses; crystallographic data for 2, 4, and 5, including atomic positional and thermal parameters (CIF). This material is available free of charge via the Internet at http://pubs.acs.org.

## REFERENCES

(1) Aldridge, A.; Downs, A. J.; Tang, C. Y.; Parsons, S.; Clarke, M. C.; Johnstone, R. D. L.; Robertson, H. E.; Rankin, D. W. H.; Wann, D. A. J. Am. Chem. Soc. **2009**, *131*, 2231–2243.

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