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CORRECTION

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Correction: A binary catalyst system of a cationic Ru–CNC pincer complex with an alkali metal salt for selective hydroboration of carbon dioxide

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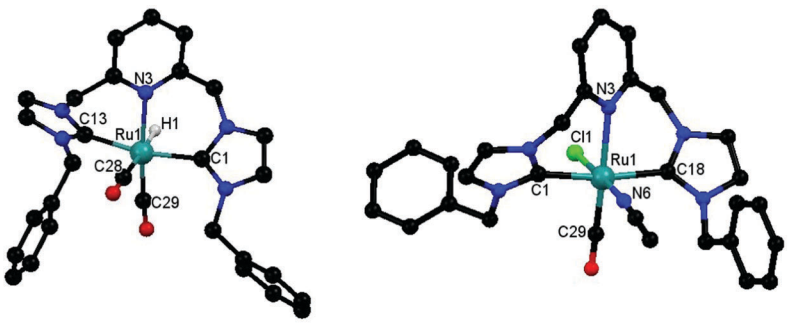
Correction for 'A binary catalyst system of a cationic Ru–CNC pincer complex with an alkali metal salt for selective hydroboration of carbon dioxide' by Chee Koon Ng *et al.*, *Chem. Commun.*, 2016, **52**, 11842–11845.

In the original article, compound **14** was reported as a PF_6^- salt. However, the presented structural data were representative of a PF_4^- salt due to unsatisfactory crystal structure refinement.

This compound has therefore been re-crystallized and the crystallographic analysis repeated. The new diffraction data have been correctly and satisfactorily refined as a PF_6^- salt. We provide below a new image and caption for Fig. 1, while corrected ESI files have been uploaded in place of those originally supplied.



12 **14**

Fig. 1 Single-crystal X-ray molecular structures of **12** and **14** (hydrogen atoms, with the exception of the hydrido ligand, solvent molecules and PF_6 anions have been omitted for clarity). Selected bond lengths (Å) and bond angles (deg): (**12**) Ru1–N3, 2.197(4); Ru1–C1, 2.087(5); Ru1–C13, 2.080(5); Ru1–C28, 1.972(5); Ru1–C29, 1.853(5); and C13–Ru1–C1, 168.78(19). (**14**) Ru1–N3, 2.228(2); Ru1–C1, 2.099(3); Ru1–C18, 2.090(3); Ru1–C29, 1.838(3); Ru1–Cl1, 2.4175(10); Ru1–N6, 2.031(3); and C1–Ru1–C18, 170.99(11).

The associated carbene bond data description on page 11843 should therefore also be amended to 'The two CNC carbene ligands are *trans* to each other (C_{NHC}–Ru–C_{NHC} = 168.8°–171.0°) and the Ru–C_{NHC} bond lengths of 2.08–2.10 Å are consistent with literature values.'

We are thankful to the CCDC for kindly notifying us of this irregularity.

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

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