

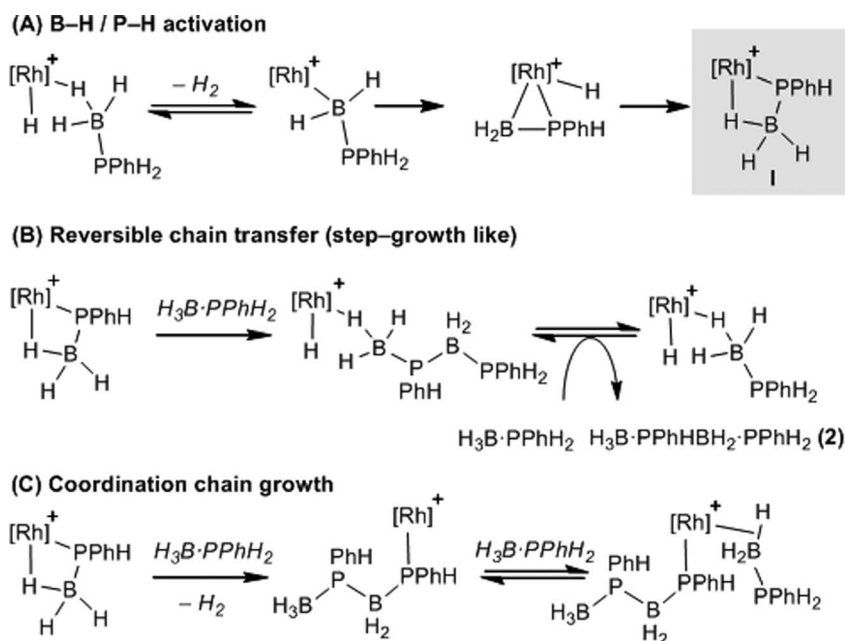
## CORRECTION

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click for updatesCite this: *Chem. Sci.*, 2016, 7, 2437**Correction: Dehydrocoupling of phosphine–boranes using the  $[\text{RhCp}^*\text{Me}(\text{PMe}_3)(\text{CH}_2\text{Cl}_2)][\text{BAR}^{\text{F}}_4]$  precatalyst: stoichiometric and catalytic studies**Thomas N. Hooper,<sup>a</sup> Andrew S. Weller,<sup>\*a</sup> Nicholas A. Beattie<sup>b</sup>  
and Stuart A. Macgregor<sup>\*b</sup>

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[www.rsc.org/chemicalscience](http://www.rsc.org/chemicalscience)Correction for 'Dehydrocoupling of phosphine–boranes using the  $[\text{RhCp}^*\text{Me}(\text{PMe}_3)(\text{CH}_2\text{Cl}_2)][\text{BAR}^{\text{F}}_4]$  precatalyst: stoichiometric and catalytic studies' by Thomas N. Hooper *et al.*, *Chem. Sci.*, 2016, DOI: 10.1039/c5sc04150c.

The authors regret that in the original article the structures of two of the compounds in Scheme 12 contained errors. A corrected version of Scheme 12 is presented herein, where a  $\text{PMe}_3$  ligand has been removed from the third compound in part A and a hydrogen atom has been removed from the  $\text{PPhH}$  group of the first compound in part C.



Scheme 12 Suggested mechanisms for dehydropolymerization.  $[\text{Rh}] = \text{Rh}(\text{PR}_3)\text{Cp}^*$  ( $\text{PR}_3 = \text{PMe}_3$  or  $\text{PPhH}_2$ ).

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

<sup>a</sup>Department of Chemistry, Chemistry Research Laboratories, University of Oxford, Mansfield Road, Oxford, OX1 3TA, UK. E-mail: [andrew.weller@chem.ox.ac.uk](mailto:andrew.weller@chem.ox.ac.uk)

<sup>b</sup>Institute of Chemical Sciences, Heriot Watt University, Edinburgh, EH14 4AS, UK. E-mail: [S.A.Macgregor@hw.ac.uk](mailto:S.A.Macgregor@hw.ac.uk)