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Addition/Correction

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Correction to Rhodium(III)-Catalyzed Isoquinolone Synthesis: The N−O Bond as a Handle for C−N Bond Formation and Catalyst Turnover

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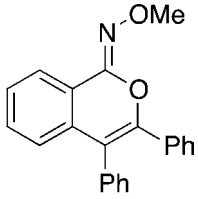
Supporting Information

Following a recent report from Huang et al.,[1] we further investigated the structure of **3a**. In accordance with the proposed structure in Huang's report, X-ray crystallographic analysis confirmed that **3a** is not an isoquinolone but rather the heterocycle resulting from C−O bond reductive elimination (see structure below). This correction does not alter the outcome of our work as it is focused on the synthesis of isoquinolones where the N−O bond has undergone cleavage.



3a

The Supporting Information was modified to match the correct structure of **3a**. An electronic crystallographic information file (CIF) was also made available.

■ ASSOCIATED CONTENT

Supporting Information

Detailed experimental procedures and characterization data for all new compounds. This material is available free of charge via the Internet at http://pubs.acs.org

■ REFERENCES

(1) Zhong, H.; Yang, D.; Wang, S.; Huang, J. *Chem. Commun.* **2012**, *48*, 3236.

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