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Addition/Correction

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Correction to "Nickel-Catalyzed Asymmetric Reductive Cross-Coupling To Access 1,1-Diarylalkanes"

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Supporting Information

Pages 5684 and 5685, Table of Contents, and Supporting Information. The stereochemistry of **L1**, depicted as the (*S*,*S*)-enantiomer in Figure 1, Table 1, the TOC graphic (identical to Figure 1), and the Supporting Information of the original publication, was incorrect. (*R*,*R*)-**L1** was used in this study. The stereochemistry of (*R*,*R*)-**L1** has been confirmed by single-crystal X-ray diffraction; the X-ray diffraction data and CIF file for (*R*,*R*)-**L1** have been added to the Supporting Information. The corrected TOC graphic/Figure 1 is shown here.

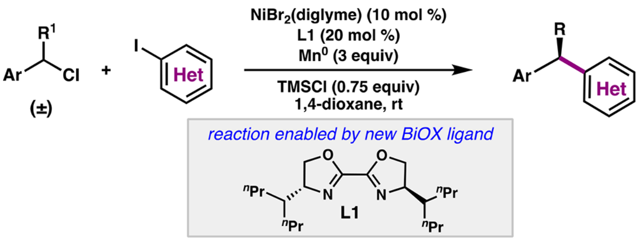
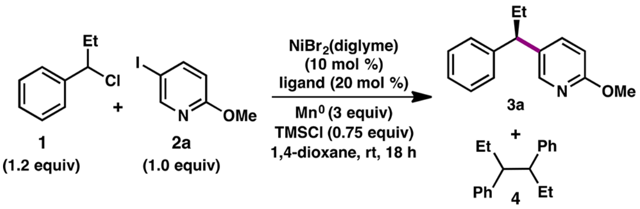


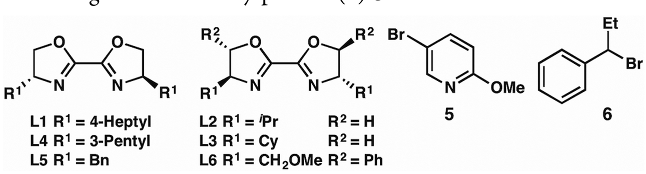
Figure 1. Ni-catalyzed enantioselective reductive cross-coupling to prepare diarylalkanes.

Table 1. Optimization of Reaction Conditions[a]



entry[a]	ligand	deviation from standard conditions	yield 3a (%)[b]	yield 4 (%)[b]	ee 3a (%)[c]
1	**L1**	none	84	8	90
2	**L2**	–	22	30	68[d]
3	**L3**	–	64	21	75[d]
4	**L4**	–	74	20	80
5	**L5**	–	4	0	60
6	**L6**	–	31	9	86[d]
7	**L1**	Zn^0 instead of Mn^0	0	26	–
8	**L1**	TDAE instead of Mn^0	3	0	66
9	**L1**	TFA instead of TMSCl	0	25	–
10	**L1**	DMA instead of dioxane	14	13	67
11	**L1**	**5** instead of **2a**	72	24	89
12	**L1**	**6** instead of **1**	8	37	81

[a]Reactions conducted under N_2 on 0.05 mmol scale for 18 h. [b]Determined by ^{1}H NMR versus an internal standard. [c]Determined by SFC using chiral stationary phase. [d](*S*)-**3a** is formed.



(*R*,*R*)-**L4** and (*R*,*R*)-**L5** were also used in Table 1 and incorrectly depicted as (*S*,*S*)-**L4** and (*S*,*S*)-**L5** in the original publication. To reflect that different enantiomeric series of catalysts were used, Table 1 has been updated to indicate that entries 2, 3, and 6 produce (*S*)-**3a**.

This correction does not change the stereochemical assignment of the diarylalkane products, or the conclusions of the Communication. The stereochemistry of the products was assigned by obtaining an X-ray structure of diarylalkane **3k**, and the rest of the compounds were assigned by analogy.

■ ASSOCIATED CONTENT

Supporting Information

The Supporting Information is available free of charge on the ACS Publications website at DOI: 10.1021/jacs.8b05247.

Detailed experimental procedures, compound characterization data (corrected) (PDF)
^{1}H and ^{13}C NMR spectra (corrected) (PDF)
Crystallographic data for (*R*,*R*)-**L1** (CIF)