JOC *Additions and Corrections*

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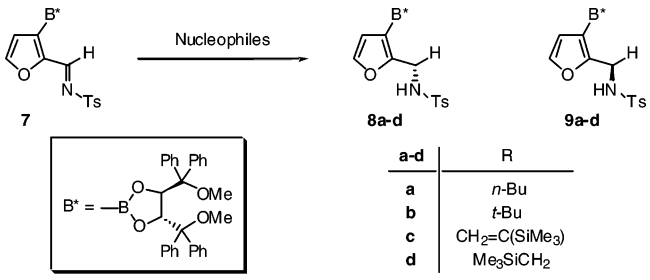
Ho-Kee Yim and Henry N. C. Wong*. Diastereoselective Addition Reactions of Furyl Sulfonylimine Using Chiral Boronates as Auxiliary: Application to the Enantioselective Synthesis of 2,3-Disubstituted Furyl Sulfonylamides.

Page 2893, left column. The text should read, "It was found that diastereomers **9a**−**d** were more polar than diastereomers **8a**−**d**", instead of "It was found that the *R* diastereomers **8a**−**d** were more polar than *S* diastereomers **9a**−**d**".

Page 2893, Table 1. Table 1 should have appeared as shown below. The description in the Supporting Information has been changed accordingly.

TABLE 1. Addition Reactions of Various Nucleophiles to 7[a]

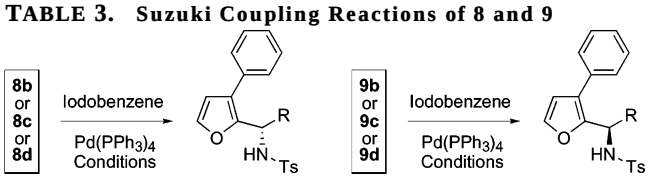


entry	nucleophile	conditions (solvent, *T* (°C))	products		yield[b] (%)	de[c] (%)
1	*n*-BuLi	THF, −78	**8a**	**9a**	45	33 (**9a**)
2	*n*-BuLi	DME, −60	**8a**	**9a**	58	33 (**9a**)
3	*n*-BuLi	PhMe, −78	**8a**	**9a**	64	33 (**9a**)
4	*t*-BuLi	THF, −78	**8b**	**9b**	35	33 (**9b**)
5	Me_3SiCH_2MgCl	THF, −40	**8c**	**9c**	94	20 (**9c**)
6	Me_3SiCH_2MgCl	DME, −30	**8c**	**9c**	81	83 (**9c**)
7	Me_3SiCH_2MgCl	THF/Et_2O 1:4	**8c**	**9c**	88	80 (**9c**)
8	Me_3SiCH_2MgCl	THF/Et_2O 1:8	**8c**	**9c**	85	86 (**9c**)
9	Me_3SiCH_2MgCl	DCM, −40	**8c**	**9c**	82	25 (**9c**)
10	Me_3SiCH_2MgCl	PhH, −40	**8c**	**9c**	78	56 (**9c**)
11	CH_2=C($SiMe_3$)MgBr	THF, −40	**8d**	**9d**	90	25 (**9d**)
12	CH_2=C($SiMe_3$)MgBr	DME, −30	**8d**	**9d**	75	74 (**9d**)
13	CH_2=C($SiMe_3$)MgBr	THF/Et_2O 1:4	**8d**	**9d**	80	78 (**9d**)
14	CH_2=C($SiMe_3$)MgBr	THF/Et_2O 1:8	**8d**	**9d**	76	80 (**9d**)

[a] All reactions were carried out by adding nucleophiles (4 equiv) to furyl sulfonylimine 7. [b] Total isolated yield of **8** and **9**. [c] Determined by ^1H NMR analysis of crude mixture. The major diastereomer is indicated in the parentheses.

Page 2895, Table 3. Table 3 should have appeared as shown below. The description in the Supporting Information has been changed accordingly.

TABLE 3. Suzuki Coupling Reactions of 8 and 9



starting	condition[a]	R	product	yield (%)
8b	A	*t*-Bu	**12a**	62
8c	B	Me_3SiCH_2	**12b**	88
8d	B	CH_2=C($SiMe_3$)	**12c**	82
9b	A	*t*-Bu	**13a**	64
9c	B	Me_3SiCH_2	**13b**	90
9d	B	CH_2=C($SiMe_3$)	**13c**	82

[a] A = $Ba(OH)_2$, PhMe/MeOH (1:1), reflux, 2 h. B = 2 M K_3PO_4, DME/H_2O (4:1), reflux, 2 h.

Page 2895, Figure 5. Figure 5 should have appeared as shown below. The description in the Supporting Information has been changed accordingly.

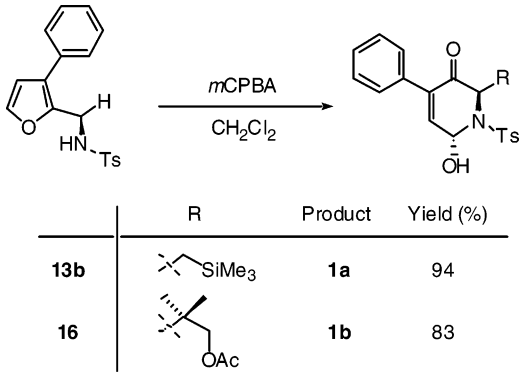


FIGURE 5. Oxidative rearrangement of **13b** and **16**.

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