

Correction to 3d Early Transition Metal Complexes Supported by a New Sterically Demanding Aryloxy Ligand

Keith Searles, Ba L. Tran, Maren Pink, Chun-Hsing Chen, and Daniel J. Mindiola*

Inorg. Chem. **2013**, *52* (19), 11126–11135 DOI: 10.1021/ic401363p

Page 11132. The solvent used to record the ^1H NMR spectrum of compound **1** was incorrectly reported as C_6D_6 . The actual solvent used when recording the reported ^1H NMR spectrum was CDCl_3 . Although this does not greatly affect the assignment of the chemical shifts, the correct referencing of the residual ^1H solvent resonance of CDCl_3 results in a downfield shift of 0.10 ppm for all previously reported chemical shifts of **1**. The correctly referenced chemical shifts are shown in a revised ^1H NMR spectrum (Figure 1). Additionally, the ^1H NMR spectrum of compound **1** recorded in C_6D_6 is shown in Figure 2. The authors apologize for the error.

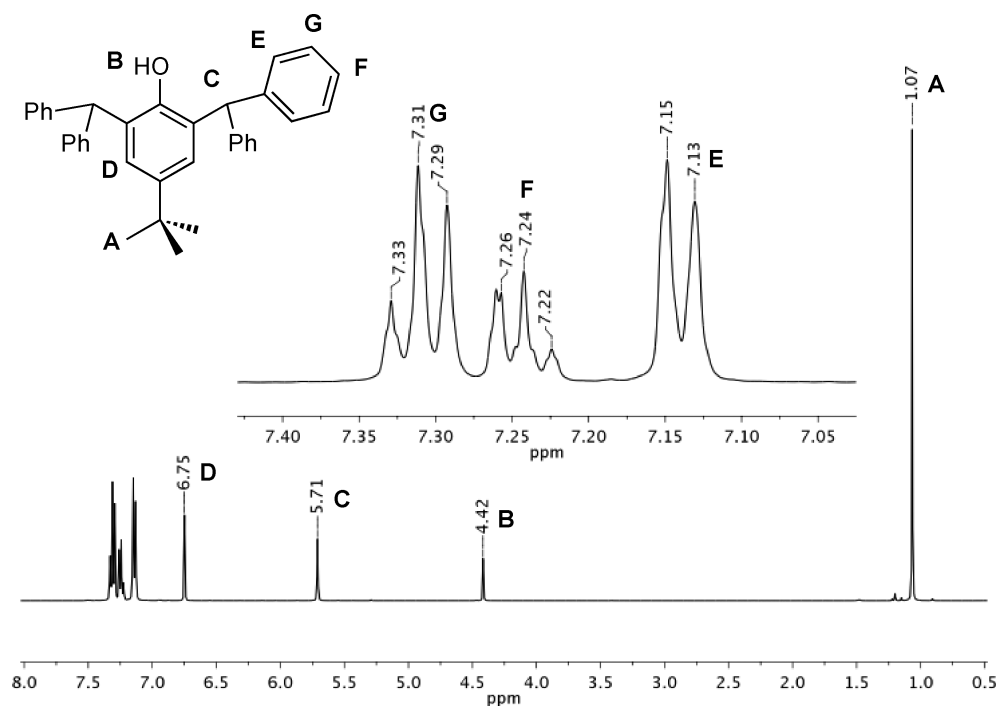


Figure 1. ^1H NMR spectrum of compound **1** (25 °C, 400.11 MHz, CDCl_3).

Published: February 28, 2014



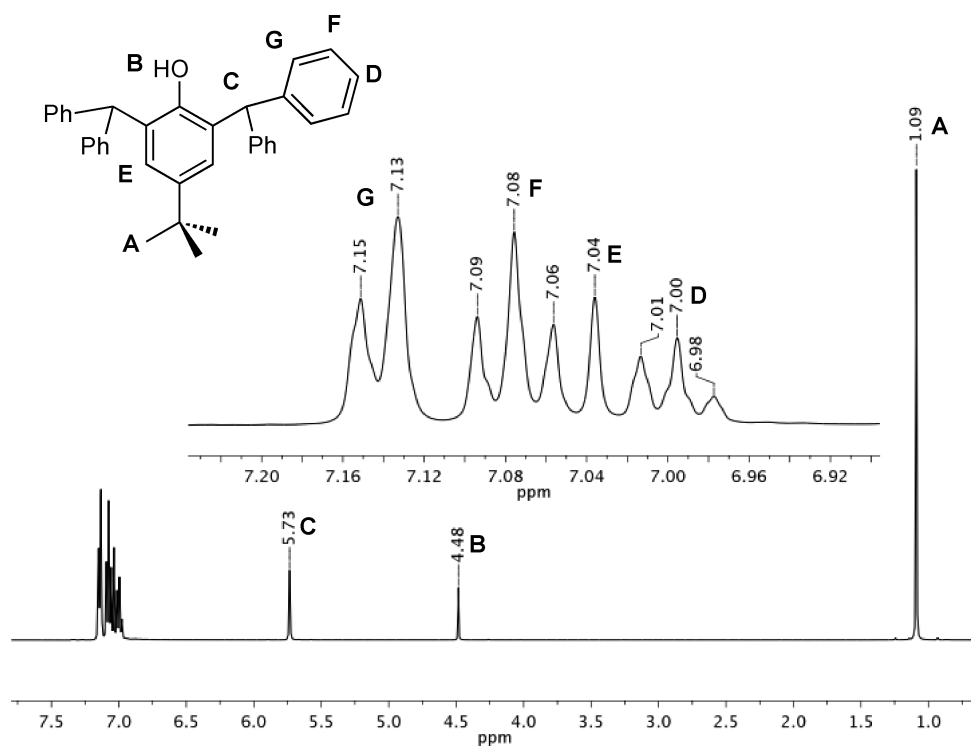


Figure 2. ¹H NMR spectrum of compound 1 (25 °C, 400.11 MHz, C₆D₆).