

Correction to Density Functional Theory Calculations on Mössbauer Parameters of Nonheme Iron Nitrosyls [*Inorganic Chemistry* **2009**, *48*, 9155–9165 DOI: 10.1021/ic9008784].
Kathrin H. Hopmann, Abhik Ghosh,* and Louis Noodleman*

Page 9163. In Table 4, the $\{\text{FeNO}\}^7$ and $\{\text{Fe}(\text{NO})_2\}^9$ labels to the rows for $[\text{Fe}(\text{NO})_2\{\text{Fe}(\text{NO})(\text{N}(\text{CH}_2\text{CH}_2\text{S})_3)\}_2\text{-S,S}']$ were inadvertently reversed. Thus, it is the $\{\text{FeNO}\}^7$ unit that possesses an experimental isomer shift of 0.37 mm s^{-1} at 77 K (and predicted values of $0.39\text{--}0.40 \text{ mm s}^{-1}$). Likewise, the $\{\text{Fe}(\text{NO})_2\}^9$ unit actually possesses an experimental isomer shift of 0.18 mm s^{-1} at 77 K (and a predicted value of 0.19 mm s^{-1}).

DOI: 10.1021/ic200575z

Published on Web 03/31/2011