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Addition/Correction

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Correction to "A Combined Mössbauer, Magnetic Circular Dichroism, and Density Functional Theory Approach for Iron Cross-Coupling Catalysis: Electronic Structure, In Situ Formation, and Reactivity of Iron-Mesityl-Bisphosphines"

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Page 9134. Due to a formatting change during production, the value of D for $FeMes_2$(SciOPP) is incorrectly given as 1191 ± 2 cm^{-1}. The correct value for $FeMes_2$(SciOPP) is D = |19| ± 2 cm^{-1}.

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