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Addition/Correction

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Correction to "Influence of Pyrazolate vs *N*-Heterocyclic Carbene Ligands on the Slow Magnetic Relaxation of Homoleptic Trischelate Lanthanide(III) and Uranium(III) Complexes"

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Page 6061. In Table 2, the values of the relaxation barrier, U_{eff}, given for $U(Bc^{Me})_3$ are incorrect. While the unit given for the entries in column 2 in the table is cm^{-1}, the numerical values given for $U(Bc^{Me})_3$ are in units of Kelvin (1 Kelvin = 0.695 cm^{-1}). Thus, the barrier for $U(Bc^{Me})_3$ is actually 22.0 cm^{-1}, and that for 12% $U(Bc^{Me})_3$ is 23.2 cm^{-1} (33.328(4) K). Please note further that the U_{eff} values reported for $U(Bc^{Me})_3$ in the text also are given in units of Kelvin. All other reported U_{eff} values are correct as presented in the published paper.

This error is an unfortunate oversight; however, it does not change any element of the original work beyond these two strict numerical U_{eff} values. For the purposes of future literature comparisons of U_{eff} etc., this correction is particularly essential.

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