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Erratum

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ERRATUM

Ismail I. Fasfous, Jamal N. Dawoud, Abdulwahab K. Sallabi and Taghreed S. Hassouneh. A density functional theory study of the $Cu^{+}\cdot(CO)_n$ (n = 1–3) complexes. J. Coord. Chem., http://dx.doi.org/10.1080/00958972.2015.1021343

Due to an editing error, when the above article was first published online, "kcal mol^{-1}" had been changed to "kcal M^{-1}" throughout. This has now been corrected in both print and online versions. Taylor & Francis apologizes for this error.