ERRATUM

Erratum to: Improved accuracy benchmarks of small molecules using correlation consistent basis sets

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In the original publication of the article an incorrect conversion factor was used to convert the enthalpy of formation for CN ($^2\Sigma^+$) at 0 K to 298.15 K. The error appeared in Table 7. The correct $\Delta H_f(298\text{ K})$ should be $105.08 \pm 0.16\text{ kcal/mol}$ ($439.65 \pm 0.04\text{ kJ/mol}$). This compares to an Active Thermochemical Tables value of $105.16 \pm 0.03\text{ kcal/mol}$ ($440.001 \pm 0.145\text{ kJ/mol}$). The authors would like to thank Dr. John Simmie from the National University of Ireland for making us aware of the problem.

The online version of the original article can be found under
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