Unpublished erratum

"Density Functional for Spectroscopy: No Long-Range Self-Interaction Error, Good Performance for Rydberg and Charge-Transfer States, and Better Performance on Average than B3LYP for Ground States," Y. Zhao and D. G. Truhlar, Journal of Physical Chemistry A Journal of Physical Chemistry A **110**, 13126-13130 (2006).

In ref. 18, volume 121 should be volume 120.