Unpublished errata for

“Performance of Recent and High-Performance Approximate Density Functionals for Time-Dependent Density Functional Theory Calculations of Valence and Rydberg Electronic Transition Energies,” by M. Isegawa, R. Peverati, and D. G. Truhlar, Journal of Chemical Physics **137**, 244104/1-17 (2012). [dx.doi.org/10.1063/1.4769078](http://dx.doi.org/10.1063/1.4769078)

Erratum: **140**, 129901 (2014). [dx.doi.org/10.1063/1.4869516](http://dx.doi.org/10.1063/1.4869516)

In the "Computational Details" section:

1. line 3: "6-311(2+,2+)G\*\*" should be "6-31(2+,2+)G\*\*" as the basis set for TDDFT calculations.

2.  line 6: "0.00108434" should be "0.0108434".

The authors are grateful to Shaohong Li for finding these typos.