Unpublished errata for

“Valence Excitation Energies of Alkenes, Carbonyl Compounds, and Azabenzenes by Time-Dependent Density Functional Theory: Linear Response of the Ground State Compared to Collinear and Noncollinear Spin-Flip TDDFT with the Tamm-Dancoff Approximation,” M. Isegawa and D. G. Truhlar, Journal of Chemical Physics **138**, 134111/1-13 (2013).

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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.