

Correction to “Eigenstates of Thiophosgene Near the Dissociation Threshold: Deviations From Ergodicity”

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In a recent paper,¹ a typographical error was made in the definition of one of the quantities. The definition of the polyad K in eq 7 of the paper should be

$$K = \nu_1 + \nu_2 + \nu_3$$

In addition, it should be noted that ref 57 in the paper,¹ an earlier work² of the author, has the identical typographical error. Note that all the calculations and results reported in the paper, and the earlier work,² use the correct definition of K . Hence, the conclusions of the paper are not changed.

■ REFERENCES

- (1) Keshavamurthy, S. Eigenstates of Thiophosgene Near the Dissociation Threshold: Deviations from Ergodicity. *J. Phys. Chem. A* **2013**, *117*, 8729–8736.
- (2) Keshavamurthy, S. On the Nature of Highly Excited Vibrational States of Thiophosgene. *J. Chem. Sci.* **2012**, *124*, 291–300.