



Erratum: Theory of vibrational equilibria and pooling at solid-diatom interfaces [J. Chem. Phys.139, 124107 (2013)]

E. T. D. Boney and R. A. Marcus

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Erratum: Theory of vibrational equilibria and pooling at solid-diatom interfaces [J. Chem. Phys. **139**, 124107 (2013)]

E. T. D. Boney and R. A. Marcus^{a)}

Noyes Laboratory of Chemical Physics, California Institute of Technology, Pasadena, California 91125-7200, USA

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We make the following corrections, occasioned by an accidental substitution of one figure by another in a recent paper.¹ On page 124107-3, first column:

- 5-7 lines before Fig. 3, delete “but varies from having slightly shorter to slightly longer effective time constants throughout the calculation.”
- Replace Fig. 3 and its caption by Fig. 3 in this Erratum.

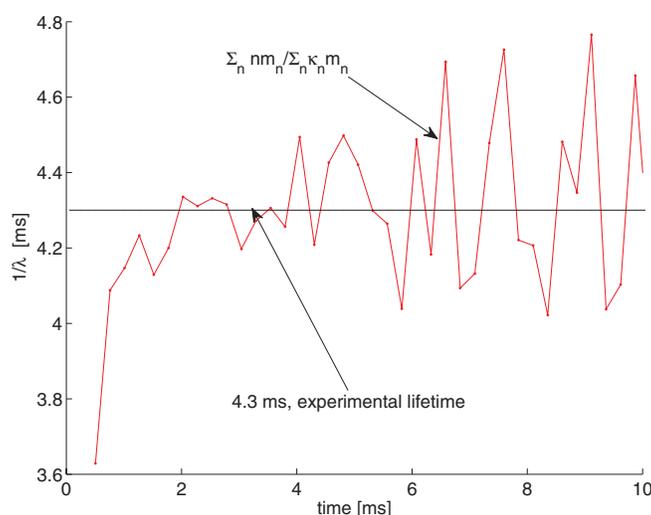


FIG. 3. Calculated $1/\lambda(t)$ versus time, compared with the single exponential observable. Like the experiment, the calculated decay is a single exponential, as in Fig. 4.

Correction 1 arises because the calculated time constant in the corrected figure is more constant than in the earlier figure, no longer showing a systematic variation.

¹E. T. D. Boney and R. A. Marcus, *J. Chem. Phys.* **139**, 124107 (2013).

^{a)}Electronic mail: ram@caltech.edu