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An error was made in the presentation of the Raman scattering cross-sections for the nitrate standard. Fig. 1 was mislabeled and the A-term fitting parameters were reported incorrectly. The corrected figure and legend are given below. The nitrate cross-sections used in the intensity analysis were correct so these changes have no effect on the rhodopsin absolute cross-sections, the intensity analysis, or the conclusions in the original paper. Also, a typographical error was made in Eq. 2.

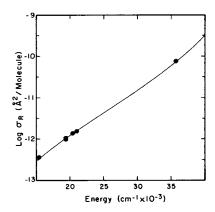


FIGURE 1 A-term fit to the nitrate cross-sections. The solid line was calculated with Eq. 2 using $K = 9.647 \times 10^{-13} \,\text{Å}^2/\text{molecule-sr}$, $\nu_e = 51,940 \,\text{cm}^{-1}$, and $C = 4.504 \times 10^{-10} \,\text{cm}^2$. The estimated errors are on the order of the point size.

$$\sigma_{R} = \frac{8\pi}{3} \left(\frac{1+2\rho}{1+\rho} \right) K \nu_{0} \nu^{3} \left[\frac{\nu_{e}^{2} + \nu_{0}^{2}}{(\nu_{e}^{2} - \nu_{0}^{2})^{2}} + C \right]^{2}. \tag{2}$$