

---

**Erratum**

---

**An Analytic Treatment of Adiabatic and Nonadiabatic Effects on the  
Vibration-rotational Energies of Diatomic Molecules****[Chinese Journal of Physics 30 (2), 177 (1992)]**F. M. Fernandez<sup>1</sup> and J. F. Ogilvie<sup>2</sup><sup>1</sup>*Programa QUINOR, Facultad de Ciencias Exactas, Universidad Nacional de La Plata,  
Calle 47 y 115, Casilla de Correo 962, 1900 La Plata, Argentina*<sup>2</sup>*Institute of Atomic and Molecular Sciences, Academia Sinica, P. O. Box 23-166,  
Taipei Taiwan 10764, R.O.C.*

An incorrect sign of an intermediate quantity has led to erroneous signs in several equations printed in this paper. Specifically, below equation 16,  $W = 1$  should read  $W = -1$ . Correspondingly, all the external signs of quantities on the right-hand side of equation 17 should be reversed; the two minus signs at the top level in equation 32 should be replaced by plus signs; the sign before each coefficient  $\Xi_j$ ,  $0 \leq j \leq 4$ , in the table and in the three equations appearing in the text on page 190 should be reversed. In no way are the discussion and the conclusions affected by these modifications.

An obvious typographical error appears in the expression for  $Z_{1,0}$  in the table:  $\gamma_2$  should be replaced by  $\eta_2$  in the second term that hence should read " $+\gamma(-3a_1\eta_1/2 + \eta_2)$ ".

We thank Professor E. Tiemann for helpful comments and the National Science Council of the Republic of China for support of this research.