Errata

Erratum: Collision-induced rotovibrational spectra of H_2-He pairs from first principles [Phys. Rev. A 35, 632 (1987)]

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In this paper, Tables I and II of the *ab initio* radial matrix elements of the induced dipole moment of H_2-He pairs were recently found to contain errors in the order of 10% or 15% that we correct in the enclosed tables. The errors are due to a sign error in one of the matrix elements $\langle 00|\rho^n|10\rangle$ used to derive the data from Table III of Ref. 1.

This unfortunate fact renders the comparisons of the computed rotovibrational collision-induced absorption spectra obsolete. New computations of these spectra at a variety of temperatures, which also reflect a new accounting for the rotation dependence of these matrix elements, will be given in Ref. 2. These differ from our previous computation mostly in the Q branch.

For the work at the University of Texas, the support of the National Science Foundation, Grant No. AST-8613085, is acknowledged.

TABLE I. Induced dipole components $B^{01}_{\lambda L}$ in 10^{-6} a.u.

R	λL=01	23	21	45
3	25 987	8284	−11 067	2370
4	7310	2306	−2429	526
5	1821	678	−500	109
6	381	231	−92	23
7	63	99	−16	6
8	5	52	−2	2
10	−2	21	0	0

TABLE II. Coefficients of Eq. (4), with R_0=5.80 bohrs.

λL	N	C_N	c_2	c_3	c_4
01	7	−80.9	0.002 46	−1.677	−0.045
23	4	1.17	0.000 217	−1.656	−0.062
21	7	−11.8	−0.000 277	−1.813	−0.051
45	6	1.19	0.000 027 3	−1.884	−0.099

[1]W. Meyer and L. Frommhold, Phys. Rev. A **34**, 2771 (1986).
[2]A. Borysow, L. Frommhold, and W. Meyer, Phys. Rev. A (to be published).

