

SELF- AND AIR-BROADENING AND SHIFT COEFFICIENTS OF CH₄ LINES IN THE 3 μm REGION

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Self- and air-broadening and shift coefficients for numerous CH₄ transitions in the 3 μm region have been determined through analysis of high-resolution absorption spectra recorded by two different Fourier transform spectrometer (FTS) systems. The first set of spectra was recorded using the 1-m FTS at the McMath-Pierce facility of the National Solar Observatory at Kitt Peak; resolution was 0.01 cm⁻¹. This data set included both self-broadened CH₄ spectra with cell path lengths of 5, 10, 25, 150, and 2500 cm and pressures ranging from 0.4 Torr to 456 Torr, and air-broadened spectra of ¹²CH₄ and ¹³CH₄ with cell path lengths of 5, 25, 50, and 150 cm and total pressures from about 100 Torr to 400 Torr. A second set of self-broadened CH₄ spectra was recorded at higher resolution (0.0027 cm⁻¹) using the FTS of the Laboratoire de Physique Moléculaire et Applications in Orsay. Path lengths for these spectra varied from 1.7 cm to 4 m, and pressures ranged from less than 1 Torr to over 500 Torr. A multispectrum nonlinear least-squares procedure^a was used to fit over 40 spectra simultaneously to determine self- and air-broadening and shift coefficients for more than 500 lines; most of these are in the ν₃ and ν₂ + ν₄ bands. Differences between air- and self-broadening and shifts of ¹²CH₄ lines will be discussed, and air-broadening and shift coefficients for ¹²CH₄ and ¹³CH₄ will be compared. We will also compare our measurements with the values on the current HITRAN compilation^b and with other available measurements.

^aD. Chris Benner, C. P. Rinsland, V. Malathy Devi, M. A. H. Smith, and D. Atkins, *JQSRT* **53**, 705-721 (1995).

^bL. S. Rothman *et al.*, *JQSRT* **48**, 469-507 (1992); L. S. Rothman *et al.*, *JQSRT*, in press (1998).