

## CHARACTERIZATION OF THE H<sub>2</sub>O+CO<sub>2</sub> CONTINUUM ABSORPTION WITHIN THE INFRARED TRANSPARENCY WINDOWS FOR PLANETARY APPLICATIONS

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Accurate knowledge of the absorption by a gas mixture of CO<sub>2</sub> and water is crucial for planetary sciences, as it allows for better modeling the atmospheres of rocky planets, e.g. improving our understanding of the early climate of Mars or why Venus and the Earth have evolved so differently. In addition to local monomer lines proportional to the density of each species, the absorption spectrum of such a gas mixture includes binary absorption features varying smoothly with frequency: self-continuum absorption proportional to the squared density, and “crossed” absorption involving both species and scaling as the density product  $\rho_{CO_2}\rho_{H_2O}$ . We used highly sensitive spectroscopy techniques (CRDS and OFCEAS) to measure the absorption by H<sub>2</sub>O+CO<sub>2</sub> gas mixtures in several spectral regions situated in transparency windows where the monomer absorption of both species is weak (1.5-1.53  $\mu\text{m}$ , 1.68-1.75  $\mu\text{m}$ , 2.06  $\mu\text{m}$ , 2.2-2.35  $\mu\text{m}$ , 3.5  $\mu\text{m}$ ). For both water and CO<sub>2</sub>, the monomer lines, modeled using HITRAN parameters, and the self-continuum absorption, calculated from literature values or measured in dedicated experiments, were subtracted from the measured absorption. The obtained “crossed absorption” coefficients are compared to the only available empirical model based on far wings of line shape profiles scaled by  $\chi$ -factors.<sup>a</sup> An additional absorption peak centered at about 6000 cm<sup>-1</sup> was attributed to a collision-induced simultaneous transition of H<sub>2</sub>O and CO<sub>2</sub> through the  $\nu_1$  and  $\nu_3$  modes, respectively. The assignment was confirmed using humidified <sup>13</sup>CO<sub>2</sub>, where a similar band was observed about 68 cm<sup>-1</sup> away corresponding to the isotopic spectral shift of the  $\nu_3$  band of CO<sub>2</sub>. Classical molecular dynamics simulations (CMDS) of the considered collision-induced absorption were conducted and are found in good agreement with the experiment.<sup>b</sup>

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<sup>a</sup>Fleurbaey H, Campargue A, Carreira Mendès Da Silva Y, Grilli R, Kassi S, Mondelain D. *J Quant Spectrosc Radiat Transf* 108119 (2022)

<sup>b</sup>Fleurbaey H, Mondelain D, Fakhardji W, Hartmann J-M, Campargue A. Submitted to *J Quant Spectrosc Radiat Transf*