

## MOLECULAR TRANSITION FREQUENCIES OF CO<sub>2</sub> NEAR 1.6 $\mu$ WITH KHZ-LEVEL UNCERTAINTIES

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We present measurements of molecular transition frequencies based on the comb-locked cavity ring-down spectroscopy technique [1], reporting vacuum transition frequencies of Doppler-broadened <sup>12</sup>C<sup>16</sup>O<sub>2</sub> in the 1.6  $\mu$ m region for the (30012)  $\leftarrow$  (00001) and (30013)  $\leftarrow$  (00001) bands with an average combined standard uncertainty of 1 kHz. A global multi-state model was fit to these data and literature values to provide spectroscopic parameters and a best-case fit precision of 4 kHz. We identified and assigned an interaction between the (30012) and (33301) states which was manifest as an observed Fermi resonance in the (30012)  $\leftarrow$  (00001) band. This interaction was accounted for in the global fit, significantly reducing uncertainties in the spectroscopic parameters of the (30012) state. We find excellent agreement with literature values, including better than 1 kHz RMS deviation with a recent saturation cavity ring-down study of the (30013)  $\leftarrow$  (00001) band [2]. Agreement is excellent with CDSD 2019 [3] and HITRAN 2016 [4] for the (30013)  $\leftarrow$  (00001) band, but much poorer for the (30012)  $\leftarrow$  (00001) band.

[1] Z. D. Reed, D. A. Long, H. Fleurbaey, and J. T. Hodges, *Optica* 7, 1209-1220 (2020).

[2] H. Wu, et al., *Phys. Chem. Chem. Phys.* 22, 2841-2848 (2020).

[3] S. A. Tashkun, et al., *J. Quant. Spectros. Rad. Trans.* 228, 124-131 (2019).

[4] I. E. Gordon, et al. *J. Quant. Spectros. Rad. Trans.* 203, 3-69 (2017).