

All data taken at Pacific Northwest National Laboratory (PNNL)  
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Composite spectrum for CO\_5T

Effective burden of composite spectrum: 1 part-per-million-meter (ppm-meter) at 296 K

#### **Sample Conditions-**

- Chemical name and CAS number: Carbon monoxide, CO : [630-08-0]
- Physical properties: M.W. 28 amu, F.P. -207C, B.P. -190C
- Supplier and stated purity: Air Products and Chemicals, Inc., 99.99%
- Sample class: I (PNNL scale).
- Temperature of sample:  $5.04 \pm 0.02$  C
- Diluent: Sample back filled with ultra high purity nitrogen to  $760 \pm 5$  Torr
- Individual samples at 0.58135, 0.99162, 1.6484, 8.4457, 145.88 and 49.33 Torr. Path length = 19.96 cm. Final data is a composite spectrum.
- Preparation: Multiple freeze-thaw cycles at 77K to remove air.

#### **Instrument Parameters-**

- Bruker-66V FTIR, temperature controlled environment, evacuated optics bench
- Spectral range: 6,500 to 600  $\text{cm}^{-1}$  (1.534 to 16.667 microns)
- Instrumental resolution (interferogram): 0.112  $\text{cm}^{-1}$
- Spectral intervals after FFT: 0.06  $\text{cm}^{-1}$
- Interferogram zero-fill: 2X
- Apodization: Boxcar
- Phase correction: Mertz
- Beam splitter: Potassium bromide (KBr)
- IR source: Carbide glowbar (22 V)
- Scanner velocity: 9 (Bruker arbitrary)
- Number of interferograms averaged per single channel spectra: 256
- Detector: Mid-band HgCdTe, photoconductive, 77K operation
- Folding limits: 15798 to 0  $\text{cm}^{-1}$

#### **Post Processing and Related Parameters-**

- Non-linearity detector correction (Bruker proprietary) applied to interferogram ( $\alpha = 0.85$ ,  $\beta = 530$ )
- Composite spectrum created from 6 individual absorbance (base-10) spectra via classical least squares fit: Intercept=0, slope is fitted, individual absorbance values weighted by  $T^2$  (transmission squared), all absorbance values  $> 1.6$  are given zero weight
- Calculated and estimated errors: Type A = 0.78%, Type B = 3%
- Frequency correction:  $V(\text{corrected}) = V(\text{instrument}) * 0.9999984669 + 0.005187$
- Axis units: X=wavenumbers ( $\text{cm}^{-1}$ ), Y=Absorbance (base-10)
- Trace carbon dioxide and water removed by spectral subtraction. Baseline straightened by subtraction of 7<sup>th</sup>-order polynomial.