

# Non-thermal desorption of complex organic molecules

## Efficient CH<sub>3</sub>OH and CH<sub>3</sub>COOCH<sub>3</sub> sputtering by cosmic rays (Corrigendum)

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A typographical error must be corrected in Table 2. The integrated band strength ( $A$  value,  $\text{cm molec}^{-1}$ ) for the CO<sub>2</sub> stretch at 2350  $\text{cm}^{-1}$  was incorrectly typeset and attributed the same

value as the CH<sub>3</sub>OH C–O stretch ( $1.8 \times 10^{-17} \text{cm molec}^{-1}$ ), and the reference D'Hendecourt & Allamandola (1986). They have been replaced in Table 2.

**Table 2.** Integrated band strengths used in the analysis.

Species	Mode	Position $\text{cm}^{-1}$	$A$ $\text{cm molec}^{-1}$	Ref.
CO	CO stretch	2140	$1.1 \times 10^{-17}$	Jiang et al. (1975)
			$1.1 \times 10^{-17}$	Gerakines et al. (1995)
			$1.12 \times 10^{-17}$	Bouilloud et al. (2015)
			$1.1 \times 10^{-17}$	Adopted for this work
CO <sub>2</sub>	CO <sub>2</sub> stretch	2350	$7.6 \times 10^{-17}$	Gerakines et al. (1995)
H <sub>2</sub> O	OH stretch	3600–2700	$2.2 \times 10^{-16}$	D'Hendecourt & Allamandola (1986)
			$2.2 \pm_{0.2}^0 \times 10^{-16}$	Adopted for this work
CH <sub>3</sub> OH	OH stretch	3600–2700	$1.1 \times 10^{-16}$	D'Hendecourt & Allamandola (1986)
			$1.28 \times 10^{-16}$	Palumbo et al. (1999)
			$1.0 \times 10^{-16}$	Bouilloud et al. (2015)
			$1.1 \pm 0.15 \times 10^{-16}$	Adopted for this work
CH <sub>3</sub> OH	C–O stretch	1030	$1.8 \times 10^{-17}$	D'Hendecourt & Allamandola (1986)
			$1.8 \times 10^{-17}$	Sandford & Allamandola (1993)
			$1.2 \times 10^{-17}$	Palumbo et al. (1999)
			$1.07 \times 10^{-17}$	Bouilloud et al. (2015)
			$1.5 \pm_{0.4}^{0.3} \times 10^{-17}$	Adopted for this work
CH <sub>3</sub> COOCH <sub>3</sub>	C–O stretch	1255	$5 \times 10^{-17} \text{ }^{(a)}$	D'Hendecourt & Allamandola (1986)
			$5 \pm 0 \times 10^{-17}$	Adopted for this work

**Notes.** <sup>(a)</sup>Band strength from the CO stretching mode of ethyl acetate measurement.