

# IUPAC Task Group on Atmospheric Chemical Kinetic Data Evaluation – Data Sheet SOx62

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## CH<sub>3</sub>SO + NO<sub>2</sub> → products

### Rate coefficient data

$k/\text{cm}^3 \text{ molecule}^{-1} \text{ s}^{-1}$	Temp./K	Reference	Technique/ Comments
<i>Absolute Rate Coefficients</i>			
$(3 \pm 2) \times 10^{-11}$	298	Mellouki et al., 1988	DF-MS
$(8 \pm 5) \times 10^{-12}$	298	Tyndall and Ravishankara, 1989	PLP-LIF
$(1.2 \pm 0.25) \times 10^{-11}$	298	Dominé et al., 1990	DF-MS
$(1.5 \pm 0.4) \times 10^{-11}$	300	Kukui et al, 2000	(a) PLP-LIF, DF-LIF

### Comment

- (a) PLP of CH<sub>3</sub>S<sub>2</sub>CH<sub>3</sub>/NO<sub>2</sub> mixtures at 351 and 248 nm relative to CH<sub>3</sub>I/NO<sub>2</sub> at 351 nm in the range 16-814 mbar He at 300 K with LIF detection of CH<sub>3</sub>O. The temporal profile of CH<sub>3</sub>O was monitored in order to obtain  $k$  by fitting which was found to be independent of pressure. Supporting measurements on pressure-dependent CH<sub>3</sub> yields in the range 243-333 K and 16-814 mbar He have been performed. Ancillary measurements of absolute SO<sub>2</sub> yields at 1.3 mbar He result in  $1.0 \pm 0.1$  using Cl + CH<sub>3</sub>SH in DF-LIF measurements.

### Preferred Values

$k = 1.2 \times 10^{-11} \text{ cm}^3 \text{ molecule}^{-1} \text{ s}^{-1}$  at 298 K.

#### Reliability

$\Delta \log k = \pm 0.2$  at 298 K.

#### Comments on Preferred Values

The measured values (Mellouki et al., 1988; Tyndall and Ravishankara, 1989; Dominé et al., 1990, Kukui et al., 2000) of  $k$  at 298 K agree within their error limits, some of which are substantial. The rate coefficient for this reaction is difficult to measure because of the lack of a clean primary source of CH<sub>3</sub>SO radicals and the complexity of the secondary chemistry. The two most recent determinations (Dominé et al., 1990; Kukui et al., 2000) agree well with each other, the latter of which being a relative rate study extended to 814 mbar (He) using CH<sub>3</sub> + NO<sub>2</sub> as a reference reaction. The preferred value is that of Dominé et al. (1990) as it agrees with all studies within the given uncertainty limits.

## References

- Dominé, F., Murrells, T. P. and Howard, C. J.: J. Phys. Chem. 94, 5839, 1990.  
Kukui, A., Bossoutrot, V., Laverdet, G. and Le Bras, G.: J. Phys. Chem. A 104, 935, 2000.  
Mellouki, A., Jourdain, J. L. and Le Bras, G.: Chem. Phys. Lett. 148, 231, 1988.  
Tyndall, G. S. and Ravishankara, A. R.: J. Phys. Chem. 93, 2426, 1989.