

IUPAC Task Group on Atmospheric Chemical Kinetic Data Evaluation – Data Sheet oClOx63

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This data sheet last evaluated: June 2015; last change in preferred values: November 2003.

HO + C(O)Cl₂ → products

Rate coefficient data

<i>k</i> /cm ³ molecule ⁻¹ s ⁻¹	Temp./K	Reference	Technique/ Comments
<i>Relative Rate Coefficients</i> <1 x 10 ⁻¹⁵	298 ± 3	Nelson et al., 1990	RR (a)

Comments

- (a) HO radicals were generated by the photolysis of CH₃ONO in CH₃ONO-NO-C(O)Cl₂-reference compound-air mixtures at ~1 bar pressure. The concentrations of C(O)Cl₂ and the reference compound were monitored during the experiments by GC. No reaction of C(O)Cl₂ was observed. However, no details concerning the reference compound used or the amount of reference compound reacted were given.

Preferred Values

Parameter	Value	T/K
<i>k</i> /cm ³ molecule ⁻¹ s ⁻¹	<5 x 10 ⁻¹⁵	298

Comments on Preferred Values

The preferred upper limit to the 298 K rate coefficient is based on the sole reported study of Nelson et al. (1990), with the preferred upper limit being increased by a factor of 5 over that cited by Nelson et al. (1990).

References

Nelson, L., Shanahan, I., Sidebottom, H. W., Treacy, J. and Nielsen, O. J.: Int. J. Chem. Kinet., 22, 577, 1990.