

## IUPAC Task Group on Atmospheric Chemical Kinetic Data Evaluation – Data Sheet oFOx108

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This data sheet last evaluated: September 2011; last change in preferred values: September 2011.



### Rate coefficient data ( $k = k_1 + k_2$ )

$k/\text{cm}^3 \text{ molecule}^{-1} \text{ s}^{-1}$	$T/\text{K}$	Reference	Technique/ Comments
<i>Relative Rate Coefficients</i>			
$(6.89 \pm 0.81) \times 10^{-15}$	297	Chen et al. (1997)	(a)

### Comments

- (a) Cl atoms were generated by the photolysis of  $\text{Cl}_2$  in  $\text{CHF}_2\text{CH}_2\text{CF}_3/\text{CF}_3\text{CHClF}/\text{Cl}_2$  mixtures in 700 Torr (933 mbar) of  $\text{N}_2$ . The decays of  $\text{CHF}_2\text{CH}_2\text{CF}_3/\text{CF}_3\text{CHClF}$  were measured by FTIR spectroscopy. The measured rate coefficient ratio of  $k(\text{Cl}+\text{CHF}_2\text{CH}_2\text{CF}_3)/k(\text{Cl}+\text{CF}_3\text{CHClF}) = 2.55 \pm 0.30$  was placed on an absolute basis using  $k(\text{Cl}+\text{CF}_3\text{CHClF}) = 2.7 \times 10^{-15} \text{ cm}^3 \text{ molecule}^{-1} \text{ s}^{-1}$  (Atkinson et al., 2008).

### Preferred Values

Parameter	Value	$T/\text{K}$
$k/\text{cm}^3 \text{ molecule}^{-1} \text{ s}^{-1}$	$6.9 \times 10^{-15}$	298
<i>Reliability</i>		
$\Delta \log k$	$\pm 0.15$	298

### Comments on Preferred Values

The recommended value is based on the study by Chen et al. (1997).

### References

Atkinson, R., Baulch, D. L., Cox, R. A., Crowley, J. N., Hampson, R. F., Hynes, R. G., Jenkin, M. E., Rossi, M. J., Troe, J., and Wallington, T. J.: Atmos. Chem. Phys., 9, 4141, 2008; IUPAC Task Group on Atmospheric Chemical Kinetic Data Evaluation, <http://iupac.pole-ether.fr>.  
Chen, J., Young, V., Niki, H., and Magid, H.: J. Phys. Chem., A, 101, 2648, 1997.