

IUPAC Task Group on Atmospheric Chemical Kinetic Data Evaluation – Data Sheet HO_x_VOC77

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This data sheet last evaluated: June 2016; last change in preferred values: June 2016.

HO + C₂H₅CH(OH)CH₂ONO₂ → products

Rate coefficient data

<i>k</i> /cm ³ molecule ⁻¹ s ⁻¹	<i>T</i> /K	Reference	Technique/ Comments
<i>Relative Rate Coefficients</i>			
(7.0 ± 1.4) × 10 ⁻¹²	296 ± 2	Treves and Rudich, 2003	RR (a, b)

Comments

- (a) The rate coefficient was measured at atmospheric pressure of dry synthetic air in a photochemical reactor (100 L Tedlar chamber) by the relative rate technique using solid-phase microextraction (SPME) coupled to gas chromatography (GC) for the detection of the organic reactants. HO radicals were generated by the photolysis of CH₃ONO (or C₂H₅ONO)-NO-air + 1-nitrooxy-2-butanol + 1-pentanol (or 1-octanol) mixtures at wavelengths > 300 nm. The measured rate coefficient ratios of *k*(HO + 1-nitrooxy-2-butanol)/*k*(HO + 1-pentanol) and *k*(HO + 1-nitrooxy-2-butanol)/*k*(HO + 1-octanol) were not given by the authors, they were placed on an absolute basis using *k*(HO + 1-pentanol) = (1.12 ± 0.15) × 10⁻¹¹ and *k*(HO + 1-octanol) = (1.44 ± 0.15) × 10⁻¹¹ cm³ molecule⁻¹ s⁻¹ (Nelson et al., 1990).
- (b) Average of the data obtained relative to HO + 1-pentanol and HO + 1-octanol.

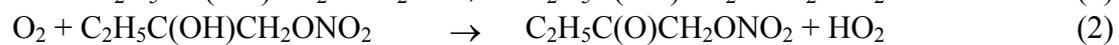
Preferred Values

Parameter	Value	<i>T</i> /K
<i>k</i> /cm ³ molecule ⁻¹ s ⁻¹	7.0 × 10 ⁻¹²	298
<i>Reliability</i>		
Δ log <i>k</i>	± 0.2	298

Comments on Preferred Values

The preferred value is an average of the relative rate coefficients of Treves and Rudich (2003) obtained with two reference compounds. The reaction of OH with C₂H₅CH(OH)CH₂ONO₂ is expected to proceed mainly via H-atom abstraction from the CH group in the α-position to the hydroxyl group leading to C₂H₅C(OH)CH₂ONO₂

radicals which reacts with O₂ leading to HO₂ and the corresponding carbonyl (C₂H₅C(O)CH₂ONO₂).



References

Nelson, L., Rattigan, O., Neavyn, R., Sidebottom, H., Treacy, J., and Nielsen, O.J.: *Int. J. Chem. Kinet.*, 22, 1111, 1990.

Treves, K., and Rudich, Y.: *J. Phys. Chem. A*, 107, 7809, 2003.