

The Chemical Mechanism of ECHAM6.3-HAM2.3-MOZ1.0

JAM version: 002b

The numbers below refer to the number of species or reactions.

| | |
|--|-----|
| Chemical species (Table S1): | 205 |
| Tropospheric photolysis reactions (Table S2): | 97 |
| Stratospheric photolysis reactions (Table S3): | 45 |
| (Tropospheric) O _x reactions (Table S4): | 7 |
| (Tropospheric) HO _x reactions (Table S5): | 17 |
| NO _x reactions (Table S6): | 25 |
| C1 oxidation reactions (Table S7): | 20 |
| C2 oxidation reactions (Table S8): | 48 |
| C3 oxidation reactions (Table S9): | 31 |
| C4 oxidation reactions (Table S10): | 63 |
| C5 oxidation reactions (Table S11): | 137 |
| C6 oxidation reactions (Table S12): | 31 |
| C7 oxidation reactions (Table S13): | 20 |
| C8 oxidation reactions (Table S14): | 13 |
| C10/C15 oxidation reactions (Table S15): | 36 |
| Tropospheric halogen reactions (Table S16): | 21 |
| Sulfur reactions (Table S17): | 11 |
| Stratospheric O(1D) reactions (Table S18): | 28 |
| Stratospheric inorganic halogen reactions (Table S19): | 48 |
| Stratospheric organic halogen reactions (Table S20): | 12 |
| (Tropospheric) heterogeneous reactions (Table S21): | 7 |
| Stratospheric sulfate aerosol reactions (Table S22): | 6 |
| Stratospheric nitric acid dilydrate reactions (Table S23): | 5 |
| Stratospheric ice aerosol reactions (Table S24): | 6 |
| Total number of reactions: | 734 |
| Henry coefficients (Table S25): | 135 |

Note: Where no specific reference is given for a rate constant or products, they are taken from previous versions of MOZART, or set to be equal to those for analogous compounds.

Table S1: Chemical species in the ECHAM-HAMMOZ JAM2 mechanism.

| Species | Name | Structure | Moleweight |
|---------|--|--|------------|
| O1D | O singlett D | O | 16 |
| OH | hydroxyl radical | OH | 17.00734 |
| HONO | nitrous acid | HONO | 47.01344 |
| HO2NO2 | pernitric acid | HO ₂ NO ₂ | 79.01 |
| NH3 | ammonia | NH ₃ | 17.03052 |
| CL | atomic chlorine | Cl | 35.45 |
| CL2 | chlorine | Cl ₂ | 70.906 |
| HCL | hydrogen chloride | HCl | 36.46094 |
| CLO | chlorine oxide | ClO | 51.4524 |
| OCLO | chlorine dioxide | OCLO | 67.4518 |
| CL2O2 | cichlorine dioxide | Cl ₂ O ₂ | 102.9048 |
| HOCL | hypochlorous acid | HOCl | 52.46034 |
| CLONO2 | chlorine nitrate | ClONO ₂ | 97.4579 |
| BR | atomic bromine | Br | 79.9 |
| BR2 | bromine | Br ₂ | 159.81 |
| HBR | hydrobromic acid | HBr | 80.91194 |
| BRO | bromine oxide | BrO | 95.9034 |
| HOBR | hypobromous acid | HOBr | 96.91134 |
| BRONO | bromo nitrite | BrONO | 125.91 |
| BRONO2 | nitryl bromide | BrONO ₂ | 141.9089 |
| BRCL | bromine chloride | BrCl | 115.357 |
| HCOOH | formic acid | HC(O)OH | 46.02538 |
| HOCH2OO | peroxy radical from formaldehyde | HOCH ₂ OO | 63.03 |
| CH3OH | methanol | CH ₃ OH | 32.04186 |
| CH3OOH | methyl peroxide | CH ₃ OOH | 48.04126 |
| HCN | hydrogen cyanide | HCN | 27.02534 |
| CH3SO3H | peroxide from methanesulfonate | CH ₃ SO ₃ H | 96.1 |
| C2H5OH | ethanol | CH ₃ CH ₂ OH | 46.06844 |
| CH3CHO | acetaldehyde | CH ₃ CHO | 44.05256 |
| C2H5OOH | ethyl hydro peroxide | CH ₃ CH ₂ OOH | 62.06784 |
| EO2 | peroxy radical from ethene | HOCH ₂ CH ₂ O ₂ | 77.06 |
| EO | product of EO2+NO | HOCH ₂ CH ₂ O | 61.06 |
| EOOH | peroxide from (2-hydroxyethyl)dioxidanyl | HOCH ₂ CH ₂ OOH | 78.06724 |
| CH3CO3 | peroxyacetyl radical | CH ₃ C(O)OO | 75.04 |

Table S1: Chemical species in JAM2 (... continued)

| Species | Name | Structure | Moleweight |
|--------------------------------------|---|--|------------|
| PAN | peroxyacetyl nitrate | $\text{CH}_3\text{C}(\text{O})\text{OONO}_2$ | 121.04892 |
| GLYOXAL | glyoxal | CHOCHO | 58.03608 |
| GLYALD | glycolaldehyde | HOCH_2CHO | 60.05196 |
| CH_3COOH | acetic acid | CH_3COOH | 60.05196 |
| CH_3COOOH | peracetic acid | $\text{CH}_3\text{C}(\text{O})\text{OOH}$ | 76.05136 |
| $\text{HOCH}_2\text{CO}_2\text{H}$ | peroxy radical from GLYALD | $\text{HOCH}_2\text{CO}_2\text{H}$ | 76.05136 |
| $\text{HOCH}_2\text{CO}_3\text{H}$ | hydroxyacetyl hydroperoxide | $\text{HOCH}_2\text{CO}_3\text{H}$ | 92.05 |
| HCOCO_2H | glyoxylate | HCOCO_2H | 74.03548 |
| HCOCO_3H | peroxide from HCOCO_3 | HCOCO_3H | 90.03488 |
| CH_3CN | methyl cyanide | CH_3CN | 41.05192 |
| DMS | dimethyl sulfide | CH_3SCH_3 | 62.13404 |
| DMSO | dimethyl sulfoxide | $\text{CH}_3\text{S}(\text{O})\text{CH}_3$ | 78.13344 |
| CH_3COCH_3 | acetone | $\text{CH}_3\text{C}(\text{O})\text{CH}_3$ | 58.07914 |
| CH_3COCHO | methyl glyoxal | $\text{CH}_3\text{C}(\text{O})\text{CHO}$ | 72.06266 |
| $\text{C}_3\text{H}_7\text{OOH}$ | lumped propyl hydro peroxide | $\text{C}_3\text{H}_8\text{O}_2$ | 76.09 |
| PO_2 | beta-hydroxy-peroxy-radical from propene | $\text{HOCH}_2\text{CH}(\text{CH}_3)\text{O}_2$ | 91.08 |
| $\text{PRONO}_3\text{BO}_2$ | nitro-peroxy radical from C_3H_6 | $\text{CH}_3\text{CH}(\text{OO})\text{CH}_2\text{ONO}_2$ | 136.08 |
| NOA | alpha-nitrooxy-acetone | $\text{CH}_3\text{C}(\text{O})\text{CH}_2\text{ONO}_2$ | 119.0761 |
| POOH | beta-hydroxy-hydroperoxide from propene | $\text{CH}_3\text{CH}(\text{OOH})\text{CH}_2\text{OH}$ | 92.09382 |
| $\text{CH}_3\text{COCH}_2\text{O}_2$ | peroxy radical from acetone | $\text{CH}_3\text{C}(\text{O})\text{CH}_2\text{O}_2$ | 89.07 |
| ROOH | hydroperoxide from $\text{CH}_3\text{COCH}_2\text{O}_2$ | $\text{CH}_3\text{C}(\text{O})\text{CH}_2\text{OOH}$ | 90.07794 |
| HYAC | 1-hydroxy acetone | $\text{CH}_3\text{C}(\text{O})\text{CH}_2\text{OH}$ | 74.07854 |
| $\text{PR}_2\text{O}_2\text{HNO}_3$ | hydroperoxy propyl nitrate | $\text{CH}_3\text{CH}(\text{OOH})\text{CH}_2\text{ONO}_2$ | 137.09 |
| BIGENE | large alkenes | C_4H_8 | 56.1 |
| MEK | methyl acetone | $\text{CH}_3\text{C}(\text{O})\text{CH}_2\text{CH}_3$ | 72.10572 |
| MACR | methacrolein | $\text{CH}_2\text{CCH}_3\text{CHO}$ | 70.08984 |
| MVK | methyl vinyl ketone | $\text{CH}_3\text{C}(\text{O})\text{CHCH}_2$ | 70.08984 |
| MPAN | peroxy methacryloyl nitrate | $\text{CH}_2\text{CCH}_3\text{CO}_3\text{NO}_2$ | 147.0862 |
| MEKO_2 | peroxy radical from MEK | $\text{C}_4\text{H}_7\text{O}_3$ | 103.09 |
| MEKOOH | peroxide from MEK | $\text{C}_4\text{H}_8\text{O}_3$ | 104.10452 |
| ENEO_2 | peroxy radical from BIGENE/OLTP | $\text{C}_4\text{H}_9\text{O}_3$ | 105.11 |
| MCO_3 | peroxy radical from methacrolein | $\text{CH}_3\text{C}(\text{CH}_2)\text{C}(\text{O})\text{OO}$ | 101.08 |
| MACRO_2 | peroxy radical from methacrolein | $\text{CH}_3\text{C}(\text{OO})(\text{CHO})\text{CH}_2\text{OH}$ | 119.09 |
| MACO_2H | methacroleic acid | $\text{CH}_3\text{C}(\text{CH}_2)\text{C}(\text{O})\text{OH}$ | 86.08924 |
| MACO_3H | methacroleic peroxy acid | $\text{CH}_3\text{C}(\text{CH}_2)\text{C}(\text{O})\text{OOH}$ | 102.08864 |

Table S1: Chemical species in JAM2 (... continued)

| Species | Name | Structure | Moleweight |
|------------|--|---|------------|
| MACROH | multifunctional aldehyde | $\text{HOCH}_2\text{C}(\text{CH}_3)(\text{OH})\text{CHO}$ | 104.10452 |
| LHMKABO2 | peroxy radicals from MVK | $\text{C}_4\text{H}_7\text{O}_4$ | 119.09 |
| BIACETOH | multifunctional aldehyde | $\text{CH}_3\text{C}(\text{O})\text{C}(\text{O})\text{CH}_2\text{OH}$ | 102.0886 |
| LHMKABOOH | hydroperoxides from LHMKABO2 | $\text{C}_4\text{H}_8\text{O}_4$ | 120.1 |
| IBUTALOH | multifunctional aldehyde | $\text{CH}_3\text{C}(\text{OH})(\text{CHO})\text{CH}_3$ | 88.10512 |
| IBUTALOH2 | peroxy radical from IBUTALOH | $\text{CH}_3\text{C}(\text{OH})(\text{C}(\text{O})\text{OO})\text{CH}_3$ | 119.09 |
| IBUTALOOH | peroxide from IBUTALOH | $\text{CH}_3\text{C}(\text{OH})(\text{C}(\text{O})\text{OOH})\text{CH}_3$ | 120.1039 |
| CO2H3CHO | aldehyde from isoprene oxidation | $\text{CH}_3\text{C}(\text{O})\text{CHOHCHO}$ | 102.08864 |
| CO2H3CO3 | peroxy radical from 2-hydroxy-3-oxobutanal | $\text{CH}_3\text{C}(\text{O})\text{CHOHC}(\text{O})\text{OO}$ | 133.08 |
| CO2H3CO3H | peroxide from CO2H3CO3 | $\text{CH}_3\text{C}(\text{O})\text{CHOHC}(\text{O})\text{OOH}$ | 134.09 |
| BIGALD1 | malealdehyde | $\text{C}_4\text{H}_4\text{O}_2$ | 84.07336 |
| MALO2 | peroxy radical from photolysis of BIGALD1 | $\text{C}_4\text{H}_3\text{O}_4$ | 115.06 |
| BIGALKANE | large alkanes | C_5H_{12} | 72.14 |
| MBO | methyl butenol | $\text{C}_5\text{H}_{10}\text{O}$ | 86.1323 |
| ALKO2 | peroxy radical from large alkanes | $\text{C}_5\text{H}_{11}\text{O}_2$ | 103.14 |
| ALKOH | alcohol from BIGALKANE | $\text{C}_5\text{H}_{12}\text{O}$ | 88.14 |
| ALKOOH | peroxide from large alkanes | $\text{C}_5\text{H}_{12}\text{O}_2$ | 104.14 |
| ALKNO3 | nitrate from BIGALKANE | $\text{C}_5\text{H}_{11}\text{NO}_3$ | 133.141 |
| MBOO2 | peroxy radical from MBO | $\text{C}_5\text{H}_{11}\text{O}_4$ | 135.13 |
| MBOOOH | peroxide from MBO | $\text{C}_5\text{H}_{12}\text{O}_4$ | 136.14 |
| MBONO3O2 | peroxy nitrate radical from MBO+NO3 | $\text{C}_5\text{H}_{10}\text{NO}_6$ | 180.13 |
| ISOPBO2 | beta-hydroperoxy-radical from isoprene (1-OH-2-OO) | $\text{HOCH}_2\text{C}(\text{OO})(\text{CH}_3)\text{CH}=\text{CH}_2$ | 117.12 |
| ISOPDO2 | beta-hydroperoxy-radical from isoprene (4-OH-3-OO) | $\text{CH}_2=\text{C}(\text{CH}_3)\text{CH}(\text{OO})\text{CH}_2\text{OH}$ | 117.12 |
| LISOPACO2 | delta-hydroperoxy-radical from isoprene | $\text{C}_5\text{H}_9\text{O}_3$ | 117.12 |
| NISOPO2 | nitro-peroxy-radical from C5H8+NO3 | $\text{O}_2\text{NOCH}_2\text{C}(\text{CH}_3)=\text{CHCH}_2\text{O}_2$ | 162.12 |
| LISOPACOOH | peroxide from LISOPACO2 | $\text{CH}_3\text{C}(\text{CH}_2\text{OH})=\text{CHCH}_2\text{OOH}$ | 118.13 |
| LISOPACNO3 | nitrate from LISOPACO2 | $\text{CH}_3\text{C}(\text{CH}_2\text{OH})=\text{CHCH}_2\text{ONO}_2$ | 147.1293 |
| LHC4ACCHO | carbonyls from LISOPACO2 | $\text{HOCH}_2\text{C}(\text{H})=\text{C}(\text{CH}_3)\text{CHO}$ | 100.11 |
| LIEPOX | epoxides from LISOPACOOH | $\text{CH}_3\text{C}(\text{CH}_2\text{OH})\text{OHCHOCH}_2$ | 118.13 |
| LIECHO | aldehyde from LIEPOX | $\text{CH}_3\text{C}(\text{CHO})\text{OHCHOCH}_2$ | 116.11 |
| LIECO3 | peroxy radical from LIECHO | $\text{CH}_3\text{C}(\text{C}(\text{O})\text{OO})\text{OHCHOCH}_2$ | 147.1 |
| IEC1O2 | peroxy radical from LIEPOX+OH | $\text{HOCH}_2\text{C}(\text{O})\text{C}(\text{CH}_3)(\text{CH}_2\text{OH})\text{OO}$ | 149.12 |
| LIECO3H | peroxide from LIECO3 | $\text{C}_5\text{H}_8\text{O}_5$ | 148.11 |
| ISOPAOH | 2-methyl-2-butene-1,4-diol | $\text{HOCH}_2\text{C}(\text{CH}_3)=\text{CHCH}_2\text{OH}$ | 102.1317 |
| ISOPBOH | 2-methyl-3-butene-1,2-diol | $\text{HOCH}_2\text{C}(\text{CH}_3)\text{OHCH}=\text{CH}_2$ | 102.1317 |

Table S1: Chemical species in JAM2 (... continued)

| Species | Name | Structure | Moleweight |
|------------|--|---|------------|
| ISOPDOH | 3-methyl-3-butene-1,2-diol | $\text{CH}_2=\text{C}(\text{CH}_3)\text{CHOHCH}_2\text{OH}$ | 102.1317 |
| ISOPBOOH | peroxide from ISOPBO2 | $\text{HOCH}_2\text{C}(\text{CH}_3)(\text{OOH})\text{CH}=\text{CH}_2$ | 118.13 |
| ISOPDOOH | peroxide from ISOPDO2 | $\text{CH}_2=\text{C}(\text{CH}_3)\text{CHOOHCH}_2\text{OH}$ | 118.13 |
| ISOPBNO3 | alkyl nitrate from ISOPBO2 | $\text{HOCH}_2\text{C}(\text{CH}_3)\text{ONO}_2\text{CH}=\text{CH}_2$ | 147.13 |
| ISOPDNO3 | alkyl nitrate from ISOPDNO3 | $\text{CH}_2=\text{C}(\text{CH}_3)\text{CHONO}_2\text{CH}_2\text{OH}$ | 147.13 |
| HCOC5 | ketone from isoprene oxidation | $\text{CH}_2=\text{C}(\text{CH}_3)\text{C}(\text{O})\text{CH}_2\text{OH}$ | 100.11 |
| NISOPOOH | nitro-hydro-peroxide | $\text{O}_2\text{NOCH}_2\text{C}(\text{CH}_3)=\text{CHCH}_2\text{OOH}$ | 163.13 |
| NC4CHO | nitro-aldehyde | $\text{O}_2\text{NOCH}_2\text{C}(\text{CH}_3)=\text{CHCHO}$ | 145.11338 |
| LNISO3 | nitro-peroxy-radical | $\text{C}_5\text{H}_6\text{NO}_6$ | 176.1 |
| LNISOOH | lumped nitro-peroxide from LNISO3 | $\text{C}_5\text{H}_7\text{NO}_6$ | 177.11 |
| LHC4ACCO3 | acyl peroxy radical from LHC4ACCHO | $\text{C}_5\text{H}_7\text{O}_4$ | 131.1 |
| LHC4ACCO2H | hydroxy-methylbutenoic acid | $\text{C}_5\text{H}_8\text{O}_3$ | 116.11 |
| LHC4ACCO3H | peroxide from LHC4ACCO3 | $\text{C}_5\text{H}_8\text{O}_4$ | 132.11 |
| LC578O2 | peroxy radicals from C5-hydroxy aldehydes | $\text{C}_5\text{H}_9\text{O}_5$ | 149.12 |
| LC578OOH | lumped hydro-peroxide from LC578O2 | $\text{C}_5\text{H}_{10}\text{O}_5$ | 150.13 |
| LC5PAN1719 | homologues of PAN from LHC4ACCO3 | $\text{C}_5\text{H}_7\text{NO}_6$ | 177.11 |
| C59O2 | alkyl peroxy radicals from HCOC5 | $\text{C}_5\text{H}_9\text{O}_5$ | 149.12 |
| C59OOH | peroxide from C59O2 | $\text{HOCH}_2\text{C}(\text{CH}_3)(\text{OOH})\text{COCH}_2\text{OH}$ | 150.13 |
| DICARBO2 | dicarbonyl from photolysis of BIGALD2 | $\text{C}_5\text{H}_5\text{O}_4$ | 129.09 |
| BIGALD2 | acetylacrolein | $\text{CHOC}(\text{H})=\text{C}(\text{H})\text{C}(=\text{O})\text{CH}_3$ | 98.09994 |
| BIGALD3 | aldehyde from toluene and xylene oxidation | $\text{C}_5\text{H}_6\text{O}_2$ | 98.09994 |
| MDIALO2 | peroxy radical from photolysis of BIGALD3 | $\text{C}_5\text{H}_5\text{O}_4$ | 129.09 |
| BENZ | benzene | C_6H_6 | 78.11184 |
| TOL | toluene | C_7H_8 | 92.13842 |
| XYL | xylene | C_8H_{10} | 106.165 |
| PHENOL | carbolic acid | $\text{C}_6\text{H}_5\text{OH}$ | 94.11124 |
| BEPOMUC | benzene eopoxy diol | $\text{C}_6\text{H}_6\text{O}_3$ | 126.11 |
| BENZO2 | peroxy radical from benzene | $\text{C}_6\text{H}_7\text{O}_5$ | 159.11 |
| CATECHOL | pyrocatechol | $\text{C}_6\text{H}_6\text{O}_2$ | 110.11064 |
| PHENO2 | peroxy radical from phenol | $\text{C}_6\text{H}_7\text{O}_6$ | 175.11 |
| PHENOOH | peroxide from PHENO2 | $\text{C}_6\text{H}_8\text{O}_6$ | 176.12 |
| CATEC1O | peroxy radical from C6H5O | $\text{C}_6\text{H}_5\text{O}_2$ | 109.1 |
| C6H5OOH | peroxide from C6H5O2 | $\text{C}_6\text{H}_5\text{OOH}$ | 110.11 |
| BENZOOH | peroxide from BENZO2 | $\text{C}_6\text{H}_8\text{O}_5$ | 160.12 |
| BIGALD4 | aldehyde from xylene oxidation | $\text{C}_6\text{H}_8\text{O}_2$ | 112.12 |

Table S1: Chemical species in JAM2 (... continued)

| Species | Name | Structure | Moleweight |
|----------|---------------------------------------|------------------------|------------|
| TOLO2 | peroxy radical from toluene | $C_7H_9O_5$ | 173.14 |
| TOLOOH | peroxide from toluene | $C_7H_{10}O_5$ | 174.15 |
| CRESOL | cresol | C_7H_8O | 108.13782 |
| BZOO | peroxy radical from toluene | $C_6H_5CH_2OO$ | 123.13 |
| BZOOH | peroxide from BZOO | $C_6H_5CH_2OOH$ | 124.14 |
| BZALD | benzaldehyde | C_6H_5CHO | 106.12194 |
| ACBZO2 | acyl peroxy radical from benzaldehyde | $C_6H_5C(O)OO$ | 137.11 |
| PBZNIT | nitrate from benzaldehyde | $C_6H_5C(O)OONO_2$ | 183.12 |
| TEPOMUC | epoxide from toluene | $C_7H_8O_3$ | 140.13 |
| XYLOL | xylol | $C_6H_3(CH_3)(CH_3)OH$ | 122.16 |
| XYLOLO2 | peroxy radical from xylol | $C_8H_{11}O_6$ | 203.17 |
| XYLOLOOH | peroxide from xylol | $C_8H_{12}O_6$ | 204.17 |
| XYLENO2 | peroxy radical from xylene | $C_8H_{11}O_5$ | 187.17 |
| XYLENOOH | peroxide from XYLENO2 | $C_8H_{12}O_5$ | 188.17 |
| TERPROD2 | terpene oxidation product C9 | $C_7H_{10}O_2$ | 126.15 |
| APIN | α -pinene | $C_{10}H_{16}$ | 136.23404 |
| BPIN | β -pinene | $C_{10}H_{16}$ | 136.23404 |
| LIMON | limonene | $C_{10}H_{16}$ | 136.23404 |
| MYRC | myrcene | $C_{10}H_{16}$ | 136.23404 |
| BCARY | β -caryophyllene | $C_{15}H_{24}$ | 204.35106 |
| TERPO2 | peroxy radical from terpenes | $C_{10}H_{17}O_3$ | 185.23 |
| TERPOOH | peroxide from terpenes | $C_{10}H_{18}O_3$ | 186.24 |
| TERPNO3 | nitrate from terpenes | $C_{10}H_{17}NO_4$ | 215.24 |
| NTERPO2 | nitro peroxy radical from terpenes | $C_{10}H_{16}NO_5$ | 230.23 |
| TERPROD1 | terpene oxidation product C10 | $C_{10}H_{16}O_2$ | 168.23 |
| TERP2O2 | peroxy radical from TERPROD1 | $C_{10}H_{15}O_4$ | 199.22 |
| TERP2OOH | peroxide from TERP2O2 | $C_{10}H_{16}O_4$ | 200.23 |
| CF2CLBR | Halon-1211 | CF_2ClBr | 165.364506 |
| CF3BR | Halon-1301 | $CBrF_3$ | 148.90991 |
| CH3BR | methyl bromide | CH_3Br | 94.93852 |
| CH2BR2 | dibromomethane | CH_2Br_2 | 173.83 |
| CHBR3 | bromoform | $CHBr_3$ | 252.73 |
| HCFC22 | HCFC-22 | $CClF_2H$ | 86.468446 |
| CH3CL | chloromethane | CH_3Cl | 50.48752 |
| CCL4 | perchloromethane | CCl_4 | 153.8227 |

Table S1: Chemical species in JAM2 (... continued)

| Species | Name | Structure | Moleweight |
|-------------|---|--|------------|
| COF2 | CF2O | CF ₂ O | 66.006906 |
| COFCL | carbonic chloride fluoride | CClFO | 82.461503 |
| CFC11 | CFC11 | CCl ₃ F | 137.368103 |
| CFC12 | CFC12 | CCl ₂ F ₂ | 120.913506 |
| CFC113 | CFC113 | CCl ₂ FCClF ₂ | 187.37561 |
| CFC114 | CFC114 | CClF ₂ CClF ₂ | 170.92 |
| CFC115 | CFC115 | CClF ₂ CF ₃ | 154.47 |
| HCFC141B | HCFC-141b | CH ₃ CCl ₂ F | 116.949623 |
| HCFC142B | HCFC-142b | CH ₃ CClF ₂ | 100.495026 |
| CH3CCL3 | methylchloroform | CH ₃ CCl ₃ | 133.40422 |
| H1202 | Halon 1202 | CBr ₂ F ₂ | 209.815506 |
| H2402 | Halon 2402 | C ₂ Br ₂ F ₄ | 259.823013 |
| BRNO2 | bromonitrite | BrNO ₂ | 125.91 |
| CATEC1O2 | peroxy radical from catechol | C ₆ H ₅ O ₃ | 125.1 |
| CATEC1OOH | peroxide from catechol | C ₆ H ₆ O ₃ | 126.1 |
| CH3O2NO2 | methyl peroxy nitrate | CH ₃ NO ₄ | 93.04 |
| ELVOC | highly oxidized, low volatility compound from terpene oxidation | C ₁₀ H ₁₆ O ₈ | 264.2 |
| HCOCO3 | peroxy radical from glyoxal oxidation | C ₂ HO ₄ | 82.03 |
| HOCH2CO3 | peroxy radical from various molecules | C ₂ H ₃ O ₄ | 91.04 |
| HPALD | hydroperoxy aldehydes from isoprene oxidation | C ₅ H ₈ O ₃ | 116.12 |
| LISOPNO3NO3 | radical product from isoprene oxidation | C ₅ H ₁₀ O ₈ N ₂ | 226.1 |
| LISOPNO3O2 | radical product from isoprene oxidation | C ₅ H ₁₀ O ₇ N | 196.1 |
| LISOPNO3OOH | peroxide from isoprene oxidation | C ₅ H ₁₁ O ₇ N | 197.1 |
| LISOPOOHO2 | radical product from isoprene oxidation | C ₅ H ₁₁ O ₆ | 167.1 |
| LISOPOOHOOH | peroxide from isoprene oxidation | C ₅ H ₁₂ O ₆ | 168.1 |
| MACRN | nitrate from methacrolein | C ₄ H ₇ NO ₅ | 149.1 |
| MACROOH | peroxide from methacrolein | C ₄ H ₈ O ₄ | 120.1 |
| MEKNO3 | nitrate from methyl ethyl ketone | C ₄ H ₇ O ₅ N | 149.1 |
| MVKN | nitrate from methyl vinyl ketone | C ₄ H ₇ NO ₅ | 149.1 |
| NTERPNO3 | nitrate from terpene oxidation | C ₁₀ H ₁₇ NO ₅ | 231.2 |
| PACALD | aldehyde from isoprene oxidation | C ₅ H ₆ O ₄ | 130.1 |

Table S2: Tropospheric photolysis reactions

| reaction | reference |
|--|-----------|
| $\text{O}_3 + h\nu \longrightarrow \text{O}_1\text{D} + \text{O}_2$ | |
| $\text{O}_3 + h\nu \longrightarrow \text{O} + \text{O}_2$ | |
| $\text{H}_2\text{O}_2 + h\nu \longrightarrow 2 \cdot \text{OH}$ | |
| $\text{N}_2\text{O} + h\nu \longrightarrow \text{N}_2 + \text{O}_1\text{D}$ | |
| $\text{NO} + h\nu \longrightarrow \text{N} + \text{O}$ | |
| $\text{NO}_2 + h\nu \longrightarrow \text{NO} + \text{O}$ | |
| $\text{NO}_3 + h\nu \longrightarrow \text{NO}_2 + \text{O}$ | |
| $\text{NO}_3 + h\nu \longrightarrow \text{NO} + \text{O}_2$ | |
| $\text{HNO}_3 + h\nu \longrightarrow \text{NO}_2 + \text{OH}$ | |
| $\text{HONO} + h\nu \longrightarrow \text{NO} + \text{OH}$ | |
| $\text{HO}_2\text{NO}_2 + h\nu \longrightarrow \text{NO}_3 + \text{OH}$ | |
| $\text{HO}_2\text{NO}_2 + h\nu \longrightarrow \text{HO}_2 + \text{NO}_2$ | |
| $\text{N}_2\text{O}_5 + h\nu \longrightarrow \text{NO}_2 + \text{NO}_3$ | |
| $\text{N}_2\text{O}_5 + h\nu \longrightarrow \text{NO} + \text{O} + \text{NO}_3$ | |
| $\text{CO}_2 + h\nu \longrightarrow \text{CO} + \text{O}$ | |
| $\text{CH}_4 + h\nu \longrightarrow \text{CH}_3\text{O}_2 + \text{H}$ | |
| $\text{CH}_4 + h\nu \longrightarrow 1.44 \cdot \text{H}_2 + 0.18 \cdot \text{CH}_2\text{O} + 0.18 \cdot \text{O} + 0.66 \cdot \text{OH} + 0.44 \cdot \text{CO}_2 + 0.38 \cdot \text{CO} + 0.05 \cdot \text{H}_2\text{O}$ | |
| $\text{CH}_2\text{O} + h\nu \longrightarrow \text{CO} + 2 \cdot \text{H}$ | |
| $\text{CH}_2\text{O} + h\nu \longrightarrow \text{CO} + \text{H}_2$ | |
| $\text{CH}_3\text{OOH} + h\nu \longrightarrow \text{CH}_2\text{O} + \text{H} + \text{OH}$ | |
| $\text{CH}_3\text{O}_2\text{NO}_2 + h\nu \longrightarrow \text{HO}_2 + \text{NO}_3 + \text{HCHO}$ | |
| $\text{CH}_3\text{O}_2\text{NO}_2 + h\nu \longrightarrow \text{CH}_3\text{O}_2 + \text{NO}_2$ | |
| $\text{CH}_3\text{CHO} + h\nu \longrightarrow \text{CH}_3\text{O}_2 + \text{CO} + \text{HO}_2$ | |
| $\text{CH}_3\text{COOOH} + h\nu \longrightarrow \text{CH}_3\text{O}_2 + \text{OH} + \text{CO}_2$ | |
| $\text{C}_2\text{H}_5\text{OOH} + h\nu \longrightarrow \text{CH}_3\text{CHO} + \text{HO}_2 + \text{OH}$ | |
| $\text{PAN} + h\nu \longrightarrow 0.6 \cdot \text{CH}_3\text{CO}_3 + 0.6 \cdot \text{NO}_2 + 0.4 \cdot \text{CH}_3\text{O}_2 + 0.4 \cdot \text{NO}_3 + 0.4 \cdot \text{CO}_2$ | |
| $\text{EOOH} + h\nu \longrightarrow \text{EO} + \text{OH}$ | |
| $\text{GLYOXAL} + h\nu \longrightarrow 2 \cdot \text{CO} + 2 \cdot \text{HO}_2$ | |
| $\text{GLYALD} + h\nu \longrightarrow 2 \cdot \text{HO}_2 + \text{CO} + \text{CH}_2\text{O}$ | |
| $\text{HOCH}_2\text{CO}_3\text{H} + h\nu \longrightarrow \text{CH}_2\text{O} + \text{HO}_2 + \text{OH} + \text{CO}_2$ | |
| $\text{HCOCO}_2\text{H} + h\nu \longrightarrow 2 \cdot \text{HO}_2 + \text{CO} + \text{CO}_2$ | |
| $\text{HCOCO}_3\text{H} + h\nu \longrightarrow \text{CO} + \text{HO}_2 + \text{OH} + \text{CO}_2$ | |
| $\text{CH}_3\text{COCH}_3 + h\nu \longrightarrow \text{CH}_3\text{CO}_3 + \text{CH}_3\text{O}_2$ | |

Table S2: Tropospheric photolysis reactions (... continued)

| reaction | reference |
|--|--|
| $\text{C}_3\text{H}_7\text{OOH} + h\nu \longrightarrow 0.82 \cdot \text{CH}_3\text{COCH}_3 + \text{OH} + \text{HO}_2 + 0.27 \cdot \text{CH}_3\text{CHO}$ | |
| $\text{POOH} + h\nu \longrightarrow \text{CH}_2\text{O} + \text{CH}_3\text{CHO} + \text{HO}_2 + \text{OH}$ | |
| $\text{HYAC} + h\nu \longrightarrow \text{CH}_3\text{CO}_3 + \text{HO}_2 + \text{CH}_2\text{O}$ | |
| $\text{CH}_3\text{COCHO} + h\nu \longrightarrow \text{CH}_3\text{CO}_3 + \text{CO} + \text{HO}_2$ | |
| $\text{ROOH} + h\nu \longrightarrow \text{CH}_2\text{O} + \text{CH}_3\text{CO}_3 + \text{OH}$ | |
| $\text{PR}_2\text{O}_2\text{HNO}_3 + h\nu \longrightarrow 0.83 \cdot \text{HO}_2 + 0.83 \cdot \text{NOA} + 0.17 \cdot \text{CH}_2\text{O} + 0.17 \cdot \text{CH}_3\text{CHO} + \text{OH}$ | |
| $\text{NOA} + h\nu \longrightarrow \text{CH}_2\text{O} + \text{CH}_3\text{CO}_3 + \text{NO}_2$ | Müller et al. (2014) |
| $\text{MEK} + h\nu \longrightarrow \text{C}_2\text{H}_5\text{O}_2 + \text{CH}_3\text{CO}_3$ | |
| $\text{MEKOOH} + h\nu \longrightarrow \text{CH}_3\text{CO}_3 + \text{OH} + \text{CH}_3\text{CHO}$ | |
| $\text{MEKNO}_3 + h\nu \longrightarrow \text{CH}_3\text{CHO} + \text{CH}_3\text{CO}_3 + \text{NO}_2$ | Müller et al. (2014) |
| $\text{MACR} + h\nu \longrightarrow \text{HO}_2 + 0.5 \cdot \text{MCO}_3 + 0.5 \cdot \text{CH}_2\text{O} + 0.5 \cdot \text{CH}_3\text{CO}_3 + 0.5 \cdot \text{CO}$ | |
| $\text{MACR} + h\nu \longrightarrow \text{HO}_2 + 0.5 \cdot \text{MCO}_3 + 0.5 \cdot \text{CH}_2\text{O} + 0.5 \cdot \text{CH}_3\text{CO}_3 + 0.5 \cdot \text{CO}$ | |
| $\text{MACROOH} + h\nu \longrightarrow \text{HO}_2 + \text{HYAC} + \text{OH} + \text{CO}$ | |
| $\text{MACROH} + h\nu \longrightarrow \text{CO} + \text{HYAC} + 2 \cdot \text{HO}_2 + \text{H}_2\text{O}$ | |
| $\text{MPAN} + h\nu \longrightarrow \text{MCO}_3 + \text{NO}_2$ | |
| $\text{MACO}_3\text{H} + h\nu \longrightarrow \text{CH}_2\text{O} + \text{CH}_3\text{CO}_3 + \text{OH} + \text{CO}_2$ | |
| $\text{MVK} + h\nu \longrightarrow 0.5 \cdot \text{C}_3\text{H}_6 + 0.5 \cdot \text{CH}_3\text{CO}_3 + 0.5 \cdot \text{CH}_2\text{O} + \text{CO} + 0.5 \cdot \text{HO}_2$ | |
| $\text{LHMKABOOH} + h\nu \longrightarrow 0.3 \cdot \text{CH}_3\text{COCHO} + \text{OH} + 0.3 \cdot \text{CH}_2\text{O} + 0.3 \cdot \text{HO}_2 + 0.7 \cdot \text{CH}_3\text{CO}_3 + 0.7 \cdot \text{GLYALD}$ | |
| $\text{MVKN} + h\nu \longrightarrow \text{CH}_3\text{CO}_3 + \text{GLYALD} + \text{NO}_2$ | Müller et al. (2014) |
| $\text{MACRN} + h\nu \longrightarrow \text{CO} + \text{HYAC} + \text{HO}_2 + \text{NO}_2$ | Müller et al. (2014) |
| $\text{CO}_2\text{H}_3\text{CHO} + h\nu \longrightarrow \text{CH}_3\text{COCHO} + \text{CO} + 2 \cdot \text{HO}_2$ | |
| $\text{CO}_2\text{H}_3\text{CO}_3\text{H} + h\nu \longrightarrow \text{CH}_3\text{COCHO} + \text{HO}_2 + \text{OH} + \text{CO}_2$ | |
| $\text{BIACETOH} + h\nu \longrightarrow \text{CH}_3\text{CO}_3 + \text{HOCH}_2\text{CO}_3$ | |
| $\text{ALKOOH} + h\nu \longrightarrow 0.4 \cdot \text{CH}_3\text{CHO} + 0.25 \cdot \text{CH}_2\text{O} + 0.25 \cdot \text{CH}_3\text{COCH}_3 + \text{HO}_2 + 0.8 \cdot \text{MEK} + \text{OH}$ | |
| $\text{ALKNO}_3 + h\nu \longrightarrow 0.4 \cdot \text{CH}_3\text{CHO} + 0.25 \cdot \text{CH}_2\text{O} + 0.25 \cdot \text{CH}_3\text{COCH}_3 + \text{HO}_2 + 0.8 \cdot \text{MEK} + \text{NO}_2$ | |
| $\text{LISOPACOOH} + h\nu \longrightarrow \text{HO}_2 + \text{LHC}_4\text{ACCHO} + \text{OH}$ | |
| $\text{LISOPACNO}_3 + h\nu \longrightarrow \text{HO}_2 + \text{LHC}_4\text{ACCHO} + \text{NO}_2$ | |
| $\text{HPALD} + h\nu \longrightarrow \text{LHC}_4\text{ACCO}_3 + \text{OH}$ | D. Taraborrelli: $J(\text{MACR})/\phi(\text{MACR}) = 2 \cdot j_{\text{macr_a}}/0.004 = 500 \cdot j_{\text{macr_a}}$ |
| $\text{PACALD} + h\nu \longrightarrow \text{OH} + 0.5 \cdot \text{HO}_2 + 0.5 \cdot \text{CO} + 0.5 \cdot \text{CH}_3\text{COCHO} + 0.5 \cdot \text{GLYOX} + 0.5 \cdot \text{CH}_3\text{CO}_3$ | D. Taraborrelli: average of product yields of C5PACALD 1 and 2 in MCMv3.3.1, $2 \cdot J(\text{MACR})/\phi(\text{MACR}) = 2 \cdot j_{\text{hpald}} = 1000 \cdot j_{\text{macr_a}}$ |

Table S2: Tropospheric photolysis reactions (... continued)

| reaction | reference |
|--|----------------------|
| $\text{LIECHO} + h\nu \longrightarrow \text{CO} + \text{HO}_2 + 0.6 \cdot \text{LHMKABO}_2 + 0.4 \cdot \text{MACRO}_2$ | |
| $\text{LIECO}_3\text{H} + h\nu \longrightarrow 0.6 \cdot \text{LHMKABO}_2 + 0.4 \cdot \text{MACRO}_2 + \text{CO}_2 + \text{OH}$ | |
| $\text{ISOPBOOH} + h\nu \longrightarrow \text{CH}_2\text{O} + \text{MVK} + \text{HO}_2 + \text{OH}$ | |
| $\text{ISOPBNO}_3 + h\nu \longrightarrow \text{CH}_2\text{O} + \text{MVK} + \text{HO}_2 + \text{NO}_2$ | |
| $\text{ISOPDOOH} + h\nu \longrightarrow \text{CH}_2\text{O} + \text{MACR} + \text{HO}_2 + \text{OH}$ | |
| $\text{ISOPDNO}_3 + h\nu \longrightarrow \text{CH}_2\text{O} + \text{MACR} + \text{HO}_2 + \text{NO}_2$ | |
| $\text{NISOPOOH} + h\nu \longrightarrow \text{HO}_2 + \text{NC}_4\text{CHO} + \text{OH}$ | |
| $\text{NC}_4\text{CHO} + h\nu \longrightarrow \text{LHC}_4\text{ACCO}_3 + \text{NO}_2$ | Müller et al. (2014) |
| $\text{LNISOOH} + h\nu \longrightarrow \text{NOA} + \text{OH} + 0.5 \cdot \text{GLYOXAL} + 0.5 \cdot \text{CO} + \text{HO}_2 + 0.5 \cdot \text{CO}_2$ | |
| $\text{LHC}_4\text{ACCHO} + h\nu \longrightarrow 0.5 \cdot \text{LHC}_4\text{ACCO}_3 + 0.25 \cdot \text{HYAC} + 0.25 \cdot \text{GLYALD} + 0.25 \cdot \text{CH}_3\text{CO}_3 + 0.75 \cdot \text{CO} + 1.25 \cdot \text{HO}_2$ | |
| $\text{LC}_{578}\text{OOH} + h\nu \longrightarrow 0.5 \cdot \text{HYAC} + 0.5 \cdot \text{CH}_3\text{COCHO} + 0.5 \cdot \text{GLYOXAL} + 0.5 \cdot \text{GLYALD} + \text{HO}_2 + \text{OH}$ | |
| $\text{LHC}_4\text{ACCO}_3\text{H} + h\nu \longrightarrow 0.5 \cdot \text{HYAC} + 0.5 \cdot \text{GLYALD} + 0.5 \cdot \text{CH}_3\text{CO}_3 + 0.5 \cdot \text{CO} + 0.5 \cdot \text{HO}_2 + \text{OH} + \text{CO}_2$ | |
| $\text{HCOC}_5 + h\nu \longrightarrow \text{CH}_2\text{O} + \text{CH}_3\text{CO}_3 + \text{HOCH}_2\text{CO}_3$ | |
| $\text{C}_{59}\text{OOH} + h\nu \longrightarrow \text{HOCH}_2\text{CO}_3 + \text{HYAC} + \text{NO}_2 + \text{OH}$ | |
| $\text{LISOPOOHOOH} + h\nu \longrightarrow 0.25 \cdot \text{CH}_3\text{COCHO} + 0.25 \cdot \text{GLYALD} + 0.25 \cdot \text{GLYOXAL} + 0.25 \cdot \text{HYAC} + 0.25 \cdot \text{CO} + 0.25 \cdot \text{MACROH} + 0.25 \cdot \text{CH}_2\text{O} + 0.25 \cdot \text{CO}_2\text{H}_3\text{CHO} + \text{HO}_2$ | |
| $\text{LISOPNO}_3\text{OOH} + h\nu \longrightarrow \text{HOCH}_2\text{CO}_3 + \text{HYAC} + \text{NO}_2 + \text{OH}$ | |
| $\text{LISOPNO}_3\text{NO}_3 + h\nu \longrightarrow \text{HOCH}_2\text{CO}_3 + \text{HYAC} + \text{NO}_2 + \text{NO}_2$ | |
| $\text{MBOOOH} + h\nu \longrightarrow \text{HO}_2 + \text{OH} + 0.67 \cdot \text{GLYALD} + 0.67 \cdot \text{CH}_3\text{COCH}_3 + 0.33 \cdot \text{IBUTALOH} + 0.33 \cdot \text{CH}_2\text{O}$ | |
| $\text{IBUTALOH} + h\nu \longrightarrow 2 \cdot \text{HO}_2 + \text{CO} + \text{CH}_3\text{COCH}_3$ | |
| $\text{IBUTALOOH} + h\nu \longrightarrow \text{CH}_3\text{COCH}_3 + \text{OH} + \text{CO}_2 + \text{HO}_2$ | |
| $\text{BEPOMUC} + h\nu \longrightarrow \text{BIGALD}_1 + 1.5 \cdot \text{HO}_2 + 1.5 \cdot \text{CO}$ | |
| $\text{BIGALD}_1 + h\nu \longrightarrow 0.6 \cdot \text{MALO}_2 + \text{HO}_2$ | |
| $\text{TOLOOH} + h\nu \longrightarrow \text{OH} + 0.6 \cdot \text{GLYOXAL} + 0.4 \cdot \text{CH}_3\text{COCHO} + \text{HO}_2 + 0.2 \cdot \text{BIGALD}_1 + 0.2 \cdot \text{BIGALD}_2 + 0.2 \cdot \text{BIGALD}_3$ | |
| $\text{TEPOMUC} + h\nu \longrightarrow 0.5 \cdot \text{CH}_3\text{CO}_3 + \text{HO}_2 + 1.5 \cdot \text{CO}$ | |
| $\text{CATEC}_1\text{OOH} + h\nu \longrightarrow \text{CATEC}_1\text{O} + \text{OH}$ | |
| $\text{BIGALD}_2 + h\nu \longrightarrow 0.6 \cdot \text{DICARBO}_2 + 0.6 \cdot \text{HO}_2$ | |
| $\text{BIGALD}_3 + h\nu \longrightarrow 0.6 \cdot \text{CO} + 0.6 \cdot \text{HO}_2 + 0.6 \cdot \text{MDIALO}_2$ | |
| $\text{BIGALD}_4 + h\nu \longrightarrow \text{CO} + \text{HO}_2 + \text{CH}_3\text{COCHO} + \text{CH}_3\text{CO}_3$ | |
| $\text{TERPOOH} + h\nu \longrightarrow 0.4 \cdot \text{CH}_2\text{O} + 0.05 \cdot \text{CH}_3\text{COCH}_3 + 0.945 \cdot \text{TERPROD}_1 + \text{HO}_2 + \text{OH}$ | |

Table S2: Tropospheric photolysis reactions (... continued)

| reaction | reference |
|---|-----------|
| $\text{TERPROD}_1 + h\nu \longrightarrow \text{CO} + \text{HO}_2 + \text{TERPROD}_2$ | |
| $\text{TERP}_2\text{OOH} + h\nu \longrightarrow \text{OH} + 0.372 \cdot \text{CH}_2\text{O} + 0.3 \cdot \text{CH}_3\text{COCH}_3 + 0.25 \cdot \text{CO} + \text{CO}_2 +$ | |
| $\text{TERPROD}_2 + \text{HO}_2 + 0.25 \cdot \text{GLYALD}$ | |
| $\text{TERPROD}_2 + h\nu \longrightarrow 0.15 \cdot \text{CH}_3\text{COCH}_2\text{O}_2 + 0.68 \cdot \text{CH}_2\text{O} + 0.8 \cdot \text{CO}_2 + 0.5 \cdot \text{CH}_3\text{COCH}_3 +$ | |
| $1.2 \cdot \text{HO}_2 + 1.7 \cdot \text{CO}$ | |
| $\text{ISOPBNO}_3 + h\nu \longrightarrow \text{HO}_2 + \text{NO}_2 + \text{TERPROD}_1$ | |
| $\text{NTERPNO}_3 + h\nu \longrightarrow \text{NO}_2 + \text{OH} + \text{TERPROD}_1$ | |
| $\text{ELVOC} + h\nu \longrightarrow \text{HO}_2 + \text{OH} + \text{TERPROD}_2$ | |

Table S3: Stratospheric photolysis reactions

| reaction | reference |
|---|--|
| $\text{O}_2 + h\nu \longrightarrow \text{O} + \text{O}_1\text{D}$ | |
| $\text{O}_2 + h\nu \longrightarrow 2 \cdot \text{O}$ | |
| $\text{H}_2\text{O} + h\nu \longrightarrow \text{H} + \text{OH}$ | |
| $\text{H}_2\text{O} + h\nu \longrightarrow \text{H}_2 + \text{O}_1\text{D}$ | |
| $\text{H}_2\text{O} + h\nu \longrightarrow 2 \cdot \text{H} + \text{O}$ | |
| $\text{CL}_2 + h\nu \longrightarrow 2 \cdot \text{CL}$ | |
| $\text{CL}_2\text{O}_2 + h\nu \longrightarrow 2 \cdot \text{CL}$ | |
| $\text{CLO} + h\nu \longrightarrow \text{CL} + \text{O}$ | |
| $\text{HCL} + h\nu \longrightarrow \text{CL} + \text{H}$ | |
| $\text{HOCL} + h\nu \longrightarrow \text{CL} + \text{OH}$ | |
| $\text{CLONO}_2 + h\nu \longrightarrow \text{CL} + \text{NO}_3$ | |
| $\text{CLONO}_2 + h\nu \longrightarrow \text{CLO} + \text{NO}_2$ | |
| $\text{OCLO} + h\nu \longrightarrow \text{CLO} + \text{O}$ | |
| $\text{BRO} + h\nu \longrightarrow \text{BR} + \text{O}$ | |
| $\text{HBR} + h\nu \longrightarrow \text{BR} + \text{H}$ | |
| $\text{HOBR} + h\nu \longrightarrow \text{BR} + \text{OH}$ | |
| $\text{BRONO} + h\nu \longrightarrow \text{BR} + \text{NO}_2$ | 50% branching ratio assigned to both possible channels. Cross-sections consistent with Burkholder and Orlando (2000) |
| $\text{BRONO} + h\nu \longrightarrow \text{BRO} + \text{NO}$ | 50% branching ratio assigned to both possible channels. Cross-sections consistent with Burkholder and Orlando (2000) |
| $\text{BRNO}_2 + h\nu \longrightarrow \text{BR} + \text{NO}_2$ | |
| $\text{BRONO}_2 + h\nu \longrightarrow \text{BR} + \text{NO}_3$ | |

Table S3: Stratospheric photolysis reactions (... continued)

| reaction | reference |
|--|-----------|
| $\text{BRONO}_2 + \text{h}\nu \longrightarrow \text{BRO} + \text{NO}_2$ | |
| $\text{BR}_2 + \text{h}\nu \longrightarrow 2 \cdot \text{BR}$ | |
| $\text{BRCL} + \text{h}\nu \longrightarrow \text{BR} + \text{CL}$ | |
| $\text{HF} + \text{h}\nu \longrightarrow \text{F} + \text{H}$ | |
| $\text{SF}_6 + \text{h}\nu \longrightarrow$ | |
| $\text{CH}_3\text{BR} + \text{h}\nu \longrightarrow \text{BR} + \text{CH}_3\text{O}_2$ | |
| $\text{CH}_2\text{BR}_2 + \text{h}\nu \longrightarrow 2 \cdot \text{BR}$ | |
| $\text{CHBR}_3 + \text{h}\nu \longrightarrow 3 \cdot \text{BR}$ | |
| $\text{CH}_3\text{CL} + \text{h}\nu \longrightarrow \text{CH}_3\text{O}_2 + \text{CL}$ | |
| $\text{CH}_3\text{CCL}_3 + \text{h}\nu \longrightarrow 3 \cdot \text{CL}$ | |
| $\text{CF}_3\text{BR} + \text{h}\nu \longrightarrow \text{BR} + \text{F} + \text{COF}_2$ | |
| $\text{CF}_2\text{CLBR} + \text{h}\nu \longrightarrow \text{BR} + \text{CL} + \text{COF}_2$ | |
| $\text{CCL}_4 + \text{h}\nu \longrightarrow 4 \cdot \text{CL} + \text{CO}_2$ | |
| $\text{CFC}_{11} + \text{h}\nu \longrightarrow 2 \cdot \text{CL} + \text{COFCL}$ | |
| $\text{CFC}_{12} + \text{h}\nu \longrightarrow 2 \cdot \text{CL} + \text{COF}_2$ | |
| $\text{CFC}_{113} + \text{h}\nu \longrightarrow 2 \cdot \text{CL} + \text{COFCL} + \text{COF}_2$ | |
| $\text{CFC}_{114} + \text{h}\nu \longrightarrow 2 \cdot \text{CL} + 2 \cdot \text{COF}_2$ | |
| $\text{CFC}_{115} + \text{h}\nu \longrightarrow \text{CL} + \text{F} + 2 \cdot \text{COF}_2$ | |
| $\text{HCFC}_{22} + \text{h}\nu \longrightarrow \text{CL} + \text{COF}_2$ | |
| $\text{HCFC}_{141}\text{B} + \text{h}\nu \longrightarrow \text{CL} + \text{COFCL}$ | |
| $\text{HCFC}_{142}\text{B} + \text{h}\nu \longrightarrow \text{CL} + \text{COF}_2$ | |
| $\text{H}_{1202} + \text{h}\nu \longrightarrow 2 \cdot \text{BR} + \text{COF}_2$ | |
| $\text{H}_{2402} + \text{h}\nu \longrightarrow 2 \cdot \text{BR} + 2 \cdot \text{COF}_2$ | |
| $\text{COF}_2 + \text{h}\nu \longrightarrow 2 \cdot \text{F}$ | |
| $\text{COFCL} + \text{h}\nu \longrightarrow \text{CL} + \text{F}$ | |

Table S4: (Tropospheric) O_x reactions

| reaction | rate coefficient | reference |
|--|---------------------------------------|-------------------|
| $\text{O} + \text{O}_2 + \text{M} \longrightarrow \text{M} + \text{O}_3$ | O_{O_2} | |
| $\text{O} + \text{O}_3 \longrightarrow 2 \cdot \text{O}_2$ | $8.000 \cdot 10^{-12} \exp(-2060./T)$ | JPL (2011) |
| $\text{O} + \text{O} + \text{M} \longrightarrow \text{M} + \text{O}_2$ | O_{O} | not in JPL (2011) |
| $\text{O}_1\text{D} + \text{N}_2 \longrightarrow \text{N}_2 + \text{O}$ | $2.150 \cdot 10^{-11} \exp(110./T)$ | |
| $\text{O}_1\text{D} + \text{O}_2 \longrightarrow \text{O} + \text{O}_2$ | $3.135 \cdot 10^{-11} \exp(55./T)$ | |

Table S4: (Tropospheric) O_x reactions (... continued)

| reaction | rate coefficient | reference |
|---|------------------------------------|------------|
| $\text{O}_1\text{D} + \text{O}_2 \longrightarrow \text{O} + \text{O}_2$ | $1.650 \cdot 10^{-12} \exp(55./T)$ | |
| $\text{O}_1\text{D} + \text{H}_2\text{O} \longrightarrow 2 \cdot \text{OH}$ | $1.630 \cdot 10^{-10} \exp(60./T)$ | JPL (2011) |

Table S5: (Tropospheric) HO_x reactions

| reaction | rate coefficient | reference |
|---|---|--------------|
| $\text{H} + \text{O}_2 + \text{M} \longrightarrow \text{HO}_2 + \text{M}$ | <i>ktroe</i> ($4.400 \cdot 10^{-32}, 1.3, 7.500 \cdot 10^{-11}, -0.2, 0.6$) | JPL (2011) |
| $\text{H} + \text{O}_3 \longrightarrow \text{O}_2 + \text{OH}$ | $1.400 \cdot 10^{-10} \exp(-470./T)$ | JPL (2011) |
| $\text{H} + \text{HO}_2 \longrightarrow 2 \cdot \text{OH}$ | $7.200 \cdot 10^{-11}$ | JPL (2011) |
| $\text{H} + \text{HO}_2 \longrightarrow \text{H}_2\text{O} + \text{O}$ | $1.600 \cdot 10^{-12}$ | JPL (2011) |
| $\text{H} + \text{HO}_2 \longrightarrow \text{H}_2 + \text{O}_2$ | $6.900 \cdot 10^{-12}$ | JPL (2011) |
| $\text{H}_2 + \text{O} \longrightarrow \text{H} + \text{OH}$ | $1.600 \cdot 10^{-11} \exp(-4570./T)$ | |
| $\text{H}_2 + \text{OH} \longrightarrow \text{H} + \text{H}_2\text{O}$ | $2.800 \cdot 10^{-12} \exp(-1800./T)$ | JPL (2011) |
| $\text{OH} + \text{O} \longrightarrow \text{H} + \text{O}_2$ | $1.800 \cdot 10^{-11} \exp(180./T)$ | JPL (2011) |
| $\text{OH} + \text{OH} \longrightarrow \text{H}_2\text{O} + \text{O}$ | $1.800 \cdot 10^{-12}$ | JPL (2011) |
| $\text{OH} + \text{OH} + \text{M} \longrightarrow \text{H}_2\text{O}_2 + \text{M}$ | <i>ktroe</i> ($6.900 \cdot 10^{-31}, 1., 2.600 \cdot 10^{-11}, 0., 0.6$) | JPL (2011) |
| $\text{OH} + \text{O}_3 \longrightarrow \text{HO}_2 + \text{O}_2$ | $1.700 \cdot 10^{-12} \exp(-940./T)$ | JPL (2011) |
| $\text{HO}_2 + \text{O} \longrightarrow \text{O}_2 + \text{OH}$ | $3.000 \cdot 10^{-11} \exp(200./T)$ | JPL (2011) |
| $\text{HO}_2 + \text{OH} \longrightarrow \text{H}_2\text{O} + \text{O}_2$ | $4.800 \cdot 10^{-11} \exp(250./T)$ | JPL (2011) |
| $\text{HO}_2 + \text{O}_3 \longrightarrow \text{OH} + 2 \cdot \text{O}_2$ | | IUPAC (2004) |
| $\text{HO}_2 + \text{HO}_2 \longrightarrow \text{H}_2\text{O}_2 + \text{O}_2$ | HO ₂ –HO ₂ | |
| $\text{H}_2\text{O}_2 + \text{O} \longrightarrow \text{HO}_2 + \text{OH}$ | $1.400 \cdot 10^{-12} \exp(-2000./T)$ | JPL (2011) |
| $\text{H}_2\text{O}_2 + \text{OH} \longrightarrow \text{H}_2\text{O} + \text{HO}_2$ | $1.800 \cdot 10^{-12}$ | JPL (2011) |

Table S6: NO_x reactions

| reaction | rate coefficient | reference |
|--|---|-----------------------------------|
| $\text{N} + \text{OH} \longrightarrow \text{H} + \text{NO}$ | $5.000 \cdot 10^{-11}$ | |
| $\text{N} + \text{O}_2 \longrightarrow \text{NO} + \text{O}$ | $1.500 \cdot 10^{-11} \exp(-3600./T)$ | JPL (2011) |
| $\text{N} + \text{NO} \longrightarrow \text{N}_2 + \text{O}$ | $2.100 \cdot 10^{-11} \exp(100./T)$ | JPL (2011) |
| $\text{N} + \text{NO}_2 \longrightarrow 0.5 \cdot \text{N}_2\text{O} + 0.5 \cdot \text{O} + 0.5 \cdot \text{NO} + 0.25 \cdot \text{N}_2 + 0.25 \cdot \text{O}_2$ | $5.800 \cdot 10^{-12} \exp(220./T)$ | JPL (2011), products: D. Kinnison |
| $\text{NO} + \text{O} + \text{M} \longrightarrow \text{M} + \text{NO}_2$ | <i>ktroe</i> ($9.000 \cdot 10^{-32}, 1.5, 3.000 \cdot 10^{-11}, 0., 0.6$) | JPL (2011) |
| $\text{NO} + \text{O}_3 \longrightarrow \text{NO}_2 + \text{O}_2$ | $3.000 \cdot 10^{-12} \exp(-1500./T)$ | JPL (2011) |

Table S6: NOx reactions (... continued)

| reaction | rate coefficient | reference |
|--|--|--|
| $\text{NO} + \text{HO}_2 \longrightarrow \text{NO}_2 + \text{OH}$ | $3.300 \cdot 10^{-12} \exp(270./T)$ | JPL (2011) |
| $\text{NO}_2 + \text{O} \longrightarrow \text{NO} + \text{O}_2$ | $5.100 \cdot 10^{-12} \exp(210./T)$ | JPL (2011) |
| $\text{NO}_2 + \text{O} + \text{M} \longrightarrow \text{M} + \text{NO}_3$ | $ktroe(2.500 \cdot 10^{-31}, 1.8, 2.200 \cdot 10^{-11}, 0.7, 0.6)$ | JPL (2011) |
| $\text{NO}_2 + \text{O}_3 \longrightarrow \text{NO}_3 + \text{O}_2$ | $1.200 \cdot 10^{-13} \exp(-2450./T)$ | JPL (2011) |
| $\text{NO}_2 + \text{H} \longrightarrow \text{NO} + \text{OH}$ | $4.000 \cdot 10^{-10} \exp(-340./T)$ | JPL (2011) |
| $\text{NO}_2 + \text{OH} + \text{M} \longrightarrow \text{HNO}_3 + \text{M}$ | $ktroe(1.800 \cdot 10^{-30}, 3., 2.800 \cdot 10^{-11}, 0., 0.6)$ | JPL (2011) |
| $\text{NO}_3 + \text{O} \longrightarrow \text{NO}_2 + \text{O}_2$ | $1.000 \cdot 10^{-11}$ | JPL (2011) |
| $\text{NO}_3 + \text{OH} \longrightarrow \text{HO}_2 + \text{NO}_2$ | $2.200 \cdot 10^{-11}$ | JPL (2011) |
| $\text{NO}_3 + \text{HO}_2 \longrightarrow \text{NO}_2 + \text{OH} + \text{O}_2$ | $3.500 \cdot 10^{-12}$ | JPL (2011) |
| $\text{NO}_3 + \text{NO} \longrightarrow 2 \cdot \text{NO}_2$ | $1.500 \cdot 10^{-11} \exp(170./T)$ | JPL (2011) |
| $\text{HNO}_3 + \text{OH} \longrightarrow \text{H}_2\text{O} + \text{NO}_3$ | | $X/(1 + X/(2.7 \cdot 10^{-17} \exp(2199/T))) + 2.4 \cdot 10^{-14} \exp(460/T); X = M6.5 \cdot 10^{-34} \exp(1335/T) - \text{JPL (2011)}$ |
| $\text{NO} + \text{OH} \longrightarrow \text{HONO}$ | $ktroe(7.000 \cdot 10^{-31}, 2.6, 3.600 \cdot 10^{-11}, 0.1, 0.6)$ | JPL (2011) |
| $\text{HONO} + \text{OH} \longrightarrow \text{H}_2\text{O} + \text{NO}_2$ | $1.800 \cdot 10^{-11} \exp(-390./T)$ | JPL (2011) |
| $\text{NO}_2 + \text{HO}_2 + \text{M} \longrightarrow \text{HO}_2\text{NO}_2 + \text{M}$ | $ktroe(2.000 \cdot 10^{-31}, 3.4, 2.900 \cdot 10^{-12}, 1.1, 0.6)$ | JPL (2011) |
| $\text{HO}_2\text{NO}_2 + \text{OH} \longrightarrow \text{H}_2\text{O} + \text{NO}_2 + \text{O}_2$ | $1.300 \cdot 10^{-12} \exp(380./T)$ | JPL (2011) |
| $\text{HO}_2\text{NO}_2 + \text{M} \longrightarrow \text{HO}_2 + \text{NO}_2 + \text{M}$ | | $[\text{NO}_2 - \text{NO}_3] * \exp(-10900./T)/2.1 \cdot 10^{-27} \text{ JPL (2011)}$ |
| $\text{NO}_2 + \text{NO}_3 + \text{M} \longrightarrow \text{M} + \text{N}_2\text{O}_5$ | $ktroe(2.000 \cdot 10^{-30}, 4.4, 1.400 \cdot 10^{-12}, 0.7, 0.6)$ | JPL (2011) |
| $\text{N}_2\text{O}_5 + \text{M} \longrightarrow \text{NO}_2 + \text{NO}_3 + \text{M}$ | $kN2O5$ | $[\text{NO}_2 - \text{NO}_3] * \exp(-11000/T)/2.7 \cdot 10^{-27} \text{ JPL (2011)}$ |
| $\text{NH}_3 + \text{OH} \longrightarrow$ | $1.700 \cdot 10^{-12} \exp(-710./T)$ | JPL (2011) |

Table S7: C1 oxidation reactions

| reaction | rate coefficient | reference |
|---|---|----------------------------|
| $\text{CO} + \text{OH} \longrightarrow \text{CO}_2 + \text{H}$ | $kactiv(1.5 \cdot 10^{-13}, -0.6, 2.1 \cdot 10^9, -6.1)$ | JPL (2011) |
| $\text{CO} + \text{OH} + \text{M} \longrightarrow \text{CO}_2 + \text{HO}_2 + \text{M}$ | $ktroe(5.900 \cdot 10^{-33}, 1.4, 1.100 \cdot 10^{-12}, -1.3, 0.6)$ | |
| $\text{CH}_4 + \text{OH} \longrightarrow \text{CH}_3\text{O}_2 + \text{H}_2\text{O}$ | $2.450 \cdot 10^{-12} \exp(-1775./T)$ | JPL (2011) |
| $\text{CH}_3\text{OH} + \text{OH} \longrightarrow \text{CH}_2\text{O} + \text{HO}_2 + \text{H}_2\text{O}$ | $2.900 \cdot 10^{-12} \exp(-345./T)$ | JPL (2011) |
| $\text{CH}_3\text{O}_2 + \text{NO} \longrightarrow \text{CH}_2\text{O} + \text{NO}_2 + \text{HO}_2$ | $1.960 \cdot 10^{-12} \exp(403./T)$ | Orlando and Tyndall (2012) |
| $\text{CH}_3\text{O}_2 + \text{HO}_2 \longrightarrow \text{CH}_3\text{OOH} + \text{O}_2$ | $3.800 \cdot 10^{-13} \exp(730./T)$ | Orlando and Tyndall (2012) |
| $\text{CH}_3\text{O}_2 + \text{CH}_3\text{O}_2 \longrightarrow 2 \cdot \text{CH}_2\text{O} + 2 \cdot \text{HO}_2$ | $7.400 \cdot 10^{-13} \exp(-520./T)$ | IUPAC (2006) |

Table S7: C1 oxidation reactions (... continued)

| reaction | rate coefficient | reference |
|--|--|-------------------------------------|
| $\text{CH}_3\text{O}_2 + \text{CH}_3\text{O}_2 \longrightarrow \text{CH}_2\text{O} + \text{CH}_3\text{OH}$ | $2.330 \cdot 10^{-14} \exp(678./T)$ | own fit for $T = 240 - 300\text{K}$ |
| $\text{CH}_2\text{O} + \text{O} \longrightarrow \text{HO}_2 + \text{OH} + \text{CO}$ | $3.400 \cdot 10^{-11} \exp(-1600./T)$ | JPL (2011) |
| $\text{CH}_2\text{O} + \text{OH} \longrightarrow \text{CO} + \text{H}_2\text{O} + \text{H}$ | $5.500 \cdot 10^{-12} \exp(125./T)$ | JPL (2011) |
| $\text{CH}_2\text{O} + \text{HO}_2 \longrightarrow \text{HOCH}_2\text{OO}$ | $9.700 \cdot 10^{-15} \exp(625./T)$ | IUPAC (2006) |
| $\text{CH}_2\text{O} + \text{NO}_3 \longrightarrow \text{CO} + \text{HO}_2 + \text{HNO}_3$ | $6.000 \cdot 10^{-13} \exp(-2058./T)$ | |
| $\text{CH}_3\text{OOH} + \text{OH} \longrightarrow 0.7 \cdot \text{CH}_3\text{O}_2 + 0.3 \cdot \text{OH} + 0.3 \cdot \text{CH}_2\text{O} + \text{H}_2\text{O}$ | $3.800 \cdot 10^{-12} \exp(200./T)$ | JPL (2011) |
| $\text{HCOOH} + \text{OH} \longrightarrow \text{CO}_2 + \text{HO}_2 + \text{H}_2\text{O}$ | $4.000 \cdot 10^{-13}$ | JPL (2011) |
| $\text{HOCH}_2\text{OO} \longrightarrow \text{CH}_2\text{O} + \text{HO}_2$ | $2.400 \cdot 10^{12} \exp(-7000./T)$ | Lamarque et al. (2012) |
| $\text{HOCH}_2\text{OO} + \text{NO} \longrightarrow \text{HCOOH} + \text{NO}_2 + \text{HO}_2$ | $2.600 \cdot 10^{-12} \exp(265./T)$ | Lamarque et al. (2012) |
| $\text{HOCH}_2\text{OO} + \text{HO}_2 \longrightarrow \text{H}_2\text{O} + \text{HCOOH}$ | $7.500 \cdot 10^{-13} \exp(700./T)$ | Lamarque et al. (2012) |
| $\text{HCN} + \text{OH} \longrightarrow \text{CO} + \text{NO} + \text{H}_2\text{O}$ | $1.200 \cdot 10^{-13} \exp(-400./T)$ | JPL (2011), products: Tyndall |
| $\text{CH}_3\text{O}_2 + \text{NO}_2 + \text{M} \longrightarrow \text{CH}_3\text{O}_2\text{NO}_2 + \text{M}$ | $ktroe(1.000 \cdot 10^{-30}, 4.8, 7.200 \cdot 10^{-12}, 2.1, 0.6)$ | JPL (2011) |
| $\text{CH}_3\text{O}_2\text{NO}_2 + \text{M} \longrightarrow \text{CH}_3\text{O}_2 + \text{NO}_2 + \text{M}$ | | |

Table S8: C2 oxidation reactions

| reaction | rate coefficient | reference |
|---|---|---|
| $\text{C}_2\text{H}_2 + \text{OH} + \text{M} \longrightarrow 0.65 \cdot \text{GLYOXAL} + 0.65 \cdot \text{OH} + 0.35 \cdot \text{HCOOH} + 0.35 \cdot \text{HO}_2 + 0.35 \cdot \text{CO} + \text{M}$ | $ktroe(5.500 \cdot 10^{-30}, 0., 8.300 \cdot 10^{-13}, -2., 0.6)$ | JPL (2011) |
| $\text{C}_2\text{H}_4 + \text{OH} + \text{M} \longrightarrow \text{EO}_2 + \text{M}$ | $ktroe(1.000 \cdot 10^{-28}, 4.5, 7.500 \cdot 10^{-12}, 0.85, 0.6)$ | JPL (2011) |
| $\text{C}_2\text{H}_4 + \text{O}_3 \longrightarrow \text{CH}_2\text{O} + 0.65 \cdot \text{CO} + 0.15 \cdot \text{OH} + 0.15 \cdot \text{HO}_2 + 0.5 \cdot \text{H}_2\text{O} + 0.35 \cdot \text{HCOOH}$ | $9.100 \cdot 10^{-15} \exp(-2580./T)$ | IUPAC (2006) |
| $\text{C}_2\text{H}_6 + \text{OH} \longrightarrow \text{C}_2\text{H}_5\text{O}_2 + \text{H}_2\text{O}$ | $7.660 \cdot 10^{-12} \exp(-1020./T)$ | JPL (2011) |
| $\text{C}_2\text{H}_5\text{OH} + \text{OH} \longrightarrow \text{CH}_3\text{CHO} + \text{HO}_2 + \text{H}_2\text{O}$ | $3.350 \cdot 10^{-12} \exp(0./T)$ | JPL (2011) |
| $\text{CH}_3\text{CHO} + \text{OH} \longrightarrow \text{CH}_3\text{CO}_3 + \text{H}_2\text{O}$ | $4.630 \cdot 10^{-12} \exp(350./T)$ | JPL (2011) |
| $\text{CH}_3\text{CHO} + \text{NO}_3 \longrightarrow \text{CH}_3\text{CO}_3 + \text{HNO}_3$ | $1.400 \cdot 10^{-12} \exp(-1900./T)$ | JPL (2011) |
| $\text{CH}_3\text{COOOH} + \text{OH} \longrightarrow 0.5 \cdot \text{CH}_3\text{CO}_3 + \text{H}_2\text{O} + 0.5 \cdot \text{CH}_2\text{O} + 0.5 \cdot \text{CO}_2 + 0.5 \cdot \text{OH}$ | $1.000 \cdot 10^{-12}$ | Orlando (p.c.) added OH |
| $\text{C}_2\text{H}_5\text{O}_2 + \text{NO} \longrightarrow \text{CH}_3\text{CHO} + \text{HO}_2 + \text{NO}_2$ | $2.620 \cdot 10^{-12} \exp(373./T)$ | Orlando and Tyndall (2012) |
| $\text{C}_2\text{H}_5\text{O}_2 + \text{HO}_2 \longrightarrow \text{C}_2\text{H}_5\text{OOH} + \text{O}_2$ | $7.400 \cdot 10^{-13} \exp(700./T)$ | Orlando and Tyndall (2012) |
| $\text{C}_2\text{H}_5\text{O}_2 + \text{CH}_3\text{O}_2 \longrightarrow 0.7 \cdot \text{CH}_2\text{O} + 0.8 \cdot \text{CH}_3\text{CHO} + \text{HO}_2 + 0.3 \cdot \text{CH}_3\text{OH} + 0.2 \cdot \text{C}_2\text{H}_5\text{OH}$ | $2.000 \cdot 10^{-13}$ | Orlando (p.c.), products: Tyndall (p.c.) |
| $\text{C}_2\text{H}_5\text{O}_2 + \text{CH}_3\text{CO}_3 \longrightarrow \text{CH}_3\text{CHO} + \text{HO}_2 + \text{CH}_3\text{O}_2 + \text{CO}_2$ | $1.800 \cdot 10^{-12} \exp(500./T)$ | 10% lower than $\text{CH}_3\text{O}_2 + \text{CH}_3\text{CO}_3$; Orlando and Tyndall (2012) only give k@298K |

Table S8: C2 oxidation reactions (... continued)

| reaction | rate coefficient | reference |
|---|--|--|
| $\text{C}_2\text{H}_5\text{O}_2 + \text{C}_2\text{H}_5\text{O}_2 \longrightarrow 1.6 \cdot \text{CH}_3\text{CHO} + 1.2 \cdot \text{HO}_2 + 0.4 \cdot \text{C}_2\text{H}_5\text{OH}$ | $7.600 \cdot 10^{-14}$ | Orlando and Tyndall (2012) |
| $\text{C}_2\text{H}_5\text{OOH} + \text{OH} \longrightarrow 0.5 \cdot \text{C}_2\text{H}_5\text{O}_2 + 0.5 \cdot \text{CH}_3\text{CHO} + 0.5 \cdot \text{OH} + \text{H}_2\text{O}$ | $3.800 \cdot 10^{-12} \exp(200./T)$ | no data, analog to $\text{CH}_3\text{OOH} + \text{OH}$ |
| $\text{CH}_3\text{CO}_3 + \text{NO} \longrightarrow \text{CH}_3\text{O}_2 + \text{CO}_2 + \text{NO}_2$ | $7.500 \cdot 10^{-12} \exp(290./T)$ | Orlando and Tyndall (2012) |
| $\text{CH}_3\text{CO}_3 + \text{NO}_2 + \text{M} \longrightarrow \text{M} + \text{PAN}$ | $ktroe(2.700 \cdot 10^{-28}, 7.1, 1.200 \cdot 10^{-11}, 0.9, 0.6)$ | |
| $\text{CH}_3\text{CO}_3 + \text{HO}_2 \longrightarrow 0.4 \cdot \text{CH}_3\text{COOOH} + 0.2 \cdot \text{CH}_3\text{COOH} + 0.2 \cdot \text{O}_3 + 0.4 \cdot \text{CH}_3\text{O}_2 + 0.4 \cdot \text{OH} + 0.4 \cdot \text{CO}_2$ | $5.200 \cdot 10^{-13} \exp(980./T)$ | Orlando and Tyndall (2012) |
| $\text{CH}_3\text{CO}_3 + \text{CH}_3\text{O}_2 \longrightarrow \text{CH}_2\text{O} + 0.9 \cdot \text{CH}_3\text{O}_2 + 0.9 \cdot \text{HO}_2 + 0.9 \cdot \text{CO}_2 + 0.1 \cdot \text{CH}_3\text{COOH}$ | $2.000 \cdot 10^{-12} \exp(500./T)$ | Orlando and Tyndall (2012) |
| $\text{CH}_3\text{CO}_3 + \text{CH}_3\text{CO}_3 \longrightarrow 2 \cdot \text{CH}_3\text{O}_2 + 2 \cdot \text{CO}_2$ | $2.900 \cdot 10^{-12} \exp(500./T)$ | Orlando and Tyndall (2012) |
| $\text{CH}_3\text{COOH} + \text{OH} \longrightarrow \text{CH}_3\text{O}_2 + \text{CO}_2 + \text{H}_2\text{O}$ | $3.150 \cdot 10^{-14} \exp(920./T)$ | JPL (2011) |
| $\text{PAN} + \text{M} \longrightarrow \text{CH}_3\text{CO}_3 + \text{NO}_2 + \text{M}$ | | |
| $\text{PAN} + \text{OH} \longrightarrow \text{CH}_2\text{O} + \text{CO}_2 + \text{NO}_3$ | $4.000 \cdot 10^{-14}$ | JPL (2011), includes implicit $\text{NO} \longrightarrow \text{NO}_2$ conversion |
| $\text{EO}_2 + \text{NO} \longrightarrow 0.75 \cdot \text{EO} + \text{NO}_2 + 0.5 \cdot \text{CH}_2\text{O} + 0.25 \cdot \text{HO}_2$ | $4.200 \cdot 10^{-12} \exp(180./T)$ | Lamarque et al. (2012) |
| $\text{EO}_2 + \text{HO}_2 \longrightarrow \text{EOOH}$ | $7.500 \cdot 10^{-13} \exp(700./T)$ | |
| $\text{EO}_2 + \text{CH}_3\text{O}_2 \longrightarrow 0.5 \cdot \text{EO} + 0.5 \cdot \text{O}_2 + 0.5 \cdot \text{CH}_2\text{O} + 0.5 \cdot \text{HO}_2 + 0.5 \cdot \text{CH}_3\text{OH} + 0.5 \cdot \text{GLYALD}$ | $4.000 \cdot 10^{-12} \exp(1000./T)$ | Tyndall (p.c.) |
| $\text{EO}_2 + \text{CH}_3\text{CO}_3 \longrightarrow \text{CH}_3\text{O}_2 + \text{EO} + \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{EO} + \text{O}_2 \longrightarrow \text{GLYALD} + \text{HO}_2$ | $1.000 \cdot 10^{-14}$ | Lamarque et al. (2012) |
| $\text{EO} \longrightarrow 2 \cdot \text{CH}_2\text{O} + \text{HO}_2$ | $1.600 \cdot 10^{11} \exp(-4150./T)$ | Lamarque et al. (2012) |
| $\text{GLYOXAL} + \text{OH} \longrightarrow 0.6 \cdot \text{HO}_2 + 1.2 \cdot \text{CO} + \text{H}_2\text{O} + 0.4 \cdot \text{HCOCO}_3$ | $3.100 \cdot 10^{-12} \exp(340./T)$ | MCM3.2 |
| $\text{GLYOXAL} + \text{NO}_3 \longrightarrow 0.6 \cdot \text{HO}_2 + 1.2 \cdot \text{CO} + 0.4 \cdot \text{HCOCO}_3 + \text{HNO}_3$ | $2.500 \cdot 10^{-12}$ | Taraborrelli et al. (2009) |
| $\text{GLYALD} + \text{OH} \longrightarrow 0.2 \cdot \text{GLYOXAL} + 0.2 \cdot \text{HO}_2 + 0.8 \cdot \text{HOCH}_2\text{CO}_3 + \text{H}_2\text{O}$ | $1.000 \cdot 10^{-11}$ | Taraborrelli et al. (2009) |
| $\text{GLYALD} + \text{NO}_3 \longrightarrow \text{HNO}_3 + \text{HOCH}_2\text{CO}_3$ | $1.440 \cdot 10^{-12} \exp(-1862./T)$ | Taraborrelli et al. (2009) |
| $\text{HOCH}_2\text{CO}_3 + \text{NO}_2 + \text{M} \longrightarrow \text{CH}_2\text{O} + \text{CO}_2 + \text{HNO}_3 + \text{M}$ | $ktroe(2.700 \cdot 10^{-28}, 7.1, 1.200 \cdot 10^{-11}, 0.9, 0.6)$ | Orlando and Tyndall (2012) |
| $\text{HOCH}_2\text{CO}_3 + \text{HO}_2 \longrightarrow 0.4 \cdot \text{HOCH}_2\text{CO}_3\text{H} + 0.2 \cdot \text{HOCH}_2\text{CO}_2\text{H} + 0.2 \cdot \text{O}_3 + 0.4 \cdot \text{CO}_2 + 0.4 \cdot \text{OH} + 0.4 \cdot \text{HO}_2 + 0.4 \cdot \text{CH}_2\text{O}$ | $4.300 \cdot 10^{-13} \exp(1040./T)$ | Orlando (p.c.) |
| $\text{HOCH}_2\text{CO}_3 + \text{CH}_3\text{O}_2 \longrightarrow 2 \cdot \text{CH}_2\text{O} + \text{CO}_2 + 2 \cdot \text{HO}_2$ | $1.000 \cdot 10^{-11}$ | Taraborrelli et al. (2009) |
| $\text{HOCH}_2\text{CO}_3 + \text{CH}_3\text{CO}_3 \longrightarrow \text{CH}_3\text{O}_2 + 2 \cdot \text{CO}_2 + \text{HO}_2 + \text{CH}_2\text{O}$ | $1.000 \cdot 10^{-11}$ | Taraborrelli et al. (2009) |

Table S8: C2 oxidation reactions (... continued)

| reaction | rate coefficient | reference |
|--|---------------------------------------|----------------------------|
| $\text{HOCH}_2\text{CO}_3 + \text{NO} \longrightarrow \text{HO}_2 + \text{NO}_2 + \text{CH}_2\text{O} + \text{CO}_2$ | $8.100 \cdot 10^{-12} \exp(270./T)$ | Taraborrelli et al. (2009) |
| $\text{HOCH}_2\text{CO}_3 + \text{NO}_3 \longrightarrow \text{HO}_2 + \text{NO}_2 + \text{CH}_2\text{O} + \text{CO}_2$ | $4.000 \cdot 10^{-12}$ | Taraborrelli et al. (2009) |
| $\text{HOCH}_2\text{CO}_2\text{H} + \text{OH} \longrightarrow \text{CH}_2\text{O} + \text{HO}_2 + \text{CO}_2 + \text{H}_2\text{O}$ | $2.730 \cdot 10^{-12}$ | Taraborrelli et al. (2009) |
| $\text{HOCH}_2\text{CO}_3\text{H} + \text{OH} \longrightarrow \text{H}_2\text{O} + \text{HOCH}_2\text{CO}_3$ | $6.190 \cdot 10^{-12}$ | Taraborrelli et al. (2009) |
| $\text{HCOCO}_3 + \text{CH}_3\text{O}_2 \longrightarrow \text{CO} + 2 \cdot \text{HO}_2 + \text{CO}_2 + \text{CH}_2\text{O}$ | $1.000 \cdot 10^{-11}$ | Taraborrelli et al. (2009) |
| $\text{HCOCO}_3 + \text{CH}_3\text{CO}_3 \longrightarrow \text{CO} + \text{HO}_2 + \text{CH}_3\text{O}_2 + 2 \cdot \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | Taraborrelli et al. (2009) |
| $\text{HCOCO}_3 + \text{HO}_2 \longrightarrow 0.7 \cdot \text{HCOCO}_3\text{H} + 0.7 \cdot \text{O}_2 + 0.3 \cdot \text{HCOCO}_2\text{H} + 0.3 \cdot \text{O}_3$ | $4.300 \cdot 10^{-13} \exp(1040./T)$ | Taraborrelli et al. (2009) |
| $\text{HCOCO}_3 + \text{NO} \longrightarrow \text{CO} + \text{HO}_2 + \text{NO}_2 + \text{CO}_2$ | $8.100 \cdot 10^{-12} \exp(270./T)$ | Taraborrelli et al. (2009) |
| $\text{HCOCO}_3 + \text{NO}_3 \longrightarrow \text{CO} + \text{HO}_2 + \text{NO}_2 + \text{CO}_2$ | $4.000 \cdot 10^{-12}$ | Taraborrelli et al. (2009) |
| $\text{HCOCO}_2\text{H} + \text{OH} \longrightarrow \text{CO} + \text{HO}_2 + \text{CO}_2 + \text{H}_2\text{O}$ | $1.230 \cdot 10^{-11}$ | Taraborrelli et al. (2009) |
| $\text{HCOCO}_3\text{H} + \text{OH} \longrightarrow \text{H}_2\text{O} + \text{HCOCO}_3$ | $1.580 \cdot 10^{-11}$ | Taraborrelli et al. (2009) |
| $\text{CH}_3\text{CN} + \text{OH} \longrightarrow \text{CH}_2\text{O} + \text{H}_2\text{O} + \text{CO} + \text{NO}_2$ | $7.800 \cdot 10^{-13} \exp(-1050./T)$ | products: Tyndall |

Table S9: C3 oxidation reactions

| reaction | rate coefficient | reference |
|---|---|---|
| $\text{C}_3\text{H}_6 + \text{OH} + \text{M} \longrightarrow \text{M} + \text{PO}_2$ | <i>ktroe</i> ($8.000 \cdot 10^{-27}, 3.5, 3.000 \cdot 10^{-11}, 0., 0.5$) | IUPAC (2006) |
| $\text{C}_3\text{H}_6 + \text{O}_3 \longrightarrow 0.28 \cdot \text{CH}_3\text{O}_2 + 0.1 \cdot \text{CH}_4 + 0.075 \cdot \text{CH}_3\text{COOH} + 0.56 \cdot \text{CO} + 0.075 \cdot \text{HCOOH} + 0.09 \cdot \text{H}_2\text{O}_2 + 0.28 \cdot \text{HO}_2 + 0.2 \cdot \text{CO}_2 + 0.545 \cdot \text{CH}_3\text{CHO} + 0.545 \cdot \text{CH}_2\text{O} + 0.36 \cdot \text{OH}$ | $5.500 \cdot 10^{-15} \exp(-1880./T)$ | IUPAC (2006) |
| $\text{C}_3\text{H}_6 + \text{NO}_3 \longrightarrow \text{PRONO}_3\text{BO}_2$ | $4.600 \cdot 10^{-13} \exp(-1156./T)$ | IUPAC (2006) |
| $\text{C}_3\text{H}_8 + \text{OH} \longrightarrow \text{C}_3\text{H}_7\text{O}_2 + \text{H}_2\text{O}$ | $7.600 \cdot 10^{-12} \exp(-585./T)$ | IUPAC (2006) |
| $\text{CH}_3\text{COCH}_3 + \text{OH} \longrightarrow \text{CH}_3\text{COCH}_2\text{O}_2 + \text{H}_2\text{O}$ | $\text{CH}_3\text{COCH}_3 - \text{OH}$ | |
| $\text{C}_3\text{H}_7\text{O}_2 + \text{NO} \longrightarrow 0.82 \cdot \text{CH}_3\text{COCH}_3 + \text{NO}_2 + \text{HO}_2 + 0.27 \cdot \text{CH}_3\text{CHO}$ | $2.900 \cdot 10^{-12} \exp(350./T)$ | Orlando and Tyndall (2012) |
| $\text{C}_3\text{H}_7\text{O}_2 + \text{HO}_2 \longrightarrow \text{C}_3\text{H}_7\text{OOH} + \text{O}_2$ | $7.500 \cdot 10^{-13} \exp(700./T)$ | Lamarque et al. (2012) |
| $\text{C}_3\text{H}_7\text{O}_2 + \text{CH}_3\text{O}_2 \longrightarrow \text{CH}_2\text{O} + 2 \cdot \text{HO}_2 + 0.82 \cdot \text{CH}_3\text{COCH}_3 + 0.27 \cdot \text{CH}_3\text{CHO}$ | $3.750 \cdot 10^{-13} \exp(-40./T)$ | Lamarque et al. (2012) |
| $\text{C}_3\text{H}_7\text{O}_2 + \text{CH}_3\text{CO}_3 \longrightarrow 0.82 \cdot \text{CH}_3\text{COCH}_3 + \text{HO}_2 + 0.27 \cdot \text{CH}_3\text{CHO} + \text{CO}_2 + \text{CH}_3\text{O}_2$ | $1.000 \cdot 10^{-11}$ | Lamarque et al. (2012) |
| $\text{C}_3\text{H}_7\text{OOH} + \text{OH} \longrightarrow 0.41 \cdot \text{CH}_3\text{COCH}_3 + 0.5 \cdot \text{OH} + 0.5 \cdot \text{C}_3\text{H}_7\text{O}_2 + 0.5 \cdot \text{H}_2\text{O} + 0.135 \cdot \text{CH}_3\text{CHO}$ | $3.800 \cdot 10^{-12} \exp(200./T)$ | Lamarque et al. (2012); $1.5 \cdot \text{CH}_3\text{CHO}$ as surrogate for propanal |
| $\text{PO}_2 + \text{NO} \longrightarrow \text{CH}_2\text{O} + \text{CH}_3\text{CHO} + \text{HO}_2 + \text{NO}_2$ | $4.200 \cdot 10^{-12} \exp(180./T)$ | Lamarque et al. (2012) |

Table S9: C3 oxidation reactions (... continued)

| reaction | rate coefficient | reference |
|--|---------------------------------------|---------------------------------|
| $\text{PO}_2 + \text{NO}_3 \longrightarrow \text{CH}_2\text{O} + \text{CH}_3\text{CHO} + \text{HO}_2 + \text{NO}_2$ | $2.500 \cdot 10^{-12}$ | Taraborrelli et al. (2009) |
| $\text{PO}_2 + \text{HO}_2 \longrightarrow \text{O}_2 + \text{POOH}$ | $7.500 \cdot 10^{-13} \exp(700./T)$ | Lamarque et al. (2012) |
| $\text{PO}_2 + \text{CH}_3\text{O}_2 \longrightarrow 0.5 \cdot \text{CH}_3\text{CHO} + 1.25 \cdot \text{CH}_2\text{O} + \text{HO}_2 + 0.5 \cdot \text{HYAC} + 0.25 \cdot \text{CH}_3\text{OH}$ | $8.300 \cdot 10^{-13}$ | products: Tyndall (p.c.) |
| $\text{PO}_2 + \text{CH}_3\text{CO}_3 \longrightarrow \text{CH}_2\text{O} + \text{CH}_3\text{CHO} + \text{HO}_2 + \text{CH}_3\text{O}_2 + \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{POOH} + \text{OH} \longrightarrow 0.5 \cdot \text{PO}_2 + 0.5 \cdot \text{HYAC} + 0.5 \cdot \text{OH} + \text{H}_2\text{O}$ | $3.800 \cdot 10^{-12} \exp(200./T)$ | Lamarque et al. (2012) |
| $\text{HYAC} + \text{OH} \longrightarrow \text{CH}_3\text{COCHO} + \text{HO}_2 + \text{H}_2\text{O}$ | $3.000 \cdot 10^{-12}$ | IUPAC (2006) |
| $\text{CH}_3\text{COCHO} + \text{OH} \longrightarrow \text{CH}_3\text{CO}_3 + \text{CO} + \text{H}_2\text{O}$ | $8.400 \cdot 10^{-13} \exp(830./T)$ | Lamarque et al. (2012) |
| $\text{CH}_3\text{COCHO} + \text{NO}_3 \longrightarrow \text{CO} + \text{HNO}_3 + \text{CH}_3\text{CO}_3$ | $1.400 \cdot 10^{-12} \exp(-1860./T)$ | Lamarque et al. (2012) |
| $\text{CH}_3\text{COCH}_2\text{O}_2 + \text{NO} \longrightarrow \text{CH}_2\text{O} + \text{CH}_3\text{CO}_3 + \text{NO}_2$ | $2.900 \cdot 10^{-12} \exp(300./T)$ | Lamarque et al. (2012) |
| $\text{CH}_3\text{COCH}_2\text{O}_2 + \text{HO}_2 \longrightarrow 0.85 \cdot \text{O}_2 + 0.85 \cdot \text{ROOH} + 0.15 \cdot \text{CH}_2\text{O} + 0.15 \cdot \text{CH}_3\text{CO}_3 + 0.15 \cdot \text{OH} + 0.15 \cdot \text{H}_2\text{O}$ | $8.600 \cdot 10^{-13} \exp(700./T)$ | Lamarque et al. (2012) + MCM3.2 |
| $\text{CH}_3\text{COCH}_2\text{O}_2 + \text{CH}_3\text{O}_2 \longrightarrow 0.3 \cdot \text{CH}_3\text{CO}_3 + 0.8 \cdot \text{CH}_2\text{O} + 0.3 \cdot \text{HO}_2 + 0.2 \cdot \text{HYAC} + 0.5 \cdot \text{CH}_3\text{COCHO} + 0.5 \cdot \text{CH}_3\text{OH}$ | $7.100 \cdot 10^{-13} \exp(500./T)$ | Lamarque et al. (2012) |
| $\text{CH}_3\text{COCH}_2\text{O}_2 + \text{CH}_3\text{CO}_3 \longrightarrow \text{CH}_2\text{O} + \text{CH}_3\text{CO}_3 + \text{CH}_3\text{O}_2 + \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{ROOH} + \text{OH} \longrightarrow \text{CH}_3\text{COCH}_2\text{O}_2 + \text{H}_2\text{O}$ | $3.800 \cdot 10^{-12} \exp(200./T)$ | Lamarque et al. (2012) |
| $\text{PRONO}_3\text{BO}_2 + \text{NO} \longrightarrow 0.83 \cdot \text{HO}_2 + 0.83 \cdot \text{NOA} + 0.17 \cdot \text{CH}_2\text{O} + 0.17 \cdot \text{CH}_3\text{CHO} + 1.17 \cdot \text{NO}_2$ | $2.540 \cdot 10^{-12} \exp(360./T)$ | Taraborrelli et al. (2009) |
| $\text{PRONO}_3\text{BO}_2 + \text{NO}_3 \longrightarrow 0.83 \cdot \text{HO}_2 + 0.83 \cdot \text{NOA} + 0.17 \cdot \text{CH}_2\text{O} + 0.17 \cdot \text{CH}_3\text{CHO} + 1.17 \cdot \text{NO}_2$ | $2.500 \cdot 10^{-12}$ | Taraborrelli et al. (2009) |
| $\text{PRONO}_3\text{BO}_2 + \text{HO}_2 \longrightarrow \text{PR}_2\text{O}_2\text{HNO}_3$ | $1.320 \cdot 10^{-12} \exp(360./T)$ | Taraborrelli et al. (2009) |
| $\text{PRONO}_3\text{BO}_2 + \text{CH}_3\text{O}_2 \longrightarrow 0.915 \cdot \text{HO}_2 + 0.915 \cdot \text{NOA} + 0.835 \cdot \text{CH}_2\text{O} + 0.085 \cdot \text{CH}_3\text{CHO} + 0.25 \cdot \text{CH}_3\text{OH}$ | $1.000 \cdot 10^{-12}$ | |
| $\text{PRONO}_3\text{BO}_2 + \text{CH}_3\text{CO}_3 \longrightarrow 0.83 \cdot \text{HO}_2 + 0.83 \cdot \text{NOA} + 0.17 \cdot \text{CH}_2\text{O} + 0.17 \cdot \text{CH}_3\text{CHO} + \text{CH}_3\text{O}_2 + \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{PR}_2\text{O}_2\text{HNO}_3 + \text{OH} \longrightarrow 0.5 \cdot \text{PRONO}_3\text{BO}_2 + 0.5 \cdot \text{NOA} + 0.5 \cdot \text{OH} + \text{H}_2\text{O}$ | $7.000 \cdot 10^{-12}$ | Taraborrelli et al. (2009) |
| $\text{NOA} + \text{OH} \longrightarrow \text{CH}_3\text{COCHO} + \text{NO}_2 + \text{H}_2\text{O}$ | $1.300 \cdot 10^{-13}$ | Taraborrelli et al. (2009) |

Table S10: C4 oxidation reactions

| reaction | rate coefficient | reference |
|---|------------------------|------------------------|
| $\text{BIGENE} + \text{OH} \longrightarrow \text{ENEO}_2$ | $5.400 \cdot 10^{-11}$ | Lamarque et al. (2012) |

Table S10: C4 oxidation reactions (... continued)

| reaction | rate coefficient | reference |
|--|---------------------------------------|---|
| $\text{MEK} + \text{OH} \longrightarrow \text{H}_2\text{O} + \text{MEKO}_2$ | $2.300 \cdot 10^{-12} \exp(-170./T)$ | Lamarque et al. (2012) |
| $\text{MEKO}_2 + \text{NO} \longrightarrow \text{CH}_3\text{CHO} + \text{CH}_3\text{CO}_3 + \text{NO}_2$ | $4.032 \cdot 10^{-12} \exp(180./T)$ | treated like MEKBO_2 from MCM3.2; $A = 4.200 \cdot 10^{-12} \cdot 0.96$; 4% nitrate yield as for MVKN |
| $\text{MEKO}_2 + \text{NO} \longrightarrow \text{MEKNO}_3$ | $1.680 \cdot 10^{-13} \exp(180./T)$ | $A = 4.200 \cdot 10^{-12} \cdot 0.04$; 4% nitrate yield as for MVKN |
| $\text{MEKO}_2 + \text{HO}_2 \longrightarrow \text{MEKOOH}$ | $7.500 \cdot 10^{-13} \exp(700./T)$ | Lamarque et al. (2012) |
| $\text{MEKO}_2 + \text{CH}_3\text{O}_2 \longrightarrow 0.3 \cdot \text{CH}_3\text{CHO} + 0.3 \cdot \text{CH}_3\text{CO}_3 + \text{CH}_2\text{O} + 0.3 \cdot \text{HO}_2 + 0.3 \cdot \text{O}_2 + 0.5 \cdot \text{BIACETOH} + 0.5 \cdot \text{CH}_3\text{OH} + 0.266 \cdot \text{HYAC}$ | $1.000 \cdot 10^{-12}$ | Tyndall (p.c.) added CH_2O to first and third channel |
| $\text{MEKO}_2 + \text{CH}_3\text{CO}_3 \longrightarrow \text{CH}_3\text{CHO} + \text{CH}_3\text{CO}_3 + \text{CH}_3\text{O}_2 + \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{MEKOOH} + \text{OH} \longrightarrow \text{H}_2\text{O} + \text{MEKO}_2$ | $3.800 \cdot 10^{-12} \exp(200./T)$ | Lamarque et al. (2012) |
| $\text{ENEO}_2 + \text{NO} \longrightarrow \text{CH}_3\text{CHO} + 0.5 \cdot \text{CH}_2\text{O} + 0.5 \cdot \text{CH}_3\text{COCH}_3 + \text{HO}_2 + \text{NO}_2$ | $4.200 \cdot 10^{-12} \exp(180./T)$ | Lamarque et al. (2012) |
| $\text{ENEO}_2 + \text{HO}_2 \longrightarrow 1.333 \cdot \text{POOH} + \text{O}_2$ | $7.500 \cdot 10^{-13} \exp(700./T)$ | factor 4/3 to preserve carbon |
| $\text{ENEO}_2 + \text{CH}_3\text{O}_2 \longrightarrow 0.665 \cdot \text{HYAC} + 0.5 \cdot \text{CH}_3\text{OH} + 0.5 \cdot \text{CH}_3\text{CHO} + 0.25 \cdot \text{CH}_3\text{COCH}_3 + 0.75 \cdot \text{CH}_2\text{O} + \text{HO}_2$ | $1.000 \cdot 10^{-12}$ | products: Tyndall (p.c.) |
| $\text{ENEO}_2 + \text{CH}_3\text{CO}_3 \longrightarrow \text{CH}_3\text{CHO} + 0.5 \cdot \text{CH}_2\text{O} + 0.5 \cdot \text{CH}_3\text{COCH}_3 + \text{HO}_2 + \text{CH}_3\text{O}_2 + \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{MACR} + \text{OH} \longrightarrow 0.45 \cdot \text{MCO}_3 + 0.55 \cdot \text{MACRO}_2$ | $1.860 \cdot 10^{-11} \exp(175./T)$ | Tnydall (p.c.) |
| $\text{MACR} + \text{O}_3 \longrightarrow 0.59 \cdot \text{CH}_3\text{COCHO} + 0.41 \cdot \text{CH}_3\text{CO}_3 + 0.82 \cdot \text{CO} + 0.41 \cdot \text{HO}_2 + 0.82 \cdot \text{OH} + 0.033750 \cdot \text{HCOOH} + 0.556250 \cdot \text{CH}_2\text{O} + 0.123750 \cdot \text{H}_2\text{O}_2$ | $1.360 \cdot 10^{-15} \exp(-2112./T)$ | |
| $\text{MACR} + \text{NO}_3 \longrightarrow \text{HNO}_3 + \text{MCO}_3$ | $2.880 \cdot 10^{-12} \exp(-1862./T)$ | |
| $\text{MCO}_3 + \text{CH}_3\text{O}_2 \longrightarrow 0.315 \cdot \text{CH}_3\text{CO}_3 + 0.585 \cdot \text{CH}_3\text{O}_2 + 0.585 \cdot \text{CO} + 1.9 \cdot \text{CH}_2\text{O} + 0.9 \cdot \text{CO}_2 + 0.9 \cdot \text{HO}_2 + 0.1 \cdot \text{MACO}_2\text{H}$ | $1.000 \cdot 10^{-11}$ | MCO_3 is an acyl radical, therefore $\text{kRO}_2\text{CH}_3\text{CO}_3$ |
| $\text{MCO}_3 + \text{CH}_3\text{CO}_3 \longrightarrow \text{CH}_2\text{O} + 0.35 \cdot \text{CH}_3\text{CO}_3 + 1.65 \cdot \text{CH}_3\text{O}_2 + 0.65 \cdot \text{CO} + 2 \cdot \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{MCO}_3 + \text{HO}_2 \longrightarrow 0.44 \cdot \text{OH} + 0.154 \cdot \text{CH}_3\text{CO}_3 + 0.286 \cdot \text{CH}_3\text{O}_2 + 0.286 \cdot \text{CO} + 0.44 \cdot \text{CH}_2\text{O} + 0.44 \cdot \text{CO}_2 + 0.15 \cdot \text{MACO}_2\text{H} + 0.15 \cdot \text{O}_3 + 0.41 \cdot \text{MACO}_3\text{H} + 0.41 \cdot \text{O}_2$ | $4.300 \cdot 10^{-13} \exp(1040./T)$ | |
| $\text{MCO}_3 + \text{NO} \longrightarrow \text{CH}_2\text{O} + 0.35 \cdot \text{CH}_3\text{CO}_3 + 0.65 \cdot \text{CH}_3\text{O}_2 + 0.65 \cdot \text{CO} + \text{NO}_2 + \text{CO}_2$ | $8.700 \cdot 10^{-12} \exp(290./T)$ | |
| $\text{MCO}_3 + \text{NO}_3 \longrightarrow \text{CH}_2\text{O} + 0.35 \cdot \text{CH}_3\text{CO}_3 + 0.65 \cdot \text{CH}_3\text{O}_2 + 0.65 \cdot \text{CO} + \text{NO}_2 + \text{CO}_2$ | $4.000 \cdot 10^{-12}$ | |

Table S10: C4 oxidation reactions (... continued)

| reaction | rate coefficient | reference |
|--|--|--|
| $\text{MCO}_3 + \text{NO}_2 + \text{M} \longrightarrow \text{M} + \text{MPAN}$ | $ktroe(2.700 \cdot 10^{-28}, 7.1, 1.200 \cdot 10^{-11}, 0.9, 0.3)$ | |
| $\text{MACRO}_2 + \text{CH}_3\text{O}_2 \longrightarrow 0.75 \cdot \text{CO} + 0.75 \cdot \text{HYAC} + \text{CH}_2\text{O} + 1.5 \cdot \text{HO}_2 + 0.25 \cdot \text{MACROH} + 0.25 \cdot \text{O}_2$ | $9.200 \cdot 10^{-14}$ | products: Tyndall (p.c.) |
| $\text{MACRO}_2 + \text{CH}_3\text{CO}_3 \longrightarrow 0.85 \cdot \text{CO} + 0.85 \cdot \text{HYAC} + 0.15 \cdot \text{CH}_2\text{O} + 0.15 \cdot \text{CH}_3\text{COCHO} + \text{HO}_2 + \text{CH}_3\text{O}_2 + \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{MACRO}_2 + \text{NO} \longrightarrow \text{NO}_2 + 0.85 \cdot \text{CO} + 0.85 \cdot \text{HYAC} + \text{HO}_2 + 0.15 \cdot \text{CH}_2\text{O} + 0.15 \cdot \text{CH}_3\text{COCHO}$ | $2.464 \cdot 10^{-12} \exp(360./T)$ | $A = 2.54 \cdot 10^{-12} \cdot 0.97$; 3% MACRN-yield inferred from HYAC 42% yield at high-NO from Crounse et al. (2011) (Tyndall, p.c.) |
| $\text{MACRO}_2 + \text{NO} \longrightarrow \text{MACRN}$ | $7.620 \cdot 10^{-14} \exp(360./T)$ | $A = 2.54 \cdot 10^{-12} \cdot 0.03$; 3% MACRN-yield inferred from HYAC 42% yield at high-NO from Crounse et al. (2011) (Tyndall, p.c.) |
| $\text{MACRO}_2 + \text{NO}_3 \longrightarrow 0.85 \cdot \text{CO} + 0.85 \cdot \text{HYAC} + 0.15 \cdot \text{CH}_2\text{O} + 0.15 \cdot \text{CH}_3\text{COCHO} + \text{HO}_2 + \text{NO}_2$ | $2.500 \cdot 10^{-12}$ | |
| $\text{MACRO}_2 + \text{HO}_2 \longrightarrow 0.6 \cdot \text{MACROOH} + 0.4 \cdot \text{CO} + 0.4 \cdot \text{HYAC} + 0.4 \cdot \text{HO}_2 + 0.4 \cdot \text{OH}$ | $1.820 \cdot 10^{-13} \exp(1300./T)$ | |
| $\text{MACRO}_2 \longrightarrow \text{CO} + \text{HYAC} + \text{OH}$ | $2.900 \cdot 10^7 \exp(-5297./T)$ | Isomerisation according to Crounse et al. (2012) |
| $\text{MACROOH} + \text{OH} \longrightarrow \text{CO} + \text{HYAC} + \text{OH} + \text{H}_2\text{O}$ | $1.800 \cdot 10^{-11}$ | |
| $\text{MACROH} + \text{OH} \longrightarrow \text{CO}_2 + \text{HYAC} + \text{HO}_2 + \text{H}_2\text{O}$ | $1.800 \cdot 10^{-11}$ | |
| $\text{MPAN} + \text{M} \longrightarrow \text{MCO}_3 + \text{NO}_2 + \text{M}$ | | |
| $\text{MPAN} + \text{OH} \longrightarrow \text{CO} + \text{HYAC} + \text{NO}_2$ | $3.200 \cdot 10^{-11}$ | rate: Orlando et al. (2002) |
| $\text{MACO}_2\text{H} + \text{OH} \longrightarrow \text{CH}_2\text{O} + \text{CH}_3\text{CO}_3 + \text{CO}_2 + \text{H}_2\text{O}$ | $1.510 \cdot 10^{-11}$ | products should be pyruvic acid + CH_2O + HO_2 |
| $\text{MACO}_3\text{H} + \text{OH} \longrightarrow \text{H}_2\text{O} + \text{MCO}_3$ | $1.870 \cdot 10^{-11}$ | |
| $\text{MVK} + \text{OH} \longrightarrow \text{LHMKABO}_2$ | $4.130 \cdot 10^{-12} \exp(452./T)$ | |
| $\text{MVK} + \text{O}_3 \longrightarrow 0.85 \cdot \text{CH}_3\text{COCHO} + 0.85 \cdot \text{HCOOH} + 0.15 \cdot \text{CH}_3\text{CO}_3 + 0.15 \cdot \text{OH} + 0.15 \cdot \text{CO} + 0.15 \cdot \text{CH}_2\text{O}$ | $7.510 \cdot 10^{-16} \exp(-1521./T)$ | products according to IUPAC (2006) |
| $\text{LHMKABO}_2 + \text{CH}_3\text{O}_2 \longrightarrow 0.9 \cdot \text{CH}_2\text{O} + 0.35 \cdot \text{GLYALD} + 0.65 \cdot \text{HO}_2 + 0.35 \cdot \text{CH}_3\text{CO}_3 + 0.175 \cdot \text{BIACETOH} + 0.25 \cdot \text{CH}_3\text{OH} + 0.25 \cdot \text{MACROH} + 0.15 \cdot \text{CH}_3\text{COCHO} + 0.075 \cdot \text{CO}_2\text{H}_3\text{CHO}$ | $1.000 \cdot 10^{-12}$ | products: Tyndall (p.c.) |
| $\text{LHMKABO}_2 + \text{CH}_3\text{CO}_3 \longrightarrow 0.3 \cdot \text{CH}_3\text{COCHO} + 0.7 \cdot \text{GLYALD} + 0.7 \cdot \text{CH}_3\text{CO}_3 + 0.3 \cdot \text{CH}_2\text{O} + 0.3 \cdot \text{HO}_2 + \text{CH}_3\text{O}_2 + \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |

Table S10: C4 oxidation reactions (... continued)

| reaction | rate coefficient | reference |
|--|--|--|
| LHMKABO ₂ + HO ₂ → 0.34 · LHMKABOOH + 0.66 · CH ₃ CO ₃ + 0.66 · OH + 0.66 · GLYALD | $1.820 \cdot 10^{-13} \exp(1300./T)$ | Praske et al. (2015) |
| LHMKABO ₂ + NO → NO ₂ + 0.3 · CH ₂ O + 0.3 · CH ₃ COCHO + 0.3 · HO ₂ + 0.7 · CH ₃ CO ₃ + 0.7 · GLYALD | $2.438 \cdot 10^{-12} \exp(360./T)$ | $A = 2.54 \cdot 10^{-12} \cdot 0.96$; 4% nitrate yield Praske et al. (2015) |
| LHMKABO ₂ + NO → MVKN | $1.020 \cdot 10^{-13} \exp(360./T)$ | $A = 2.54 \cdot 10^{-12} \cdot 0.04$; 4% nitrate yield Praske et al. (2015) |
| LHMKABO ₂ + NO ₃ → 0.3 · CH ₃ COCHO + 0.7 · GLYALD + 0.7 · CH ₃ CO ₃ + 0.3 · CH ₂ O + 0.3 · HO ₂ + NO ₂ | $2.500 \cdot 10^{-12}$ | |
| MVKN + OH → CH ₃ COCHO + HO ₂ + CO ₂ + NO ₂ + H ₂ O | $5.600 \cdot 10^{-12}$ | k from Paulot et al. (2009); simplified products |
| MACRN + OH → CO ₂ + HYAC + NO ₂ + H ₂ O | $5.000 \cdot 10^{-11}$ | k from Paulot et al. (2009); simplified products |
| LHMKABOOH + OH → 0.3 · CO ₂ H ₃ CHO + OH + H ₂ O + 0.7 · BIACETOH | $4.500 \cdot 10^{-12}$ | |
| CO ₂ H ₃ CHO + OH → CO ₂ H ₃ CO ₃ + H ₂ O | $2.450 \cdot 10^{-11}$ | |
| CO ₂ H ₃ CHO + NO ₃ → CO ₂ H ₃ CO ₃ + HNO ₃ | $5.760 \cdot 10^{-12} \exp(-1862./T)$ | |
| CO ₂ H ₃ CO ₃ + CH ₃ O ₂ → CH ₃ COCHO + 2 · HO ₂ + CO ₂ + CH ₂ O | $1.000 \cdot 10^{-11}$ | products: Tyndall (p.c.) |
| CO ₂ H ₃ CO ₃ + CH ₃ CO ₃ → CH ₃ COCHO + HO ₂ + CH ₃ O ₂ + 2 · CO ₂ | $1.000 \cdot 10^{-11}$ | |
| CO ₂ H ₃ CO ₃ + HO ₂ → 0.6 · CO ₂ H ₃ CO ₃ H + 0.4 · CO ₂ + 0.4 · OH + 0.4 · HO ₂ + 0.4 · CH ₃ COCHO | $4.300 \cdot 10^{-13} \exp(1040./T)$ | |
| CO ₂ H ₃ CO ₃ + NO → CH ₃ COCHO + HO ₂ + NO ₂ + CO ₂ | $8.100 \cdot 10^{-12} \exp(270./T)$ | |
| CO ₂ H ₃ CO ₃ + NO ₃ → CH ₃ COCHO + HO ₂ + NO ₂ + CO ₂ | $4.000 \cdot 10^{-12}$ | |
| CO ₂ H ₃ CO ₃ H + OH → CO ₂ H ₃ CO ₃ + H ₂ O | $1.000 \cdot 10^{-12}$ | Orlando (p.c.) |
| MALO ₂ + NO ₂ + M → 0.8 · LC ₅ PAN ₁₇₁₉ + M | $ktroe(2.700 \cdot 10^{-28}, 7.1, 1.200 \cdot 10^{-11}, 0.9, 0.6)$ | Orlando and Tyndall (2012) - Same as k(CH ₃ CO ₃ + NO ₂); LC ₅ PAN as a surrogate |
| MALO ₂ + NO → 0.4 · GLYOXAL + HO ₂ + 0.4 · CO + 0.4 · CO ₂ + NO ₂ + 0.6 · CO ₂ H ₃ CHO | $7.500 \cdot 10^{-12} \exp(290./T)$ | products: Tyndall (p.c.) |
| MALO ₂ + HO ₂ → 0.16 · GLYOXAL + HO ₂ + 0.16 · CO + 0.16 · CO ₂ + 0.16 · OH + 0.84 · CO ₂ H ₃ CHO | $4.300 \cdot 10^{-13} \exp(1040./T)$ | |
| MALO ₂ + CH ₃ O ₂ → 0.4 · GLYOXAL + 2 · HO ₂ + 0.4 · CO + 0.4 · CO ₂ + 0.6 · CO ₂ H ₃ CHO + CH ₂ O | $1.000 \cdot 10^{-12}$ | products: Tyndall (p.c.) |
| MALO ₂ + CH ₃ CO ₃ → 0.4 · GLYOXAL + HO ₂ + 0.4 · CO + 1.4 · CO ₂ + 0.6 · CO ₂ H ₃ CHO + CH ₃ O ₂ | $1.000 \cdot 10^{-11}$ | |

Table S10: C4 oxidation reactions (... continued)

| reaction | rate coefficient | reference |
|--|--|--|
| MDIALO ₂ + HO ₂ \longrightarrow 0.4 · OH + 0.332 · HO ₂ + 0.068 · CH ₃ COCHO + 0.136 · CO + 0.068 · CH ₃ O ₂ + 0.068 · GLYOXAL | $4.300 \cdot 10^{-13} \exp(1040./T)$ | |
| MDIALO ₂ + NO \longrightarrow NO ₂ + 0.83 · HO ₂ + 0.17 · CH ₃ COCHO + 0.34 · CO + 0.17 · CH ₃ O ₂ + 0.17 · GLYOXAL | $7.500 \cdot 10^{-12} \exp(290./T)$ | |
| MDIALO ₂ + NO ₂ + M \longrightarrow M | $ktroe(2.700 \cdot 10^{-28}, 7.1, 1.200 \cdot 10^{-11}, 0.9, 0.6)$ | Orlando and Tyndall (2012); Same as k(CH ₃ CO ₃ + NO ₂) |
| MDIALO ₂ + CH ₃ O ₂ \longrightarrow 1.83 · HO ₂ + 0.17 · CH ₃ COCHO + 0.34 · CO + 0.17 · CH ₃ O ₂ + 0.17 · GLYOXAL + CH ₂ O | $1.000 \cdot 10^{-12}$ | |
| MDIALO ₂ + CH ₃ CO ₃ \longrightarrow 0.83 · HO ₂ + 0.17 · CH ₃ COCHO + 0.34 · CO + 1.17 · CH ₃ O ₂ + 0.17 · GLYOXAL + CO ₂ | $1.000 \cdot 10^{-11}$ | |

Table S11: C5 oxidation reactions

| reaction | rate coefficient | reference |
|---|---------------------------------------|--|
| BIGALKANE + OH \longrightarrow ALKO ₂ + H ₂ O | $3.500 \cdot 10^{-12}$ | |
| C ₅ H ₈ + OH \longrightarrow 0.4 · LISOPACO ₂ + 0.35 · ISOPBO ₂ + 0.25 · ISOPDO ₂ | $2.700 \cdot 10^{-11} \exp(390./T)$ | Tyndall (p.c.); MCM3.2 has yields .25, .5, .25 |
| C ₅ H ₈ + O ₃ \longrightarrow 0.051 · CH ₃ O ₂ + 0.1575 · CH ₃ CO ₃ + 0.054 · LHMVKABO ₂ + 0.522 · CO + 0.068750 · HCOOH + 0.11 · H ₂ O ₂ + 0.324750 · MACR + 0.1275 · C ₃ H ₆ + 0.2625 · HO ₂ + 0.255 · CO ₂ + 0.749750 · CH ₂ O + 0.041250 · MACO ₂ H + 0.27 · OH + 0.244 · MVK | $7.860 \cdot 10^{-15} \exp(-1913./T)$ | |
| C ₅ H ₈ + NO ₃ \longrightarrow NISOPO ₂ | $3.030 \cdot 10^{-12} \exp(-446./T)$ | |
| MBO + OH \longrightarrow MBOO ₂ | $8.100 \cdot 10^{-12} \exp(610./T)$ | |
| MBO + O ₃ \longrightarrow 0.35 · CO + 0.5 · CH ₂ O + 0.1 · CH ₃ COCH ₃ + 0.9 · IBUTALOH + 0.25 · HCOOH + 0.06 · HO ₂ + 0.06 · OH | $1.000 \cdot 10^{-17}$ | |
| MBO + NO ₃ \longrightarrow MBONO ₃ O ₂ | $4.600 \cdot 10^{-14} \exp(-400./T)$ | |
| ALKO ₂ + NO \longrightarrow 0.4 · CH ₃ CHO + 0.25 · CH ₂ O + 0.25 · CH ₃ COCH ₃ + HO ₂ + 0.8 · MEK + NO ₂ | $3.780 \cdot 10^{-12} \exp(180./T)$ | $A = 4.200 \cdot 10^{-12} \cdot 0.9$; 10% ALKNO ₃ -yield |
| ALKO ₂ + NO \longrightarrow ALKNO ₃ | $4.200 \cdot 10^{-13} \exp(180./T)$ | 10% ALKNO ₃ -yield |
| ALKO ₂ + HO ₂ \longrightarrow ALKOOH | $7.500 \cdot 10^{-13} \exp(700./T)$ | |
| ALKO ₂ + CH ₃ O ₂ \longrightarrow 0.3 · CH ₃ CHO + 1.1875 · CH ₂ O + 0.1875 · CH ₃ COCH ₃ + 0.75 · HO ₂ + 0.6 · MEK + 0.25 · ALKOH | $1.000 \cdot 10^{-12}$ | |

Table S11: C5 oxidation reactions (... continued)

| reaction | rate coefficient | reference |
|---|---------------------------------------|--|
| $\text{ALKO}_2 + \text{CH}_3\text{CO}_3 \longrightarrow 0.4 \cdot \text{CH}_3\text{CHO} + 0.25 \cdot \text{CH}_2\text{O} + 0.25 \cdot \text{CH}_3\text{COCH}_3 + \text{HO}_2 + 0.8 \cdot \text{MEK} + \text{CH}_3\text{O}_2 + \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{ALKOOH} + \text{OH} \longrightarrow \text{ALKO}_2 + \text{H}_2\text{O}$ | $3.800 \cdot 10^{-12} \exp(200./T)$ | |
| $\text{ALKOH} + \text{OH} \longrightarrow 1.25 \cdot \text{MEK} + \text{HO}_2 + \text{H}_2\text{O}$ | $5.000 \cdot 10^{-12}$ | Tyndall (p.c.), MEK yield to account for C |
| $\text{ALKNO}_3 + \text{OH} \longrightarrow 0.4 \cdot \text{CH}_3\text{CHO} + 0.25 \cdot \text{CH}_2\text{O} + 0.25 \cdot \text{CH}_3\text{COCH}_3 + \text{HO}_2 + 0.8 \cdot \text{MEK} + \text{NO}_2$ | $2.000 \cdot 10^{-12}$ | |
| $\text{LISOPACO}_2 + \text{HO}_2 \longrightarrow \text{LISOPACOOH}$ | $2.050 \cdot 10^{-13} \exp(1300./T)$ | add OH channel |
| $\text{LISOPACO}_2 + \text{NO} \longrightarrow \text{HO}_2 + 0.977 \cdot \text{LHC}_4\text{ACCHO} + \text{NO}_2 + 0.0277 \cdot \text{CH}_3\text{COCHO} + 0.0277 \cdot \text{GLYOXAL} + 0.0277 \cdot \text{HYAC} + 0.0277 \cdot \text{GLYALD}$ | $2.235 \cdot 10^{-12} \exp(360./T)$ | $A = 2.54 \cdot 10^{-12} \cdot 0.88$; average 12% nitrate yield from Paulot et al. (2009); Tyndall (p.c.); direct GLYOXAL channel from lab meas. |
| $\text{LISOPACO}_2 + \text{NO} \longrightarrow \text{LISOPACNO}_3$ | $3.050 \cdot 10^{-13} \exp(360./T)$ | $A = 2.54 \cdot 10^{-12} \cdot 0.12$; average 12% nitrate yield from Paulot et al. (2009) |
| $\text{LISOPACO}_2 + \text{NO}_3 \longrightarrow \text{HO}_2 + \text{LHC}_4\text{ACCHO} + \text{NO}_2$ | $2.500 \cdot 10^{-12}$ | |
| $\text{LISOPACO}_2 + \text{CH}_3\text{O}_2 \longrightarrow 0.75 \cdot \text{CH}_2\text{O} + 0.75 \cdot \text{LHC}_4\text{ACCHO} + 0.25 \cdot \text{CH}_3\text{OH} + 0.25 \cdot \text{ISOPAOH} + \text{HO}_2$ | $2.400 \cdot 10^{-12}$ | products: Tyndall (p.c.) |
| $\text{LISOPACO}_2 + \text{CH}_3\text{CO}_3 \longrightarrow \text{HO}_2 + \text{LHC}_4\text{ACCHO} + \text{CH}_3\text{O}_2 + \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{LISOPACO}_2 \longrightarrow \text{HO}_2 + \text{HPALD}$ | $4.1 \cdot 10^8 \exp(-7700./T)$ | Bulk isomerization rate constant for all ISOPO2 by Crounse et al. (2011) |
| $\text{ISOPBO}_2 \longrightarrow \text{HO}_2 + \text{HPALD}$ | $4.1 \cdot 10^8 \exp(-7700./T)$ | see note of [LISOPACO2] |
| $\text{ISOPDO}_2 \longrightarrow \text{HO}_2 + \text{HPALD}$ | $4.1 \cdot 10^8 \exp(-7700./T)$ | see note of [LISOPACO2] |
| $\text{LISOPACOOH} + \text{OH} \longrightarrow 0.415 \cdot \text{LIEPOX} + 0.415 \cdot \text{OH} + 0.415 \cdot \text{LISOPOOHO}_2 + 0.14 \cdot \text{LHC}_4\text{ACCHO} + 0.03 \cdot \text{H}_2\text{O} + 0.03 \cdot \text{LISOPACO}_2$ | $1.540 \cdot 10^{-10}$ | k from MCMv3.3.1 and OH-addition branching ratios estimated with site-specific SAR by Peeters et al. (2007) and H-abstraction channel assumed to be like the one for $\text{CH}_3\text{OOH} + \text{OH}$ reaction and abstraction from the alpha-hydroperoxyl allyl hydrogen estimated by SAR of MOM (Taraborrelli in prep.) being $2.12 \cdot 10^{-11}$ |
| $\text{ISOPAOH} + \text{OH} \longrightarrow \text{LISOPOOHO}_2$ | $9.300 \cdot 10^{-11}$ | OH-addition to double bond and products approximated with the one from $\text{ISOPOOH} + \text{OH}$ reaction leading to similar SOA precursors. |
| $\text{LISOPACNO}_3 + \text{OH} \longrightarrow \text{LISOPNO}_3\text{O}_2$ | $6.000 \cdot 10^{-11}$ | |
| $\text{LIEPOX} + \text{OH} \longrightarrow 0.29 \cdot \text{IEC}_1\text{O}_2 + 0.71 \cdot \text{LIECHO} + 0.71 \cdot \text{HO}_2 + \text{H}_2\text{O}$ | $1.500 \cdot 10^{-11}$ | MCM3.2 |
| $\text{LIECHO} + \text{OH} \longrightarrow \text{H}_2\text{O} + \text{LIECO}_3$ | $1.760 \cdot 10^{-11}$ | MCM3.2 |
| $\text{LIECHO} + \text{NO}_3 \longrightarrow \text{HNO}_3 + \text{LIECO}_3$ | $1.050 \cdot 10^{-11} \exp(-1860./T)$ | MCM3.2 |

Table S11: C5 oxidation reactions (... continued)

| reaction | rate coefficient | reference |
|---|--------------------------------------|---|
| $\text{LIECO}_3 + \text{HO}_2 \longrightarrow 0.6 \cdot \text{LIECO}_3\text{H} + 0.4 \cdot \text{CO}_2 + 0.4 \cdot \text{OH} + 0.25 \cdot \text{LHMKABO}_2 + 0.15 \cdot \text{MACRO}_2$ | $5.200 \cdot 10^{-13} \exp(980./T)$ | MCM3.2 |
| $\text{LIECO}_3 + \text{NO} \longrightarrow 0.6 \cdot \text{LHMKABO}_2 + 0.4 \cdot \text{MACRO}_2 + \text{NO}_2 + \text{CO}_2$ | $7.500 \cdot 10^{-12} \exp(290./T)$ | MCM3.2 |
| $\text{LIECO}_3 + \text{NO}_3 \longrightarrow 0.6 \cdot \text{LHMKABO}_2 + 0.4 \cdot \text{MACRO}_2 + \text{NO}_2 + \text{CO}_2$ | $4.000 \cdot 10^{-12}$ | MCM3.2 |
| $\text{LIECO}_3\text{H} + \text{OH} \longrightarrow \text{H}_2\text{O} + \text{LIECO}_3$ | $1.040 \cdot 10^{-11}$ | MCM3.2 |
| $\text{IEC}_1\text{O}_2 + \text{HO}_2 \longrightarrow \text{LIECO}_3\text{H}$ | $2.050 \cdot 10^{-13} \exp(1300./T)$ | MCM3.2 plus shortcut |
| $\text{IEC}_1\text{O}_2 + \text{NO} \longrightarrow \text{BIACETOH} + \text{NO}_2 + \text{CH}_2\text{O} + \text{HO}_2$ | $2.700 \cdot 10^{-12} \exp(360./T)$ | MCM3.2 |
| $\text{IEC}_1\text{O}_2 + \text{NO}_3 \longrightarrow \text{BIACETOH} + \text{NO}_2 + \text{CH}_2\text{O} + \text{HO}_2$ | $2.300 \cdot 10^{-12}$ | MCM3.2 |
| $\text{ISOPBO}_2 + \text{HO}_2 \longrightarrow \text{ISOPBOOH}$ | $2.050 \cdot 10^{-13} \exp(1300./T)$ | |
| $\text{ISOPBO}_2 + \text{NO} \longrightarrow \text{CH}_2\text{O} + \text{MVK} + \text{HO}_2 + \text{NO}_2$ | $2.235 \cdot 10^{-12} \exp(360./T)$ | $A = 2.54 \cdot 10^{-12} \cdot 0.88$; average 12% nitrate yield from Paulot et al. (2009); |
| $\text{ISOPBO}_2 + \text{NO} \longrightarrow \text{ISOPBNO}_3$ | $3.050 \cdot 10^{-13} \exp(360./T)$ | $A = 2.54 \cdot 10^{-12} \cdot 0.12$; average 12% nitrate yield from Paulot et al. (2009); |
| $\text{ISOPBO}_2 + \text{NO}_3 \longrightarrow \text{CH}_2\text{O} + \text{MVK} + \text{HO}_2 + \text{NO}_2$ | $2.500 \cdot 10^{-12}$ | |
| $\text{ISOPBO}_2 + \text{CH}_3\text{O}_2 \longrightarrow 0.75 \cdot \text{MVK} + 1.75 \cdot \text{CH}_2\text{O} + 1.5 \cdot \text{HO}_2 + 0.25 \cdot \text{ISOPBOH}$ | $8.000 \cdot 10^{-13}$ | products: Tyndall (p.c.) |
| $\text{ISOPBO}_2 + \text{CH}_3\text{CO}_3 \longrightarrow \text{CH}_2\text{O} + \text{MVK} + \text{HO}_2 + \text{CH}_3\text{O}_2 + \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{ISOPBOOH} + \text{OH} \longrightarrow 0.63 \cdot \text{LIEPOX} + 0.63 \cdot \text{OH} + 0.12 \cdot \text{LISOPOOHO}_2 + 0.15 \cdot \text{H}_2\text{O} + 0.15 \cdot \text{ISOPBO}_2$ | $7.500 \cdot 10^{-11}$ | St. Clair et al. (2015) |
| $\text{ISOPBOH} + \text{OH} \longrightarrow \text{LISOPOOHO}_2$ | $3.850 \cdot 10^{-11}$ | OH-addition to double bond and products approximated with the one from $\text{ISOPOOH} + \text{OH}$ reaction leading to similar SOA precursors. |
| $\text{ISOPBNO}_3 + \text{OH} \longrightarrow \text{LISOPNO}_3\text{O}_2$ | $1.360 \cdot 10^{-11}$ | |
| $\text{ISOPDO}_2 + \text{HO}_2 \longrightarrow \text{ISOPDOOH}$ | $2.050 \cdot 10^{-13} \exp(1300./T)$ | |
| $\text{ISOPDO}_2 + \text{NO} \longrightarrow \text{CH}_2\text{O} + \text{MACR} + \text{HO}_2 + \text{NO}_2$ | $2.235 \cdot 10^{-12} \exp(360./T)$ | $A = 2.54 \cdot 10^{-12} \cdot 0.88$; average 12% nitrate yield from Paulot et al. (2009); |
| $\text{ISOPDO}_2 + \text{NO} \longrightarrow \text{ISOPDNO}_3$ | $3.050 \cdot 10^{-13} \exp(360./T)$ | $A = 2.54 \cdot 10^{-12} \cdot 0.12$; average 12% nitrate yield from Paulot et al. (2009); |
| $\text{ISOPDO}_2 + \text{NO}_3 \longrightarrow \text{CH}_2\text{O} + \text{MACR} + \text{HO}_2 + \text{NO}_2$ | $2.500 \cdot 10^{-12}$ | |
| $\text{ISOPDO}_2 + \text{CH}_3\text{O}_2 \longrightarrow 0.5 \cdot \text{MACR} + 1.25 \cdot \text{CH}_2\text{O} + \text{HO}_2 + 0.25 \cdot \text{CH}_3\text{OH} + 0.25 \cdot \text{HCOC}_5 + 0.25 \cdot \text{ISOPDOH}$ | $2.900 \cdot 10^{-12}$ | products: Tyndall (p.c.) |
| $\text{ISOPDO}_2 + \text{CH}_3\text{CO}_3 \longrightarrow \text{CH}_2\text{O} + \text{MACR} + \text{HO}_2 + \text{CH}_3\text{O}_2 + \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |

Table S11: C5 oxidation reactions (... continued)

| reaction | rate coefficient | reference |
|--|---------------------------------------|---|
| ISOPDOOH + OH \longrightarrow 0.79 · LIEPOX + 0.79 · OH + 0.14 · LISOPOOHO ₂ + 0.07 · H ₂ O + 0.07 · ISOPDO ₂ | $1.180 \cdot 10^{-10}$ | St. Clair et al. (2015) |
| ISOPDOH + OH \longrightarrow LISOPOOHO ₂ | $7.380 \cdot 10^{-11}$ | OH-addition to double bond and products approximated with the one from ISOPOOH + OH reaction leading to similar SOA precursors. |
| ISOPDNO ₃ + OH \longrightarrow LISOPNO ₃ O ₂ | $6.100 \cdot 10^{-11}$ | OH-addition to double bond |
| NISOPO ₂ + HO ₂ \longrightarrow NISOPOOH | $2.050 \cdot 10^{-13} \exp(1300./T)$ | |
| NISOPO ₂ + NO \longrightarrow HO ₂ + NC ₄ CHO + NO ₂ | $2.540 \cdot 10^{-12} \exp(360./T)$ | |
| NISOPO ₂ + NO ₃ \longrightarrow HO ₂ + NC ₄ CHO + NO ₂ | $2.500 \cdot 10^{-12}$ | |
| NISOPO ₂ + CH ₃ O ₂ \longrightarrow 0.75 · CH ₂ O + 0.75 · NC ₄ CHO + HO ₂ + 0.25 · CH ₃ OH + 0.25 · LISOPACNO ₃ | $1.300 \cdot 10^{-12}$ | products: Tyndall (p.c.) |
| NISOPO ₂ + CH ₃ CO ₃ \longrightarrow HO ₂ + NC ₄ CHO + CH ₃ O ₂ + CO ₂ | $1.000 \cdot 10^{-11}$ | |
| NISOPOOH + OH \longrightarrow NC ₄ CHO + OH + H ₂ O | $1.030 \cdot 10^{-10}$ | |
| NC ₄ CHO + OH \longrightarrow H ₂ O + LNISO ₃ | $4.160 \cdot 10^{-11}$ | |
| NC ₄ CHO + O ₃ \longrightarrow 0.445 · NO ₂ + 0.89 · CO + 0.075625 · H ₂ O ₂ + 0.034375 · HCOCO ₂ H + 0.555 · NOA + 0.445 · HO ₂ + 0.520625 · GLYOXAL + 0.89 · OH + 0.445 · CH ₃ COCHO | $2.400 \cdot 10^{-17}$ | |
| NC ₄ CHO + NO ₃ \longrightarrow HNO ₃ + LNISO ₃ | $6.120 \cdot 10^{-12} \exp(-1862./T)$ | |
| LNISO ₃ + HO ₂ \longrightarrow 0.8 · LNISOOH + 0.2 · NOA + 0.2 · OH + 0.2 · CO ₂ + 0.2 · CO + 0.2 · HO ₂ | $1.930 \cdot 10^{-13} \exp(1300./T)$ | products: Tyndall (p.c.) |
| LNISO ₃ + NO \longrightarrow NOA + 0.5 · GLYOXAL + 0.5 · CO + HO ₂ + NO ₂ + 0.5 · CO ₂ | $4.270 \cdot 10^{-12} \exp(360./T)$ | |
| LNISO ₃ + NO ₃ \longrightarrow NOA + 0.5 · GLYOXAL + 0.5 · CO + HO ₂ + NO ₂ + 0.5 · CO ₂ | $3.302 \cdot 10^{-12} \exp(360./T)$ | |
| LNISO ₃ + CH ₃ O ₂ \longrightarrow 0.375 · GLYOXAL + 0.875 · NOA + CH ₂ O + 1.75 · HO ₂ + 0.625 · CO ₂ + 0.0625 · MACRN + 0.0625 · MVKN + 0.5 · CO | $1.000 \cdot 10^{-12}$ | products: Tyndall (p.c.) |
| LNISO ₃ + CH ₃ CO ₃ \longrightarrow NOA + 0.5 · GLYOXAL + 0.5 · CO + HO ₂ + CH ₃ O ₂ + 1.5 · CO ₂ | $1.000 \cdot 10^{-11}$ | |
| LNISOOH + OH \longrightarrow H ₂ O + LNISO ₃ | $2.650 \cdot 10^{-11}$ | |
| LHC ₄ ACCHO + OH \longrightarrow 0.52 · LC ₅₇₈ O ₂ + 0.48 · LHC ₄ ACCO ₃ + H ₂ O | $4.520 \cdot 10^{-11}$ | |

Table S11: C5 oxidation reactions (... continued)

| reaction | rate coefficient | reference |
|---|---------------------------------------|--------------------------|
| $\text{LHC}_4\text{ACCHO} + \text{O}_3 \longrightarrow 0.2225 \cdot \text{CH}_3\text{CO}_3 +$ $0.89 \cdot \text{CO} + 0.017188 \cdot \text{HOCH}_2\text{CO}_2\text{H} + 0.075625 \cdot \text{H}_2\text{O}_2 +$ $0.017188 \cdot \text{HCOCO}_2\text{H} + 0.2775 \cdot \text{HYAC} + 0.6675 \cdot \text{HO}_2 +$ $0.260313 \cdot \text{GLYOXAL} + 0.2225 \cdot \text{CH}_2\text{O} + 0.89 \cdot \text{OH} +$ $0.260313 \cdot \text{GLYALD} + 0.5 \cdot \text{CH}_3\text{COCHO}$ | $2.400 \cdot 10^{-17}$ | |
| $\text{LHC}_4\text{ACCHO} + \text{NO}_3 \longrightarrow \text{HNO}_3 + \text{LHC}_4\text{ACCO}_3$ | $6.120 \cdot 10^{-12} \exp(-1862./T)$ | |
| $\text{LC}_{578}\text{O}_2 + \text{NO} \longrightarrow 0.25 \cdot \text{CH}_3\text{COCHO} + 0.25 \cdot \text{GLYALD} +$ $0.25 \cdot \text{GLYOXAL} + 0.25 \cdot \text{HYAC} + 0.25 \cdot \text{CO} + 0.25 \cdot \text{MACROH} +$ $0.25 \cdot \text{CH}_2\text{O} + 0.25 \cdot \text{CO}_2\text{H}_3\text{CHO} + \text{HO}_2 + \text{NO}_2$ | $2.540 \cdot 10^{-12} \exp(360./T)$ | products: Tyndall (p.c.) |
| $\text{LC}_{578}\text{O}_2 + \text{NO}_3 \longrightarrow 0.25 \cdot \text{CH}_3\text{COCHO} + 0.25 \cdot \text{GLYALD} +$ $0.25 \cdot \text{GLYOXAL} + 0.25 \cdot \text{HYAC} + 0.25 \cdot \text{CO} + 0.25 \cdot \text{MACROH} +$ $0.25 \cdot \text{CH}_2\text{O} + 0.25 \cdot \text{CO}_2\text{H}_3\text{CHO} + \text{HO}_2 + \text{NO}_2$ | $2.500 \cdot 10^{-12}$ | products: Tyndall (p.c.) |
| $\text{LC}_{578}\text{O}_2 + \text{HO}_2 \longrightarrow \text{LC}_{578}\text{OOH}$ | $2.050 \cdot 10^{-13} \exp(1300./T)$ | add OH channel |
| $\text{LC}_{578}\text{O}_2 + \text{CH}_3\text{O}_2 \longrightarrow 0.156250 \cdot \text{CH}_3\text{COCHO} +$ $0.156250 \cdot \text{GLYALD} + 0.156250 \cdot \text{GLYOXAL} + 0.156250 \cdot \text{HYAC} +$ $0.156250 \cdot \text{CO} + 0.468750 \cdot \text{MACROH} + 1.031250 \cdot \text{CH}_2\text{O} +$ $0.3125 \cdot \text{CO}_2\text{H}_3\text{CHO} + 1.25 \cdot \text{HO}_2 + 0.125 \cdot \text{CH}_3\text{OH}$ | $1.000 \cdot 10^{-12}$ | products: Tyndall (p.c.) |
| $\text{LC}_{578}\text{O}_2 + \text{CH}_3\text{CO}_3 \longrightarrow 0.25 \cdot \text{CH}_3\text{COCHO} + 0.25 \cdot \text{GLYALD} +$ $0.25 \cdot \text{GLYOXAL} + 0.25 \cdot \text{HYAC} + 0.25 \cdot \text{CO} + 0.25 \cdot \text{MACROH} +$ $0.25 \cdot \text{CH}_2\text{O} + 0.25 \cdot \text{CO}_2\text{H}_3\text{CHO} + \text{CH}_3\text{O}_2 + \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{LC}_{578}\text{OOH} + \text{OH} \longrightarrow \text{H}_2\text{O} + \text{LC}_{578}\text{O}_2$ | $3.160 \cdot 10^{-11}$ | |
| $\text{LHC}_4\text{ACCO}_3 \longrightarrow \text{HO}_2 + \text{PACALD}$ | $4.1 \cdot 10^8 \exp(-7700./T)$ | see note of [LISOPACO2] |
| $\text{LHC}_4\text{ACCO}_3 + \text{CH}_3\text{O}_2 \longrightarrow \text{CH}_2\text{O} + 0.1 \cdot \text{LHC}_4\text{ACCO}_2\text{H} +$ $0.45 \cdot \text{GLYALD} + 0.45 \cdot \text{HYAC} + 0.45 \cdot \text{CH}_3\text{CO}_3 + 0.45 \cdot \text{CO} +$ $0.45 \cdot \text{HO}_2 + 0.9 \cdot \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | products: Tyndall (p.c.) |
| $\text{LHC}_4\text{ACCO}_3 + \text{CH}_3\text{CO}_3 \longrightarrow 0.5 \cdot \text{HYAC} + 0.5 \cdot \text{GLYALD} +$ $0.5 \cdot \text{CH}_3\text{CO}_3 + 0.5 \cdot \text{CO} + 0.5 \cdot \text{HO}_2 + \text{CH}_3\text{O}_2 + 2 \cdot \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{LHC}_4\text{ACCO}_3 + \text{HO}_2 \longrightarrow 0.4 \cdot \text{LHC}_4\text{ACCO}_3\text{H} +$ $0.2 \cdot \text{LHC}_4\text{ACCO}_2\text{H} + 0.2 \cdot \text{O}_3 + 0.4 \cdot \text{CO}_2 + 0.4 \cdot \text{OH} +$ $0.2 \cdot \text{HYAC} + 0.2 \cdot \text{GLYALD} + 0.2 \cdot \text{CH}_3\text{CO}_3 + 0.2 \cdot \text{CO} +$ $0.2 \cdot \text{HO}_2$ | $4.300 \cdot 10^{-13} \exp(1040./T)$ | |
| $\text{LHC}_4\text{ACCO}_3 + \text{NO} \longrightarrow 0.5 \cdot \text{HYAC} + 0.5 \cdot \text{GLYALD} +$ $0.5 \cdot \text{CH}_3\text{CO}_3 + 0.5 \cdot \text{CO} + 0.5 \cdot \text{HO}_2 + \text{NO}_2 + \text{CO}_2$ | $8.100 \cdot 10^{-12} \exp(270./T)$ | |
| $\text{LHC}_4\text{ACCO}_3 + \text{NO}_3 \longrightarrow 0.5 \cdot \text{HYAC} + 0.5 \cdot \text{GLYALD} +$ $0.5 \cdot \text{CH}_3\text{CO}_3 + 0.5 \cdot \text{CO} + 0.5 \cdot \text{HO}_2 + \text{NO}_2 + \text{CO}_2$ | $4.000 \cdot 10^{-12}$ | |

Table S11: C5 oxidation reactions (... continued)

| reaction | rate coefficient | reference |
|--|--|---|
| $\text{LHC}_4\text{ACCO}_3 + \text{NO}_2 + \text{M} \longrightarrow \text{LC}_5\text{PAN}_{1719} + \text{M}$ | $ktroe(2.700 \cdot 10^{-28}, 7.1, 1.200 \cdot 10^{-11}, 0.9, 0.3)$ | |
| $\text{LHC}_4\text{ACCO}_2\text{H} + \text{OH} \longrightarrow 0.5 \cdot \text{HYAC} + 0.5 \cdot \text{GLYALD} + 0.5 \cdot \text{CH}_3\text{CO}_3 + 0.5 \cdot \text{CO} + 0.5 \cdot \text{HO}_2 + \text{CO}_2 + \text{H}_2\text{O}$ | $2.520 \cdot 10^{-11}$ | |
| $\text{LHC}_4\text{ACCO}_3\text{H} + \text{OH} \longrightarrow \text{H}_2\text{O} + \text{LHC}_4\text{ACCO}_3$ | $2.880 \cdot 10^{-11}$ | |
| $\text{LC}_5\text{PAN}_{1719} + \text{M} \longrightarrow \text{LHC}_4\text{ACCO}_3 + \text{NO}_2 + \text{M}$ | | $k_0 = 4.9 \cdot 10^{-3} \exp(-12100./T) * M, k_{inf} = 5.4 \cdot 10^{16} \exp(-13830./T), f_c = 0.3$ |
| $\text{LC}_5\text{PAN}_{1719} + \text{OH} \longrightarrow \text{CO} + \text{MACROH} + \text{NO}_2$ | $2.520 \cdot 10^{-11}$ | |
| $\text{HCOC}_5 + \text{OH} \longrightarrow \text{C}_{59}\text{O}_2$ | $3.810 \cdot 10^{-11}$ | |
| $\text{C}_{59}\text{O}_2 + \text{CH}_3\text{O}_2 \longrightarrow 0.75 \cdot \text{HOCH}_2\text{CO}_3 + 0.75 \cdot \text{HYAC} + \text{CH}_2\text{O} + 0.75 \cdot \text{HO}_2 + 0.3125 \cdot \text{MACROH}$ | $1.000 \cdot 10^{-12}$ | Tyndall (p.c.) |
| $\text{C}_{59}\text{O}_2 + \text{CH}_3\text{CO}_3 \longrightarrow \text{HOCH}_2\text{CO}_3 + \text{HYAC} + \text{CH}_3\text{O}_2 + \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{C}_{59}\text{O}_2 + \text{NO} \longrightarrow \text{HOCH}_2\text{CO}_3 + \text{HYAC} + \text{NO}_2$ | $2.540 \cdot 10^{-12} \exp(360./T)$ | |
| $\text{C}_{59}\text{O}_2 + \text{NO}_3 \longrightarrow \text{HOCH}_2\text{CO}_3 + \text{HYAC} + \text{NO}_2$ | $2.500 \cdot 10^{-12}$ | |
| $\text{C}_{59}\text{O}_2 + \text{HO}_2 \longrightarrow \text{C}_{59}\text{OOH}$ | $2.050 \cdot 10^{-13} \exp(1300./T)$ | |
| $\text{C}_{59}\text{OOH} + \text{OH} \longrightarrow \text{C}_{59}\text{O}_2 + \text{H}_2\text{O}$ | $9.700 \cdot 10^{-12}$ | |
| $\text{MBOO}_2 + \text{NO} \longrightarrow \text{HO}_2 + 0.67 \cdot \text{CH}_3\text{COCH}_3 + 0.67 \cdot \text{GLYALD} + 0.33 \cdot \text{CH}_2\text{O} + 0.33 \cdot \text{IBUTALOH} + \text{NO}_2$ | $2.600 \cdot 10^{-12} \exp(365./T)$ | |
| $\text{MBOO}_2 + \text{CH}_3\text{O}_2 \longrightarrow 0.9165 \cdot \text{CH}_2\text{O} + 0.625 \cdot \text{MACROH} + 0.25 \cdot \text{CH}_3\text{OH} + \text{HO}_2 + 0.3335 \cdot \text{CH}_3\text{COCH}_3 + 0.3335 \cdot \text{GLYALD} + 0.1665 \cdot \text{IBUTALOH}$ | $3.750 \cdot 10^{-13} \exp(-40./T)$ | |
| $\text{MBOO}_2 + \text{CH}_3\text{CO}_3 \longrightarrow 0.67 \cdot \text{CH}_3\text{COCH}_3 + 0.67 \cdot \text{GLYALD} + 0.33 \cdot \text{CH}_2\text{O} + 0.33 \cdot \text{IBUTALOH} + \text{HO}_2 + \text{CH}_3\text{O}_2 + \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{MBOO}_2 + \text{HO}_2 \longrightarrow \text{MBOOOH}$ | $7.500 \cdot 10^{-13} \exp(700./T)$ | |
| $\text{MBOOOH} + \text{OH} \longrightarrow 0.5 \cdot \text{MBOO}_2 + 0.625 \cdot \text{MACROH} + 0.5 \cdot \text{OH} + \text{H}_2\text{O}$ | $3.800 \cdot 10^{-12} \exp(200./T)$ | |
| $\text{IBUTALOH} + \text{OH} \longrightarrow \text{H}_2\text{O} + \text{IBUTALOHO}_2$ | $1.400 \cdot 10^{-11}$ | |
| $\text{IBUTALOHO}_2 + \text{NO} \longrightarrow \text{CO}_2 + \text{NO}_2 + \text{HO}_2 + \text{CH}_3\text{COCH}_3$ | $2.600 \cdot 10^{-12} \exp(365./T)$ | |
| $\text{IBUTALOHO}_2 + \text{HO}_2 \longrightarrow 0.6 \cdot \text{IBUTALOHOOH} + 0.4 \cdot \text{HO}_2 + 0.4 \cdot \text{OH} + 0.4 \cdot \text{CH}_3\text{COCH}_3 + 0.4 \cdot \text{CO}_2$ | $4.300 \cdot 10^{-13} \exp(1040./T)$ | |
| $\text{IBUTALOHO}_2 + \text{CH}_3\text{O}_2 \longrightarrow \text{CH}_3\text{COCH}_3 + 2 \cdot \text{HO}_2 + \text{CO}_2 + \text{CH}_2\text{O}$ | $1.000 \cdot 10^{-12}$ | products: Tyndall (p.c.) |
| $\text{IBUTALOHO}_2 + \text{CH}_3\text{CO}_3 \longrightarrow \text{CH}_3\text{COCH}_3 + \text{HO}_2 + \text{CH}_3\text{O}_2 + 2 \cdot \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{IBUTALOHOOH} + \text{OH} \longrightarrow \text{H}_2\text{O} + \text{IBUTALOHO}_2$ | $1.000 \cdot 10^{-12}$ | Tyndall (p.c.) |

Table S11: C5 oxidation reactions (... continued)

| reaction | rate coefficient | reference |
|---|--|---|
| $\text{MBONO}_3\text{O}_2 + \text{HO}_2 \longrightarrow$ | $4.300 \cdot 10^{-13} \exp(1040./T)$ | |
| $\text{MBONO}_3\text{O}_2 + \text{NO} \longrightarrow 0.25 \cdot \text{CH}_2\text{O} + 0.25 \cdot \text{IBUTALOH} + 1.25 \cdot \text{NO}_2 + 0.500250 \cdot \text{NOA} + 0.75 \cdot \text{CH}_3\text{COCH}_3 + 0.75 \cdot \text{HO}_2$ | $2.600 \cdot 10^{-12} \exp(365./T)$ | |
| $\text{MBONO}_3\text{O}_2 + \text{NO}_3 \longrightarrow 0.25 \cdot \text{CH}_2\text{O} + 0.25 \cdot \text{IBUTALOH} + 1.25 \cdot \text{NO}_2 + 0.500250 \cdot \text{NOA} + 0.75 \cdot \text{CH}_3\text{COCH}_3 + 0.75 \cdot \text{HO}_2$ | $2.400 \cdot 10^{-12}$ | |
| $\text{MBONO}_3\text{O}_2 + \text{CH}_3\text{O}_2 \longrightarrow 0.875 \cdot \text{CH}_2\text{O} + 0.125 \cdot \text{IBUTALOH} + 0.125 \cdot \text{NO}_2 + 0.250125 \cdot \text{NOA} + 0.375 \cdot \text{CH}_3\text{COCH}_3 + 0.875 \cdot \text{HO}_2 + 0.25 \cdot \text{CH}_3\text{OH} + 0.625 \cdot \text{MACROH}$ | $1.000 \cdot 10^{-12}$ | Tyndall (p.c.) |
| $\text{MBONO}_3\text{O}_2 + \text{CH}_3\text{CO}_3 \longrightarrow 0.25 \cdot \text{CH}_2\text{O} + 0.25 \cdot \text{IBUTALOH} + 0.25 \cdot \text{NO}_2 + 0.500250 \cdot \text{NOA} + 0.75 \cdot \text{CH}_3\text{COCH}_3 + 0.75 \cdot \text{HO}_2 + \text{CH}_3\text{O}_2 + \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{DICARBO}_2 + \text{HO}_2 \longrightarrow 0.4 \cdot \text{OH} + 0.068 \cdot \text{CH}_3\text{COCHO} + 0.068 \cdot \text{HO}_2 + 0.068 \cdot \text{CO} + 0.4 \cdot \text{CO}_2 + 0.332 \cdot \text{CH}_3\text{CO}_3 + 0.332 \cdot \text{GLYOXAL}$ | $4.300 \cdot 10^{-13} \exp(1040./T)$ | |
| $\text{DICARBO}_2 + \text{NO} \longrightarrow 0.17 \cdot \text{CH}_3\text{COCHO} + 0.17 \cdot \text{HO}_2 + 0.17 \cdot \text{CO} + \text{CO}_2 + 0.83 \cdot \text{CH}_3\text{CO}_3 + 0.83 \cdot \text{GLYOXAL} + \text{NO}_2$ | $7.500 \cdot 10^{-12} \exp(290./T)$ | |
| $\text{DICARBO}_2 + \text{NO}_2 + \text{M} \longrightarrow \text{M}$ | $ktroe(2.700 \cdot 10^{-28}, 7.1, 1.200 \cdot 10^{-11}, 0.9, 0.6)$ | Orlando and Tyndall (2012), Same as $k(\text{CH}_3\text{CO}_3 + \text{NO}_2)$ |
| $\text{DICARBO}_2 + \text{CH}_3\text{O}_2 \longrightarrow 0.17 \cdot \text{CH}_3\text{COCHO} + 1.17 \cdot \text{HO}_2 + 0.17 \cdot \text{CO} + \text{CO}_2 + 0.83 \cdot \text{CH}_3\text{CO}_3 + 0.83 \cdot \text{GLYOXAL} + \text{CH}_2\text{O}$ | $1.000 \cdot 10^{-12}$ | |
| $\text{DICARBO}_2 + \text{CH}_3\text{CO}_3 \longrightarrow 0.17 \cdot \text{CH}_3\text{COCHO} + 0.17 \cdot \text{HO}_2 + 0.17 \cdot \text{CO} + 2 \cdot \text{CO}_2 + 0.83 \cdot \text{CH}_3\text{CO}_3 + 0.83 \cdot \text{GLYOXAL} + \text{CH}_3\text{O}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{HPALD} + \text{OH} \longrightarrow 0.641 \cdot \text{OH} + 0.385 \cdot \text{PACALD} + 0.256 \cdot \text{BIGALD}_3 + 0.359 \cdot \text{CH}_3\text{COCHO} + 0.359 \cdot \text{GLYOX} + 0.359 \cdot \text{HO}_2$ | $5.200 \cdot 10^{-11}$ | simplification of chemistry in MCMv3.3.1 |
| $\text{PACALD} + \text{OH} \longrightarrow \text{CH}_3\text{COCHO} + \text{HCOCO}_3\text{H} + \text{HO}_2$ | $4.720 \cdot 10^{-11}$ | k and products for $\text{C}_5\text{PACALD}_2$ from MCMv3.3.1 assuming an implicit $\text{RO}_2 \longrightarrow \text{RO}$ conversion |
| $\text{LISOPOOHO}_2 + \text{HO}_2 \longrightarrow \text{LISOPOOHOOH}$ | $2.050 \cdot 10^{-13} \exp(1300./T)$ | |
| $\text{LISOPOOHO}_2 + \text{NO} \longrightarrow \text{CH}_2\text{O} + \text{HO}_2 + 0.5 \cdot \text{MACROOH} + 0.5 \cdot \text{LHMKABOOH} + \text{NO}_2$ | $2.540 \cdot 10^{-12} \exp(360./T)$ | $A = 2.54 \cdot 10^{-12} \cdot 0.12$; average 12% nitrate yield from Paulot et al. (2009); nitrate yield left for the moment equal to the one of the simple ISOPO2 (it should be higher) |
| $\text{LISOPOOHO}_2 + \text{NO} \longrightarrow \text{LISOPNO}_3\text{OOH}$ | $3.050 \cdot 10^{-13} \exp(360./T)$ | $A = 2.54 \cdot 10^{-12} \cdot 0.12$; average 12% nitrate yield from Paulot et al. (2009); nitrate yield left for the moment equal to the one of the simple ISOPO2 (it should be higher) |

Table S11: C5 oxidation reactions (... continued)

| reaction | rate coefficient | reference |
|--|--------------------------------------|---|
| $\text{LISOPOOHO}_2 + \text{NO}_3 \longrightarrow \text{CH}_2\text{O} + \text{HO}_2 + 0.5 \cdot \text{MACROOH} + 0.5 \cdot \text{LHMKABOOH} + \text{NO}_2$ | $2.500 \cdot 10^{-12}$ | |
| $\text{LISOPOOHO}_2 + \text{CH}_3\text{O}_2 \longrightarrow 1.5 \cdot \text{CH}_2\text{O} + 0.75 \cdot \text{HO}_2 + 0.375 \cdot \text{MACROOH} + 0.375 \cdot \text{LHMKABOOH} + 0.25 \cdot \text{LISOPOOHOOH}$ | $8.000 \cdot 10^{-13}$ | |
| $\text{LISOPOOHO}_2 + \text{CH}_3\text{CO}_3 \longrightarrow \text{CH}_2\text{O} + \text{HO}_2 + 0.5 \cdot \text{MACROOH} + 0.5 \cdot \text{LHMKABOOH} + \text{CH}_3\text{O}_2 + \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{LISOPNO}_3\text{O}_2 + \text{HO}_2 \longrightarrow \text{LISOPNO}_3\text{OOH}$ | $2.050 \cdot 10^{-13} \exp(1300./T)$ | |
| $\text{LISOPNO}_3\text{O}_2 + \text{NO} \longrightarrow \text{CH}_2\text{O} + 0.5 \cdot \text{MACRN} + 0.5 \cdot \text{MVKN} + \text{HO}_2 + \text{NO}_2$ | $2.235 \cdot 10^{-12} \exp(360./T)$ | $A = 2.54 \cdot 10^{-12} \cdot 0.88$; average 12% nitrate yield from Paulot et al. (2009); nitrate yield left for the moment equal to the one of the simple ISOPO2 (it should be higher); previous RONO_2 split into 50% MACRN + 50% MVKN |
| $\text{LISOPNO}_3\text{O}_2 + \text{NO} \longrightarrow \text{LISOPNO}_3\text{NO}_3$ | $3.050 \cdot 10^{-13} \exp(360./T)$ | $A = 2.54 \cdot 10^{-12} \cdot 0.12$; average 12% nitrate yield from Paulot et al. (2009); nitrate yield left for the moment equal to the one of the simple ISOPO2 (it should be higher) |
| $\text{LISOPNO}_3\text{O}_2 + \text{NO}_3 \longrightarrow \text{CH}_2\text{O} + 0.5 \cdot \text{MACRN} + 0.5 \cdot \text{MVKN} + \text{HO}_2 + \text{NO}_2$ | $2.500 \cdot 10^{-12}$ | |
| $\text{LISOPNO}_3\text{O}_2 + \text{CH}_3\text{O}_2 \longrightarrow 0.3525 \cdot \text{MACRN} + 0.3525 \cdot \text{MVKN} + 1.75 \cdot \text{CH}_2\text{O} + 1.5 \cdot \text{HO}_2 + 0.25 \cdot \text{LISOPNO}_3\text{OOH}$ | $8.000 \cdot 10^{-13}$ | |
| $\text{LISOPNO}_3\text{O}_2 + \text{CH}_3\text{CO}_3 \longrightarrow \text{CH}_2\text{O} + 0.5 \cdot \text{MACRN} + 0.5 \cdot \text{MVKN} + \text{HO}_2 + \text{CH}_3\text{O}_2 + \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{LISOPOOHOOH} + \text{OH} \longrightarrow \text{H}_2\text{O} + \text{LISOPOOHO}_2$ | $7.600 \cdot 10^{-12} \exp(200./T)$ | twice the $k(\text{CH}_3\text{OOH} + \text{OH} \longrightarrow \text{CH}_3\text{O}_2)$ |
| $\text{LISOPOOHOOH} + \text{OH} \longrightarrow \text{LC}_{578}\text{OOH} + \text{OH}$ | $2.104 \cdot 10^{-11}$ | k for H-abstractions from SAR in MOM by a secondary carbon bearing a -OH group and a secondary and a tertiary carbon atoms bearing an -OOH group: $8.42 \cdot 10^{-13} \cdot 3.44 + (8.42 \cdot 10^{-13} + 1.75 \cdot 10^{-12}) \cdot 7$ |
| $\text{LISOPNO}_3\text{OOH} + \text{OH} \longrightarrow \text{H}_2\text{O} + \text{LISOPNO}_3\text{O}_2$ | $3.800 \cdot 10^{-12} \exp(200./T)$ | |
| $\text{LISOPNO}_3\text{OOH} + \text{OH} \longrightarrow \text{C}_{59}\text{OOH} + \text{OH}$ | $1.515 \cdot 10^{-11}$ | k for H-abstractions from SAR in MOM by a secondary carbon bearing a -OH group and a tertiary carbon atom bearing an -OOH group: $8.42 \cdot 10^{-13} \cdot 3.44 + 1.75 \cdot 10^{-12} \cdot 7$ |
| $\text{LISOPNO}_3\text{NO}_3 + \text{OH} \longrightarrow 0.5 \cdot \text{MACRN} + 0.5 \cdot \text{MVKN} + \text{NO}_2 + \text{CH}_2\text{O} + \text{HO}_2$ | $8.916 \cdot 10^{-12}$ | k for H-abstractions from SAR in MOM by a secondary and a tertiary carbon bearing a -OH group: $(8.42 \cdot 10^{-13} + 1.75 \cdot 10^{-12}) \cdot 3.44$; previous RONO_2 split into 50% MACRN + 50% MVKN |

Table S12: C6 oxidation reactions

| reaction | rate coefficient | reference |
|---|--------------------------------------|---|
| $\text{BENZ} + \text{OH} \longrightarrow 0.53 \cdot \text{PHENOL} + 0.12 \cdot \text{BEPOMUC} + 0.65 \cdot \text{HO}_2 + 0.35 \cdot \text{BENZO}_2$ | $2.300 \cdot 10^{-12} \exp(-193./T)$ | |
| $\text{PHENOL} + \text{OH} \longrightarrow 0.14 \cdot \text{PHENO}_2 + 0.8 \cdot \text{HO}_2 + 0.8 \cdot \text{CATECHOL} + 0.06 \cdot \text{C}_6\text{H}_5\text{O}$ | $4.700 \cdot 10^{-13} \exp(1220./T)$ | |
| $\text{PHENOL} + \text{NO}_3 \longrightarrow 0.26 \cdot \text{PHENO}_2 + 0.74 \cdot \text{C}_6\text{H}_5\text{O} + 0.74 \cdot \text{HNO}_3$ | $3.800 \cdot 10^{-12}$ | NPHEO2 approximated with PHENO ₂ |
| $\text{PHENO}_2 + \text{NO} \longrightarrow \text{HO}_2 + 0.7 \cdot \text{GLYOXAL} + \text{NO}_2$ | $2.600 \cdot 10^{-12} \exp(365./T)$ | |
| $\text{PHENO}_2 + \text{HO}_2 \longrightarrow \text{PHENOOH}$ | $7.500 \cdot 10^{-13} \exp(700./T)$ | |
| $\text{PHENO}_2 + \text{CH}_3\text{O}_2 \longrightarrow 2 \cdot \text{HO}_2 + 0.7 \cdot \text{GLYOXAL} + \text{CH}_2\text{O}$ | $1.000 \cdot 10^{-12}$ | |
| $\text{PHENO}_2 + \text{CH}_3\text{CO}_3 \longrightarrow \text{HO}_2 + 0.7 \cdot \text{GLYOXAL} + \text{CH}_3\text{O}_2 + \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{PHENOOH} + \text{OH} \longrightarrow \text{H}_2\text{O} + \text{PHENO}_2$ | $3.800 \cdot 10^{-12} \exp(200./T)$ | |
| $\text{C}_6\text{H}_5\text{O} + \text{NO}_2 \longrightarrow$ | $2.100 \cdot 10^{-12}$ | |
| $\text{C}_6\text{H}_5\text{O} + \text{O}_3 \longrightarrow \text{C}_6\text{H}_5\text{O}_2$ | $2.800 \cdot 10^{-13}$ | |
| $\text{C}_6\text{H}_5\text{O}_2 + \text{NO} \longrightarrow \text{C}_6\text{H}_5\text{O} + \text{NO}_2$ | $2.600 \cdot 10^{-12} \exp(365./T)$ | |
| $\text{C}_6\text{H}_5\text{O}_2 + \text{NO}_3 \longrightarrow \text{C}_6\text{H}_5\text{O} + \text{NO}_2$ | $2.300 \cdot 10^{-12}$ | MCM3.2 |
| $\text{C}_6\text{H}_5\text{O}_2 + \text{HO}_2 \longrightarrow \text{C}_6\text{H}_5\text{OOH}$ | $7.500 \cdot 10^{-13} \exp(700./T)$ | |
| $\text{C}_6\text{H}_5\text{O}_2 + \text{CH}_3\text{O}_2 \longrightarrow \text{C}_6\text{H}_5\text{O} + \text{CH}_2\text{O} + \text{HO}_2$ | $1.000 \cdot 10^{-12}$ | |
| $\text{C}_6\text{H}_5\text{O}_2 + \text{CH}_3\text{CO}_3 \longrightarrow \text{C}_6\text{H}_5\text{O} + \text{CH}_3\text{O}_2 + \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{C}_6\text{H}_5\text{OOH} + \text{OH} \longrightarrow \text{C}_6\text{H}_5\text{O}_2$ | $3.800 \cdot 10^{-12} \exp(200./T)$ | |
| $\text{BENZO}_2 + \text{NO} \longrightarrow \text{GLYOXAL} + \text{NO}_2 + 0.5 \cdot \text{BIGALD}_1 + \text{HO}_2$ | $2.600 \cdot 10^{-12} \exp(365./T)$ | MCM3.2 |
| $\text{BENZO}_2 + \text{HO}_2 \longrightarrow \text{BENZOOH}$ | $7.500 \cdot 10^{-13} \exp(700./T)$ | |
| $\text{BENZO}_2 + \text{CH}_3\text{O}_2 \longrightarrow \text{GLYOXAL} + 0.5 \cdot \text{BIGALD}_1 + 2 \cdot \text{HO}_2 + \text{CH}_2\text{O}$ | $1.000 \cdot 10^{-12}$ | |
| $\text{BENZO}_2 + \text{CH}_3\text{CO}_3 \longrightarrow \text{GLYOXAL} + 0.5 \cdot \text{BIGALD}_1 + \text{HO}_2 + \text{CH}_3\text{O}_2 + \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{BENZOOH} + \text{OH} \longrightarrow \text{BENZO}_2$ | $3.800 \cdot 10^{-12} \exp(200./T)$ | |
| $\text{CATECHOL} + \text{OH} \longrightarrow \text{CATEC}_1\text{O}$ | $1.000 \cdot 10^{-10}$ | |
| $\text{CATECHOL} + \text{NO}_3 \longrightarrow \text{CATEC}_1\text{O} + \text{HNO}_3$ | $9.900 \cdot 10^{-11}$ | |
| $\text{CATEC}_1\text{O} + \text{NO}_2 \longrightarrow$ | $2.100 \cdot 10^{-12}$ | |
| $\text{CATEC}_1\text{O} + \text{O}_3 \longrightarrow \text{CATEC}_1\text{O}_2$ | $2.800 \cdot 10^{-13}$ | |
| $\text{CATEC}_1\text{O}_2 + \text{HO}_2 \longrightarrow \text{CATEC}_1\text{OOH}$ | $7.500 \cdot 10^{-13} \exp(700./T)$ | |
| $\text{CATEC}_1\text{O}_2 + \text{NO} \longrightarrow \text{CATEC}_1\text{O} + \text{NO}_2$ | $2.600 \cdot 10^{-12} \exp(365./T)$ | |
| $\text{CATEC}_1\text{O}_2 + \text{NO}_3 \longrightarrow \text{CATEC}_1\text{O} + \text{NO}_2$ | $2.300 \cdot 10^{-12}$ | MCM3.2 |
| $\text{CATEC}_1\text{O}_2 + \text{CH}_3\text{O}_2 \longrightarrow \text{CATEC}_1\text{O} + \text{CH}_2\text{O} + \text{HO}_2$ | $1.000 \cdot 10^{-12}$ | |
| $\text{CATEC}_1\text{O}_2 + \text{CH}_3\text{CO}_3 \longrightarrow \text{CATEC}_1\text{OCH}_3\text{O}_2 + \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{CATEC}_1\text{OOH} + \text{OH} \longrightarrow \text{CATEC}_1\text{O}_2$ | $1.900 \cdot 10^{-12} \exp(190./T)$ | |

Table S13: C7 oxidation reactions

| reaction | rate coefficient | reference |
|--|---|--|
| TOL + OH \longrightarrow 0.18 · CRESOL + 0.1 · TEPOMUC + 0.07 · BZOO + 0.65 · TOLO ₂ + 0.28 · HO ₂ | $1.700 \cdot 10^{-12} \exp(352./T)$ | |
| CRESOL + OH \longrightarrow 0.2 · PHENO ₂ + 0.73 · HO ₂ + 0.73 · CATECHOL + 0.07 · C ₆ H ₅ O | $4.700 \cdot 10^{-11}$ | CATECHOL and PHENO ₂ omits one CH ₃ group of MCATECHOL and CRESO ₂ |
| CRESOL + NO ₃ \longrightarrow 0.61 · PHENO ₂ + 0.39 · C ₆ H ₅ O + 0.49 · HNO ₃ | $1.400 \cdot 10^{-11}$ | CRESO ₂ and NCRESO ₂ approximated with PHENO ₂ ; TOL ₁ O with C ₆ H ₅ O (see MCM3.2 for details) |
| TOLO ₂ + HO ₂ \longrightarrow TOLOOH | $7.500 \cdot 10^{-13} \exp(700./T)$ | |
| TOLO ₂ + NO \longrightarrow NO ₂ + 0.6 · GLYOXAL + 0.4 · CH ₃ COCHO + HO ₂ + 0.2 · BIGALD ₁ + 0.2 · BIGALD ₂ + 0.2 · BIGALD ₃ | $2.600 \cdot 10^{-12} \exp(365./T)$ | |
| TOLO ₂ + CH ₃ O ₂ \longrightarrow 0.6 · GLYOXAL + 0.4 · CH ₃ COCHO + 2 · HO ₂ + CH ₂ O | $1.000 \cdot 10^{-12}$ | |
| TOLO ₂ + CH ₃ CO ₃ \longrightarrow 0.6 · GLYOXAL + 0.4 · CH ₃ COCHO + HO ₂ + CH ₃ O ₂ + CO ₂ + 0.2 · BIGALD ₁ + 0.2 · BIGALD ₂ + 0.2 · BIGALD ₃ | $1.000 \cdot 10^{-11}$ | |
| TOLOOH + OH \longrightarrow TOLO ₂ | $3.800 \cdot 10^{-12} \exp(200./T)$ | |
| BZOO + HO ₂ \longrightarrow BZOOH | $7.500 \cdot 10^{-13} \exp(700./T)$ | |
| BZOO + NO \longrightarrow BZALD + NO ₂ + HO ₂ | $2.600 \cdot 10^{-12} \exp(365./T)$ | MCM3.2 forms 10% nitrate |
| BZOO + CH ₃ O ₂ \longrightarrow BZALD + 2 · HO ₂ + CH ₂ O | $1.000 \cdot 10^{-12}$ | |
| BZOO + CH ₃ CO ₃ \longrightarrow BZALD + HO ₂ + CH ₃ O ₂ + CO ₂ | $1.000 \cdot 10^{-11}$ | |
| BZOOH + OH \longrightarrow BZOO | $3.800 \cdot 10^{-12} \exp(200./T)$ | |
| BZALD + OH \longrightarrow ACBZO ₂ | $5.900 \cdot 10^{-12} \exp(225./T)$ | |
| ACBZO ₂ + NO ₂ + M \longrightarrow M + PBZNIT | <i>ktroe</i> ($2.700 \cdot 10^{-28}$, 7.1, $1.200 \cdot 10^{-11}$, 0.9, 0.6) | Orlando and Tyndall (2012) - Same as k(CH ₃ CO ₃ + NO ₂) |
| PBZNIT + M \longrightarrow ACBZO ₂ + NO ₂ + M | | |
| ACBZO ₂ + NO \longrightarrow C ₆ H ₅ O ₂ + NO ₂ | $7.500 \cdot 10^{-12} \exp(290./T)$ | |
| ACBZO ₂ + HO ₂ \longrightarrow 0.4 · C ₆ H ₅ O ₂ + 0.4 · OH | $4.300 \cdot 10^{-13} \exp(1040./T)$ | |
| ACBZO ₂ + CH ₃ O ₂ \longrightarrow C ₆ H ₅ O ₂ + CH ₂ O + HO ₂ | $1.000 \cdot 10^{-12}$ | |
| ACBZO ₂ + CH ₃ CO ₃ \longrightarrow C ₆ H ₅ O ₂ + CH ₃ O ₂ + CO ₂ | $1.000 \cdot 10^{-11}$ | |

Table S14: C8 oxidation reactions

| reaction | rate coefficient | reference |
|--|------------------------|-----------|
| XYL + OH \longrightarrow 0.15 · XYLOL + 0.23 · TEPOMUC + 0.06 · BZOO + 0.56 · XYLENO ₂ + 0.38 · HO ₂ | $1.700 \cdot 10^{-11}$ | |

Table S14: C8 oxidation reactions (... continued)

| reaction | rate coefficient | reference |
|--|-------------------------------------|--|
| $\text{XYLOL} + \text{OH} \longrightarrow 0.3 \cdot \text{XYLOLO}_2 + 0.63 \cdot \text{HO}_2 + 0.63 \cdot \text{CATECHOL} + 0.07 \cdot \text{C}_6\text{H}_5\text{O}$ | $8.400 \cdot 10^{-11}$ | CATECHOL omits two CH ₃ groups of O-, M- and P-XYCATECH |
| $\text{XYLOL} + \text{NO}_3 \longrightarrow 0.61 \cdot \text{PHENO}_2 + 0.39 \cdot \text{C}_6\text{H}_5\text{O} + 0.49 \cdot \text{HNO}_3$ | $3.200 \cdot 10^{-11}$ | XYLOLO ₂ and NXYLOLO ₂ approximated with PHENO ₂ ; XY1O with C ₆ H ₅ O (see MCM3.2 for details) |
| $\text{XYLOLO}_2 + \text{NO} \longrightarrow \text{HO}_2 + \text{NO}_2 + 0.17 \cdot \text{GLYOXAL} + 0.51 \cdot \text{CH}_3\text{COCHO}$ | $2.600 \cdot 10^{-12} \exp(365./T)$ | |
| $\text{XYLOLO}_2 + \text{HO}_2 \longrightarrow \text{XYLOLOOH}$ | $7.500 \cdot 10^{-13} \exp(700./T)$ | |
| $\text{XYLOLO}_2 + \text{CH}_3\text{O}_2 \longrightarrow 0.17 \cdot \text{GLYOXAL} + 0.51 \cdot \text{CH}_3\text{COCHO} + 2 \cdot \text{HO}_2 + \text{CH}_2\text{O}$ | $1.000 \cdot 10^{-12}$ | |
| $\text{XYLOLO}_2 + \text{CH}_3\text{CO}_3 \longrightarrow \text{HO}_2 + 0.17 \cdot \text{GLYOXAL} + 0.51 \cdot \text{CH}_3\text{COCHO} + \text{CH}_3\text{O}_2 + \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{XYLOLOOH} + \text{OH} \longrightarrow \text{XYLOLO}_2$ | $3.800 \cdot 10^{-12} \exp(200./T)$ | |
| $\text{XYLENO}_2 + \text{HO}_2 \longrightarrow \text{XYLENOOH}$ | $7.500 \cdot 10^{-13} \exp(700./T)$ | |
| $\text{XYLENO}_2 + \text{NO} \longrightarrow \text{HO}_2 + \text{NO}_2 + 0.34 \cdot \text{GLYOXAL} + 0.54 \cdot \text{CH}_3\text{COCHO} + 0.06 \cdot \text{BIGALD}_1 + 0.2 \cdot \text{BIGALD}_2 + 0.15 \cdot \text{BIGALD}_3 + 0.21 \cdot \text{BIGALD}_4$ | $2.600 \cdot 10^{-12} \exp(365./T)$ | |
| $\text{XYLENO}_2 + \text{CH}_3\text{O}_2 \longrightarrow 0.34 \cdot \text{GLYOXAL} + 0.54 \cdot \text{CH}_3\text{COCHO} + 2 \cdot \text{HO}_2 + \text{CH}_2\text{O} + 0.06 \cdot \text{BIGALD}_1 + 0.2 \cdot \text{BIGALD}_2 + 0.15 \cdot \text{BIGALD}_3 + 0.21 \cdot \text{BIGALD}_4$ | $1.000 \cdot 10^{-12}$ | |
| $\text{XYLENO}_2 + \text{CH}_3\text{CO}_3 \longrightarrow \text{HO}_2 + 0.34 \cdot \text{GLYOXAL} + 0.54 \cdot \text{CH}_3\text{COCHO} + 0.06 \cdot \text{BIGALD}_1 + 0.2 \cdot \text{BIGALD}_2 + 0.15 \cdot \text{BIGALD}_3 + 0.21 \cdot \text{BIGALD}_4 + \text{CH}_3\text{O}_2 + \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{XYLENOOH} + \text{OH} \longrightarrow \text{XYLENO}_2$ | $3.800 \cdot 10^{-12} \exp(200./T)$ | |

Table S15: C10/C15 oxidation reactions

| reaction | rate coefficient | reference |
|---|--------------------------------------|--|
| $\text{APIN} + \text{OH} \longrightarrow \text{TERPO}_2$ | $1.200 \cdot 10^{-11} \exp(440./T)$ | |
| $\text{BPIN} + \text{OH} \longrightarrow \text{TERPO}_2$ | $1.600 \cdot 10^{-11} \exp(470./T)$ | |
| $\text{LIMON} + \text{OH} \longrightarrow \text{TERPO}_2$ | $4.200 \cdot 10^{-11} \exp(400./T)$ | |
| $\text{MYRC} + \text{OH} \longrightarrow \text{TERPO}_2$ | $2.100 \cdot 10^{-10}$ | |
| $\text{BCARY} + \text{OH} \longrightarrow \text{TERPO}_2$ | $2.000 \cdot 10^{-10}$ | |
| $\text{APIN} + \text{O}_3 \longrightarrow 0.07 \cdot \text{ELVOC} + 0.39 \cdot \text{TERPROD}_1 + 0.27 \cdot \text{TERPROD}_2 + 0.63 \cdot \text{OH} + 0.57 \cdot \text{HO}_2 + 0.23 \cdot \text{CO} + 0.27 \cdot \text{CO}_2 + 0.52 \cdot \text{CH}_3\text{COCH}_3 + 0.34 \cdot \text{CH}_2\text{O} + 0.05 \cdot \text{HCOOH} + 0.05 \cdot \text{BIGALKANE} + 0.06 \cdot \text{CH}_3\text{CO}_3 + 0.06 \cdot \text{CH}_3\text{COCH}_2\text{O}_2$ | $6.300 \cdot 10^{-16} \exp(-580./T)$ | 7% ELVOC-yield according to Ehn et al. (2014) for endocyclic alkenes |
| $\text{BPIN} + \text{O}_3 \longrightarrow 0.43 \cdot \text{TERPROD}_1 + 0.3 \cdot \text{TERPROD}_2 + 0.63 \cdot \text{OH} + 0.57 \cdot \text{HO}_2 + 0.23 \cdot \text{CO} + 0.27 \cdot \text{CO}_2 + 0.52 \cdot \text{CH}_3\text{COCH}_3 + 0.34 \cdot \text{CH}_2\text{O} + 0.05 \cdot \text{HCOOH} + 0.05 \cdot \text{BIGALKANE} + 0.06 \cdot \text{CH}_3\text{CO}_3 + 0.06 \cdot \text{CH}_3\text{COCH}_2\text{O}_2$ | 1.700 $10^{-15} \exp(-1300./T)$ | . |

Table S15: C10/C15 oxidation reactions (... continued)

| reaction | rate coefficient | reference |
|--|--------------------------------------|---|
| $\text{LIMON} + \text{O}_3 \longrightarrow 0.07 \cdot \text{ELVOC} + 0.39 \cdot \text{TERPROD}_1 + 0.27 \cdot \text{TERPROD}_2 + 0.63 \cdot \text{OH} + 0.57 \cdot \text{HO}_2 + 0.23 \cdot \text{CO} + 0.27 \cdot \text{CO}_2 + 0.52 \cdot \text{CH}_3\text{COCH}_3 + 0.34 \cdot \text{CH}_2\text{O} + 0.05 \cdot \text{HCOOH} + 0.05 \cdot \text{BIGALKANE} + 0.06 \cdot \text{CH}_3\text{CO}_3 + 0.06 \cdot \text{CH}_3\text{COCH}_2\text{O}_2$ | $3.000 \cdot 10^{-15} \exp(-780./T)$ | 7% ELVOC-yield according to Ehn et al. (2014) for endocyclic alkenes |
| $\text{MYRC} + \text{O}_3 \longrightarrow 0.43 \cdot \text{TERPROD}_1 + 0.3 \cdot \text{TERPROD}_2 + 0.63 \cdot \text{OH} + 0.57 \cdot \text{HO}_2 + 0.23 \cdot \text{CO} + 0.27 \cdot \text{CO}_2 + 0.52 \cdot \text{CH}_3\text{COCH}_3 + 0.34 \cdot \text{CH}_2\text{O} + 0.05 \cdot \text{HCOOH} + 0.05 \cdot \text{BIGALKANE} + 0.06 \cdot \text{CH}_3\text{CO}_3 + 0.06 \cdot \text{CH}_3\text{COCH}_2\text{O}_2$ | $4.700 \cdot 10^{-16}$ | |
| $\text{BCARY} + \text{O}_3 \longrightarrow 0.645 \cdot \text{TERPROD}_1 + 0.45 \cdot \text{TERPROD}_2 + 0.63 \cdot \text{OH} + 0.57 \cdot \text{HO}_2 + 0.23 \cdot \text{CO} + 0.27 \cdot \text{CO}_2 + 0.52 \cdot \text{CH}_3\text{COCH}_3 + 0.34 \cdot \text{CH}_2\text{O} + 0.05 \cdot \text{HCOOH} + 0.05 \cdot \text{BIGALKANE} + 0.06 \cdot \text{CH}_3\text{CO}_3 + 0.06 \cdot \text{CH}_3\text{COCH}_2\text{O}_2$ | $1.200 \cdot 10^{-14}$ | |
| $\text{APIN} + \text{NO}_3 \longrightarrow \text{NTERPO}_2$ | $1.200 \cdot 10^{-12} \exp(490./T)$ | |
| $\text{BPIN} + \text{NO}_3 \longrightarrow \text{NTERPO}_2$ | $2.500 \cdot 10^{-12}$ | |
| $\text{LIMON} + \text{NO}_3 \longrightarrow \text{NTERPO}_2$ | $1.100 \cdot 10^{-11}$ | |
| $\text{MYRC} + \text{NO}_3 \longrightarrow \text{NTERPO}_2$ | $1.200 \cdot 10^{-11}$ | |
| $\text{BCARY} + \text{NO}_3 \longrightarrow \text{NTERPO}_2 + 0.5 \cdot \text{TERPROD}_1$ | $1.900 \cdot 10^{-11}$ | |
| $\text{TERPO}_2 + \text{NO} \longrightarrow 0.26 \cdot \text{TERPNO}_3 + 0.74 \cdot \text{NO}_2 + 0.36 \cdot \text{CH}_2\text{O} + 0.045 \cdot \text{CH}_3\text{COCH}_3 + 0.695 \cdot \text{TERPROD}_1 + 0.74 \cdot \text{HO}_2$ | $4.200 \cdot 10^{-12} \exp(180./T)$ | alkyl nitrate yield according to Rindelaub et al. (2015) for alpha-pinene |
| $\text{TERPO}_2 + \text{HO}_2 \longrightarrow \text{TERPOOH}$ | $7.500 \cdot 10^{-13} \exp(700./T)$ | |
| $\text{TERPO}_2 + \text{CH}_3\text{O}_2 \longrightarrow 1.15 \cdot \text{CH}_2\text{O} + 0.05 \cdot \text{CH}_3\text{COCH}_3 + 0.945 \cdot \text{TERPROD}_1 + \text{HO}_2 + 0.25 \cdot \text{CH}_3\text{OH}$ | $2.000 \cdot 10^{-12} \exp(500./T)$ | |
| $\text{TERPO}_2 + \text{CH}_3\text{CO}_3 \longrightarrow 0.4 \cdot \text{CH}_2\text{O} + 0.05 \cdot \text{CH}_3\text{COCH}_3 + 0.945 \cdot \text{TERPROD}_1 + \text{HO}_2 + \text{CH}_3\text{O}_2 + \text{CO}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{TERPOOH} + \text{OH} \longrightarrow \text{H}_2\text{O} + \text{TERPO}_2$ | $3.300 \cdot 10^{-11}$ | |
| $\text{TERPROD}_1 + \text{OH} \longrightarrow \text{TERP}_2\text{O}_2$ | $5.700 \cdot 10^{-11}$ | |
| $\text{TERPROD}_1 + \text{NO}_3 \longrightarrow 0.5 \cdot \text{TERP}_2\text{O}_2 + 0.5 \cdot \text{NTERPO}_2 + 0.5 \cdot \text{NO}_2$ | $1.000 \cdot 10^{-12}$ | |
| $\text{TERPNO}_3 + \text{OH} \longrightarrow \text{NO}_2 + \text{TERPROD}_1 + \text{H}_2\text{O}$ | $3.500 \cdot 10^{-12}$ | |
| $\text{TERP}_2\text{O}_2 + \text{NO} \longrightarrow 0.1 \cdot \text{TERPNO}_3 + 0.9 \cdot \text{NO}_2 + 0.34 \cdot \text{CH}_2\text{O} + 0.27 \cdot \text{CH}_3\text{COCH}_3 + 0.225 \cdot \text{CO} + 0.9 \cdot \text{CO}_2 + 0.9 \cdot \text{TERPROD}_2 + 0.9 \cdot \text{HO}_2 + 0.225 \cdot \text{GLYALD}$ | $4.200 \cdot 10^{-12} \exp(180./T)$ | |
| $\text{TERP}_2\text{O}_2 + \text{HO}_2 \longrightarrow \text{TERP}_2\text{OOH}$ | $7.500 \cdot 10^{-13} \exp(700./T)$ | |
| $\text{TERP}_2\text{O}_2 + \text{CH}_3\text{O}_2 \longrightarrow \text{TERPROD}_2 + 0.93 \cdot \text{CH}_2\text{O} + 0.25 \cdot \text{CH}_3\text{OH} + \text{HO}_2 + 0.5 \cdot \text{CO}_2 + 0.125 \cdot \text{CO} + 0.125 \cdot \text{GLYALD} + 0.15 \cdot \text{CH}_3\text{COCH}_3$ | $2.000 \cdot 10^{-12} \exp(500./T)$ | |
| $\text{TERP}_2\text{O}_2 + \text{CH}_3\text{CO}_3 \longrightarrow 0.34 \cdot \text{CH}_2\text{O} + 0.27 \cdot \text{CH}_3\text{COCH}_3 + 0.225 \cdot \text{CO} + 2 \cdot \text{CO}_2 + \text{TERPROD}_2 + \text{HO}_2 + 0.225 \cdot \text{GLYALD} + \text{CH}_3\text{O}_2$ | $1.000 \cdot 10^{-11}$ | |
| $\text{TERP}_2\text{OOH} + \text{OH} \longrightarrow \text{H}_2\text{O} + \text{TERP}_2\text{O}_2$ | $2.300 \cdot 10^{-11}$ | |

Table S15: C10/C15 oxidation reactions (... continued)

| reaction | rate coefficient | reference |
|--|-------------------------------------|---|
| TERPROD ₂ + OH \longrightarrow 0.15 · CH ₃ COCH ₂ O ₂ + 0.68 · CH ₂ O + 1.8 · CO ₂ + 0.5 · CH ₃ COCH ₃ + 0.65 · CH ₃ CO ₃ + 0.2 · HO ₂ + 0.7 · CO | $3.400 \cdot 10^{-11}$ | |
| NTERPO ₂ + NO \longrightarrow 0.26 · NTERPNO ₃ + 1.48 · NO ₂ + 0.74 · TERPROD ₁ | $4.200 \cdot 10^{-12} \exp(180./T)$ | alkyl nitrate yield according to Rindelaub et al. (2015) for alpha-pinene |
| NTERPO ₂ + HO ₂ \longrightarrow NTERPNO ₃ | $7.500 \cdot 10^{-13} \exp(700./T)$ | |
| NTERPO ₂ + CH ₃ O ₂ \longrightarrow 0.5 · NTERPNO ₃ + 0.75 · CH ₂ O + 0.25 · CH ₃ OH + 0.5 · HO ₂ + 0.5 · TERPROD ₁ + 0.5 · NO ₂ | $2.000 \cdot 10^{-12} \exp(500./T)$ | |
| NTERPO ₂ + CH ₃ CO ₃ \longrightarrow CH ₃ O ₂ + TERPROD ₁ + CO ₂ + NO ₂ | $1.000 \cdot 10^{-11}$ | |
| NTERPO ₂ + NO ₃ \longrightarrow 2 · NO ₂ + TERPROD ₁ | $2.400 \cdot 10^{-12}$ | |
| NTERPNO ₃ + OH \longrightarrow NO ₂ + TERPROD ₁ + H ₂ O | $3.500 \cdot 10^{-12}$ | |
| ELVOC + OH \longrightarrow HO ₂ + TERPROD ₁ | $1.000 \cdot 10^{-11}$ | a general rate coefficient for oxygenated VOC |

Table S16: Tropospheric halogen reactions

| reaction | rate coefficient | reference |
|---|--|--|
| CL + CH ₂ O \longrightarrow HCL + HO ₂ + CO | $8.100 \cdot 10^{-11} \exp(-30./T)$ | |
| CL + CH ₄ \longrightarrow CH ₃ O ₂ + HCL | $7.300 \cdot 10^{-12} \exp(-1280./T)$ | |
| CL + CH ₃ CN \longrightarrow CH ₂ O + HCL + CO + NO | $1.600 \cdot 10^{-11} \exp(-2140./T)$ | JPL (2011), products: Tyndall |
| CL + C ₂ H ₂ + M \longrightarrow 0.1 · CL + 0.1 · GLYOXAL + 0.9 · HCL + 0.9 · HO ₂ + 1.8 · CO + M | $ktroe(5.200 \cdot 10^{-30}, 2.4, 2.200 \cdot 10^{-10}, 0.7, 0.6)$ | |
| CL + C ₂ H ₄ + M \longrightarrow HO ₂ + 2 · CO + HCL + M | $ktroe(1.600e-29, 3.3, 3.100 \cdot 10^{-10}, 1., 0.6)$ | |
| CL + C ₂ H ₆ \longrightarrow C ₂ H ₅ O ₂ + HCL | $7.200 \cdot 10^{-11} \exp(-70./T)$ | |
| CL + CH ₃ O ₂ \longrightarrow 0.5 · CLO + 0.5 · CH ₂ O + 0.5 · HCL + 0.5 · OH + 0.5 · CO + HO ₂ | $1.600 \cdot 10^{-10}$ | 50% CH ₂ OO is produced and assumed to be produced mostly in dry environments and thus decomposing |
| CL + CH ₃ OH \longrightarrow CH ₂ O + HO ₂ + HCL | $7.100 \cdot 10^{-11} \exp(-75./T)$ | IUPAC (2008) |
| CL + CH ₃ OOH \longrightarrow CH ₂ O + OH + HCL | $5.900 \cdot 10^{-11}$ | IUPAC (2008) |
| CL + CH ₃ CHO \longrightarrow CH ₃ CO ₃ + HCL | $8.000 \cdot 10^{-11}$ | IUPAC (2008) |
| CL + GLYALD \longrightarrow 0.35 · GLYOXAL + 0.35 · HO ₂ + 0.65 · HOCH ₂ CO ₃ + HCL | $7.600 \cdot 10^{-11}$ | k by Bacher et al. (2001); products by Niki et al. (1987) |
| CL + GLYOXAL \longrightarrow HCL + HCOCO ₃ | $3.441 \cdot 10^{-11}$ | k(298K) from Niki et al. (1985); k(CH ₂ O + Cl) has been adjusted for the k(CH ₂ O + Cl) by JPL (2011) used here |
| CL + C ₃ H ₈ \longrightarrow C ₃ H ₇ O ₂ + HCL | $1.400 \cdot 10^{-10}$ | IUPAC (2008) |

Table S16: Tropospheric halogen reactions (... continued)

| reaction | rate coefficient | reference |
|--|--------------------------------------|--|
| $\text{CL} + \text{CH}_3\text{COCH}_3 \longrightarrow \text{CH}_3\text{COCH}_2\text{O}_2 + \text{HCL}$ | $1.500 \cdot 10^{-11} \exp(-590./T)$ | IUPAC (2008) |
| $\text{CL} + \text{HYAC} \longrightarrow \text{CH}_3\text{COCHO} + \text{HO}_2 + \text{HCL}$ | $5.400 \cdot 10^{-11}$ | Calvert et al. (2008a) |
| $\text{CL} + \text{BIGALKANE} \longrightarrow \text{ALKO}_2 + \text{HCL}$ | $1.935e - 10$ | from the reaction with OH, BIGALKANE seems to be methyl-butane whose k is an average of k given by Qian et al. (2002) and Anderson et al. (2007) |
| $\text{CL} + \text{MEK} \longrightarrow \text{HCL} + \text{MEKO}_2$ | $3.800 \cdot 10^{-11} \exp(16./T)$ | Calvert et al. (2008b) |
| $\text{CLO} + \text{CH}_3\text{O}_2 \longrightarrow \text{CL} + \text{HO}_2 + \text{CH}_2\text{O}$ | $3.300 \cdot 10^{-12} \exp(-115./T)$ | |
| $\text{BR} + \text{CH}_2\text{O} \longrightarrow \text{HBR} + \text{HO}_2 + \text{CO}$ | $1.700 \cdot 10^{-11} \exp(-800./T)$ | |
| $\text{BR} + \text{CH}_3\text{CHO} \longrightarrow \text{CH}_3\text{CO}_3 + \text{HBR}$ | $1.800 \cdot 10^{-11} \exp(-460./T)$ | IUPAC (2008) |
| $\text{BRO} + \text{CH}_3\text{O}_2 \longrightarrow \text{HOBR} + \text{OH} + \text{HO}_2 + \text{CO}$ | $2.420 \cdot 10^{-14} \exp(1617./T)$ | Shallcross et al. (2015); CH_2OO assumed to be produced mostly in dry environments and thus decomposing |

Table S17: Sulfur reactions

| reaction | rate coefficient | reference |
|--|--|------------|
| $\text{SO}_2 + \text{OH} + \text{M} \longrightarrow \text{H}_2\text{SO}_4 + \text{HO}_2$ | $ktroe(3.300 \cdot 10^{-31}, 4.3, 1.600 \cdot 10^{-12}, 0., 0.6)$ | JPL (2011) |
| $\text{DMS} + \text{OH} \longrightarrow \text{CH}_3\text{SO}_2 + \text{HCHO}$ | $1.130 \cdot 10^{-11} \exp(-253./T)$ | |
| $\text{DMS} + \text{OH} \longrightarrow \text{DMSO} + \text{HO}_2$ | $1. \cdot 10^{-9} \exp(5820./T) \cdot [\text{O}_2] / (1. \cdot 10^{30} + 5. \cdot \exp(6280./T) \cdot [\text{O}_2])$ | |
| $\text{DMS} + \text{NO}_3 \longrightarrow \text{CH}_3\text{SO}_2 + \text{HNO}_3 + \text{HCHO}$ | $1.900 \cdot 10^{-13} \exp(520./T)$ | |
| $\text{DMS} + \text{CL} \longrightarrow \text{CH}_3\text{SO}_2 + \text{HCL} + \text{HCHO}$ | $3.300 \cdot 10^{-10}$ | |
| $\text{DMS} + \text{BR} \longrightarrow \text{CH}_3\text{SO}_2 + \text{HBR} + \text{HCHO}$ | $9.000 \cdot 10^{-11} \exp(-2386./T)$ | |
| $\text{DMS} + \text{BRO} \longrightarrow \text{BR} + \text{DMSO}$ | $4.400 \cdot 10^{-13}$ | |
| $\text{DMSO} + \text{OH} \longrightarrow 0.6 \cdot \text{SO}_2 + \text{HCHO} + 0.6 \cdot \text{CH}_3\text{O}_2 + 0.4 \cdot \text{HO}_2 + 0.4 \cdot \text{CH}_3\text{SO}_3\text{H}$ | $1.000 \cdot 10^{-10}$ | |
| $\text{CH}_3\text{SO}_2 \longrightarrow \text{CH}_3\text{O}_2 + \text{SO}_2$ | $1.800 \cdot 10^{13} \exp(-8661./T)$ | |
| $\text{CH}_3\text{SO}_2 + \text{O}_3 \longrightarrow \text{CH}_3\text{SO}_3$ | $3.000 \cdot 10^{-13}$ | |
| $\text{CH}_3\text{SO}_3 + \text{HO}_2 \longrightarrow \text{CH}_3\text{SO}_3\text{H}$ | $5.000 \cdot 10^{-11}$ | |

Table S18: Stratospheric O(1D) reactions (... continued)

| reaction | rate coefficient | reference |
|---|-------------------------------------|---|
| Table S18: Stratospheric O(1D) reactions | | |
| reaction | rate coefficient | reference |
| $\text{O}_1\text{D} + \text{N}_2\text{O} \longrightarrow 2 \cdot \text{NO}$ | $7.250 \cdot 10^{-11} \exp(20./T)$ | |
| $\text{O}_1\text{D} + \text{N}_2\text{O} \longrightarrow \text{N}_2 + \text{O}_2$ | $4.630 \cdot 10^{-11} \exp(20./T)$ | |
| $\text{O}_1\text{D} + \text{O}_3 \longrightarrow 2 \cdot \text{O}_2$ | $1.200 \cdot 10^{-10}$ | |
| $\text{O}_1\text{D} + \text{CFC}_{11} \longrightarrow 2 \cdot \text{CL} + \text{COFCL}$ | $2.020 \cdot 10^{-10}$ | |
| $\text{O}_1\text{D} + \text{CFC}_{12} \longrightarrow 2 \cdot \text{CL} + \text{COF}_2$ | $1.204 \cdot 10^{-10}$ | |
| $\text{O}_1\text{D} + \text{CFC}_{113} \longrightarrow 2 \cdot \text{CL} + \text{COFCL} + \text{COF}_2$ | $1.500 \cdot 10^{-10}$ | |
| $\text{O}_1\text{D} + \text{CFC}_{114} \longrightarrow 2 \cdot \text{CL} + 2 \cdot \text{COF}_2$ | $9.750 \cdot 10^{-11}$ | |
| $\text{O}_1\text{D} + \text{CFC}_{115} \longrightarrow \text{CL} + \text{F} + 2 \cdot \text{COF}_2$ | $1.500 \cdot 10^{-11}$ | |
| $\text{O}_1\text{D} + \text{HCFC}_{22} \longrightarrow \text{CL} + \text{COF}_2$ | $7.200 \cdot 10^{-11}$ | |
| $\text{O}_1\text{D} + \text{HCFC}_{141}\text{B} \longrightarrow \text{CL} + \text{COFCL}$ | $1.794 \cdot 10^{-10}$ | |
| $\text{O}_1\text{D} + \text{HCFC}_{142}\text{B} \longrightarrow \text{CL} + \text{COF}_2$ | $1.628 \cdot 10^{-10}$ | |
| $\text{O}_1\text{D} + \text{CCL}_4 \longrightarrow 4 \cdot \text{CL}$ | $2.840 \cdot 10^{-10}$ | |
| $\text{O}_1\text{D} + \text{CH}_3\text{BR} \longrightarrow \text{BR}$ | $1.674 \cdot 10^{-10}$ | |
| $\text{O}_1\text{D} + \text{CF}_2\text{CLBR} \longrightarrow \text{BR} + \text{CL} + \text{COF}_2$ | $9.600 \cdot 10^{-11}$ | |
| $\text{O}_1\text{D} + \text{CF}_3\text{BR} \longrightarrow \text{BR} + \text{F} + \text{COF}_2$ | $4.100 \cdot 10^{-11}$ | |
| $\text{O}_1\text{D} + \text{H}_{1202} \longrightarrow 2 \cdot \text{BR} + \text{COF}_2$ | $1.012 \cdot 10^{-10}$ | |
| $\text{O}_1\text{D} + \text{H}_{2402} \longrightarrow 2 \cdot \text{BR} + 2 \cdot \text{COF}_2$ | $1.200 \cdot 10^{-10}$ | |
| $\text{O}_1\text{D} + \text{CHBR}_3 \longrightarrow 3 \cdot \text{BR}$ | $4.490 \cdot 10^{-10}$ | |
| $\text{O}_1\text{D} + \text{CH}_2\text{BR}_2 \longrightarrow 2 \cdot \text{BR}$ | $2.570 \cdot 10^{-10}$ | |
| $\text{O}_1\text{D} + \text{COF}_2 \longrightarrow 2 \cdot \text{F}$ | $2.140 \cdot 10^{-11}$ | |
| $\text{O}_1\text{D} + \text{COFCL} \longrightarrow \text{CL} + \text{F}$ | $1.900 \cdot 10^{-10}$ | |
| $\text{O}_1\text{D} + \text{CH}_4 \longrightarrow \text{CH}_3\text{O}_2 + \text{OH}$ | $1.310 \cdot 10^{-10}$ | |
| $\text{O}_1\text{D} + \text{CH}_4 \longrightarrow \text{CH}_2\text{O} + \text{H} + \text{HO}_2$ | $3.500 \cdot 10^{-11}$ | |
| $\text{O}_1\text{D} + \text{CH}_4 \longrightarrow \text{CH}_2\text{O} + \text{H}_2$ | $9.000 \cdot 10^{-12}$ | |
| $\text{O}_1\text{D} + \text{H}_2 \longrightarrow \text{H} + \text{OH}$ | $1.200 \cdot 10^{-10}$ | |
| $\text{O}_1\text{D} + \text{HCL} \longrightarrow \text{CL} + \text{OH}$ | $1.500 \cdot 10^{-10}$ | |
| $\text{O}_1\text{D} + \text{HBR} \longrightarrow \text{BR} + \text{OH}$ | $1.200 \cdot 10^{-10}$ | |
| $\text{O}_1\text{D} + \text{HCN} \longrightarrow \text{CO} + \text{OH} + \text{NO}$ | $7.700 \cdot 10^{-11} \exp(100./T)$ | Strekowski et al. (2001), products: Tyndall |

Table S19: Stratospheric inorganic halogen reactions

| reaction | rate coefficient | reference |
|--|--|---|
| $\text{CL} + \text{O}_3 \longrightarrow \text{CLO} + \text{O}_2$ | $2.300 \cdot 10^{-11} \exp(-200./T)$ | |
| $\text{CL} + \text{H}_2 \longrightarrow \text{H} + \text{HCL}$ | $3.050 \cdot 10^{-11} \exp(-2270./T)$ | |
| $\text{CL} + \text{H}_2\text{O}_2 \longrightarrow \text{HCL} + \text{HO}_2$ | $1.100 \cdot 10^{-11} \exp(-980./T)$ | |
| $\text{CL} + \text{HO}_2 \longrightarrow \text{HCL} + \text{O}_2$ | $1.400 \cdot 10^{-11} \exp(270./T)$ | |
| $\text{CL} + \text{HO}_2 \longrightarrow \text{CLO} + \text{OH}$ | $3.600 \cdot 10^{-11} \exp(-375./T)$ | |
| $\text{CL}_2\text{O}_2 + \text{M} \longrightarrow 2 \cdot \text{CLO} + \text{M}$ | k_{CL2O2} | |
| $\text{CLO} + \text{O} \longrightarrow \text{CL} + \text{O}_2$ | $2.800 \cdot 10^{-11} \exp(85./T)$ | |
| $\text{CLO} + \text{OH} \longrightarrow \text{CL} + \text{HO}_2$ | $7.400 \cdot 10^{-12} \exp(270./T)$ | |
| $\text{CLO} + \text{OH} \longrightarrow \text{HCL} + \text{O}_2$ | $6.000 \cdot 10^{-13} \exp(230./T)$ | |
| $\text{CLO} + \text{HO}_2 \longrightarrow \text{HOCL} + \text{O}_2$ | $2.600 \cdot 10^{-12} \exp(290./T)$ | |
| $\text{CLO} + \text{NO} \longrightarrow \text{CL} + \text{NO}_2$ | $6.400 \cdot 10^{-12} \exp(290./T)$ | |
| $\text{CLO} + \text{NO}_2 + \text{M} \longrightarrow \text{CLONO}_2 + \text{M}$ | $ktroe(1.800 \cdot 10^{-31}, 3.4, 1.500 \cdot 10^{-11}, 1.9, 0.6)$ | |
| $\text{CLO} + \text{CLO} \longrightarrow 2 \cdot \text{CL} + \text{O}_2$ | $3.000 \cdot 10^{-11} \exp(-2450./T)$ | |
| $\text{CLO} + \text{CLO} \longrightarrow \text{CL}_2 + \text{O}_2$ | $1.000 \cdot 10^{-12} \exp(-1590./T)$ | |
| $\text{CLO} + \text{CLO} \longrightarrow \text{CL} + \text{OCLO}$ | $3.500 \cdot 10^{-13} \exp(-1370./T)$ | |
| $\text{CLO} + \text{CLO} + \text{M} \longrightarrow \text{CL}_2\text{O}_2 + \text{M}$ | $ktroe(1.600 \cdot 10^{-32}, 4.5, 3.000 \cdot 10^{-12}, 2., 0.6)$ | |
| $\text{HCL} + \text{OH} \longrightarrow \text{CL} + \text{H}_2\text{O}$ | $1.800 \cdot 10^{-12} \exp(-250./T)$ | |
| $\text{HCL} + \text{O} \longrightarrow \text{CL} + \text{OH}$ | $1.000 \cdot 10^{-11} \exp(-3300./T)$ | |
| $\text{HOCL} + \text{O} \longrightarrow \text{CLO} + \text{OH}$ | $1.700 \cdot 10^{-13}$ | |
| $\text{HOCL} + \text{CL} \longrightarrow \text{CLO} + \text{HCL}$ | $3.400 \cdot 10^{-12} \exp(-130./T)$ | |
| $\text{HOCL} + \text{OH} \longrightarrow \text{CLO} + \text{H}_2\text{O}$ | $3.000 \cdot 10^{-12} \exp(-500./T)$ | |
| $\text{CLONO}_2 + \text{O} \longrightarrow \text{CLO} + \text{NO}_3$ | $3.600 \cdot 10^{-12} \exp(-840./T)$ | |
| $\text{CLONO}_2 + \text{OH} \longrightarrow \text{HOCL} + \text{NO}_3$ | $1.200 \cdot 10^{-12} \exp(-330./T)$ | |
| $\text{CLONO}_2 + \text{CL} \longrightarrow \text{CL}_2 + \text{NO}_3$ | $6.500 \cdot 10^{-12} \exp(135./T)$ | |
| $\text{BR} + \text{O}_3 \longrightarrow \text{BRO} + \text{O}_2$ | $1.600 \cdot 10^{-11} \exp(-780./T)$ | |
| $\text{BR} + \text{HO}_2 \longrightarrow \text{HBR} + \text{O}_2$ | $4.800 \cdot 10^{-12} \exp(-310./T)$ | |
| $\text{BR} + \text{NO}_2 + \text{M} \longrightarrow 0.85 \cdot \text{BRONO} + 0.15 \cdot \text{BRNO}_2 + \text{M}$ | $ktroe(4.200 \cdot 10^{-31}, 2.4, 2.700 \cdot 10^{-11}, 0., 0.6)$ | |
| $\text{BRONO} + \text{M} \longrightarrow \text{BR} + \text{NO}_2 + \text{M}$ | $1.648 \cdot 10^{11} \exp(-7399./T)$ | fit to upper limit by Wine et al. (1993) and scaled 1.5 factor to match data by Orlando and Burkholder (2000) |
| $\text{BRO} + \text{O} \longrightarrow \text{BR} + \text{O}_2$ | $1.900 \cdot 10^{-11} \exp(230./T)$ | |
| $\text{BRO} + \text{OH} \longrightarrow \text{BR} + \text{HO}_2$ | $1.700 \cdot 10^{-11} \exp(250./T)$ | |
| $\text{BRO} + \text{HO}_2 \longrightarrow \text{HOBR} + \text{O}_2$ | $4.500 \cdot 10^{-12} \exp(460./T)$ | |

Table S19: Stratospheric inorganic halogen reactions (... continued)

| reaction | rate coefficient | reference |
|---|--|--|
| $\text{BRO} + \text{NO} \longrightarrow \text{BR} + \text{NO}_2$ | $8.800 \cdot 10^{-12} \exp(260./T)$ | |
| $\text{BRO} + \text{NO}_2 + \text{M} \longrightarrow \text{BRONO}_2 + \text{M}$ | $ktroe(5.200 \cdot 10^{-31}, 3.2, 6.900 \cdot 10^{-12}, 2.9, 0.6)$ | |
| $\text{BRO} + \text{CLO} \longrightarrow \text{BR} + \text{OCLO}$ | $9.500 \cdot 10^{-13} \exp(550./T)$ | |
| $\text{BRO} + \text{CLO} \longrightarrow \text{BR} + \text{CL} + \text{O}_2$ | $2.300 \cdot 10^{-12} \exp(260./T)$ | |
| $\text{BRO} + \text{CLO} \longrightarrow \text{BRCL} + \text{O}_2$ | $4.100 \cdot 10^{-13} \exp(290./T)$ | |
| $\text{BRO} + \text{BRO} \longrightarrow 2 \cdot \text{BR} + \text{O}_2$ | $2.400 \cdot 10^{-12} \exp(40./T)$ | |
| $\text{BRO} + \text{BRO} \longrightarrow \text{BR}_2 + \text{O}_2$ | $2.800 \cdot 10^{-14} \exp(860./T)$ | |
| $\text{HBR} + \text{OH} \longrightarrow \text{BR} + \text{H}_2\text{O}$ | $5.500 \cdot 10^{-12} \exp(200./T)$ | |
| $\text{HBR} + \text{O} \longrightarrow \text{BR} + \text{OH}$ | $5.800 \cdot 10^{-12} \exp(-1500./T)$ | |
| $\text{HOBR} + \text{O} \longrightarrow \text{BRO} + \text{OH}$ | $1.200 \cdot 10^{-10} \exp(-430./T)$ | |
| $\text{BRONO}_2 + \text{O} \longrightarrow \text{BRO} + \text{NO}_3$ | $1.900 \cdot 10^{-11} \exp(215./T)$ | |
| $\text{BRONO}_2 + \text{BR} \longrightarrow \text{BR}_2 + \text{NO}_3$ | $5.805 \cdot 10^{-11}$ | Average of k at 298K by Orlando and Tyndall (1996) and Harwood et al. (1998) |
| $\text{BR}_2 + \text{OH} \longrightarrow \text{BR} + \text{HOBR}$ | $2.100 \cdot 10^{-11} \exp(240./T)$ | |
| $\text{F} + \text{H}_2\text{O} \longrightarrow \text{HF} + \text{OH}$ | $1.400 \cdot 10^{-11}$ | |
| $\text{F} + \text{H}_2 \longrightarrow \text{H} + \text{HF}$ | $1.400 \cdot 10^{-10} \exp(-500./T)$ | |
| $\text{F} + \text{CH}_4 \longrightarrow \text{CH}_3\text{O}_2 + \text{HF}$ | $1.600 \cdot 10^{-10} \exp(-260./T)$ | |
| $\text{F} + \text{HNO}_3 \longrightarrow \text{HF} + \text{NO}_3$ | $6.000 \cdot 10^{-12} \exp(400./T)$ | |

Table S20: Stratospheric organic halogen reactions

| reaction | rate coefficient | reference |
|--|---------------------------------------|--|
| $\text{CH}_3\text{BR} + \text{OH} \longrightarrow \text{BR} + \text{H}_2\text{O} + \text{HO}_2$ | $2.350 \cdot 10^{-12} \exp(-1300./T)$ | |
| $\text{CH}_3\text{BR} + \text{CL} \longrightarrow \text{HCL} + \text{HO}_2 + \text{BR}$ | $1.400 \cdot 10^{-11} \exp(-1030./T)$ | |
| $\text{CH}_2\text{BR}_2 + \text{OH} \longrightarrow 2 \cdot \text{BR} + \text{H}_2\text{O}$ | $2.000 \cdot 10^{-12} \exp(-840./T)$ | |
| $\text{CHBR}_3 + \text{OH} \longrightarrow 3 \cdot \text{BR}$ | $1.350 \cdot 10^{-12} \exp(-600./T)$ | |
| $\text{CH}_2\text{BR}_2 + \text{CL} \longrightarrow 2 \cdot \text{BR} + \text{HCL}$ | $6.300 \cdot 10^{-12} \exp(-800./T)$ | |
| $\text{CHBR}_3 + \text{CL} \longrightarrow 3 \cdot \text{BR} + \text{HCL}$ | $4.850 \cdot 10^{-12} \exp(-850./T)$ | |
| $\text{CH}_3\text{CL} + \text{CL} \longrightarrow \text{CO} + \text{HO}_2 + 2 \cdot \text{HCL}$ | $2.170 \cdot 10^{-11} \exp(-1130./T)$ | |
| $\text{CH}_3\text{CL} + \text{OH} \longrightarrow \text{CO} + \text{HO}_2 + \text{HCL} + \text{H}_2\text{O}$ | $2.400 \cdot 10^{-12} \exp(-1250./T)$ | products: Tyndall(p.c.), implicitly includes $\text{NO} \longrightarrow \text{NO}_2$ conversion with CH_2ClO_2 |
| $\text{CH}_3\text{CCL}_3 + \text{OH} \longrightarrow \text{H}_2\text{O} + 3 \cdot \text{CL}$ | $1.640 \cdot 10^{-12} \exp(-1520./T)$ | |
| $\text{HCFC}_{22} + \text{OH} \longrightarrow \text{CL} + \text{H}_2\text{O} + \text{COF}_2$ | $1.050 \cdot 10^{-12} \exp(-1600./T)$ | |
| $\text{HCFC}_{141}\text{B} + \text{OH} \longrightarrow \text{CL} + \text{COFCL}$ | $1.250 \cdot 10^{-12} \exp(-1600./T)$ | |

Table S20: Stratospheric organic halogen reactions (... continued)

| reaction | rate coefficient | reference |
|--|---------------------------------------|-----------|
| $\text{HCFC}_{142}\text{B} + \text{OH} \longrightarrow \text{CL} + \text{COF}_2$ | $1.300 \cdot 10^{-12} \exp(-1770./T)$ | |

Table S21: (Tropospheric) heterogeneous reactions

| reaction | reaction probability | reference |
|--|--|------------------------|
| $\text{O}_3 \longrightarrow \text{HO}_2$ | $\gamma = 10^{-6}$ | Stadtler et al. (2017) |
| $\text{HO}_2 \longrightarrow 0.5 \cdot \text{H}_2\text{O}_2$ | $\gamma = 0.2$ | Stadtler et al. (2017) |
| $\text{NO}_3 \longrightarrow$ | $\gamma = 0.001$ | Stadtler et al. (2017) |
| $\text{NO}_2 \longrightarrow 0.5 \cdot \text{HONO}$ | $\gamma = 10^{-4}$ | Stadtler et al. (2017) |
| $\text{HNO}_3 \longrightarrow$ | $\gamma_{SS} = 0.01$, $\gamma_{DU} = 0.1$ | Stadtler et al. (2017) |
| $\text{N}_2\text{O}_5 \longrightarrow$ | $\gamma = f(T, RH)$ | Stadtler et al. (2017) |

Table S22: Stratospheric sulfate aerosol reactions

| reaction | reaction probability | reference |
|---|--|------------------------|
| $\text{N}_2\text{O}_5 \longrightarrow 2 \cdot \text{HNO}_3$ | $\gamma = 0.04$ | Lamarque et al. (2012) |
| $\text{CLONO}_2 \longrightarrow \text{HOCL} + \text{HNO}_3$ | $f(\text{sulfuricacidwt}\%)$ | Lamarque et al. (2012) |
| $\text{BRONO}_2 \longrightarrow \text{HOBR} + \text{HNO}_3$ | $f(T, P, HCl, \text{H}_2\text{O}, r)$ | Lamarque et al. (2012) |
| $\text{CLONO}_2 + \text{HCL} \longrightarrow \text{CL}_2 + \text{HNO}_3$ | $f(T, P, \text{H}_2\text{O}, r)$ | Lamarque et al. (2012) |
| $\text{HOCL} + \text{HCL} \longrightarrow \text{CL}_2 + \text{H}_2\text{O}$ | $f(T, P, HCl, \text{H}_2\text{O}, r)$ | Lamarque et al. (2012) |
| $\text{HOBR} + \text{HCL} \longrightarrow \text{BRCL} + \text{H}_2\text{O}$ | $f(T, P, HCl, \text{HOBr}, \text{H}_2\text{O}, r)$ | Lamarque et al. (2012) |

Table S23: Stratospheric nitric acid dihydrate reactions

| reaction | reaction probability | reference |
|---|----------------------|------------------------|
| $\text{N}_2\text{O}_5 \longrightarrow 2 \cdot \text{HNO}_3$ | $\gamma = 0.0004$ | Lamarque et al. (2012) |
| $\text{CLONO}_2 \longrightarrow \text{HOCL} + \text{HNO}_3$ | $\gamma = 0.004$ | Lamarque et al. (2012) |
| $\text{CLONO}_2 + \text{HCL} \longrightarrow \text{CL}_2 + \text{HNO}_3$ | $\gamma = 0.2$ | Lamarque et al. (2012) |
| $\text{HOCL} + \text{HCL} \longrightarrow \text{CL}_2 + \text{H}_2\text{O}$ | $\gamma = 0.1$ | Lamarque et al. (2012) |
| $\text{BRONO}_2 \longrightarrow \text{HOBR} + \text{HNO}_3$ | $\gamma = 0.3$ | Lamarque et al. (2012) |

Table S24: Stratospheric ice aerosol reactions

| reaction | reaction probability | reference |
|---|----------------------|------------------------|
| $\text{N}_2\text{O}_5 \longrightarrow 2 \cdot \text{HNO}_3$ | $\gamma = 0.02$ | Lamarque et al. (2012) |
| $\text{CLONO}_2 \longrightarrow \text{HOCL} + \text{HNO}_3$ | $\gamma = 0.3$ | Lamarque et al. (2012) |
| $\text{BRONO}_2 \longrightarrow \text{HOBR} + \text{HNO}_3$ | $\gamma = 0.3$ | Lamarque et al. (2012) |
| $\text{CLONO}_2 + \text{HCL} \longrightarrow \text{CL}_2 + \text{HNO}_3$ | $\gamma = 0.3$ | Lamarque et al. (2012) |
| $\text{HOCL} + \text{HCL} \longrightarrow \text{CL}_2 + \text{H}_2\text{O}$ | $\gamma = 0.2$ | Lamarque et al. (2012) |
| $\text{HOBR} + \text{HCL} \longrightarrow \text{BRCL} + \text{H}_2\text{O}$ | $\gamma = 0.3$ | Lamarque et al. (2012) |

Table S25: Henry coefficients (H_0 and temperature factor) and dry deposition reactivities for gas-phase species in ECHAM-HAMMOZ.

| Species | H | Reactivity coefficient | Source |
|--|------------------------------|------------------------|---|
| CO | $9.81 \cdot 10^{-4}$, 1720. | 0. | JPL (2011) |
| H ₂ | $7.8 \cdot 10^{-4}$, 500. | 0. | Sander (1999) |
| H ₂ O ₂ | $8.44 \cdot 10^4$, 7600. | 1 | JPL (2011) |
| HCN | 12., 5000. | 1 | Sander (1999) |
| HNO ₃ | $3.2 \cdot 10^{11}$, 8700. | 1 | $H = 2.1 \cdot 10^5$ M/atm and $Ka = 15.4$ M (Schwartz and White, 1981). At an average cloud droplet pH=5 $\Rightarrow H^* = H \cdot (1 + Ka/[H+]) = 2.1 \cdot 10^5 \cdot (1 + 15.4/1 \cdot 10^{-5}) = 3.2 \cdot 10^{11}$ |
| HO ₂ | 690., 0. | 1 | JPL (2006) |
| HO ₂ NO ₂ | $1.2 \cdot 10^4$, 6900. | 1 | Sander (1999) |
| HONO | $5.05 \cdot 10^3$, 4800. | 0.1 | Sander (1999): $H=49$ and $Ka = 5.1 \cdot 10^{-4} \Rightarrow H^* = 49(1 + 5.1 \cdot 10^{-4} * 1 \cdot 10^7) = 5.05 \cdot 10^3$ |
| N ₂ O ₅ | 2.1, 3400. | 1 | Sander (1999): 2 ref. give it as infinite and one as 2.1 M/atm |
| N ₂ O ₅ | $3.2 \cdot 10^{11}$, 8700. | 1 | HNO ₃ as proxy as done in GEOS-Chem |
| NH ₃ | $1.02 \cdot 10^4$, 4200. | 0. | Sander (2015): $H(298) = 58$ M/atm and $H^*(298)$ calculated at pH=7 with $Kb = 1.75 \cdot 10^5$ M |
| NO | $1.92 \cdot 10^{-3}$, 1790. | 1 | JPL (2006) |
| NO ₂ | $12. \cdot 10^{-2}$, 2360. | 1 | JPL (2011) |
| NO ₃ | $3.8 \cdot 10^{-2}$, 0. | 1 | JPL (2006) |
| O ₃ | $1.03 \cdot 10^{-2}$, 2830. | 1 | JPL (2006) |
| SO ₂ | $2.45 \cdot 10^5$, 3100. | 0. | Sander (2015): $H(298) = 1.2$ M/atm and $H^*(298)$ calculated at pH=7 with $K1a = 1.23 \cdot 10^{-2}$ M and $K2a = 6.61 \cdot 10^{-8}$ M |
| H ₂ SO ₄ | $1.3 \cdot 10^{15}$, 20000. | 0. | Sander (2015) |
| CH ₄ | $1.41 \cdot 10^{-3}$, 2040. | 0. | JPL (2011) |
| CH ₂ O | $3.23 \cdot 10^3$, 7100. | 0.1 | JPL (2011) |
| CH ₃ OH | 203., 5640. | 0. | JPL (2011) |
| CH ₃ OOH | 300., 5280. | 0.1 | JPL (2011) |
| HCOOH | $8.9 \cdot 10^3$, 6100. | 0. | Sander (1999): measured value by Johnson et al. (1996)??, surface reactivity as in Nguyen et al. (2015) |
| CH ₃ O ₂ NO ₂ | 2.0, 4700. | 0.1 | Sander (1999): methyl nitrate as proxy species |
| C ₂ H ₅ OH | 190., 6660. | 0. | JPL (2011) |
| C ₂ H ₅ OOH | 336., 5995. | 1 | JPL (2011) |
| CH ₃ CHO | 12.9, 5890. | 0.1 | JPL (2011) |
| CH ₃ CN | 52.8, 3970. | 0. | JPL (2011) |
| PAN | 2.8, 5730. | 0.1 | JPL (2011) |
| CH ₃ COOH | $4.1 \cdot 10^3$, 6200. | 0. | JPL (2011) |

Table S25: Henry coefficients (... continued)

| Species | H | Reactivity coefficient | Source |
|--|---------------------------|------------------------|--|
| CH ₃ COOOH | 837., 5310. | 0.1 | JPL (2011) |
| GLYALD | $4.1 \cdot 10^4$, 4600. | 0.1 | Sander (1999) |
| GLYOXAL | $4.19 \cdot 10^5$, 7480. | 0.1 | JPL (2011) |
| HCOCO ₂ H | $1.1 \cdot 10^4$, 4800. | 0.1 | Sander (2015) |
| HCOCO ₃ H | $1.1 \cdot 10^4$, 4800. | 0.1 | Sander (2015) |
| HOCH ₂ CO ₂ H | $2.8 \cdot 10^4$, 4000. | 0.1 | Sander (2015) |
| HOCH ₂ CO ₃ H | $2.8 \cdot 10^4$, 4000. | 0.1 | Sander (2015) |
| C ₃ H ₇ OOH | 336., 5995. | 1 | |
| CH ₃ COCH ₃ | 27.8, 5530. | 0. | JPL (2011) |
| CH ₃ COCHO | $3.7 \cdot 10^3$, 7500. | 0.1 | Sander (1999) |
| HYAC | $7.7 \cdot 10^3$, 0. | 0.1 | Sander (2015) |
| NOA | $1. \cdot 10^3$, 0. | 0.1 | Sander (2015) |
| MACR | 6.5, 5300. | 0.1 | Sander (1999) |
| MACROOH | $2.1 \cdot 10^5$, 9900. | 1 | Sander (2015) |
| MACROH | $2.1 \cdot 10^5$, 9900. | 0. | Sander (2015) |
| MEK | 37., 8200. | 0.1 | Sander (2015) |
| MEKOOH | $1.1 \cdot 10^3$, 7300. | 1 | Sander (2015) |
| MPAN | 1.7, 0. | 0.1 | Sander (1999) |
| MVK | $4.1 \cdot 10^1$, 7800. | 0.1 | Sander (1999) |
| MACO ₂ H | $1.5 \cdot 10^3$, 6800. | 0.1 | Sander (2015) |
| MACO ₃ H | $1.5 \cdot 10^3$, 6800. | 1 | Sander (2015) |
| BIGALD ₁ | $2.5 \cdot 10^5$, 0. | 0.1 | Sander (2015) |
| CO ₂ H ₃ CHO | $4.1 \cdot 10^4$, 4600. | 0.1 | Sander (1999) |
| CO ₂ H ₃ CO ₃ H | $4.1 \cdot 10^4$, 4600. | 1 | Sander (1999) |
| BIACETOH | $4.1 \cdot 10^4$, 4600. | 0.1 | Sander (1999) |
| IBUTALOH | $4.1 \cdot 10^4$, 4600. | 0.1 | Sander (1999) |
| IBUTALOHOOH | $4.1 \cdot 10^4$, 4600. | 1 | Sander (1999) |
| LHMKABOOH | $2.1 \cdot 10^5$, 9900. | 1 | Sander (2015) |
| MVKN | $5. \cdot 10^3$, 0. | 0.1 | like for ISOPNO3 from Nguyen et al. (2015) |
| MACRN | $5. \cdot 10^3$, 0. | 0.1 | like for ISOPNO3 from Nguyen et al. (2015) |
| MBO | 64., 0. | 0.1 | Sander (2015) |
| HPALD | $4.1 \cdot 10^4$, 4600. | 1 | Sander (1999) |
| PACALD | $4.1 \cdot 10^4$, 4600. | 1 | Sander (1999) |
| ISOPBNO ₃ | $5. \cdot 10^3$, 0. | 0.1 | Nguyen et al. (2015) |

Table S25: Henry coefficients (... continued)

| Species | H | Reactivity coefficient | Source |
|--------------------------------------|--------------------------|------------------------|----------------------|
| ISOPDNO ₃ | $5. \cdot 10^3$, 0. | 0.1 | Nguyen et al. (2015) |
| LISOPACNO ₃ | $5. \cdot 10^3$, 0. | 0.1 | Nguyen et al. (2015) |
| LC ₅ PAN ₁₇₁₉ | $5. \cdot 10^3$, 0. | 0.1 | Nguyen et al. (2015) |
| LHC ₄ ACCO ₂ H | $2.1 \cdot 10^5$, 9900. | 0.1 | Sander (2015) |
| LHC ₄ ACCO ₃ H | $2.1 \cdot 10^5$, 9900. | 1 | Sander (2015) |
| LIECHO | $4.1 \cdot 10^4$, 4600. | 0.1 | Sander (1999) |
| LIECO ₃ H | $2.1 \cdot 10^5$, 9900. | 1 | Sander (2015) |
| LIEPOX | $2.1 \cdot 10^5$, 9900. | 1 | Sander (2015) |
| LNISOOH | $2.1 \cdot 10^5$, 9900. | 1 | Sander (2015) |
| MBOOOH | $3. \cdot 10^{11}$, 0. | 1 | Sander (2015) |
| NC ₄ CHO | 2.3, 0. | 0.1 | Sander (2015) |
| NISOPOOH | $3.6 \cdot 10^4$, 0. | 1 | Sander (2015) |
| LHC ₄ ACCHO | $4.1 \cdot 10^4$, 4600. | 0.1 | Sander (1999) |
| HCOC ₅ | $4.1 \cdot 10^4$, 4600. | 1 | Sander (1999) |
| ISOPBOOH | $2.1 \cdot 10^5$, 9900. | 1 | Sander (2015) |
| ISOPDOOH | $2.1 \cdot 10^5$, 9900. | 1 | Sander (2015) |
| LISOPACOOH | $2.1 \cdot 10^5$, 9900. | 1 | Sander (2015) |
| ISOPBOH | $2.1 \cdot 10^5$, 9900. | 1 | Sander (2015) |
| ISOPDOH | $2.1 \cdot 10^5$, 9900. | 1 | Sander (2015) |
| ISOPAOH | $2.1 \cdot 10^5$, 9900. | 1 | Sander (2015) |
| LISOPOOHOOH | $2. \cdot 10^{16}$, 0. | 1 | Sander (2015) |
| LISOPNO ₃ OOH | $3. \cdot 10^{11}$, 0. | 1 | Sander (2015) |
| LISOPNO ₃ NO ₃ | $2.1 \cdot 10^5$, 9900. | 1 | Sander (2015) |
| LC ₅₇₈ OOH | $3. \cdot 10^{11}$, 0. | 1 | Sander (2015) |
| C ₅₉ OOH | $3. \cdot 10^{11}$, 0. | 1 | Sander (2015) |
| BIGALD ₂ | $2.5 \cdot 10^5$, 0. | 0.1 | Sander (2015) |
| BIGALD ₃ | $2.5 \cdot 10^5$, 0. | 0.1 | Sander (2015) |
| BZALD | 38., 5500. | 0.1 | Sander (2015) |
| BZOOH | $2.9 \cdot 10^3$, 0. | 1 | Sander (2015) |
| PHENOOH | $3. \cdot 10^{11}$, 0. | 1 | Sander (2015) |
| C ₆ H ₅ OOH | $3. \cdot 10^3$, 5900. | 1 | Sander (2015) |
| PHENOL | $3. \cdot 10^3$, 5900. | 0.1 | Sander (2015) |
| CATECHOL | $4.6 \cdot 10^3$, 0. | 0.1 | Sander (2015) |
| CATEC ₁ OOH | $4.6 \cdot 10^3$, 0. | 1 | Sander (2015) |

Table S25: Henry coefficients (... continued)

| Species | H | Reactivity coefficient | Source |
|-----------------------|-----------------------------|------------------------|---|
| BEPOMUC | $2.5 \cdot 10^5$, 0. | 0.1 | Sander (2015) |
| BENZOOH | $2.1 \cdot 10^5$, 9900. | 1 | Sander (2015) |
| PBZNIT | 2.8, 5730. | 0.1 | PAN as proxy MCM3.2 name PBZN |
| BIGALD ₄ | $2.5 \cdot 10^5$, 0. | 0.1 | Sander (2015) |
| TOLOOH | $2.1 \cdot 10^5$, 9900. | 1 | Sander (2015) |
| CRESOL | $1.1 \cdot 10^3$, 6700. | 0.1 | Sander (2015) |
| TEPOMUC | $2.5 \cdot 10^5$, 0. | 0.1 | Sander (2015) |
| TERPROD ₂ | $2.5 \cdot 10^5$, 0. | 0.1 | Sander (2015) |
| XYLENOOH | $2.1 \cdot 10^5$, 9900. | 1 | Sander (2015) |
| XYLOL | $1.1 \cdot 10^3$, 6700. | 0.1 | cresol as proxy MCM3.2 name OXYLOL |
| XYLOLOOH | $2.1 \cdot 10^5$, 9900. | 1 | Sander (2015) |
| TERPOOH | $2.1 \cdot 10^5$, 9900. | 1 | Sander (2015) |
| TERP ₂ OOH | $2.1 \cdot 10^5$, 9900. | 1 | Sander (2015) |
| TERPNO ₃ | $5. \cdot 10^3$, 0. | 0.1 | like an isoprene nitrate MCM3.2 name APINANO ₃ |
| TERPROD ₁ | $2.5 \cdot 10^5$, 0. | 1 | Sander (2015) |
| NTERPNO ₃ | $5. \cdot 10^3$, 0. | 0.1 | like an isoprene nitrate MCM3.2 name NAPINAOOH |
| MTHOM | $2. \cdot 10^9$, 0. | 1 | Sander (2015) |
| ALKOH | 336., 5995. | 0.1 | |
| ALKOOH | 336., 5995. | 1 | |
| POOH | 300., 5280. | 0.1 | |
| ROOH | 300., 5280. | 1 | |
| EOOH | 300., 5280. | 0.1 | |
| ALKNO ₃ | $5. \cdot 10^3$, 0. | 0.1 | like for ISOPNO ₃ from Nguyen et al. (2015) MCM3.2 |
| HF | $1.3 \cdot 10^4$, 0. | 0. | Sander (2015) |
| HCL | $2.0 \cdot 10^{11}$, 600. | 0. | Fernandez et al. (2014) |
| HOCL | $6.6 \cdot 10^2$, 5900. | 0. | Sander (1999) |
| CL ₂ | $9.2 \cdot 10^{-2}$, 2000. | 0. | Sander (1999) |
| CLONO ₂ | $2. \cdot 10^{13}$, 0. | 1 | Sander (1999) : listed as "infinite" |
| BR ₂ | $7.2 \cdot 10^{-1}$, 0. | 0. | Sander (2015) |
| HBR | $7.2 \cdot 10^{13}$, 6100. | 0. | Fernandez et al. (2014) |
| HOBR | $6.1 \cdot 10^3$, 0. | 0. | Sander (1999) |
| BRONO ₂ | $2. \cdot 10^{13}$, 0. | 1 | Sander (1999) : listed as "infinite" |
| BRONO | $3. \cdot 10^{-1}$, 0. | 0. | Sander (2015) |
| BRNO ₂ | $3. \cdot 10^{-1}$, 0. | 0. | Sander (2015) |

Table S25: Henry coefficients (... continued)

| Species | H | Reactivity coefficient | Source |
|-----------------------------------|-----------------------------|---------------------------|----------------------|
| BRCL | $9.7 \cdot 10^{-1}$, 5600. | 0. | Sander (2015) |
| DMS | 0.54, 3460. | 0. | JPL (2011) |
| DMSO | $9.8 \cdot 10^4$, 0. | 0.1 | JPL (2011) |
| CH ₃ SO ₃ H | $1. \cdot 10^{30}$, 0. | 1 | Jöckel et al. (2006) |

References

- Anderson, J. L., Dixon, J. K., and Brennecke, J. F.: Solubility of CO₂, CH₄, C₂H₆, C₂H₄, O₂, and N₂ in 1-Hexyl-3-methylpyridinium Bis(trifluoromethylsulfonyl)imide: Comparison to Other Ionic Liquids, *Accounts of Chemical Research*, 40, 1208–1216, doi:10.1021/ar7001649, PMID: 17970599, 2007.
- Bacher, C., Tyndall, G. S., and Orlando, J. J.: The Atmospheric Chemistry of Glycolaldehyde, *Journal of Atmospheric Chemistry*, 39, 171–189, doi:10.1023/A:1010689706869, 2001.
- Burkholder, J. and Orlando, J.: UV absorption cross-sections of cis-BrONO, *CHEMICAL PHYSICS LETTERS*, 317, 603–608, doi:10.1016/S0009-2614(99)01412-8, 2000.
- Calvert, J. G., Derwent, R. G., Orlando, J. J., Tyndall, G. S., and Wallington, T. J.: *Mechanisms of Atmospheric Oxidation of the Alkanes*, Oxford University Press, Oxford, UK, 2008a.
- Calvert, J. G., Mellouki, A., Orlando, J. J., Pilling, M., and Wallington, T. J.: *The Mechanisms of Atmospheric Oxidation of the Oxygenates*, Oxford University Press, Oxford, UK, 2008b.
- Crounse, J. D., Paulot, F., Kjaergaard, H. G., and Wennberg, P. O.: Peroxy radical isomerization in the oxidation of isoprene, *PHYSICAL CHEMISTRY CHEMICAL PHYSICS*, 13, 13 607–13 613, doi:10.1039/c1cp21330j, 2011.
- Crounse, J. D., Knap, H. C., Ornsø, K. B., Jorgensen, S., Paulot, F., Kjaergaard, H. G., and Wennberg, P. O.: Atmospheric Fate of Methacrolein. 1. Peroxy Radical Isomerization Following Addition of OH and O₂, *JOURNAL OF PHYSICAL CHEMISTRY A*, 116, 5756–5762, doi:10.1021/jp211560u, 2012.
- Ehn, M., Thornton, J. A., Kleist, E., Sipila, M., Junninen, H., Pullinen, I., Springer, M., Rubach, F., Tillmann, R., Lee, B., Lopez-Hilfiker, F., Andres, S., Acir, I.-H., Rissanen, M., Jokinen, T., Schobesberger, S., Kangasluoma, J., Kontkanen, J., Nieminen, T., Kurten, T., Nielsen, L. B., Jorgensen, S., Kjaergaard, H. G., Canagaratna, M., Dal Maso, M., Berndt, T., Petaja, T., Wahner, A., Kerminen, V.-M., Kulmala, M., Worsnop, D. R., Wildt, J., and Mentel, T. F.: A large source of low-volatility secondary organic aerosol, *NATURE*, 506, 476+, doi:10.1038/nature13032, 2014.
- Fernandez, R. P., Salawitch, R. J., Kinnison, D. E., Lamarque, J.-F., and Saiz-Lopez, A.: Bromine partitioning in the tropical tropopause layer: implications for stratospheric injection, *Atmospheric Chemistry and Physics*, 14, 13 391–13 410, doi:10.5194/acp-14-13391-2014, 2014.
- Harwood, M., Burkholder, J., and Ravishankara, A.: Photodissociation of BrONO₂ and N₂O₅: Quantum yields for NO₃ production at 248, 308, and 352.5 nm, *JOURNAL OF PHYSICAL CHEMISTRY A*, 102, 1309–1317, doi:10.1021/jp9729829, 1998.
- IUPAC, Atkinson, R. and Baulch, D. L. and Cox, R. A. and Crowley, J. N. and Hampson, R. F. and Hynes, R. G. and Jenkin, M. E. and Rossi, M. J. and Troe, J.: Evaluated kinetic and photochemical data for atmospheric chemistry: Volume I - gas phase reactions of O_x, HO_x, NO_x and SO_x species, *Atmospheric Chemistry and Physics*, 4, 1461–1738, doi:10.5194/acp-4-1461-2004, 2004.
- IUPAC, Atkinson, R. and Baulch, D. L. and Cox, R. A. and Crowley, J. N. and Hampson, R. F. and Hynes, R. G. and Jenkin, M. E. and Rossi, M. J. and Troe, J. and IUPAC Subcommittee.: Evaluated kinetic and photochemical data for atmospheric chemistry: Volume II - gas phase reactions of organic species, *Atmospheric Chemistry and Physics*, 6, 3625–4055, doi:10.5194/acp-6-3625-2006, 2006.

- IUPAC, Atkinson, R. and Baulch, D. L. and Cox, R. A. and Crowley, J. N. and Hampson, R. F. and Hynes, R. G. and Jenkin, M. E. and Rossi, M. J. and Troe, J. and Wallington, T. J.: Evaluated kinetic and photochemical data for atmospheric chemistry: Volume IV gas phase reactions of organic halogen species, *Atmospheric Chemistry and Physics*, 8, 4141–4496, doi:10.5194/acp-8-4141-2008, 2008.
- Jöckel, P., Tost, H., Pozzer, A., Brühl, C., Buchholz, J., Ganzeveld, L., Hoor, P., Kerkweg, A., Lawrence, M., Sander, R., et al.: The atmospheric chemistry general circulation model ECHAM5/MESSy1: consistent simulation of ozone from the surface to the mesosphere, *Atmospheric Chemistry and Physics Discussions*, 6, 6957–7050, doi:10.5194/acp-6-5067-2006, 2006.
- Johnson, M., Kuwata, K., Wong, C., and Okumura, M.: Vibrational spectrum of I-(H₂O), *CHEMICAL PHYSICS LETTERS*, 260, 551–557, doi:10.1016/0009-2614(96)00911-6, 1996.
- JPL, S. P. Sander and J. Abbatt and J. R. Barker and J. B. Burkholder and R. R. Friedl and D. M. Golden and R. E. Huie and C. E. Kolb and M. J. Kurylo and G. K. Moortgat and V. L. Orkin and P. H. Wine.: Chemical Kinetics and Photochemical Data for Use in Atmospheric Studies, Evaluation No. 17, JPL Publication 10-6, Jet Propulsion Laboratory, Pasadena, URL <http://jpldataeval.jpl.nasa.gov>, 2011.
- JPL, S. P. Sander and R. R. Friedl and D. M. Golden and M. J. Kurylo and G. K. Moortgat and H. Keller-Rudek and P. H. Wine and A. R. Ravishankara and C. E. Kolb and M. J. Molina and B. J. Finlayson-Pitts and R. E. Huie and V. L. Orkin.: Chemical Kinetics and Photochemical Data for Use in Atmospheric Studies, Evaluation Number 15, JPL Publication 06-2, Jet Propulsion Laboratory, Pasadena, CA, URL <http://jpldataeval.jpl.nasa.gov>, 2006.
- Lamarque, J. F., Emmons, L. K., Hess, P. G., Kinnison, D. E., Tilmes, S., Vitt, F., Heald, C. L., Holland, E. A., Lauritzen, P. H., Neu, J., Orlando, J. J., Rasch, P. J., and Tyndall, G. K.: CAM-chem: description and evaluation of interactive atmospheric chemistry in the Community Earth System Model, *Geoscientific Model Development*, 5, 369–411, doi:10.5194/gmd-5-369-2012, 2012.
- Müller, J.-F., Peeters, J., and Stavrakou, T.: Fast photolysis of carbonyl nitrates from isoprene, *Atmospheric Chemistry and Physics*, 14, 2497–2508, doi:10.5194/acp-14-2497-2014, 2014.
- Nguyen, T. L., Lee, H., Matthews, D. A., McCarthy, M. C., and Stanton, J. F.: Stabilization of the Simplest Criegee Intermediate from the Reaction between Ozone and Ethylene: A High-Level Quantum Chemical and Kinetic Analysis of Ozonolysis, *The Journal of Physical Chemistry A*, 119, 5524–5533, doi:10.1021/acs.jpca.5b02088, PMID: 25945650, 2015.
- Niki, H., Maker, P., Savage, C., and Breitenbach, L.: AN FTIR STUDY OF THE CL-ATOM-INITIATED REACTION OF GLYOXAL, *INTERNATIONAL JOURNAL OF CHEMICAL KINETICS*, 17, 547–558, doi:10.1002/kin.550170507, 1985.
- Niki, H., Maker, P., Savage, C., and Hurley, M.: FOURIER-TRANSFORM INFRARED STUDY OF THE KINETICS AND MECHANISMS FOR THE CL-ATOM-INITIATED AND HO-RADICAL-INITIATED OXIDATION OF GLYCOLALDEHYDE, *JOURNAL OF PHYSICAL CHEMISTRY*, 91, 2174–2178, doi:10.1021/j100292a038, 1987.
- Orlando, J. and Burkholder, J.: Identification of BrONO as the major product in the gas-phase reaction of Br with NO₂, *JOURNAL OF PHYSICAL CHEMISTRY A*, 104, 2048–2053, doi:10.1021/jp993713g, 2000.
- Orlando, J., Tyndall, G., Bertman, S., Chen, W., and Burkholder, J.: Rate coefficient for the reaction of OH with CH₂=C(CH₃)C(O)OONO₂ (MPAN), *ATMOSPHERIC ENVIRONMENT*, 36, 1895–1900, doi:10.1016/S1352-2310(02)00090-0, 2002.

- Orlando, J. J. and Tyndall, G. S.: Rate coefficients for the thermal decomposition of BrONO_2 and the heat of formation of BrONO_2 , *J. Phys. Chem.*, 100, 19 398–19 405, 1996.
- Orlando, J. J. and Tyndall, G. S.: Laboratory studies of organic peroxy radical chemistry: an overview with emphasis on recent issues of atmospheric significance, *CHEMICAL SOCIETY REVIEWS*, 41, 6294–6317, doi:10.1039/c2cs35166h, 2012.
- Paulot, F., Crounse, J. D., Kjaergaard, H. G., Kurten, A., St Clair, J. M., Seinfeld, J. H., and Wennberg, P. O.: Unexpected Epoxide Formation in the Gas-Phase Photooxidation of Isoprene, *Science*, 325, 730–733, doi:10.1126/science.1172910, 2009.
- Peeters, J., Boullart, W., Pultau, V., Vandenberg, S., and Vereecken, L.: Structure-activity relationship for the addition of OH to (poly)alkenes: Site-specific and total rate constants, *JOURNAL OF PHYSICAL CHEMISTRY A*, 111, 1618–1631, doi:10.1021/jp066973o, 2007.
- Praske, E., Crounse, J. D., Bates, K. H., Kurten, T., Kjaergaard, H. G., and Wennberg, P. O.: Atmospheric Fate of Methyl Vinyl Ketone: Peroxy Radical Reactions with NO and HO₂, *JOURNAL OF PHYSICAL CHEMISTRY A*, 119, 4562–4572, doi:10.1021/jp5107058, 2015.
- Qian, H.-B., Turton, D., Seakins, P. W., and Pilling, M. J.: A laser flash photolysis/IR diode laser absorption study of the reaction of chlorine atoms with selected alkanes, *International Journal of Chemical Kinetics*, 34, 86–94, doi:10.1002/kin.10025, 2002.
- Rindelaub, J. D., McAvey, K. M., and Shepson, P. B.: The photochemical production of organic nitrates from alpha-pinene and loss via acid-dependent particle phase hydrolysis, *ATMOSPHERIC ENVIRONMENT*, 100, 193–201, doi:10.1016/j.atmosenv.2014.10.010, 2015.
- Sander, R.: Modeling atmospheric chemistry: Interactions between gas-phase species and liquid cloud/aerosol particles, *SURVEYS IN GEOPHYSICS*, 20, 1–31, doi:10.1023/A:1006501706704, 1999.
- Sander, R.: Compilation of Henry’s law constants (version 4.0) for water as solvent, *ATMOSPHERIC CHEMISTRY AND PHYSICS*, 15, 4399–4981, doi:10.5194/acp-15-4399-2015, 2015.
- Schwartz, S. E. and White, W. H.: Solubility equilibria of the nitrogen oxides and oxyacids in dilute aqueous solution, in: *Advances in Environmental Science and Engineering*, edited by Pfafflin, J. R. and Ziegler, E. N., vol. 4, pp. 1–45, Gordon and Breach Science Publishers, NY, 1981.
- Shallcross, D. E., Leather, K. E., Bacak, A., Xiao, P., Lee, E. P. F., Ng, M., Mok, D. K. W., Dyke, J. M., Hossaini, R., Chipperfield, M. P., Khan, M. A. H., and Percival, C. J.: Reaction between CH_3O_2 and BrO radicals: a new source of upper troposphere lower stratosphere hydroxyl radicals, *J. Phys. Chem. A*, 119, 4618–4632, 2015.
- St. Clair, J., Moon, S., Holbrook, W. S., Perron, J. T., Riebe, C. S., Martel, S. J., Carr, B., Harman, C., Singha, K., and Richter, D. d.: Geophysical imaging reveals topographic stress control of bedrock weathering, *Science*, 350, 534–538, doi:10.1126/science.aab2210, 2015.
- Stadtler, S., Simpson, D., Schröder, S., Taraborrelli, D., Bott, A., and Schultz, M.: Ozone Impacts of Gas-Aerosol Uptake in Global Chemistry Transport Models, *Atmos. Chem. Phys. Discuss.*, 2017, 1–35, doi:10.5194/acp-2017-566, URL <https://www.atmos-chem-phys-discuss.net/acp-2017-566/>, 2017.
- Strekowski, R., Nicovich, I., and Wine, P.: Kinetic and mechanistic study of the reaction of O(D-1) with CF_2HBr , *INTERNATIONAL JOURNAL OF CHEMICAL KINETICS*, 33, 262–270, doi:10.1002/kin.1019, 2001.

- Taraborrelli, D., Lawrence, M. G., Butler, T. M., Sander, R., and Lelieveld, J.: Mainz Isoprene Mechanism 2 (MIM2): an isoprene oxidation mechanism for regional and global atmospheric modelling, *Atmospheric Chemistry and Physics*, 9, 2751–2777, 2009.
- Wine, P. H., Nicovich, J. M., Stickel, R. E., Zhao, Z., Shackelford, C. J., Kreutter, K. D., Daykin, E. P., and Wang, S.: Halogen and Sulphur Reactions Relevant to Polar Chemistry, *The Tropospheric Chemistry of Ozone in the Polar Regions*, 1993.