

Corrections to “Exploring DMS oxidation and implications for global aerosol radiative forcing”

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Table 3. The three DMS oxidation reactions in the standard CAM6-chem.

Gas-phase Reactions	Rate Constants* ($\text{cm}^3 \text{ molecule}^{-1} \text{ s}^{-1}$)	References
DMS + OH \rightarrow SO ₂ (H-abstraction)	$9.60 \times 10^{-12} \exp(-234/T)$	(Emmons et al., 2010)
DMS + OH \rightarrow 0.5SO ₂ + 0.5HO ₂ (OH-addition)	$1.7 \times 10^{-42} \exp(7810/T) \times 0.21[M] / \{1 +$ $5.5 \times 10^{-31} \exp(7460/T) \times 0.21[M]\}$	(Emmons et al., 2010)
DMS + NO ₃ \rightarrow SO ₂ + HNO ₃	$1.90 \times 10^{-13} \exp(520/T)$	(Emmons et al., 2010)

* T is air temperature in K; $[M]$ is air density in molecules cm^{-3} .

Table 4. Summary of the MSA-producing branch of the H-abstraction pathway in the DMS chemistry implemented into CAM6-chem.

Gas-phase Reactions	Rate Constants* ($\text{cm}^3 \text{ molecule}^{-1} \text{ s}^{-1}$)	References
DMS + OH \rightarrow MSP ($\text{CH}_3\text{SCH}_2\text{OO}$)	$1.12 \times 10^{-11} \exp(-250/T)$	(Saunders et al., 2003)
DMS + Cl \rightarrow 0.45MSP + 0.55(CH_3) ₂ S(Cl) + 0.45HCl	3.40×10^{-10}	IUPAC
(CH_3) ₂ S(Cl) \rightarrow DMS + Cl	9.00×10^1	(Enami et al., 2004)
MSP + NO \rightarrow $\text{CH}_3\text{SCH}_2(\text{O})$ + NO ₂	$4.90 \times 10^{-12} \exp(260/T)$	(Saunders et al., 2003)
MSP + RO ₂ \rightarrow $\text{CH}_3\text{SCH}_2(\text{O})$ + O ₂	3.74×10^{-12}	(Saunders et al., 2003)
$\text{CH}_3\text{SCH}_2(\text{O}) \rightarrow \text{CH}_3\text{S} + \text{CH}_2\text{O}$	1.00×10^6	(Saunders et al., 2003)
$\text{CH}_3\text{S} + \text{O}_3 \rightarrow \text{CH}_3\text{S}(\text{O}) + \text{O}_2$	$1.15 \times 10^{-12} \exp(430/T)$	(Saunders et al., 2003)
$\text{CH}_3\text{S} + \text{O}_2 \rightarrow \text{CH}_3\text{S}(\text{OO})$	$1.20 \times 10^{-16} \exp(1580/T)$	(Saunders et al., 2003)
$\text{CH}_3\text{S}(\text{O}) + \text{O}_3 \rightarrow \text{CH}_3(\text{O}_2) + \text{SO}_2$	4.00×10^{-13}	(Saunders et al., 2003)
$\text{CH}_3\text{S}(\text{OO}) \rightarrow \text{CH}_3(\text{O}_2) + \text{SO}_2$	$5.60 \times 10^{16} \exp(-10870/T)$	(Saunders et al., 2003)
$\text{CH}_3\text{S}(\text{OO}) \rightarrow \text{CH}_3\text{SO}_2$	1.00	(Saunders et al., 2003)
$\text{CH}_3\text{SO}_2 + \text{O}_3 \rightarrow \text{CH}_3\text{SO}_3 + \text{O}_2$	3.00×10^{-13}	(Saunders et al., 2003)
$\text{CH}_3\text{SO}_2 \rightarrow \text{CH}_3(\text{O}_2) + \text{SO}_2$	$5.00 \times 10^{13} \exp(-9673/T)$	(Saunders et al., 2003)
$\text{CH}_3\text{SO}_3 + \text{HO}_2 \rightarrow \text{MSA} + \text{O}_2$	5.00×10^{-11}	(Saunders et al., 2003)
$\text{CH}_3\text{SO}_3 \rightarrow \text{CH}_3(\text{O}_2) + \text{H}_2\text{SO}_4$	$5.00 \times 10^{13} \exp(-9946/T)$	(Saunders et al., 2003)

* T is air temperature in K.

Table 5. Summary of the isomerization branch of the H-abstraction pathway in the DMS chemistry implemented into CAM6-chem.

Gas-phase Reactions	Rate Constants* ($\text{cm}^3 \text{ molecule}^{-1} \text{ s}^{-1}$)	References
MSP \rightarrow O ₂ CH ₂ SCH ₂ OOH	$2.24 \times 10^{11} \exp(-9.8 \times 10^3/T) \exp(1.03 \times 10^8/T^3)$	(Veres et al., 2020)
OOCH ₂ SCH ₂ OOH \rightarrow HPMTF (HOOCH ₂ SCHO) + OH	$6.09 \times 10^{11} \exp(-9.5 \times 10^3/T) \exp(1.1 \times 10^8/T^3)$	(Veres et al., 2020)
OOCH ₂ SCH ₂ OOH + NO \rightarrow HOOCH ₂ SCH ₂ O + NO ₂	$4.90 \times 10^{-12} \exp(260/T)$	(Saunders et al., 2003)
HOOCH ₂ SCH ₂ O \rightarrow HOOCH ₂ S + CH ₂ O	1.00×10^6	(Saunders et al., 2003)
OOCH ₂ SCH ₂ OOH + HO ₂ \rightarrow HOOCH ₂ SCH ₂ OOH + O ₂	$1.13 \times 10^{-13} \exp(1300/T)$	(Saunders et al., 2003)
HPMTF + OH \rightarrow HOOCH ₂ SCO + H ₂ O	1.11×10^{-11}	(Patroescu et al., 1996; Vermeuel et al., 2020)
HOOCH ₂ SCO \rightarrow HOOCH ₂ S + CO	$9.20 \times 10^9 \exp(-505.4/T)$	(Wu et al., 2015)
HOOCH ₂ SCO \rightarrow OH + CH ₂ O + OCS	$1.60 \times 10^7 \exp(-1468.6/T)$	(Wu et al., 2015)
HOOCH ₂ S + O ₃ \rightarrow HOOCH ₂ SO + O ₂	$1.15 \times 10^{-12} \exp(430/T)$	(Saunders et al., 2003)
HOOCH ₂ S + NO ₂ \rightarrow HOOCH ₂ SO + NO	$6.00 \times 10^{-11} \exp(240/T)$	(Saunders et al., 2003)
HOOCH ₂ SO + O ₃ \rightarrow SO ₂ + CH ₂ O + OH + O ₂	4.00×10^{-13}	(Saunders et al., 2003)
HOOCH ₂ SO + NO ₂ \rightarrow SO ₂ + CH ₂ O + OH + NO	1.20×10^{-11}	(Saunders et al., 2003)

* T is air temperature in K.

Table 6. Gas-phase DMS oxidation (OH-addition pathway) implemented into CAM6-chem in this study.

Gas-phase Reactions	Rate Constants* ($\text{cm}^3 \text{ molecule}^{-1} \text{ s}^{-1}$)	References
DMS + OH \rightarrow 0.6SO ₂ + 0.4DMSO + CH ₃ O ₂	$8.20 \times 10^{-39} [\text{O}_2] \exp(5376/T) / \{1 + 1.05 \times$ $10^{-5}([\text{O}_2]/[\text{M}]) \exp(3644/T)\}$	(Burkholder et al., 2015; Pham et al., 1995)
DMS + NO ₃ \rightarrow SO ₂ + HNO ₃ + CH ₃ O ₂ + CH ₂ O	$1.13 \times 10^{-12} \exp(530/T)$	(Burkholder et al., 2015)
DMS + BrO \rightarrow DMSO + Br	$3.39 \times 10^{-13} \exp(950/T)$	(Burkholder et al., 2015)
DMS + O ₃ \rightarrow SO ₂	1.00×10^{-19}	(Burkholder et al., 2015)
DMS + Cl \rightarrow 0.5SO ₂ + 0.5DMSO + 0.5HCl + 0.5ClO	3.40×10^{-10}	(Burkholder et al., 2015; Chen et al., 2018)
DMSO + OH \rightarrow 0.95MSIA + 0.05SO ₂	$8.94 \times 10^{-11} \exp(800/T)$	(Burkholder et al., 2015)
MSIA + OH \rightarrow 0.9SO ₂ + 0.1MSA	9.00×10^{-11}	(Burkholder et al., 2015)
MSIA + O ₃ \rightarrow MSA	2.00×10^{-18}	(Lucas, 2002)

* T is air temperature in K; $[\text{M}]$ and $[\text{O}_2]$ are density of air and O₂ in molecules cm^{-3} , respectively.

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