The Chemical Mechanism of MAFOR v2.1

KPP version: `2.2.3_rs3`

MECCA version: `4.0`

Date: November 14, 2022

Batch file: `mafor.bat`

Integrator: `rosenbrock_posdef`

Gas equation file: `gas.eqn`

Replacement file: `maforchem`

Selected reactions:
“`Tr && (G || Aa) && !Br && !Hg`”

Number of aerosol phases: 1

Number of species in selected mechanism:

Gas phase:	781
Aqueous phase:	153
All species:	934

Number of reactions in selected mechanism:

Gas phase (Gnnn):	1869
Aqueous phase (Annn):	237
Henry (Hnnn):	122
Photolysis (Jnnn):	351
Aqueous phase photolysis (PHnnn):	13
Heterogeneous (HETnnn):	0
Equilibria (EQnn):	100
Isotope exchange (IEXnnn):	0
Tagging equations (TAGnnn):	0
Dummy (Dnn):	1
All equations:	2693

Table 1: Gas phase reactions

#	labels	reaction	rate coefficient	reference
G1000	UpStTrG	$O_2 + O(^1D) \rightarrow O(^3P) + O_2$	3.3E-11*EXP(55./temp)	Burkholder et al. (2015)
G1001	UpStTrG	$O_2 + O(^3P) \rightarrow O_3$	6.0E-34*((temp/300.)**(-2.4)) *cair	Burkholder et al. (2015)
G2100	UpStTrG	$H + O_2 \rightarrow HO_2$	k_3rd(temp,cair,4.4E-32,1.3, 7.5E-11,-0.2,0.6)	Burkholder et al. (2015)
G2104	UpStTrG	$OH + O_3 \rightarrow HO_2 + O_2$	1.7E-12*EXP(-940./temp)	Burkholder et al. (2015)
G2105	UpStTrG	$OH + H_2 \rightarrow H_2O + H$	2.8E-12*EXP(-1800./temp)	Burkholder et al. (2015)
G2107	UpStTrG	$HO_2 + O_3 \rightarrow OH + 2\ O_2$	1.E-14*EXP(-490./temp)	Burkholder et al. (2015)
G2109	UpStTrG	$HO_2 + OH \rightarrow H_2O + O_2$	4.8E-11*EXP(250./temp)	Burkholder et al. (2015)
G2110	UpStTrG	$HO_2 + HO_2 \rightarrow H_2O_2 + O_2$	k_HO2_HO2	Burkholder et al. (2015)*
G2111	UpStTrG	$H_2O + O(^1D) \rightarrow 2\ OH$	1.63E-10*EXP(60./temp)	Burkholder et al. (2015)
G2112	UpStTrG	$H_2O_2 + OH \rightarrow H_2O + HO_2$	1.8E-12	Burkholder et al. (2015)
G2117	UpStTrG	$H_2O + H_2O \rightarrow (H_2O)_2$	6.521E-26*temp*EXP(1851.09/temp) *EXP(-5.10485E-3*temp)	Scribano et al. (2006)*
G2118	UpStTrG	$(H_2O)_2 \rightarrow H_2O + H_2O$	1.E0	see note*
G3101	UpStTrGN	$N_2 + O(^1D) \rightarrow O(^3P) + N_2$	2.15E-11*EXP(110./temp)	Burkholder et al. (2015)
G3103	UpStTrGN	$NO + O_3 \rightarrow NO_2 + O_2$	3.0E-12*EXP(-1500./temp)	Burkholder et al. (2015)
G3106	StTrGN	$NO_2 + O_3 \rightarrow NO_3 + O_2$	1.2E-13*EXP(-2450./temp)	Burkholder et al. (2015)
G3108	StTrGN	$NO_3 + NO \rightarrow 2\ NO_2$	1.5E-11*EXP(170./temp)	Burkholder et al. (2015)
G3109	UpStTrGN	$NO_3 + NO_2 \rightarrow N_2O_5$	k_NO3_NO2	Burkholder et al. (2015)*
G3110	StTrGN	$N_2O_5 \rightarrow NO_2 + NO_3$	k_NO3_NO2/(5.8E-27*EXP(10840./ temp))	Burkholder et al. (2015)*
G3200	TrGN	$NO + OH \rightarrow HONO$	k_3rd(temp,cair,7.0E-31,2.6, 3.6E-11,0.1,0.6)	Burkholder et al. (2015)
G3201	UpStTrGN	$NO + HO_2 \rightarrow NO_2 + OH$	3.3E-12*EXP(270./temp)	Burkholder et al. (2015)
G3202	UpStTrGN	$NO_2 + OH \rightarrow HNO_3$	k_3rd(temp,cair,1.8E-30,3.0, 2.8E-11,0.,0.6)	Burkholder et al. (2015)
G3203	StTrGN	$NO_2 + HO_2 \rightarrow HNO_4$	k_NO2_HO2	Burkholder et al. (2015)*
G3204	TrGN	$NO_3 + HO_2 \rightarrow NO_2 + OH + O_2$	3.5E-12	Burkholder et al. (2015)
G3205	TrGN	$HONO + OH \rightarrow NO_2 + H_2O$	1.8E-11*EXP(-390./temp)	Burkholder et al. (2015)
G3206	StTrGN	$HNO_3 + OH \rightarrow H_2O + NO_3$	k_HNO3_OH	Dulitz et al. (2018)*
G3207	StTrGN	$HNO_4 \rightarrow NO_2 + HO_2$	k_NO2_HO2/(2.1E-27*EXP(10900./ temp))	Burkholder et al. (2015)*
G3208	StTrGN	$HNO_4 + OH \rightarrow NO_2 + H_2O$	1.3E-12*EXP(380./temp)	Burkholder et al. (2015)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G3209	TrGN	NH_3 + OH → NH_2 + H_2O	1.7E-12*EXP(-710./temp)	Kohlmann and Poppe (1999)
G3210	TrGN	NH_2 + O_3 → NH_2O + O_2	4.3E-12*EXP(-930./temp)	Kohlmann and Poppe (1999)
G3211	TrGN	NH_2 + HO_2 → NH_2O + OH	4.8E-07*EXP(-628./temp) *temp**(-1.32)	Kohlmann and Poppe (1999)
G3212	TrGN	NH_2 + HO_2 → HNO + H_2O	9.4E-09*EXP(-356./temp) *temp**(-1.12)	Kohlmann and Poppe (1999)
G3213	TrGN	NH_2 + NO → HO_2 + OH + N_2	1.92E-12*((temp/298.)**(-1.5))	Kohlmann and Poppe (1999)
G3214	TrGN	NH_2 + NO → N_2 + H_2O	1.41E-11*((temp/298.)**(-1.5))	Kohlmann and Poppe (1999)
G3215	TrGN	NH_2 + NO_2 → N_2O + H_2O	1.2E-11*((temp/298.)**(-2.0))	Kohlmann and Poppe (1999)
G3216	TrGN	NH_2 + NO_2 → NH_2O + NO	0.8E-11*((temp/298.)**(-2.0))	Kohlmann and Poppe (1999)
G3217	TrGN	NH_2O + O_3 → NH_2 + O_2	1.2E-14	Kohlmann and Poppe (1999)
G3218	TrGN	NH_2O → NHOH	1.3E3	Kohlmann and Poppe (1999)
G3219	TrGN	HNO + OH → NO + H_2O	8.0E-11*EXP(-500./temp)	Kohlmann and Poppe (1999)
G3220	TrGN	HNO + NHOH → NH_2OH + NO	1.66E-12*EXP(-1500./temp)	Kohlmann and Poppe (1999)
G3221	TrGN	HNO + NO_2 → HONO + NO	1.0E-12*EXP(-1000./temp)	Kohlmann and Poppe (1999)
G3222	TrGN	NHOH + OH → HNO + H_2O	1.66E-12	Kohlmann and Poppe (1999)
G3223	TrGN	NH_2OH + OH → NHOH + H_2O	4.13E-11*EXP(-2138./temp)	Kohlmann and Poppe (1999)
G3224	TrGN	HNO + O_2 → HO_2 + NO	3.65E-14*EXP(-4600./temp)	Kohlmann and Poppe (1999)
G4101	StTrG	CH_4 + OH → CH_3 + H_2O	1.85E-20*EXP(2.82*LOG(temp) -987./temp)	Atkinson (2003)
G4102	TrG	CH_3OH + OH → .85 HCHO + .85 HO_2 + .15 CH_3O + H_2O	6.38E-18*(temp**2)*EXP(144./temp)	Atkinson et al. (2006)
G4103a	StTrG	CH_3O_2 + HO_2 → CH_3OOH + O_2	3.8E-13*EXP(780./temp)/(1.+1./ 498.*EXP(1160./temp))	Atkinson et al. (2006)
G4103b	StTrG	CH_3O_2 + HO_2 → HCHO + H_2O + O_2	3.8E-13*EXP(780./temp)/(1.+ 498.*EXP(-1160./temp))	Atkinson et al. (2006)
G4104a	StTrGN	CH_3O_2 + NO → CH_3O + NO_2	2.3E-12*EXP(360./temp)*(1.-beta_ CH3NO3)	Atkinson et al. (2006), Butkovskaya et al. (2012), Flocke et al. (1998)
G4104b	StTrGN	CH_3O_2 + NO → CH_3ONO_2	2.3E-12*EXP(360./temp)*beta_ CH3NO3	Atkinson et al. (2006), Butkovskaya et al. (2012), Flocke et al. (1998)*
G4105	TrGN	CH_3O_2 + NO_3 → CH_3O + NO_2 + O_2	1.2E-12	Atkinson et al. (2006)
G4106a	StTrG	CH_3O_2 → CH_3O + .5 O_2	7.4E-13*EXP(-520./temp)*RO2*2.	Atkinson et al. (2006)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G4106b	StTrG	$CH_3O_2 \rightarrow .5\ HCHO + .5\ CH_3OH + .5\ O_2$	`(k_CH3O2-7.4E-13*EXP(-520./temp))*RO2*2.`	Atkinson et al. (2006)
G4107	StTrG	$CH_3OOH + OH \rightarrow .6\ CH_3O_2 + .4\ HCHO + .4\ OH + H_2O$	`k_CH3OOH_OH`	Wallington et al. (2018)
G4108	StTrG	$HCHO + OH \rightarrow CO + H_2O + HO_2$	`9.52E-18*EXP(2.03*LOG(temp)+636./temp)`	Sivakumaran et al. (2003)
G4109	TrGN	$HCHO + NO_3 \rightarrow HNO_3 + CO + HO_2$	`3.4E-13*EXP(-1900./temp)`	Burkholder et al. (2015)*
G4110	UpStTrG	$CO + OH \rightarrow H + CO_2$	`(1.57E-13+cair*3.54E-33)`	McCabe et al. (2001)
G4111	TrG	$HCOOH + OH \rightarrow CO_2 + HO_2 + H_2O$	`2.94E-14*exp(786./temp)+9.85E-13*EXP(-1036./temp)`	Paulot et al. (2011)
G4114	StTrGN	$CH_3O_2 + NO_2 \rightarrow CH_3O_2NO_2$	`k_NO2_CH3O2`	Burkholder et al. (2015)
G4115	StTrGN	$CH_3O_2NO_2 \rightarrow CH_3O_2 + NO_2$	`k_NO2_CH3O2/(9.5E-29*EXP(11234./temp))`	Burkholder et al. (2015)*
G4116	StTrGN	$CH_3O_2NO_2 + OH \rightarrow HCHO + NO_3 + H_2O$	`3.00E-14`	see note*
G4117	StTrGN	$CH_3ONO_2 + OH \rightarrow H_2O + HCHO + NO_2$	`4.0E-13*EXP(-845./temp)`	Atkinson et al. (2006)
G4118	StTrG	$CH_3O \rightarrow HO_2 + HCHO$	`1.3E-14*exp(-663./temp)*c(ind_O2)`	Chai et al. (2014)
G4119a	StTrGN	$CH_3O + NO_2 \rightarrow CH_3ONO_2$	`k_3rd_iupac(temp,cair,8.1E-29,4.5,2.1E-11,0.,0.44)`	Atkinson et al. (2006)
G4119b	StTrGN	$CH_3O + NO_2 \rightarrow HCHO + HONO$	`9.6E-12*EXP(-1150./temp)`	Atkinson et al. (2006)
G4120a	StTrGN	$CH_3O + NO \rightarrow CH_3ONO$	`k_3rd_iupac(temp,cair,2.6E-29,2.8,3.3E-11,0.6,REAL(EXP(-temp/900.),SP))`	Atkinson et al. (2006)
G4120b	StTrGN	$CH_3O + NO \rightarrow HCHO + HNO$	`2.3E-12*(temp/300.)**0.7`	Atkinson et al. (2006)
G4121	StTrG	$CH_3O_2 + O_3 \rightarrow CH_3O + 2\ O_2$	`2.9E-16*exp(-1000./temp)`	Burkholder et al. (2015)
G4122	StTrGN	$CH_3ONO + OH \rightarrow H_2O + HCHO + NO$	`1.E-10*exp(-1764./temp)`	Nielsen et al. (1991)
G4123	StTrG	$HCHO + HO_2 \rightarrow HOCH_2O_2$	`9.7E-15*EXP(625./temp)`	Atkinson et al. (2006)
G4124	StTrG	$HOCH_2O_2 \rightarrow HCHO + HO_2$	`2.4E12*EXP(-7000./temp)`	Atkinson et al. (2006)
G4125	StTrG	$HOCH_2O_2 + HO_2 \rightarrow .5\ HOCH_2OOH + .5\ HCOOH + .2\ OH + .2\ HO_2 + .3\ H_2O + .8\ O_2$	`5.6E-15*EXP(2300./temp)`	Atkinson et al. (2006)
G4126	StTrGN	$HOCH_2O_2 + NO \rightarrow NO_2 + HO_2 + HCOOH$	`0.7275*2.3E-12*EXP(360./temp)`	Atkinson et al. (2006)*
G4127	StTrGN	$HOCH_2O_2 + NO_3 \rightarrow NO_2 + HO_2 + HCOOH$	`1.2E-12`	see note*
G4129a	StTrG	$HOCH_2O_2 \rightarrow HCOOH + HO_2$	`(k_CH3O2*5.5E-12)**0.5*RO2*2.`	Atkinson et al. (2006)
G4129b	StTrG	$HOCH_2O_2 \rightarrow .5\ HCOOH + .5\ HOCH_2OH + .5\ O_2$	`(k_CH3O2*5.7E-14*EXP(750./temp))**0.5*RO2*2.`	Atkinson et al. (2006)
G4130a	StTrG	$HOCH_2OOH + OH \rightarrow HOCH_2O_2 + H_2O$	`k_roohro`	Taraborrelli (2010)*
G4130b	StTrG	$HOCH_2OOH + OH \rightarrow HCOOH + H_2O + OH$	`k_rohro + k_s*f_sooh*f_soh`	Taraborrelli (2010)*

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G4132	StTrG	$HOCH_2OH + OH \rightarrow HO_2 + HCOOH + H_2O$	`2.*k_rohro + k_s*f_soh*f_soh`	Taraborrelli (2010)*
G4133	StTrG	$CH_3O_2 + OH \rightarrow CH_3O + HO_2$	`1.4E-10`	Bossolasco et al. (2014)*
G4134	StTrG	$CH_2OO \rightarrow CO + HO_2 + OH$	`1.124E+14*EXP(-10000/temp)`	see note*
G4135	StTrG	$CH_2OO + H_2O \rightarrow HOCH_2OOH$	`k_CH2OO_NO2*3.6E-6`	Ouyang et al. (2013)*
G4136	StTrG	$CH_2OO + (H_2O)_2 \rightarrow HOCH_2OOH + H_2O$	`5.2E-12`	Chao et al. (2015), Lewis et al. (2015)*
G4137	StTrGN	$CH_2OO + NO \rightarrow HCHO + NO_2$	`6.E-14`	Welz et al. (2012)*
G4138	StTrGN	$CH_2OO + NO_2 \rightarrow HCHO + NO_3$	`k_CH2OO_NO2`	Welz et al. (2012), Stone et al. (2014)*
G4140	StTrG	$CH_2OO + CO \rightarrow HCHO + CO_2$	`3.6E-14`	Vereecken et al. (2012)
G4141	StTrG	$CH_2OO + HCOOH \rightarrow 2\ HCOOH$	`1.E-10`	Welz et al. (2014)*
G4142	StTrG	$CH_2OO + HCHO \rightarrow 2$ LCARBON	`1.7E-12`	Stone et al. (2014)*
G4143	StTrG	$CH_2OO + CH_3OH \rightarrow 2$ LCARBON	`5.E-12`	Vereecken et al. (2012)*
G4144	StTrG	$CH_2OO + CH_3O_2 \rightarrow 2$ LCARBON	`5.E-12`	Vereecken et al. (2012)*
G4145	StTrG	$CH_2OO + HO_2 \rightarrow$ LCARBON	`5.E-12`	Vereecken et al. (2012)
G4146	StTrG	$CH_2OO + O_3 \rightarrow HCHO + 2\ O_2$	`1.E-12`	Vereecken et al. (2014)
G4147	StTrG	$CH_2OO + CH_2OO \rightarrow 2\ HCHO + O_2$	`6.E-11`	Buras et al. (2014)
G4148	StTrGN	$HOCH_2O_2 + NO_2 \rightarrow HOCH_2O_2NO_2$	`k_NO2_CH3O2`	see note*
G4149	StTrGN	$HOCH_2O_2NO_2 \rightarrow HOCH_2O_2 + NO_2$	`k_NO2_CH3O2/(9.5E-29*EXP(11234./temp))`	Barnes et al. (1985)*
G4150	StTrGN	$HOCH_2O_2NO_2 + OH \rightarrow HCOOH + NO_3 + H_2O$	`9.50E-13*EXP(-650./temp)*f_soh`	see note*
G4151	StTrG	$CH_3 + O_2 \rightarrow CH_3O_2$	`k_3rd_iupac(temp,cair,7.0E-31,3.,1.8E-12,-1.1,0.33)`	Atkinson et al. (2006)
G4152	StTrG	$CH_3 + O_3 \rightarrow .956\ HCHO + .956\ H + .044\ CH_3O + O_2$	`5.1E-12*exp(-210./temp)`	Albaladejo et al. (2002), Ogryzlo et al. (1981)
G4153	StTrG	$CH_3 + O(^3P) \rightarrow .83\ HCHO + .83\ H + .17\ CO + .17\ H_2 + .17\ H$	`1.3E-10`	Atkinson et al. (2006)
G4154	StTrG	$CH_3O + O_3 \rightarrow CH_3O_2 + O_2$	`2.53E-14`	Albaladejo et al. (2002)*
G4155	StTrG	$CH_3O + O(^3P) \rightarrow .75\ CH_3 + .75\ O_2 + .25\ HCHO + .25\ OH$	`2.5E-11`	Baulch et al. (2005)
G4156	StTrG	$CH_3O_2 + O(^3P) \rightarrow CH_3O + O_2$	`4.3E-11`	Zellner et al. (1988)
G4157	StTrG	$HCHO + O(^3P) \rightarrow .7\ OH + .7\ CO + .3\ H + .3\ CO_2 + HO_2$	`3.4E-11*EXP(-1600./temp)`	Burkholder et al. (2015)
G4158	TrG	$CH_2OO^* \rightarrow .37\ CH_2OO + .47\ CO + .47\ H_2O + .16\ HO_2 + .16\ CO + .16\ OH$	`KDEC`	Atkinson et al. (2006)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G4159	TrGN	$HCN + OH \rightarrow H_2O + CN$	k_3rd(temp,cair,4.28E-33,1.0, REAL(4.25E-13*EXP(-1150./temp) ,SP),1.0,0.8)	Kleinböhl et al. (2006)
G4160a	TrGN	$HCN + O(^1D) \rightarrow O(^3P) + HCN$	1.08E-10*EXP(105./temp) *0.15*EXP(200/temp)	Strekowski et al. (2010)
G4160b	TrGN	$HCN + O(^1D) \rightarrow H + NCO$	1.08E-10*EXP(105./temp)*0.68/2.	Strekowski et al. (2010)*
G4160c	TrGN	$HCN + O(^1D) \rightarrow OH + CN$	1.08E-10*EXP(105./temp)*(1.-(0.68/ 2.+0.15*EXP(200/temp)))	Strekowski et al. (2010)*
G4161	TrGN	$HCN + O(^3P) \rightarrow H + NCO$	1.0E-11*EXP(-4000./temp)	Burkholder et al. (2015)*
G4162	TrGN	$CN + O_2 \rightarrow NCO + O(^3P)$	1.2E-11*EXP(210./temp)*0.75	Baulch et al. (2005)
G4163	TrGN	$CN + O_2 \rightarrow CO + NO$	1.2E-11*EXP(210./temp)*0.25	Baulch et al. (2005)
G4164	TrGN	$NCO + O_2 \rightarrow CO_2 + NO$	7.E-15	Becker et al. (2000)*
G42000	TrGC	$C_2H_6 + OH \rightarrow C_2H_5O_2 + H_2O$	1.49E-17*temp*temp*EXP(-499./ temp)	Atkinson et al. (2006)
G42001	TrGC	$C_2H_4 + O_3 \rightarrow HCHO + CH_2OO^*$	9.1E-15*EXP(-2580./temp)	Atkinson et al. (2006)*
G42002	TrGC	$C_2H_4 + OH \rightarrow HOCH_2CH_2O_2$	k_3rd_iupac(temp,cair,8.6E-29, 3.1,9.E-12,0.85,0.48)	Atkinson et al. (2006), Rickard and Pascoe (2009)
G42003	TrGC	$C_2H_5O_2 + HO_2 \rightarrow C_2H_5OOH$	7.5E-13*EXP(700./temp)	Burkholder et al. (2015)
G42004a	TrGCN	$C_2H_5O_2 + NO \rightarrow CH_3CHO + HO_2 + NO_2$	2.55E-12*EXP(380./temp)*(1.-beta_ C2H5NO3)	Atkinson et al. (2006), Butkovskaya et al. (2010)
G42004b	TrGCN	$C_2H_5O_2 + NO \rightarrow C_2H_5ONO_2$	2.55E-12*EXP(380./temp)*beta_ C2H5NO3	Atkinson et al. (2006), Butkovskaya et al. (2010)
G42005	TrGCN	$C_2H_5O_2 + NO_3 \rightarrow CH_3CHO + HO_2 + NO_2$	2.3E-12	Wallington et al. (2018)
G42006	TrGC	$C_2H_5O_2 \rightarrow .8\ CH_3CHO + .6\ HO_2 + .2\ C_2H_5OH$	2.*(7.6E-14*k_CH3O2)**(.5)*RO2	Sander et al. (2018), Atkinson et al. (2006)
G42007a	TrGC	$C_2H_5OOH + OH \rightarrow C_2H_5O_2 + H_2O$	k_roohro	Sander et al. (2018)
G42007b	TrGC	$C_2H_5OOH + OH \rightarrow CH_3CHO + OH$	k_s*f_sooh	Sander et al. (2018)
G42008a	TrGC	$CH_3CHO + OH \rightarrow CH_3C(O) + H_2O$	4.4E-12*EXP(365./temp)*0.95	Atkinson et al. (2006)
G42008b	TrGC	$CH_3CHO + OH \rightarrow HCOCH_2O_2 + H_2O$	4.4E-12*EXP(365./temp)*0.05	Atkinson et al. (2006)
G42009	TrGCN	$CH_3CHO + NO_3 \rightarrow CH_3C(O) + HNO_3$	KNO3AL	Rickard and Pascoe (2009)
G42010	TrGC	$CH_3COOH + OH \rightarrow CH_3 + CO_2 + H_2O$	k_CH3CO2H_OH	Atkinson et al. (2006)*
G42011a	TrGC	$CH_3C(O)OO + HO_2 \rightarrow OH + CH_3 + CO_2$	5.20E-13*EXP(980./temp)*1.507*0.61	Groß et al. (2014)
G42011b	TrGC	$CH_3C(O)OO + HO_2 \rightarrow CH_3C(O)OOH$	5.20E-13*EXP(980./temp)*1.507*0.23	Groß et al. (2014)
G42011c	TrGC	$CH_3C(O)OO + HO_2 \rightarrow CH_3COOH + O_3$	5.20E-13*EXP(980./temp)*1.507*0.16	Groß et al. (2014)
G42012	TrGCN	$CH_3C(O)OO + NO \rightarrow CH_3 + CO_2 + NO_2$	8.1E-12*EXP(270./temp)	Tyndall et al. (2001a)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G42013	TrGCN	$CH_3C(O)OO + NO_2 \rightarrow PAN$	`k_CH3CO3_NO2`	Burkholder et al. (2015)*
G42014	TrGCN	$CH_3C(O)OO + NO_3 \rightarrow CH_3 + NO_2 + CO_2$	`4.E-12`	Canosa-Mas et al. (1996)
G42017a	TrGC	$CH_3C(O)OO \rightarrow CH_3 + CO_2$	`k1_RO2RCO3*0.9`	Sander et al. (2018)
G42017b	TrGC	$CH_3C(O)OO \rightarrow CH_3COOH$	`k1_RO2RCO3*0.1`	Sander et al. (2018)
G42018	TrGC	$CH_3C(O)OOH + OH \rightarrow CH_3C(O)OO + H_2O$	`k_roohro`	Rickard and Pascoe (2009)*
G42020	TrGCN	$PAN + OH \rightarrow HCHO + CO + NO_2 + H_2O$	`3.00E-14`	Rickard and Pascoe (2009)
G42021	TrGCN	$PAN \rightarrow CH_3C(O)OO + NO_2$	`k_PAN_M`	Burkholder et al. (2015)*
G42022a	TrGC	$C_2H_2 + OH \rightarrow GLYOX + OH$	`k_3rd(temp,cair,5.5e-30,0.0, 8.3e-13,-2.,0.6)*0.71`	Burkholder et al. (2015)*
G42022b	TrGC	$C_2H_2 + OH \rightarrow HCOOH + CO + HO_2$	`k_3rd(temp,cair,5.5e-30,0.0, 8.3e-13,-2.,0.6)*0.29`	Burkholder et al. (2015)*
G42023a	TrGC	$HOCH_2CHO + OH \rightarrow HOCH2CO + H_2O$	`8.00E-12*0.80`	Atkinson et al. (2006)
G42023b	TrGC	$HOCH_2CHO + OH \rightarrow HOCHCHO + H_2O$	`8.00E-12*0.20`	Atkinson et al. (2006)
G42024a	TrGC	$HOCH2CO + O_2 \rightarrow HOCH_2CO_3$	`5.1E-12*(1.-1./(1+1.85E-18*cair))`	Atkinson et al. (2006), Beyersdorf et al. (2010)*
G42024b	TrGC	$HOCH2CO + O_2 \rightarrow OH + HCHO + CO_2$	`5.1E-12*1./(1+1.85E-18*cair)`	Atkinson et al. (2006), Beyersdorf et al. (2010)*
G42025	TrGC	$HOCHCHO \rightarrow GLYOX + HO_2$	`KDEC`	Sander et al. (2018)
G42026	TrGCN	$HOCH_2CHO + NO_3 \rightarrow HOCH2CO + HNO_3$	`KNO3AL`	Rickard and Pascoe (2009)
G42027a	TrGC	$HOCH_2CO_3 \rightarrow HCHO + CO_2 + HO_2$	`k1_RO2RCO3*0.9`	Sander et al. (2018)
G42027b	TrGC	$HOCH_2CO_3 \rightarrow HOCH_2CO_2H$	`k1_RO2RCO3*0.1`	Sander et al. (2018)
G42028a	TrGC	$HOCH_2CO_3 + HO_2 \rightarrow HCHO + HO_2 + OH + CO_2$	`KAPHO2*rco3_oh`	Sander et al. (2018), Groß et al. (2014)
G42028b	TrGC	$HOCH_2CO_3 + HO_2 \rightarrow HOCH_2CO_3H$	`KAPHO2*rco3_ooh`	Sander et al. (2018), Groß et al. (2014)
G42028c	TrGC	$HOCH_2CO_3 + HO_2 \rightarrow HOCH_2CO_2H + O_3$	`KAPHO2*rco3_o3`	Sander et al. (2018), Groß et al. (2014)
G42029	TrGCN	$HOCH_2CO_3 + NO \rightarrow NO_2 + HO_2 + HCHO + CO_2$	`KAPNO`	Rickard and Pascoe (2009)
G42030	TrGCN	$HOCH_2CO_3 + NO_2 \rightarrow PHAN$	`k_CH3CO3_NO2`	Rickard and Pascoe (2009)
G42031	TrGCN	$HOCH_2CO_3 + NO_3 \rightarrow NO_2 + HO_2 + HCHO + CO_2$	`KRO2NO3*1.74`	Rickard and Pascoe (2009)
G42032	TrGC	$HOCH_2CO_2H + OH \rightarrow .09\ HCHO + .09\ CO_2 + .91\ HCOCO_2H + HO_2 + H_2O$	`k_co2h+k_s*f_soh*f_co2h`	Sander et al. (2018)
G42033a	TrGC	$HOCH_2CO_3H + OH \rightarrow HOCH_2CO_3 + H_2O$	`k_roohro`	Sander et al. (2018)
G42033b	TrGC	$HOCH_2CO_3H + OH \rightarrow HCOCO_3H + HO_2$	`k_s*f_soh*f_co2h`	Sander et al. (2018)
G42034	TrGCN	$PHAN \rightarrow HOCH_2CO_3 + NO_2$	`k_PAN_M`	Rickard and Pascoe (2009)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G42035	TrGCN	PHAN + OH → HCHO + CO + NO_2 + H_2O	`k_s*f_soh*f_cpan+k_rohro`	Sander et al. (2018)
G42036	TrGC	GLYOX + OH → HCOCO + H_2O	`3.1E-12*EXP(340./temp)`	Atkinson et al. (2006), Orlando and Tyndall (2001), Lockhart et al. (2013)
G42037	TrGCN	GLYOX + NO_3 → HCOCO + HNO_3	`KNO3AL`	Rickard and Pascoe (2009)
G42038a	TrGC	HCOCO → CO + CO + HO_2	`7.E11*EXP(-3160./temp) +5.E-12*c(ind_O2)`	Orlando and Tyndall (2001), Lockhart et al. (2013), Rickard and Pascoe (2009)
G42037b	TrGC	HCOCO → $HCOCO_3$	`5.E-12*c(ind_O2)*3.2*exp(-550./ temp)`	Lockhart et al. (2013), Rickard and Pascoe (2009)
G42037c	TrGC	HCOCO → OH + CO + CO_2	`5.E-12*c(ind_O2) *(1.-3.2*exp(-550./temp))`	Lockhart et al. (2013), Rickard and Pascoe (2009)
G42039a	TrGC	$HCOCO_3$ → CO + HO_2 + CO_2	`k1_RO2RCO3*0.9`	Sander et al. (2018)
G42039b	TrGC	$HCOCO_3$ → $HCOCO_2H$	`k1_RO2RCO3*0.1`	Sander et al. (2018)
G42040	TrGC	$HCOCO_3$ + HO_2 → HO_2 + CO + CO_2 + OH	`KAPHO2`	Feierabend et al. (2008), Sander et al. (2018)
G42041	TrGCN	$HCOCO_3$ + NO → HO_2 + CO + NO_2 + CO_2	`KAPNO`	Rickard and Pascoe (2009)
G42042	TrGCN	$HCOCO_3$ + NO_3 → HO_2 + CO + NO_2 + CO_2	`KRO2NO3*1.74`	Rickard and Pascoe (2009)
G42043	TrGCN	$HCOCO_3$ + NO_2 → HO_2 + CO + NO_3 + CO_2	`k_CH3CO3_NO2`	Orlando and Tyndall (2001), Sander et al. (2018)
G42044	TrGC	$HCOCO_2H$ + OH → CO + HO_2 + CO_2 + H_2O	`k_co2h+k_t*f_o*f_co2h`	Sander et al. (2018)
G42045a	TrGC	$HCOCO_3H$ + OH → $HCOCO_3$ + H_2O	`k_roohro`	Sander et al. (2018)
G42045b	TrGC	$HCOCO_3H$ + OH → CO + CO_2 + H_2O + OH	`k_t*f_o*f_co2h`	Sander et al. (2018)
G42046	TrGC	$HOCH_2CH_2O_2$ → .6 $HOCH_2CH_2O$ + .2 $HOCH_2CHO$ + .2 ETHGLY	`2.*(7.8E-14*EXP(1000./temp) *k_CH3O2)**(.5)*RO2`	Atkinson et al. (2006), Rickard and Pascoe (2009)
G42047	TrGCN	$HOCH_2CH_2O_2$ + NO → .25 HO_2 + .5 HCHO + .75 $HOCH_2CH_2O$ + NO_2	`KRO2NO*(1.-alpha_AN(3,1,0,0,0, temp,cair))`	Rickard and Pascoe (2009)*
G42048	TrGCN	$HOCH_2CH_2O_2$ + NO → ETHOHNO3	`KRO2NO*alpha_AN(3,1,0,0,0,temp, cair)`	Sander et al. (2018)
G42049a	TrGC	$HOCH_2CH_2O_2$ + HO_2 → HYETHO2H	`1.53E-13*EXP(1300./temp) *(1.-rchohch2o2_oh)`	Rickard and Pascoe (2009)
G42049b	TrGC	$HOCH_2CH_2O_2$ + HO_2 → $HOCH_2CH_2O$ + OH	`1.53E-13*EXP(1300./temp) *rchohch2o2_oh`	Rickard and Pascoe (2009)
G42050	TrGCN	ETHOHNO3 + OH → .93 NO_3CH2CHO + .93 HO_2 + .07 $HOCH_2CHO$ + .07 NO_2 + H_2O	`k_s*(f_soh*f_ch2ono2+f_ono2*f_ pch2oh)+k_rohro`	Sander et al. (2018)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G42051a	TrGC	HYETHO2H + OH → $HOCH_2CH_2O_2$ + H_2O	k_roohro	Rickard and Pascoe (2009)*
G42051b	TrGC	HYETHO2H + OH → $HOCH_2CHO$ + OH + H_2O	k_s*f_sooh*f_pch2oh	Sander et al. (2018)
G42051c	TrGC	HYETHO2H + OH → HOOCH2CHO + HO_2 + H_2O	k_s*f_soh*f_pch2oh+k_rohro	Sander et al. (2018)
G42052a	TrGC	$HOCH_2CH_2O$ → HO_2 + $HOCH_2CHO$	6.00E-14*EXP(-550./temp) *C(ind_O2)	Rickard and Pascoe (2009)
G42052b	TrGC	$HOCH_2CH_2O$ → HO_2 + HCHO + HCHO	9.50E13*EXP(-5988./temp)	Rickard and Pascoe (2009)
G42053	TrGC	ETHGLY + OH → $HOCH_2CHO$ + HO_2 + H_2O	2*k_s*f_soh*f_pch2oh+2*k_rohro	Sander et al. (2018)
G42054	TrGC	$HCOCH_2O_2$ → .6 HCHO + .6 CO + .6 HO_2 + .2 GLYOX + .2 $HOCH_2CHO$	k1_RO2pORO2	Sander et al. (2018)
G42055a	TrGC	$HCOCH_2O_2$ + HO_2 → HOOCH2CHO	KRO2HO2(2)*rcoch2o2_ooh	Sander et al. (2018)
G42055b	TrGC	$HCOCH_2O_2$ + HO_2 → HCHO + CO + HO_2 + OH	KRO2HO2(2)*rcoch2o2_oh	Sander et al. (2018)
G42056a	TrGCN	$HCOCH_2O_2$ + NO → NO_2 + HCHO + CO + HO_2	KRO2NO*(1.-alpha_AN(3,1,1,0,0, temp,cair))	Sander et al. (2018)
G42056b	TrGCN	$HCOCH_2O_2$ + NO → NO_3CH2CHO	KRO2NO*alpha_AN(3,1,1,0,0,temp, cair)	Sander et al. (2018)
G42057	TrGCN	$HCOCH_2O_2$ + NO_3 → HCHO + CO + HO_2 + NO_2	KRO2NO3	Sander et al. (2018)
G42058a	TrGC	HOOCH2CHO + OH → $HCOCH_2O_2$	k_roohro	Sander et al. (2018)
G42058b	TrGC	HOOCH2CHO + OH → HCHO + CO + OH	0.8*8.E-12	Sander et al. (2018)*
G42058c	TrGC	HOOCH2CHO + OH → GLYOX + OH	k_s*f_sooh*f_cho	Sander et al. (2018)
G42059	TrGCN	HOOCH2CHO + NO_3 → OH + HCHO + CO + HNO_3	KNO3AL	Rickard and Pascoe (2009)
G42060	TrGCN	$HOOCH_2CO_3$ + NO → NO_2 + OH + HCHO + CO_2	KAPNO	Sander et al. (2018)
G42061	TrGCN	$HOOCH_2CO_3$ + NO_3 → NO_2 + OH + HCHO + CO_2	KRO2NO3*1.74	Sander et al. (2018)
G42062a	TrGC	$HOOCH_2CO_3$ + HO_2 → 2 OH + HCHO + CO_2	KAPHO2*rco3_oh	Sander et al. (2018)
G42062b	TrGC	$HOOCH_2CO_3$ + HO_2 → HOOCH2CO3H	KAPHO2*rco3_ooh	Sander et al. (2018)
G42062c	TrGC	$HOOCH_2CO_3$ + HO_2 → HOOCH2CO2H + O_3	KAPHO2*rco3_o3	Sander et al. (2018)
G42063a	TrGC	$HOOCH_2CO_3$ → OH + HCHO + CO_2	k1_RO2RCO3*0.9	Sander et al. (2018)
G42063b	TrGC	$HOOCH_2CO_3$ → HOOCH2CO2H	k1_RO2RCO3*0.1	Sander et al. (2018)
G42064a	TrGC	HOOCH2CO3H + OH → $HOOCH_2CO_3$ + H_2O	2.*k_roohro	Sander et al. (2018)
G42064b	TrGC	HOOCH2CO3H + OH → $HCOCO_3H$ + OH + H_2O	k_s*f_sooh*f_co2h	Sander et al. (2018)
G42065	TrGC	HOOCH2CO2H + OH → $HCOCO_2H$ + OH + H_2O	k_s*f_sooh*f_co2h+k_co2h	Sander et al. (2018)
G42066	TrGC	CH2CO + OH → .6 HCHO + .6 HO_2 + .6 CO + .4 HOOCH2CO2H	2.8E-12*exp(510./temp)	Baulch et al. (2005), Sander et al. (2018)
G42067a	TrGC	CH3CHOHOOH + OH → CH_3COOH + OH	(k_t*f_tooh*f_toh + k_rohro)	Sander et al. (2018)
G42067b	TrGC	CH3CHOHOOH + OH → CH3CHOHO2	k_roohro	Sander et al. (2018)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G42068	TrGC	CH3CHOHO2 $\rightarrow$ CH_3CHO + HO_2	`3.46E12*EXP(-12500./(1.98*temp))`	Hermans et al. (2005), Sander et al. (2018)
G42069	TrGC	CH_3CHO + HO_2 $\rightarrow$ CH3CHOHO2	`3.46E12*EXP(-12500./(1.98*temp)) /(6.34E26*EXP(-14700./ (1.98*temp)))`	Hermans et al. (2005), Sander et al. (2018)
G42070	TrGC	CH3CHOHO2 + HO_2 $\rightarrow$ .5 CH3CHOHOOH + .3 CH_3COOH + .2 CH_3 + .2 HCOOH + .2 OH	`5.6E-15*EXP(2300./temp)`	Sander et al. (2018)
G42071	TrGC	CH3CHOHO2 $\rightarrow$ CH_3 + HCOOH + OH	`k1_RO2sORO2`	Sander et al. (2018)
G42072	TrGCN	CH3CHOHO2 + NO $\rightarrow$ CH_3 + HCOOH + OH + NO_2	`KRO2NO`	Sander et al. (2018)
G42073	TrGCN	$C_2H_5ONO_2$ + OH $\rightarrow$ CH_3CHO + H_2O + NO_2	`6.7E-13*EXP(-395./temp)`	Atkinson et al. (2006)
G42074a	TrGCN	NO_3CH2CHO + OH $\rightarrow$ GLYOX + NO_2 + H_2O	`k_s*f_ch2ono2*f_cho`	Paulot et al. (2009a), Sander et al. (2018)*
G42074b	TrGCN	NO_3CH2CHO + OH $\rightarrow$ NO_3CH2CO_3 + H_2O	`k_t*f_o*f_ch2ono2*3.`	Paulot et al. (2009a), Sander et al. (2018)*
G42075	TrGCN	NO_3CH2CO_3 + HO_2 $\rightarrow$ HCHO + NO_2 + CO_2 + OH	`KAPHO2`	Rickard and Pascoe (2009)*
G42076	TrGCN	NO_3CH2CO_3 + NO $\rightarrow$ HCHO + NO_2 + CO_2 + NO_2	`KAPNO`	Rickard and Pascoe (2009)
G42077	TrGCN	NO_3CH2CO_3 + NO_2 $\rightarrow$ NO_3CH2CHO	`k_CH3CO3_NO2`	Rickard and Pascoe (2009)
G42078	TrGCN	NO_3CH2CO_3 $\rightarrow$ HCHO + NO_2 + CO_2	`k1_RO2RCO3`	Rickard and Pascoe (2009)*
G42079	TrGCN	NO_3CH2CHO $\rightarrow$ NO_3CH2CO_3 + NO_2	`k_PAN_M`	Rickard and Pascoe (2009)
G42080	StTrGCN	$C_2H_5O_2$ + NO_2 $\rightarrow$ $C_2H_5O_2NO_2$	`k_3rd_iupac(temp,cair,1.3E-29, 6.2,8.8E-12,0.0,0.31)`	Atkinson et al. (2006)
G42081	StTrGCN	$C_2H_5O_2NO_2$ $\rightarrow$ $C_2H_5O_2$ + NO_2	`k_3rd_iupac(temp,cair, REAL(4.8E-4*EXP(-9285./temp) ,SP),0.0,REAL(8.8E15*EXP(-10440./ temp),SP),0.0,0.31)`	Atkinson et al. (2006)
G42082	StTrGCN	$C_2H_5O_2NO_2$ + OH $\rightarrow$ CH_3CHO + NO_3 + H_2O	`9.50E-13*EXP(-650./temp)`	Sander et al. (2018)*
G42083a	TrGC	$CH_3C(O)$ + O_2 $\rightarrow$ $CH_3C(O)OO$	`5.1E-12*(1. - 1./(1.+ 9.4E-18*cair))`	Atkinson et al. (2006), Beyersdorf et al. (2010)*
G42083b	TrGC	$CH_3C(O)$ + O_2 $\rightarrow$ OH + HCHO + CO	`5.1E-12*1./(1.+9.4E-18*cair)`	Atkinson et al. (2006), Beyersdorf et al. (2010)*
G42084	TrGC	C_2H_5OH + OH $\rightarrow$ .95 $C_2H_5O_2$ + .95 HO_2 + .05 $HOCH_2CH_2O_2$ + H_2O	`3.0E-12*EXP(20./temp)`	Sander et al. (2018), Atkinson et al. (2006)
G42085a	TrGCN	CH_3CN + OH $\rightarrow$ $NCCH_2O_2$ + H_2O	`8.1E-13*EXP(-1080./temp)*0.40`	Atkinson et al. (2006), Tyndall et al. (2001b)*

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G42085b	TrGCN	$CH_3CN + OH \rightarrow OH + CH_3C(O) + NO$	`8.1E-13*EXP(-1080./temp)*(1.-0.40)`	Atkinson et al. (2006), Tyndall et al. (2001b)*
G42086a	TrGCN	$CH_3CN + O(^1D) \rightarrow O(^3P) + CH_3CN$	`2.54E-10*EXP(-24./temp) *0.0269*EXP(137./temp)`	Strekowski et al. (2010)
G42086b	TrGCN	$CH_3CN + O(^1D) \rightarrow 2\ H + CO + HCN$	`2.54E-10*EXP(-24./temp)*0.16`	Strekowski et al. (2010)*
G42086c	TrGCN	$CH_3CN + O(^1D) \rightarrow .5\ CH_3 + .5\ NCO + .5\ NCCH_2O_2 + .5\ OH$	`2.54E-10*EXP(-24./temp)*(1.-(0.16+ 0.0269*EXP(137./temp)))`	Strekowski et al. (2010)*
G42087	TrGCN	$NCCH_2O_2 + NO \rightarrow HCN + CO_2 + HO_2 + NO_2$	`KRO2NO`	see note*
G42088	TrGCN	$NCCH_2O_2 + HO_2 \rightarrow HCN + CO_2 + HO_2$	`KRO2HO2(2)`	see note*
G42089a	TrGC	$CH_2CHOH + OH \rightarrow HCOOH + OH + HCHO$	`k_CH2CHOH_OH_HCOOH`	Sander et al. (2018), So et al. (2014)*
G42089b	TrGC	$CH_2CHOH + OH \rightarrow HOCH_2CHO + HO_2$	`k_CH2CHOH_OH_ALD`	Sander et al. (2018), So et al. (2014)
G42090	TrGC	$CH_2CHOH + HCOOH \rightarrow CH_3CHO + HCOOH$	`k_CH2CHOH_HCOOH`	Sander et al. (2018), da Silva (2010)*
G42091	TrGC	$CH_3CHO + HCOOH \rightarrow CH_2CHOH + HCOOH$	`k_ALD_HCOOH`	Sander et al. (2018), da Silva (2010)*
G43000a	TrGC	$C_3H_8 + OH \rightarrow iC_3H_7O_2 + H_2O$	`k_s`	Sander et al. (2018)
G43000b	TrGC	$C_3H_8 + OH \rightarrow C_3H_7O_2 + H_2O$	`2.*k_p`	Sander et al. (2018)
G43001a	TrGC	$C_3H_6 + O_3 \rightarrow HCHO + .16$ CH3CHOHOOH $+ .50\ OH + .50\ HCOCH_2O_2 + .05$ CH2CO $+ .09\ CH_3OH + .09\ CO + .2\ CH_4 + .2\ CO_2$	`5.5E-15*EXP(-1880./temp)*.57`	Atkinson et al. (2006)*
G43001b	TrGC	$C_3H_6 + O_3 \rightarrow CH_3CHO + CH_2OO^*$	`5.5E-15*EXP(-1880./temp)*.43`	Atkinson et al. (2006)*
G43002	TrGC	$C_3H_6 + OH \rightarrow$ HYPROPO2	`k_3rd_iupac(temp,cair,8.6E-27, 3.5,3.E-11,1.,0.5)`	Atkinson et al. (2006), Rickard and Pascoe (2009)
G43003	TrGCN	$C_3H_6 + NO_3 \rightarrow$ PRONO3BO2	`4.6E-13*EXP(-1155./temp)`	Wallington et al. (2018)
G43004	TrGC	$iC_3H_7O_2 + HO_2 \rightarrow iC_3H_7OOH$	`1.9E-13*EXP(1300./temp)`	Atkinson (1997)*
G43005a	TrGCN	$iC_3H_7O_2 + NO \rightarrow CH_3COCH_3 + HO_2 + NO_2$	`2.7E-12*EXP(360./temp)*(1.-alpha_ AN(3,2,0,0,0,temp,cair))`	Wallington et al. (2018)
G43005b	TrGCN	$iC_3H_7O_2 + NO \rightarrow iC_3H_7ONO_2$	`2.7E-12*EXP(360./temp)*alpha_ AN(3,2,0,0,0,temp,cair)`	Wallington et al. (2018)
G43006	TrGC	$iC_3H_7O_2 \rightarrow .8\ CH_3COCH_3 + .2$ IPROPOL $+ .6\ HO_2$	`2.*(1.6E-12*EXP(-2200./temp) *k_CH3O2)**(.5)*RO2`	Rickard and Pascoe (2009), Atkinson et al. (2006)
G43007a	TrGC	$iC_3H_7OOH + OH \rightarrow iC_3H_7O_2 + H_2O$	`k_roohro`	Sander et al. (2018)
G43007b	TrGC	$iC_3H_7OOH + OH \rightarrow CH_3COCH_3 + H_2O + OH$	`k_t*f_tooh`	Sander et al. (2018)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G43008	TrGC	$C_3H_7O_2 + HO_2 \rightarrow C_3H_7OOH$	`1.9E-13*EXP(1300./temp)`	Atkinson (1997)*
G43009a	TrGCN	$C_3H_7O_2 + NO \rightarrow C_2H_5CHO + HO_2 + NO_2$	`2.7E-12*EXP(360./temp)*(1.-alpha_AN(3,1,0,0,0,temp,cair))`	Wallington et al. (2018)
G43009b	TrGCN	$C_3H_7O_2 + NO \rightarrow C_3H_7ONO_2$	`2.7E-12*EXP(360./temp)*alpha_AN(3,1,0,0,0,temp,cair)`	Wallington et al. (2018)
G43010	TrGC	$C_3H_7O_2 \rightarrow .8\ CH_3COCH_3 + .2$ NPROPOL $+ .6\ HO_2$	`2.*(k_CH3O2*3.E-13)**(.5)*RO2`	Rickard and Pascoe (2009), Atkinson et al. (2006)
G43011	TrGC	$CH_3COCH_3 + OH \rightarrow CH_3COCH_2O_2 + H_2O$	`(8.8E-12*EXP(-1320./temp) +1.7E-14*EXP(423./temp))`	Atkinson et al. (2006)*
G43012a	TrGC	$CH_3COCH_2O_2 + HO_2 \rightarrow CH_3COCH_2O_2H$	`8.6E-13*EXP(700./temp)*rcoch2o2_ooh`	Tyndall et al. (2001a), Sander et al. (2018)
G43012b	TrGC	$CH_3COCH_2O_2 + HO_2 \rightarrow OH + CH_3C(O) + HCHO$	`8.6E-13*EXP(700./temp)*rcoch2o2_oh`	Tyndall et al. (2001a), Sander et al. (2018)
G43013a	TrGCN	$CH_3COCH_2O_2 + NO \rightarrow CH_3C(O) + HCHO + NO_2$	`2.9E-12*EXP(300./temp)*(1.-alpha_AN(4,1,1,0,0,temp,cair))`	Burkholder et al. (2015)
G43013b	TrGCN	$CH_3COCH_2O_2 + NO \rightarrow$ NOA	`2.9E-12*EXP(300./temp)*alpha_AN(4,1,1,0,0,temp,cair)`	Burkholder et al. (2015)
G43014	TrGC	$CH_3COCH_2O_2 \rightarrow .3\ CH_3C(O) + .3\ HCHO + .5$ MGLYOX $+ .2\ CH_3COCH_2OH$	`k1_RO2pORO2`	Orlando and Tyndall (2012)
G43015a	TrGC	$CH_3COCH_2O_2H + OH \rightarrow CH_3COCH_2O_2 + H_2O$	`k_roohro`	see note*
G43015b	TrGC	$CH_3COCH_2O_2H + OH \rightarrow$ MGLYOX $+ OH + H_2O$	`k_s*f_sooh*f_co`	Sander et al. (2018)
G43016	TrGC	$CH_3COCH_2OH + OH \rightarrow$ MGLYOX $+ HO_2 + H_2O$	`1.6E-12*EXP(305./temp)`	Atkinson et al. (2006)
G43017	TrGC	MGLYOX $+ OH \rightarrow .4\ CH_3 + .6\ CH_3C(O) + 1.4\ CO + H_2O$	`1.9E-12*EXP(575./temp)`	Baeza-Romero et al. (2007), Atkinson et al. (2006)
G43020	TrGCN	$iC_3H_7ONO_2 + OH \rightarrow CH_3COCH_3 + NO_2$	`6.2E-13*EXP(-230./temp)`	Wallington et al. (2018)
G43021	TrGCN	$CH_3COCH_2O_2 + NO_3 \rightarrow CH_3C(O) + HCHO + NO_2$	`KRO2NO3`	Rickard and Pascoe (2009)
G43022	TrGC	HYPROPO2 $\rightarrow CH_3CHO + HCHO + HO_2$	`k1_RO2sORO2`	Rickard and Pascoe (2009)
G43023a	TrGC	HYPROPO2 $+ HO_2 \rightarrow$ HYPROPO2H	`KRO2HO2(3)*(1.-rchohch2o2_oh)`	Rickard and Pascoe (2009)
G43023b	TrGC	HYPROPO2 $+ HO_2 \rightarrow CH_3CHO + HCHO + HO_2 + OH$	`KRO2HO2(3)*rchohch2o2_oh`	Rickard and Pascoe (2009)
G43024a	TrGCN	HYPROPO2 $+ NO \rightarrow CH_3CHO + HCHO + HO_2 + NO_2$	`KRO2NO*(1.-alpha_AN(4,1,0,0,0,temp,cair))`	Rickard and Pascoe (2009)
G43024b	TrGCN	HYPROPO2 $+ NO \rightarrow$ PROPOLNO3	`KRO2NO*alpha_AN(4,1,0,0,0,temp,cair)`	Rickard and Pascoe (2009)
G43025	TrGCN	HYPROPO2 $+ NO_3 \rightarrow CH_3CHO + HCHO + HO_2 + NO_2$	`KRO2NO3`	Rickard and Pascoe (2009)
G43026a	TrGC	HYPROPO2H $+ OH \rightarrow$ HYPROPO2	`k_roohro`	Rickard and Pascoe (2009)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G43026b	TrGC	HYPROPO2H + OH → CH_3COCH_2OH + OH	(k_s*f_soh*f_pch2oh+k_t*f_tooh*f_pch2oh)	Sander et al. (2018)
G43027	TrGCN	PRONO3BO2 + HO_2 → PR2O2HNO3	KRO2HO2(3)	Rickard and Pascoe (2009)
G43028	TrGCN	PRONO3BO2 + NO → NOA + HO_2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G43029	TrGCN	PRONO3BO2 + NO_3 → NOA + HO_2 + NO_2	KRO2NO3	Rickard and Pascoe (2009)
G43030a	TrGCN	PR2O2HNO3 + OH → PRONO3BO2	k_roohro	Rickard and Pascoe (2009)
G43030b	TrGCN	PR2O2HNO3 + OH → NOA + OH	k_t*f_tooh*f_ch2ono2	Sander et al. (2018)
G43031	TrGCN	MGLYOX + NO_3 → $CH_3C(O)$ + CO + HNO_3	KNO3AL*2.4	Rickard and Pascoe (2009)
G43032	TrGCN	NOA + OH → MGLYOX + NO_2	(k_s*f_co*f_ono2+k_p*f_co)	Sander et al. (2018)
G43033	TrGC	HOCH2COCHO + OH → .8609 HOCH2CO + .8609 CO + .1391 HCOCOCHO + .1391 HO_2	(1.9E-12*EXP(575./temp)+k_s*f_soh*f_co)	Sander et al. (2018)
G43034	TrGCN	HOCH2COCHO + NO_3 → HOCH2CO + CO + HNO_3	KNO3AL*2.4	Sander et al. (2018)
G43035	TrGC	CH_3COCO_2H + OH → $CH_3C(O)$ + H_2O + CO_2	4.9E-14*EXP(276./temp)	Mellouki and Mu (2003), Sander et al. (2018)
G43036	TrGC	$HCOCOCH_2O_2$ → .6 HCOCO + .6 HCHO + .2 HCOCOCHO + .2 HOCH2COCHO	k1_RO2pORO2	Sander et al. (2018)
G43037	TrGCN	$HCOCOCH_2O_2$ + NO → HCOCO + HCHO + NO_2	KRO2NO	Sander et al. (2018)*
G43038a	TrGC	$HCOCOCH_2O_2$ + HO_2 → $HCOCOCH_2OOH$	KRO2HO2(3)*rcoch2o2_ooh	Sander et al. (2018)
G43038b	TrGC	$HCOCOCH_2O_2$ + HO_2 → HCOCO + HCHO + OH	KRO2HO2(3)*rcoch2o2_oh	Sander et al. (2018)
G43039	TrGCN	$HCOCOCH_2O_2$ + NO_3 → HCOCO + HCHO + NO_2	KRO2NO3	Sander et al. (2018)
G43040a	TrGC	$HCOCOCH_2OOH$ + OH → $HOOCH_2CO_3$ + CO + H_2O	k_t*f_co*f_o	Sander et al. (2018)*
G43040b	TrGC	$HCOCOCH_2OOH$ + OH → HCOCOCHO + H_2O + OH	k_s*f_sooh*f_co	Sander et al. (2018)*
G43040c	TrGC	$HCOCOCH_2OOH$ + OH → $HCOCOCH_2O_2$ + H_2O	k_roohro	Sander et al. (2018)
G43041	TrGCN	$HCOCOCH_2OOH$ + NO_3 → $HOOCH_2CO_3$ + CO + HNO_3	KNO3AL*2.4	Sander et al. (2018)
G43042	TrGC	HOCH2COCH2O2 → HCHO + HOCH2CO	k1_RO2pORO2	Sander et al. (2018)
G43043a	TrGC	HOCH2COCH2O2 + HO_2 → HOCH2COCH2OOH	KRO2HO2(3)*rcoch2o2_ooh	Sander et al. (2018)
G43043b	TrGC	HOCH2COCH2O2 + HO_2 → HCHO + HOCH2CO + OH	KRO2HO2(3)*rcoch2o2_oh	Sander et al. (2018)
G43044	TrGCN	HOCH2COCH2O2 + NO → HCHO + HOCH2CO + NO_2	KRO2NO	Sander et al. (2018)*
G43045a	TrGC	HOCH2COCH2OOH + OH → HOCH2COCHO + OH	k_s*f_sooh*f_co	Sander et al. (2018)
G43045b	TrGC	HOCH2COCH2OOH + OH → HOCH2COCH2O2	k_roohro	Sander et al. (2018)
G43045c	TrGC	HOCH2COCH2OOH + OH → $HCOCOCH_2OOH$ + HO_2	1.60E-12*EXP(305./temp)	Sander et al. (2018)*
G43046	TrGC	CH3CHCO + OH → .72 CO + .72 CH_3CHO + .72 HO_2 + .21 CH_3COCO_2H + .07 CH_3CHO + .07 HO_2 + .07 CO_2	7.6E-11	Hatakeyama et al. (1985), Sander et al. (2018)
G43047	TrGCN	PROPOLNO3 + OH → CH_3COCH_2OH + NO_2	k_t*f_ono2*f_pch2oh+k_s*f_soh*f_ch2ono2	Sander et al. (2018)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G43048	TrGCN	$CH_3COCH_2O_2 + NO_2 \rightarrow CH_3COCH_2OONO_2$	2.3E-12*EXP(300./temp)	Tyndall et al. (2001a)*
G43049	TrGCN	$CH_3COCH_2OONO_2 \rightarrow CH_3COCH_2O_2 + NO_2$	1.9E16*EXP(-10830./temp)	Sehested et al. (1998)*
G43050	TrGCN	$CH_3COCH_2OONO_2 + OH \rightarrow MGLYOX + NO_3 + H_2O$	9.50E-13*EXP(-650./temp)*f_co	Sander et al. (2018)*
G43051a	TrGC	$C_3H_7OOH + OH \rightarrow C_3H_7O_2 + H_2O$	k_roohro	Sander et al. (2018)
G43051b	TrGC	$C_3H_7OOH + OH \rightarrow C_2H_5CHO + H_2O + OH$	k_s*f_sooh	Sander et al. (2018)
G43051c	TrGC	$C_3H_7OOH + OH \rightarrow C_2H_5CHO + HO_2 + H_2O$	k_s*f_pch2oh	Sander et al. (2018)*
G43052	TrGC	$C_2H_5CHO + OH \rightarrow C_2H_5CO_3 + H_2O$	4.9E-12*EXP(405./temp)	Atkinson et al. (2006)*
G43053	TrGCN	$C_2H_5CHO + NO_3 \rightarrow C_2H_5CO_3 + HNO_3$	6.3E-15	Atkinson et al. (2006)
G43054a	TrGC	$C_2H_5CO_3 \rightarrow C_2H_5O_2 + CO_2$	k1_RO2RCO3*0.9	Sander et al. (2018)
G43054b	TrGC	$C_2H_5CO_3 \rightarrow C_2H_5CO_2H$	k1_RO2RCO3*0.1	Sander et al. (2018)
G43055a	TrGC	$C_2H_5CO_3 + HO_2 \rightarrow C_2H_5O_2 + CO_2 + OH$	KAPHO2*rco3_oh	Sander et al. (2018), Groß et al. (2014)
G43055b	TrGC	$C_2H_5CO_3 + HO_2 \rightarrow C_2H_5CO_3H$	KAPHO2*rco3_ooh	Sander et al. (2018), Groß et al. (2014)
G43055c	TrGC	$C_2H_5CO_3 + HO_2 \rightarrow C_2H_5CO_2H + O_3$	KAPHO2*rco3_o3	Sander et al. (2018), Groß et al. (2014)
G43056	TrGCN	$C_2H_5CO_3 + NO \rightarrow NO_2 + C_2H_5O_2 + CO_2$	KAPNO	Rickard and Pascoe (2009)
G43057	TrGCN	$C_2H_5CO_3 + NO_2 \rightarrow PPN$	k_CH3CO3_NO2	Rickard and Pascoe (2009)
G43058	TrGCN	$PPN \rightarrow C_2H_5CO_3 + NO_2$	k_PAN_M	Rickard and Pascoe (2009)
G43059	TrGC	$C_2H_5CO_2H + OH \rightarrow CH_3CHO + CO_2 + H_2O$	k_co2h+k_p+k_s*f_co2h	Sander et al. (2018)*
G43060a	TrGC	$C_2H_5CO_3H + OH \rightarrow C_2H_5CO_3 + H_2O$	k_roohro	Sander et al. (2018)
G43060b	TrGC	$C_2H_5CO_3H + OH \rightarrow CH_3CHO + CO_2 + H_2O$	k_s*f_co2h+k_p	Sander et al. (2018)*
G43061	TrGCN	$PPN + OH \rightarrow CH_3CHO + CO_2 + NO_2 + H_2O$	k_s*f_cpan+k_p	Sander et al. (2018)*
G43062	TrGC	$CH_3COCO_3H + OH \rightarrow CH_3COCO_3 + H_2O$	k_roohro	Sander et al. (2018)
G43063a	TrGC	$CH_3COCO_3 + HO_2 \rightarrow CH_3C(O) + CO_2 + OH$	KAPHO2*rco3_oh	Sander et al. (2018)
G43063b	TrGC	$CH_3COCO_3 + HO_2 \rightarrow CH_3COCO_3H$	KAPHO2*(rco3_ooh+rco3_o3)	Sander et al. (2018)
G43064	TrGCN	$CH_3COCO_3 + NO \rightarrow CH_3C(O) + CO_2 + NO_2$	KAPNO	Sander et al. (2018)
G43065	TrGCN	$CH_3COCO_3 + NO_2 \rightarrow CH_3C(O) + CO_2 + NO_3$	k_CH3CO3_NO2	Sander et al. (2018)*
G43066	TrGCN	$CH_3COCO_3 + NO_3 \rightarrow CH_3C(O)OO + CO_2 + NO_2$	KRO2NO3*1.74	Sander et al. (2018)
G43067	TrGC	$CH_3COCO_3 \rightarrow CH_3C(O)OO + CO_2$	k1_RO2RCO3	Sander et al. (2018)
G43068	TrGC	$HCOCOCHO + OH \rightarrow 3\ CO + HO_2$	2.*k_t*f_co*f_o	Sander et al. (2018)
G43069	TrGC	$IPROPOL + OH \rightarrow CH_3COCH_3 + HO_2 + H_2O$	2.6E-12*EXP(200./temp)	Atkinson et al. (2006)
G43070a	TrGC	$NPROPOL + OH \rightarrow C_2H_5CHO + HO_2 + H_2O$	4.6E-12*EXP(70./temp)*(k_s*f_soh/ (k_p+k_s*f_pch2oh+k_s*f_soh))	Atkinson et al. (2006), Sander et al. (2018)*

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G43070b	TrGC	NPROPOL + OH → HYPROPO2 + H_2O	4.6E-12*EXP(70./temp)*((k_p+k_s*f_pch2oh)/(k_p+k_s*f_pch2oh+k_s*f_soh))	Atkinson et al. (2006), Sander et al. (2018)*
G43071a	TrGC	CH_2CHCH_2OH + OH → HCOOH + OH + CH_3CHO	k_CH2CHOH_OH_HCOOH	Sander et al. (2018), So et al. (2014)*
G43072	TrGC	CH_2CHCH_2OH + HCOOH → C_2H_5CHO + HCOOH	k_CH2CHOH_HCOOH	Sander et al. (2018), da Silva (2010)*
G43073	TrGC	C_2H_5CHO + HCOOH → CH_2CHCH_2OH + HCOOH	k_ALD_HCOOH	Sander et al. (2018), da Silva (2010)*
G43074	TrGC	$HCOCOCH_2OOH$ + OH → HCOCO + CO + HO_2 + OH	k_s*f_sooh*f_co+k_roohro	Sander et al. (2018)*
G43202	TrGTerC	HCOCH2CHO + OH → HCOCH2CO3	4.29E-11	Rickard and Pascoe (2009)
G43203	TrGTerCN	HCOCH2CHO + NO_3 → HCOCH2CO3 + HNO_3	2.*KNO3AL*2.4	Rickard and Pascoe (2009)
G43204a	TrGTerC	HCOCH2CO3 → $HCOCH_2O_2$ + CO_2	k1_RO2RCO3*0.9	Sander et al. (2018)
G43204b	TrGTerC	HCOCH2CO3 → HCOCH2CO2H	k1_RO2RCO3*0.1	Sander et al. (2018)
G43205	TrGTerCN	HCOCH2CO3 + NO → $HCOCH_2O_2$ + CO_2 + NO_2	KAPNO	Rickard and Pascoe (2009)
G43206	TrGTerCN	HCOCH2CO3 + NO_2 → C_3PAN2	k_CH3CO3_NO2	Rickard and Pascoe (2009)
G43207a	TrGTerC	HCOCH2CO3 + HO_2 → HCOCH2CO3H	KAPHO2*rco3_ooh	Rickard and Pascoe (2009)
G43207b	TrGTerC	HCOCH2CO3 + HO_2 → HCOCH2CO2H + O_3	KAPHO2*rco3_o3	Rickard and Pascoe (2009)
G43207c	TrGTerC	HCOCH2CO3 + HO_2 → $HCOCH_2O_2$ + CO_2 + OH	KAPHO2*rco3_oh	Rickard and Pascoe (2009)
G43210	TrGTerCN	C_3PAN2 → HCOCH2CO3 + NO_2	k_PAN_M	Rickard and Pascoe (2009)
G43211	TrGTerCN	C_3PAN2 + OH → GLYOX + CO + NO_2	2.10E-11	Rickard and Pascoe (2009)
G43212	TrGTerC	HCOCH2CO2H + OH → $HCOCH_2O_2$ + CO_2	2.14E-11	Rickard and Pascoe (2009)
G43213a	TrGTerC	$HOC_2H_4CO_3$ → $HOCH_2CH_2O_2$ + CO_2	k1_RO2RCO3*0.9	Sander et al. (2018)
G43213b	TrGTerC	$HOC_2H_4CO_3$ → HOC2H4CO2H	k1_RO2RCO3*0.1	Sander et al. (2018)
G43214	TrGTerCN	$HOC_2H_4CO_3$ + NO → $HOCH_2CH_2O_2$ + CO_2 + NO_2	KAPNO	Rickard and Pascoe (2009)
G43215a	TrGTerC	$HOC_2H_4CO_3$ + HO_2 → HOC2H4CO3H	KAPHO2*rco3_ooh	Rickard and Pascoe (2009)
G43215b	TrGTerC	$HOC_2H_4CO_3$ + HO_2 → $HOCH_2CH_2O_2$ + CO_2 + OH	KAPHO2*rco3_oh	Rickard and Pascoe (2009)
G43215c	TrGTerC	$HOC_2H_4CO_3$ + HO_2 → HOC2H4CO2H + O_3	KAPHO2*rco3_o3	Rickard and Pascoe (2009)
G43218	TrGTerCN	$HOC_2H_4CO_3$ + NO_2 → C_3PAN1	k_CH3CO3_NO2	Rickard and Pascoe (2009)
G43219	TrGTerC	HOC2H4CO2H + OH → $HOCH_2CH_2O_2$ + CO_2	1.39E-11	Rickard and Pascoe (2009)
G43220	TrGTerC	HOC2H4CO3H + OH → $HOC_2H_4CO_3$	1.73E-11	Rickard and Pascoe (2009)
G43221	TrGTerCN	C_3PAN1 → $HOC_2H_4CO_3$ + NO_2	k_PAN_M	Rickard and Pascoe (2009)
G43222	TrGTerCN	C_3PAN1 + OH → $HOCH_2CHO$ + CO + NO_2	4.51E-12	Rickard and Pascoe (2009)
G43223	TrGTerC	HCOCH2CO3H + OH → $HCOCH_2O_2$ + CO_2 + H_2O	2.49E-11	Rickard and Pascoe (2009)*
G43415	TrGAroC	C3DIALOOH + OH → HCOCOCHO + OH	1.44E-10	Rickard and Pascoe (2009)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G43418a	TrGAroC	C3DIALO2 + HO_2 → C3DIALOOH	KRO2HO2(3)*(rco3_ooh+rco3_o3)	Rickard and Pascoe (2009)
G43418b	TrGAroC	C3DIALO2 + HO_2 → GLYOX + CO + HO_2 + OH	KRO2HO2(3)*rco3_oh	Rickard and Pascoe (2009)
G43419	TrGAroCN	C3DIALO2 + NO → GLYOX + CO + HO_2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G43420	TrGAroCN	C3DIALO2 + NO_3 → GLYOX + CO + HO_2 + NO_2	KRO2NO3	Rickard and Pascoe (2009)*
G43421	TrGAroC	C3DIALO2 → GLYOX + CO + HO_2	k1_RO2sORO2	Rickard and Pascoe (2009)*
G43422a	TrGAroC	HCOCOHCO3 + HO_2 → GLYOX + CO_2 + HO_2 + OH	KAPHO2*rco3_oh	Rickard and Pascoe (2009)
G43422b	TrGAroC	HCOCOHCO3 + HO_2 → HCOCOHCO3H	KAPHO2*(rco3_ooh+rco3_o3)	Rickard and Pascoe (2009)
G43424	TrGAroCN	HCOCOHCO3 + NO → GLYOX + CO_2 + HO_2 + NO_2	KAPNO	Rickard and Pascoe (2009)
G43425	TrGAroCN	HCOCOHCO3 + NO_2 → HCOCOHPAN	k_CH3CO3_NO2	Rickard and Pascoe (2009)
G43426	TrGAroCN	HCOCOHCO3 + NO_3 → GLYOX + CO_2 + HO_2 + NO_2	KRO2NO3*1.74	Rickard and Pascoe (2009)
G43427	TrGAroC	HCOCOHCO3 → GLYOX + CO_2 + HO_2	k1_RO2RCO3	Rickard and Pascoe (2009)
G43428	TrGAroC	METACETHO + OH → $CH_3C(O)$ + CO_2	9.82E-11	Rickard and Pascoe (2009)
G43442	TrGAroCN	HCOCOHPAN + OH → GLYOX + CO + NO_2	6.97E-11	Rickard and Pascoe (2009)
G43443	TrGAroCN	HCOCOHPAN → HCOCOHCO3 + NO_2	k_PAN_M	Rickard and Pascoe (2009)
G43444	TrGAroC	C32OH13CO + OH → HCOCOHCO3	1.36E-10	Rickard and Pascoe (2009)
G43446	TrGAroC	HCOCOHCO3H + OH → HCOCOHCO3	7.33E-11	Rickard and Pascoe (2009)
G44000	TrGC	C_4H_{10} + OH → $LC_4H_9O_2$ + H_2O	2.03E-17*temp*temp*EXP(78./temp)	Atkinson et al. (2006)*
G44001a	TrGC	$LC_4H_9O_2$ → C_3H_7CHO + HO_2	(k1_RO2pRO2*0.1273+k1_RO2sRO2*0.8727)*0.1273	Rickard and Pascoe (2009), Sander et al. (2018)
G44001b	TrGC	$LC_4H_9O_2$ → .636 MEK + .636 HO_2 + .364 CH_3CHO + .364 $C_2H_5O_2$	(k1_RO2pRO2*0.1273+k1_RO2sRO2*0.8727)*0.8727	Rickard and Pascoe (2009), Sander et al. (2018)*
G44002	TrGC	$LC_4H_9O_2$ + HO_2 → LC_4H_9OOH	KRO2HO2(4)	Rickard and Pascoe (2009)
G44003a	TrGCN	$LC_4H_9O_2$ + NO → NO_2 + C_3H_7CHO + HO_2	KRO2NO*(1.-(0.1273*alpha_AN(4,1,0,0,0,temp,cair)+0.8727*alpha_AN(4,2,0,0,0,temp,cair)))*0.1273	Rickard and Pascoe (2009), Sander et al. (2018)
G44003b	TrGCN	$LC_4H_9O_2$ + NO → NO_2 + .636 MEK + .636 HO_2 + .364 CH_3CHO + .364 $C_2H_5O_2$	KRO2NO*(1.-(0.1273*alpha_AN(4,1,0,0,0,temp,cair)+0.8727*alpha_AN(4,2,0,0,0,temp,cair)))*0.8727	Rickard and Pascoe (2009), Sander et al. (2018)
G44003c	TrGCN	$LC_4H_9O_2$ + NO → LC4H9NO3	KRO2NO*(0.1273*alpha_AN(4,1,0,0,0,temp,cair)+0.8727*alpha_AN(4,2,0,0,0,temp,cair))	Rickard and Pascoe (2009)*
G44004a	TrGCN	$LC_4H_9O_2$ + NO_3 → NO_2 + C_3H_7CHO + HO_2	KRO2NO3*0.1273	Rickard and Pascoe (2009), Sander et al. (2018)
G44004b	TrGCN	$LC_4H_9O_2$ + NO_3 → NO_2 + .636 MEK + .636 HO_2 + .364 CH_3CHO + .364 $C_2H_5O_2$	KRO2NO3*0.8727	Rickard and Pascoe (2009), Sander et al. (2018)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G44005a	TrGC	$LC_4H_9OOH + OH \rightarrow LC_4H_9O_2 + H_2O$	`k_roohro`	Sander et al. (2018)
G44005b	TrGC	$LC_4H_9OOH + OH \rightarrow C_3H_7CHO + H_2O + OH$	`k_s*f_tooh*f_alk*(k_p/(k_p+k_s))`	Sander et al. (2018)
G44005c	TrGC	$LC_4H_9OOH + OH \rightarrow MEK + H_2O + OH$	`k_t*f_tooh*f_alk*(k_s/(k_p+k_s))`	Sander et al. (2018)
G44006a	TrGC	$iC_4H_{10} + OH \rightarrow TC_4H_9O_2 + H_2O$	`1.17E-17*temp*temp*EXP(213./temp) *k_t/(3.*k_p+k_t)`	Atkinson (2003)
G44006b	TrGC	$iC_4H_{10} + OH \rightarrow IC_4H_9O_2 + H_2O$	`1.17E-17*temp*temp*EXP(213./temp) *3.*k_p/(3.*k_p+k_t)`	Atkinson (2003)
G44007	TrGC	$TC_4H_9O_2 \rightarrow CH_3COCH_3 + CH_3$	`k1_RO2tRO2`	Rickard and Pascoe (2009), Sander et al. (2018)
G44008	TrGC	$TC_4H_9O_2 + HO_2 \rightarrow TC_4H_9OOH$	`KRO2HO2(4)`	Rickard and Pascoe (2009)
G44009a	TrGCN	$TC_4H_9O_2 + NO \rightarrow NO_2 + CH_3COCH_3 + CH_3$	`KRO2NO*(1.-alpha_AN(4,3,0,0,0, temp,cair))`	Rickard and Pascoe (2009), Sander et al. (2018)
G44009b	TrGCN	$TC_4H_9O_2 + NO \rightarrow$ TC4H9NO3	`KRO2NO*alpha_AN(4,3,0,0,0,temp, cair)`	Rickard and Pascoe (2009)
G44010a	TrGC	$TC_4H_9OOH + OH \rightarrow TC_4H_9O_2 + H_2O$	`k_roohro`	Sander et al. (2018)
G44010b	TrGC	$TC_4H_9OOH + OH \rightarrow CH_3COCH_3 + HCHO + OH + H_2O$	`3.*k_p*f_tch2oh`	Sander et al. (2018)*
G44011	TrGCN	TC4H9NO3 $+ OH \rightarrow CH_3COCH_3 + HCHO + NO_2 + H_2O$	`3.*k_p*f_ch2ono2`	Sander et al. (2018)*
G44012	TrGC	$IC_4H_9O_2 \rightarrow$ IPRCHO	`k1_RO2sRO2`	Rickard and Pascoe (2009), Sander et al. (2018)
G44013	TrGC	$IC_4H_9O_2 + HO_2 \rightarrow IC_4H_9OOH$	`KRO2HO2(4)`	Rickard and Pascoe (2009)
G44014a	TrGCN	$IC_4H_9O_2 + NO \rightarrow NO_2 +$ IPRCHO	`KRO2NO*(1.-alpha_AN(4,2,0,0,0, temp,cair))`	Rickard and Pascoe (2009), Sander et al. (2018)
G44014b	TrGCN	$IC_4H_9O_2 + NO \rightarrow$ IC4H9NO3	`KRO2NO*alpha_AN(4,2,0,0,0,temp, cair)`	Rickard and Pascoe (2009)
G44015a	TrGC	$IC_4H_9OOH + OH \rightarrow IC_4H_9O_2 + H_2O$	`k_roohro`	Sander et al. (2018)
G44015b	TrGC	$IC_4H_9OOH + OH \rightarrow$ IPRCHO $+ OH + H_2O$	`k_s*f_sooh+2.*k_s+k_t*f_pch2oh`	Sander et al. (2018)*
G44016	TrGCN	IC4H9NO3 $+ OH \rightarrow$ IPRCHO $+ NO_2 + H_2O$	`k_s*f_ono2+2.*k_p+k_t*f_ch2ono2`	Sander et al. (2018)*
G44017	TrGC	$MVK + O_3 \rightarrow .87\ MGLYOX + .5481\ CO + .1392\ HO_2 + .1392\ OH + .3219\ CH_2OO + .13\ HCHO + .04680\ OH + .04680\ CO + .07280\ CH_3C(O) + .026\ CH_3CHO + .026\ CO_2 + .026\ HCHO + .026\ HO_2 + .02402\ MGLYOX + .02402\ H_2O_2 + .00718\ CH_3COCO_2H$	`8.5E-16*EXP(-1520./temp)`	Sander et al. (2018)
G44018	TrGC	$MVK + OH \rightarrow$ LHMVKABO2	`2.6E-12*EXP(610./temp)`	Sander et al. (2018), Atkinson et al. (2006)*

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G44019	TrGC	MEK + OH → LMEKO2 + H_2O	1.5E-12*EXP(-90./temp)	Atkinson et al. (2006), Sander et al. (2018)*
G44020	TrGC	LMEKO2 + HO_2 → LMEKOOH	KRO2HO2(4)	Sander et al. (2018)
G44021a	TrGCN	LMEKO2 + NO → .62 CH_3CHO + .62 $CH_3C(O)$ + .38 HCHO + .38 CO_2 + .38 $HOCH_2CH_2O_2$ + NO_2	KRO2NO*(1.-(.62*alpha_AN(4,2,1,0,0,temp,cair)+.38*alpha_AN(4,1,0,1,0,temp,cair)))	Sander et al. (2018)*
G44021b	TrGCN	LMEKO2 + NO → LMEKNO3	KRO2NO*(.62*alpha_AN(4,2,1,0,0,temp,cair)+.38*alpha_AN(4,1,0,1,0,temp,cair))	Sander et al. (2018)
G44022a	TrGC	LMEKOOH + OH → LMEKO2 + H_2O	k_roohro	Sander et al. (2018)
G44022b	TrGC	LMEKOOH + OH → .62 BIACET + .38 HCHO + .38 CO_2 + .38 $HOCH_2CH_2O_2$ + H_2O + OH	(.62*k_t*f_tooh*f_co+.38*k_s*f_sooh)	Sander et al. (2018)
G44023a	TrGCN	LC4H9NO3 + OH → MEK + NO_2 + H_2O	(k_t*f_ono2*f_alk+k_p*f_alk+k_s*f_ch2ono2+k_p)*(k_s/(k_p+k_s))	Sander et al. (2018)*
G44023b	TrGCN	LC4H9NO3 + OH → C_3H_7CHO + NO_2 + H_2O	(k_p+k_s*(1+f_ch2ono2+f_ono2)*f_alk)*(k_p/(k_p+k_s))	Sander et al. (2018)*
G44024	TrGCN	MPAN + OH → CH_3COCH_2OH + CO + NO_2	3.2E-11	Orlando et al. (2002)
G44025	TrGCN	MPAN → MACO3 + NO_2	k_PAN_M	see note*
G44026	TrGC	LMEKO2 → .538 HCHO + .538 CO_2 + .459 $HOCH_2CH_2O_2$ + .079 $C_2H_5O_2$ + .462 $CH_3C(O)$ + .462 CH_3CHO	(.62*k1_RO2sOR02+.38*k1_RO2pOR02)	Rickard and Pascoe (2009)*
G44027	TrGC	MACR + OH → .45 MACO3 + .55 MACRO2	8.E-12*EXP(380./temp)	Orlando et al. (1999b), Sander et al. (2018)
G44028	TrGC	MACR + O_3 → .5481 CO + .1392 HO_2 + .1392 OH + .3219 CH_2OO + .87 MGLYOX + .13 HCHO + .13 OH + .065 $HCOCOCH_2O_2$ + .065 CO + .065 $CH_3C(O)$	1.36E-15*EXP(-2112./temp)	Sander et al. (2018)
G44029	TrGCN	MACR + NO_3 → MACO3 + HNO_3	KNO3AL*2.0	Rickard and Pascoe (2009)
G44030a	TrGC	MACO3 → $CH_3C(O)$ + HCHO + CO_2	k1_RO2RCO3*0.9	Sander et al. (2018)
G44030b	TrGC	MACO3 → MACO2H	k1_RO2RCO3*0.1	Sander et al. (2018)
G44031a	TrGC	MACO3 + HO_2 → MACO2 + OH	KAPHO2*rco3_oh	Sander et al. (2018)
G44031b	TrGC	MACO3 + HO_2 → MACO3H	KAPHO2*rco3_ooh	Sander et al. (2018)
G44031c	TrGC	MACO3 + HO_2 → MACO2H + O_3	KAPHO2*rco3_o3	Sander et al. (2018)
G44032	TrGCN	MACO3 + NO → MACO2 + NO_2	8.70E-12*EXP(290./temp)	Sander et al. (2018)
G44033	TrGCN	MACO3 + NO_2 → MPAN	k_CH3CO3_NO2	Rickard and Pascoe (2009)
G44034	TrGCN	MACO3 + NO_3 → MACO2 + NO_2	KRO2NO3*1.74	Sander et al. (2018)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G44035	TrGC	MACRO2 $\rightarrow$ .7 CH_3COCH_2OH + .7 HCHO + .7 HO_2 + .3 MACROH	k1_RO2tORO2	Rickard and Pascoe (2009)*
G44036a	TrGC	MACRO2 + HO_2 $\rightarrow$ MACRO + OH	KRO2HO2(4)*rcoch2o2_oh	Sander et al. (2018)
G44036b	TrGC	MACRO2 + HO_2 $\rightarrow$ MACROOH	KRO2HO2(4)*rcoch2o2_ooh	Sander et al. (2018)
G44037a	TrGCN	MACRO2 + NO $\rightarrow$ MACRO + NO_2	KRO2NO*(1.-alpha_AN(6,3,1,0,0, temp,cair))	Sander et al. (2018)
G44037b	TrGCN	MACRO2 + NO $\rightarrow$ MACRNO3	KRO2NO*alpha_AN(6,3,1,0,0,temp, cair)	Sander et al. (2018)
G44038	TrGCN	MACRO2 + NO_3 $\rightarrow$ MACRO + NO_2	KRO2NO3	Sander et al. (2018)
G44039a	TrGC	MACROOH + OH $\rightarrow$ MACRO2	k_roohro	Sander et al. (2018)
G44039b	TrGC	MACROOH + OH $\rightarrow$ CO + CH_3COCH_2OH + OH	k_t*f_o*f_tch2oh*f_alk	Sander et al. (2018)
G44039c	TrGC	MACROOH + OH $\rightarrow$ CO + MGLYOX + HO_2	(k_s*f_soh*f_pch2oh + k_rohro)	Sander et al. (2018)
G44040	TrGC	MACROH + OH $\rightarrow$ CH_3COCH_2OH + CO + HO_2	k_t*f_o*f_tch2oh*f_alk	Sander et al. (2018)
G44041	TrGC	MACRO $\rightarrow$ .885 CH_3COCH_2OH + .885 CO + .115 MGLYOX + .115 HCHO + HO_2	KDEC	Sander et al. (2018)
G44042	TrGC	MACO2H + OH $\rightarrow$ CH_3COCH_2OH + HO_2 + CO_2	((k_adt+k_adp)*a_co2h+k_co2h)	Sander et al. (2018)
G44043a	TrGC	MACO3H + OH $\rightarrow$ CH_3COCH_2OH + CO_2 + OH	(k_adt+k_adp)*a_co2h	Sander et al. (2018)
G44043b	TrGC	MACO3H + OH $\rightarrow$ MACO3	k_roohro	Sander et al. (2018)
G44044	TrGC	LHMVKABO2 $\rightarrow$ .024 CO2H3CHO + .072 MGLYOX + .072 HO_2 + .072 HCHO + .5280 $CH_3C(O)$ + .5280 $HOCH_2CHO$ + .176 BIACETOH + .2 HO12CO3C4	(.12*k1_RO2pORO2+.88*k1_RO2sORO2)	Sander et al. (2018)
G44045a	TrGC	LHMVKABO2 + HO_2 $\rightarrow$ OH + $HOCH_2CHO$ + $CH_3C(O)$	KRO2HO2(4)*.88*rcoch2o2_oh	Sander et al. (2018)
G44045b	TrGC	LHMVKABO2 + HO_2 $\rightarrow$ LHMVKABOOH	KRO2HO2(4)*(.12+.88*rcoch2o2_ooh)	Sander et al. (2018)
G44046a	TrGCN	LHMVKABO2 + NO $\rightarrow$ .12 MGLYOX + .12 HO_2 + .88 $HOCH_2CHO$ + .88 $CH_3C(O)$ + .12 HCHO + NO_2	KRO2NO*(1.-(.12*alpha_AN(6,1,0, 1,0,temp,cair)+.88*alpha_AN(6,2, 1,0,0,temp,cair)))	Sander et al. (2018)
G44046b	TrGCN	LHMVKABO2 + NO $\rightarrow$ MVKNO3	KRO2NO*(.12*alpha_AN(6,1,0,1,0, temp,cair)+.88*alpha_AN(6,2,1,0, 0,temp,cair))	Sander et al. (2018)*
G44047	TrGCN	LHMVKABO2 + NO_3 $\rightarrow$ .12 MGLYOX + .12 HO_2 + .88 $HOCH_2CHO$ + .88 $CH_3C(O)$ + .12 HCHO + .12 HO_2 + NO_2	KRO2NO3	Sander et al. (2018)
G44048a	TrGC	LHMVKABOOH + OH $\rightarrow$ LHMVKABO2	k_roohro	Sander et al. (2018)
G44048b	TrGC	LHMVKABOOH + OH $\rightarrow$ .12 CO2H3CHO + .88 BIACETOH + OH	(.12*k_s*f_sooh*f_pch2oh+.88*k_ t*f_tooh*f_pch2oh*f_co)	Sander et al. (2018)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G44049a	TrGC	CO2H3CHO + OH → CO2H3CO3	k_t*f_o*f_alk	Sander et al. (2018)
G44049b	TrGC	CO2H3CHO + OH → CH_3COCOCHO + HO_2 + H_2O	k_t*f_co*f_toh*f_cho	Sander et al. (2018)
G44050	TrGCN	CO2H3CHO + NO_3 → CO2H3CO3 + HNO_3	KNO3AL*4.0	Rickard and Pascoe (2009)
G44051	TrGC	CO2H3CO3 → MGLYOX + HO_2 + CO_2	k1_RO2RCO3	Sander et al. (2018)
G44052a	TrGC	CO2H3CO3 + HO_2 → OH + MGLYOX + HO_2 + CO_2	KAPHO2*rco3_oh	Sander et al. (2018)
G44052b	TrGC	CO2H3CO3 + HO_2 → CO2H3CO2H + O_3	KAPHO2*rco3_o3	Sander et al. (2018)
G44052c	TrGC	CO2H3CO3 + HO_2 → CO2H3CO3H	KAPHO2*rco3_ooh	Sander et al. (2018)
G44053	TrGCN	CO2H3CO3 + NO → MGLYOX + HO_2 + NO_2 + CO_2	KAPNO	Sander et al. (2018)
G44054	TrGCN	CO2H3CO3 + NO_3 → MGLYOX + HO_2 + NO_2 + CO_2	KRO2NO3*1.74	Sander et al. (2018)
G44055a	TrGC	CO2H3CO3H + OH → CO2H3CO3	k_roohro	Sander et al. (2018)
G44055b	TrGC	CO2H3CO3H + OH → $CH_3C(O)$ + CO + CO_2 + OH	(k_t*f_co2h*f_co*f_toh)	Sander et al. (2018)
G44056	TrGC	CO2H3CO2H + OH → CH3COCOCO2H + HO_2	k_t*f_co2h*f_co*f_toh+k_co2h	Sander et al. (2018)
G44057a	TrGC	HO12CO3C4 + OH → BIACETOH + HO_2	k_t*f_toh*f_alk*f_co	Sander et al. (2018)
G44057b	TrGC	HO12CO3C4 + OH → CO2H3CHO + HO_2	k_s*f_soh*f_alk	Sander et al. (2018)
G44058	TrGC	MACO2 → .65 CH_3 + .65 CO + .65 HCHO + .35 OH + .35 $CH_3COCH_2O_2$ + CO_2	KDEC	Sander et al. (2018)
G44059	TrGC	LHMVKABO2 → .88 MGLYOX + .88 HCHO + .12 HOOCH2CHO + .12 $CH_3C(O)$ + OH	KHSD	Sander et al. (2018)
G44060	TrGC	MACRO2 → MGLYOX + HCHO + OH	KHSB	Sander et al. (2018)
G44061a	TrGCN	MVKNO3 + OH → MGLYOX + CO_2 + HO_2 + NO_2 + H_2O	k_s*f_sooh*f_ch2ono2+k_rohro	Sander et al. (2018)*
G44061b	TrGCN	MVKNO3 + OH → BIACETOH + NO_2 + H_2O	k_t*f_ono2*f_co*f_pch2oh	Sander et al. (2018)*
G44062a	TrGCN	MACRNO3 + OH → CH_3COCH_2OH + CO_2 + NO_2 + H_2O	k_t*f_o*f_ch2ono2	Sander et al. (2018)*
G44062b	TrGCN	MACRNO3 + OH → MGLYOX + CO + NO_2 + H_2O	k_rohro+k_s*f_sooh*f_ch2ono2	Sander et al. (2018)*
G44063	TrGC	MACRO2 → CH_3COCH_2OH + OH + CO	K14HSAL	Sander et al. (2018)
G44064	TrGC	EZCH3CO2CHCHO → .9 CH_3COCHCO + .1 $CH_3C(O)$ + .01 GLYOX + .18 CO + .09 HO_2 + OH	K15HS24VYNAL	Sander et al. (2018)
G44065	TrGC	EZCH3CO2CHCHO + HO_2→ CH_3COOHCHCHO	KRO2HO2(4)	Sander et al. (2018)
G44066	TrGCN	EZCH3CO2CHCHO + NO → CH_3COCHO_2CHO + NO_2	KRO2NO	Sander et al. (2018)*
G44067	TrGCN	EZCH3CO2CHCHO + NO_3→ CH_3COCHO_2CHO + NO_2	kRO2NO3	Sander et al. (2018)
G44068	TrGC	EZCH3CO2CHCHO → CH_3COCHO_2CHO	k1_RO2sORO2	Sander et al. (2018)
G44069	TrGC	EZCHOCCH3CHO2 → HCOCCH_3CO + OH	K15HS24VYNAL	Sander et al. (2018)
G44070	TrGCN	EZCHOCCH3CHO2 + NO → $HCOCO_2CH_3CHO$ + NO_2	KRO2NO	Sander et al. (2018)*
G44071	TrGC	EZCHOCCH3CHO2 + HO_2→ HCOCCH_3CHOOH	KRO2HO2(4)	Sander et al. (2018)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G44072	TrGCN	EZCHOCCH3CHO2 + NO_3→ $HCOCO_2CH_3CHO$ + NO_2	KRO2NO3	Sander et al. (2018)
G44073	TrGC	EZCHOCCH3CHO2 → $HCOCO_2CH_3CHO$	k1_RO2pORO2	Sander et al. (2018)
G44074	TrGC	$CH_3COOHCHCHO$ → CH_3COCHO_2CHO + OH	KHYDEC	Sander et al. (2018)
G44075	TrGC	$HCOCCH_3CHOOH$ → $HCOCO_2CH_3CHO$ + OH	KHYDEC	Sander et al. (2018)
G44076	TrGCN	CH_3COCHO_2CHO + NO → $CH_3C(O)$ + GLYOX + NO_2	KRO2NO	Sander et al. (2018)*
G44077	TrGCN	CH_3COCHO_2CHO + NO_3 → $CH_3C(O)$ + GLYOX + NO_2	KRO2NO3	Sander et al. (2018)
G44078	TrGC	CH_3COCHO_2CHO + HO_2 → $CH_3C(O)$ + GLYOX + OH	KRO2HO2(4)	Sander et al. (2018)*
G44079	TrGC	CH_3COCHO_2CHO → $CH_3C(O)$ + GLYOX	k1_RO2sORO2	Sander et al. (2018)
G44080	TrGC	$HCOCO_2CH_3CHO$ → MGLYOX + CO + HO_2	k1_RO2tORO2	Sander et al. (2018)
G44081	TrGCN	$HCOCO_2CH_3CHO$ + NO → MGLYOX + CO + HO_2 + NO_2	KRO2NO	Sander et al. (2018)*
G44082	TrGC	$HCOCO_2CH_3CHO$ + HO_2 → MGLYOX + CO + HO_2 + OH	KRO2HO2(4)	Sander et al. (2018)*
G44083	TrGCN	$HCOCO_2CH_3CHO$ + NO_3 → MGLYOX + CO + HO_2 + NO_2	KRO2NO3	Sander et al. (2018)
G44084	TrGC	$HCOCCH_3CO$ + OH → CO + MGLYOX + HO_2	1E-10*a_cho	Hatakeyama et al. (1985), Sander et al. (2018)
G44085	TrGC	$CH_3COCHCO$ + OH → CO + MGLYOX + HO_2	7.6E-11*a_coch3	Hatakeyama et al. (1985), Sander et al. (2018)*
G44086	TrGCN	LMEKNO3 + OH → .62 MGLYOX + .62 HCHO + .62 HO_2 + .62 NO_2 + .38 $CH_3C(O)$ + .38 NO_3CH2CHO	.62*(k_p*(f_co+f_ch2ono2)) +.38*(k_s*f_ch2ono2*f_co)	Sander et al. (2018)*
G44087	TrGC	MEPROPENE + OH → IBUTOLBO2	9.4E-12*EXP(505./temp)	Atkinson et al. (2006)
G44088a	TrGC	MEPROPENE + O_3 → CH_3COCH_3 + CH_2OO^*	2.7E-15*EXP(-1630./temp)*0.33	Atkinson et al. (2006), Sander et al. (2018)
G44088b	TrGC	MEPROPENE + O_3 → $CH_3COCH_2O_2$ + OH + HCHO	2.7E-15*EXP(-1630./temp)*0.67	Atkinson et al. (2006), Sander et al. (2018)
G44089	TrGCN	MEPROPENE + NO_3 → CH_3COCH_3 + HCHO + NO_2	3.4E-13	Atkinson et al. (2006), Sander et al. (2018)*
G44090	TrGC	IBUTOLBO2 → CH_3COCH_3 + HCHO + HO_2	k1_RO2tORO2	Sander et al. (2018)
G44091a	TrGC	IBUTOLBO2 + HO_2 → IBUTOLBOOH	KRO2HO2(4)*rcoch2o2_ooh	Sander et al. (2018)
G44091b	TrGC	IBUTOLBO2 + HO_2 → CH_3COCH_3 + HCHO + HO_2 + OH	KRO2HO2(4)*rcoch2o2_oh	Sander et al. (2018)
G44092a	TrGCN	IBUTOLBO2 + NO → CH_3COCH_3 + HCHO + HO_2 + NO_2	KRO2NO*(1.-alpha_AN(5,3,0,0,0, temp,cair))	Sander et al. (2018)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G44092b	TrGCN	IBUTOLBO2 + NO → IBUTOLBNO3	KRO2NO*alpha_AN(5,3,0,0,0,temp, cair)	Sander et al. (2018)
G44093	TrGCN	IBUTOLBO2 + NO_3 → CH_3COCH_3 + HCHO + HO_2 + NO_2	KRO2NO3	Sander et al. (2018)
G44094a	TrGC	IBUTOLBOOH + OH → IBUTOLBO2	k_roohro	Sander et al. (2018)
G44094b	TrGC	IBUTOLBOOH + OH → CH_3COCH_3 + HCHO + HO_2	k_s*f_sooh*f_pch2oh	Sander et al. (2018)
G44095	TrGCN	IBUTOLBNO3 + OH → CH_3COCH_3 + HCHO + HO_2 + NO_2	3.*k_p	Sander et al. (2018)
G44096	TrGC	BUT1ENE + OH → LBUT1ENO2	6.6E-12*EXP(465./temp)	Atkinson et al. (2006)*
G44097a	TrGC	BUT1ENE + O_3 → HCHO + .5 C_2H_5CHO + .5 H_2O_2 + .5 CH_3CHO + .5 CO + .5 HO_2	3.35E-15*EXP(-1745./temp)*.57	Atkinson et al. (2006), Sander et al. (2018)*
G44097b	TrGC	BUT1ENE + O_3 → C_2H_5CHO + CH_2OO^*	3.35E-15*EXP(-1745./temp)*.43	Atkinson et al. (2006), Sander et al. (2018)*
G44098	TrGCN	BUT1ENE + NO_3 → C_2H_5CHO + HCHO + NO_2	3.2E-13*EXP(-950./temp)	Atkinson et al. (2006), Sander et al. (2018)*
G44099	TrGC	LBUT1ENO2 → C_2H_5CHO + HCHO + HO_2	k1_RO2sORO2	Sander et al. (2018)
G44100a	TrGC	LBUT1ENO2 + HO_2 → LBUT1ENOOH	KRO2HO2(4)*rcoch2o2_ooh	Sander et al. (2018)
G44100b	TrGC	LBUT1ENO2 + HO_2 → C_2H_5CHO + HCHO + HO_2 + OH	KRO2HO2(4)*rcoch2o2_oh	Sander et al. (2018)
G44101a	TrGCN	LBUT1ENO2 + NO → C_2H_5CHO + HCHO + HO_2 + NO_2	KRO2NO*(1.-alpha_AN(5,2,0,0,0, temp,cair))	Sander et al. (2018)
G44101b	TrGCN	LBUT1ENO2 + NO → LBUT1ENNO3	KRO2NO*alpha_AN(5,2,0,0,0,temp, cair)	Sander et al. (2018)
G44102	TrGCN	LBUT1ENO2 + NO_3 → C_2H_5CHO + HCHO + HO_2 + NO_2	KRO2NO3	Sander et al. (2018)
G44103a	TrGC	LBUT1ENOOH + OH → LBUT1ENO2	k_roohro	Sander et al. (2018)
G44103b	TrGC	LBUT1ENOOH + OH → $C_2H_5CO_3$ + HCHO + HO_2	k_t*f_tooh*f_pch2oh	Sander et al. (2018)*
G44104	TrGCN	LBUT1ENNO3 + OH → C_2H_5CHO + CO + HO_2 + NO_2	k_s*f_soh*f_ch2ono2	Sander et al. (2018)*
G44105	TrGC	CBUT2ENE + OH → BUT2OLO2	1.1E-11*EXP(485./temp)	Atkinson et al. (2006)
G44106	TrGC	CBUT2ENE + O_3 → CH_3CHO + .16 CH3CHOHOOH + .50 OH + .50 $HCOCH_2O_2$ + .05 CH2CO + .09 CH_3OH + .09 CO + .2 CH_4 + .2 CO_2	3.2E-15*EXP(-965./temp)	Atkinson et al. (2006), Sander et al. (2018)*
G44107	TrGCN	CBUT2ENE + NO_3 → 2 CH_3CHO + NO_2	3.5E-13	Atkinson et al. (2006), Sander et al. (2018)*
G44108	TrGC	TBUT2ENE + OH → BUT2OLO2	1.0E-11*EXP(553./temp)	Atkinson et al. (2006)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G44109	TrGC	TBUT2ENE + O_3 → CH_3CHO + .16 CH3CHOHOOH + .50 OH + .50 $HCOCH_2O_2$ + .05 CH2CO + .09 CH_3OH + .09 CO + .2 CH_4 + .2 CO_2	6.6E-15*EXP(-1060./temp)	Atkinson et al. (2006), Sander et al. (2018)
G44110	TrGCN	TBUT2ENE + NO_3 → 2 CH_3CHO + NO_2	1.78E-12*EXP(-530./temp) +1.28E-14*EXP(570./temp)	Atkinson et al. (2006), Sander et al. (2018)*
G44111	TrGC	BUT2OLO2 → C_2H_5CHO + HCHO + HO_2	k1_RO2sOR02	Sander et al. (2018)
G44112a	TrGC	BUT2OLO2 + HO_2 → BUT2OLOOH	KRO2HO2(4)*rcoch2o2_ooh	Sander et al. (2018)
G44112b	TrGC	BUT2OLO2 + HO_2 → 2 CH_3CHO + HO_2 + OH	KRO2HO2(4)*rcoch2o2_oh	Sander et al. (2018)
G44113a	TrGCN	BUT2OLO2 + NO → 2 CH_3CHO + HO_2 + NO_2	KRO2NO*(1.-alpha_AN(5,2,0,0,0, temp,cair))	Sander et al. (2018)
G44113b	TrGCN	BUT2OLO2 + NO → BUT2OLNO3	KRO2NO*alpha_AN(5,2,0,0,0,temp, cair)	Sander et al. (2018)
G44114	TrGCN	BUT2OLO2 + NO_3 → 2 CH_3CHO + HO_2 + NO_2	KRO2NO3	Sander et al. (2018)
G44115a	TrGC	BUT2OLOOH + OH → BUT2OLO2	k_roohro	Sander et al. (2018)
G44115b	TrGC	BUT2OLOOH + OH → LMEKOOH + HO_2	k_t*f_toh*f_pch2oh	Sander et al. (2018)
G44115c	TrGC	BUT2OLOOH + OH → BUT2OLO + OH	k_t*f_tooh*f_pch2oh	Sander et al. (2018)
G44116	TrGCN	BUT2OLNO3 + OH → LMEKNO3 + HO_2	k_t*f_toh*f_ch2ono2	Sander et al. (2018)
G44117	TrGC	BUT2OLO + OH → BIACET + HO_2	k_t*f_toh*f_co	Sander et al. (2018)
G44118	TrGC	IPRCHO + OH → IPRCO3 + H_2O	6.8E-12*EXP(410./temp)	Atkinson et al. (2006)
G44119	TrGCN	IPRCHO + NO_3 → IPRCO3 + HNO_3	1.67E-12*EXP(-1460./temp)	Atkinson et al. (2006)
G44120	TrGC	IPRCO3 → $iC_3H_7O_2$ + CO_2	k1_RO2RCO3	Rickard and Pascoe (2009)
G44121a	TrGC	IPRCO3 + HO_2 → PERIBUACID	KAPHO2*rco3_ooh	Rickard and Pascoe (2009), Sander et al. (2018)
G44121b	TrGC	IPRCO3 + HO_2 → $iC_3H_7O_2$ + CO_2 + OH	KAPHO2*(1-rco3_ooh)	Rickard and Pascoe (2009), Sander et al. (2018)
G44122	TrGCN	IPRCO3 + NO_2 → PIPN	k_CH3CO3_NO2	Rickard and Pascoe (2009)
G44123	TrGCN	IPRCO3 + NO → $iC_3H_7O_2$ + CO_2 + NO_2	KAPNO	Rickard and Pascoe (2009)
G44124a	TrGC	PERIBUACID + OH → IPRCO3 + H_2O	k_roohro	Rickard and Pascoe (2009)
G44124b	TrGC	PERIBUACID + OH → CH_3COCH_3 + H_2O + CO_2	k_s*f_co2h	Sander et al. (2018)*
G44125	TrGCN	PIPN → IPRCO3 + NO_2	k_PAN_M	Rickard and Pascoe (2009)
G44126	TrGCN	PIPN + OH → CH_3COCH_3 + CO_2 + NO_2	k_s*f_cpan	Sander et al. (2018)*
G44127	TrGC	MPROPENOL + OH → HCOOH + OH + CH_3COCH_3	k_CH2CHOH_OH_HCOOH	Sander et al. (2018), So et al. (2014)*
G44128	TrGC	MPROPENOL + HCOOH → IPRCHO + HCOOH	k_CH2CHOH_HCOOH	Sander et al. (2018), da Silva (2010)*

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G44129	TrGC	IPRCHO + HCOOH → MPROPENOL + HCOOH	k_ALD_HCOOH	Sander et al. (2018), da Silva (2010)*
G44130	TrGC	BUTENOL + OH → HCOOH + OH + C_2H_5CHO	k_CH2CHOH_OH_HCOOH	Sander et al. (2018), So et al. (2014)*
G44131	TrGC	BUTENOL + HCOOH → C_3H_7CHO + HCOOH	k_CH2CHOH_HCOOH	Sander et al. (2018), da Silva (2010)*
G44132	TrGC	C_3H_7CHO + HCOOH → BUTENOL + HCOOH	k_ALD_HCOOH	Sander et al. (2018), da Silva (2010)*
G44133	TrGC	HVMK + OH → HCOOH + OH + MGLYOX	8.8E-11	Sander et al. (2018), So et al. (2014), Messaadia et al. (2015)*
G44134	TrGC	HVMK + HCOOH → CO2C3CHO + HCOOH	k_CH2CHOH_HCOOH	Sander et al. (2018), da Silva (2010)*
G44135	TrGC	CO2C3CHO + HCOOH → HVMK + HCOOH	k_ALD_HCOOH	Sander et al. (2018), da Silva (2010)*
G44136	TrGC	HMAC + OH → HCOOH + OH + MGLYOX	8.8E-11	Sander et al. (2018), So et al. (2014), Messaadia et al. (2015)*
G44137	TrGC	HMAC + HCOOH → IBUTDIAL + HCOOH	k_CH2CHOH_HCOOH	Sander et al. (2018), da Silva (2010)*
G44138	TrGC	IBUTDIAL + HCOOH → HMAC + HCOOH	k_ALD_HCOOH	Sander et al. (2018), da Silva (2010)*
G44139	TrGC	CO2C3CHO + OH → $CH_3COCH_2O_2$ + CO_2 + H_2O	k_t*f_o*f_alk+k_s*f_cho*f_co	Sander et al. (2018)*
G44140	TrGCN	CO2C3CHO + NO_3 → $CH_3COCH_2O_2$ + CO_2 + HNO_3	KNO3AL*4.0	Sander et al. (2018)*
G44141	TrGC	IBUTDIAL + OH → CH_3CHO + CO + HO_2 + CO_2 + H_2O	2.*k_t*f_o*f_alk+k_t*f_cho*f_cho	Sander et al. (2018)*
G44142	TrGCN	IBUTDIAL + NO_3 → CH_3CHO + CO + HO_2 + CO_2 + HNO_3	2.*KNO3AL*4.0	Sander et al. (2018)*
G44200	TrGTerC	$CH_3COCOCH_2O_2$ → $CH_3C(O)$ + HCHO + CO	k1_RO2pORO2	Rickard and Pascoe (2009)
G44201	TrGTerC	$CH_3COCOCH_2O_2$ + HO_2 → $CH_3COCOCH_2OOH$	KRO2HO2(4)	Rickard and Pascoe (2009)
G44202	TrGTerCN	$CH_3COCOCH_2O_2$ + NO → $CH_3C(O)$ + HCHO + CO + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G44203a	TrGTerC	$CH_3COCOCH_2OOH$ + OH → $CH_3COCOCHO$ + OH	k_s*f_co*f_sooh	Rickard and Pascoe (2009)*
G44203b	TrGTerC	$CH_3COCOCH_2OOH$ + OH → $CH_3COCOCH_2O_2$	k_roohro	Rickard and Pascoe (2009)
G44204	TrGTerC	C44O2 + HO_2 → C44OOH	KRO2HO2(4)	Rickard and Pascoe (2009)
G44205	TrGTerCN	C44O2 + NO → HCOCH2CHO + CO_2 + HO_2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G44206	TrGTerC	C44O2 → HCOCH2CHO + CO_2 + HO_2	k1_RO2sORO2	Rickard and Pascoe (2009)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G44207	TrGTerC	C44OOH + OH → C44O2	7.46E-11	Rickard and Pascoe (2009)
G44208	TrGTerC	CHOC3COO2 → HCOCH2CO3 + HCHO	k1_RO2pORO2	Rickard and Pascoe (2009)
G44209	TrGTerC	CHOC3COO2 + HO_2 → C413COOOH	KRO2HO2(4)	Rickard and Pascoe (2009)
G44210	TrGTerCN	CHOC3COO2 + NO → HCOCH2CO3 + HCHO + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G44211	TrGTerC	C413COOOH + OH → CHOC3COO2	8.33E-11	Rickard and Pascoe (2009)
G44212	TrGTerC	C4CODIAL + OH → C312COCO3	3.39E-11	Rickard and Pascoe (2009)
G44213	TrGTerCN	C4CODIAL + NO_3 → C312COCO3 + HNO_3	2.*KNO3AL*4.0	Rickard and Pascoe (2009)
G44214	TrGTerC	C312COCO3 → $HCOCOCH_2O_2$ + CO_2	k1_RO2RCO3	Rickard and Pascoe (2009)
G44215a	TrGTerC	C312COCO3 + HO_2 → C312COCO3H	KAPHO2*rco3_ooh	Rickard and Pascoe (2009)
G44215b	TrGTerC	C312COCO3 + HO_2 → $HCOCOCH_2O_2$ + CO_2 + OH	KAPHO2*(1-rco3_ooh)	Rickard and Pascoe (2009)
G44216	TrGTerCN	C312COCO3 + NO_2 → C312COPAN	k_CH3CO3_NO2	Rickard and Pascoe (2009)
G44217	TrGTerCN	C312COCO3 + NO → $HCOCOCH_2O_2$ + CO_2 + NO_2	KAPNO	Rickard and Pascoe (2009)
G44218	TrGTerC	C312COCO3H + OH → C312COCO3	1.63E-11	Rickard and Pascoe (2009)
G44219	TrGTerCN	C312COPAN → C312COCO3 + NO_2	k_PAN_M	Rickard and Pascoe (2009)
G44220	TrGTerCN	C312COPAN + OH → HCOCOCHO + CO + NO_2	1.27E-11	Rickard and Pascoe (2009)
G44221	TrGTerC	$CH_3COCOCHO$ + OH → $CH_3C(O)$ + 2 CO	8.4E-13*EXP(830./temp)	Sander et al. (2018)*
G44222	TrGTerCN	$CH_3COCOCHO$ + NO_3 → $CH_3C(O)$ + 2 CO + HNO_3	KNO3AL*4.0	Rickard and Pascoe (2009)
G44223	TrGTerC	IBUTALOH + OH → IPRHOCO3	1.4E-11	Rickard and Pascoe (2009)
G44224a	TrGTerC	IPRHOCO3 + HO_2 → CH_3COCH_3 + CO_2 + HO_2 + OH	KAPHO2*rco3_oh	Rickard and Pascoe (2009), Sander et al. (2018)
G44224b	TrGTerC	IPRHOCO3 + HO_2 → IPRHOCO2H + O_3	KAPHO2*rco3_o3	Rickard and Pascoe (2009), Sander et al. (2018)
G44224c	TrGTerC	IPRHOCO3 + HO_2 → IPRHOCO3H	KAPHO2*rco3_ooh	Rickard and Pascoe (2009), Sander et al. (2018)
G44225	TrGTerCN	IPRHOCO3 + NO → CH_3COCH_3 + CO_2 + HO_2 + NO_2	KAPNO	Rickard and Pascoe (2009)
G44226	TrGTerCN	IPRHOCO3 + NO_2 → C4PAN5	k_CH3CO3_NO2	Rickard and Pascoe (2009)
G44227	TrGTerCN	IPRHOCO3 + NO_3 → CH_3COCH_3 + CO_2 + HO_2 + NO_2	KRO2NO3*1.74	Rickard and Pascoe (2009)
G44228a	TrGTerC	IPRHOCO3 → CH_3COCH_3 + CO_2 + HO_2	k1_RO2RCO3*0.7	Rickard and Pascoe (2009)
G44228b	TrGTerC	IPRHOCO3 → IPRHOCO2H	k1_RO2RCO3*0.3	Rickard and Pascoe (2009)
G44229	TrGTerC	IPRHOCO2H + OH → CH_3COCH_3 + CO_2 + HO_2 + H_2O	1.72E-12	Rickard and Pascoe (2009)
G44230	TrGTerC	OH + IPRHOCO3H → IPRHOCO3	4.80E-12	Rickard and Pascoe (2009)
G44231	TrGTerCN	C4PAN5 → IPRHOCO3 + NO_2	K_PAN_M	Rickard and Pascoe (2009)
G44232	TrGTerCN	C4PAN5 + OH → CH_3COCH_3 + CO + NO_2	4.75E-13	Rickard and Pascoe (2009)
G44233a	TrGTerC	MBOOO → IPRHOCO2H	1.60E-17*C(ind_H2O)*(0.08+0.15)	Rickard and Pascoe (2009), Sander et al. (2018)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G44233b	TrGTerC	MBOOO → IBUTALOH + H_2O_2	1.60E-17*C(ind_H2O)*0.77	Rickard and Pascoe (2009), Sander et al. (2018)
G44234	TrGTerC	MBOOO + CO → IBUTALOH + CO_2	1.20E-15	Rickard and Pascoe (2009)
G44235	TrGTerCN	MBOOO + NO → IBUTALOH + NO_2	1.00E-14	Rickard and Pascoe (2009)
G44236	TrGTerCN	MBOOO + NO_2 → IBUTALOH + NO_3	1.00E-15	Rickard and Pascoe (2009)
G44400	TrGAroC	MALANHY + OH → MALANHYO2	1.4E-12	Rickard and Pascoe (2009)
G44401a	TrGAroC	MALDIALOOH + OH → HOCOC4DIAL + OH	1.22E-10	Rickard and Pascoe (2009)
G44401b	TrGAroC	MALDIALOOH + OH → MALDIALO2	k_roohro	Rickard and Pascoe (2009)
G44402	TrGAroCN	NC4DCO2H + OH → MALANHY + NO_2	k_roohro	Rickard and Pascoe (2009)*
G44403	TrGAroC	CO14O3CO2H + OH → $HCOCH_2O_2$ + 2 CO_2	2.19E-11	Rickard and Pascoe (2009)
G44404	TrGAroC	BZFUOOH + OH → BZFUO2	3.68E-11	Rickard and Pascoe (2009)
G44405	TrGAroC	HOCOC4DIAL + OH → CO2C4DIAL + HO_2	3.67E-11	Rickard and Pascoe (2009)
G44406a	TrGAroC	MALDIALCO3 + HO_2 → MALDALCO2H + O_3	KAPHO2*rco3_o3	Rickard and Pascoe (2009)
G44406b	TrGAroC	MALDIALCO3 + HO_2 → MALDALCO3H	KAPHO2*rco3_ooh	Rickard and Pascoe (2009)
G44406c	TrGAroC	MALDIALCO3 + HO_2 → .6 MALANHY + HO_2 + .4 GLYOX + .4 CO + .4 CO_2 + OH	KAPHO2*rco3_oh	Rickard and Pascoe (2009)*
G44407	TrGAroCN	MALDIALCO3 + NO → .6 MALANHY + HO_2 + .4 GLYOX + .4 CO + .4 CO_2 + NO_2	KAPNO	Rickard and Pascoe (2009)*
G44408	TrGAroCN	MALDIALCO3 + NO_2 → MALDIALPAN	k_CH3CO3_NO2	Rickard and Pascoe (2009)
G44409	TrGAroCN	MALDIALCO3 + NO_3 → .6 MALANHY + HO_2 + .4 GLYOX + .4 CO + .4 CO_2 + NO_2	KRO2NO3*1.74	Rickard and Pascoe (2009)*
G44410	TrGAroC	MALDIALCO3 → .6 MALANHY + HO_2 + .4 GLYOX + .4 CO + .4 CO_2	k1_RO2RCO3	Rickard and Pascoe (2009)*
G44411	TrGAroCN	BZFUONE + NO_3 → NBZFUO2	3.00E-13	Rickard and Pascoe (2009)
G44412	TrGAroC	BZFUONE + O_3 → .3125 CO14O3CO2H + .1875 CO14O3CHO + .1875 H_2O_2 + .5 CO + .5 CO_2 + .5 $HCOCH_2O_2$ + .5 OH	2.20E-19	see note*
G44413	TrGAroC	BZFUONE + OH → BZFUO2	4.45E-11	Rickard and Pascoe (2009)
G44414	TrGAroCN	NBZFUOOH + OH → NBZFUO2	6.18E-12	Rickard and Pascoe (2009)
G44415	TrGAroC	MALDALCO3H + OH → MALDIALCO3	4.00E-11	Rickard and Pascoe (2009)
G44416	TrGAroC	EPXDLCO2H + OH → C3DIALO2 + CO_2	2.31E-11	Rickard and Pascoe (2009)
G44417a	TrGAroC	EPXDLCO3 + HO_2 → C3DIALO2 + CO_2 + OH	KAPHO2*rco3_oh	Rickard and Pascoe (2009)
G44417b	TrGAroC	EPXDLCO3 + HO_2 → EPXDLCO2H + O_3	KAPHO2*rco3_o3	Rickard and Pascoe (2009)
G44417c	TrGAroC	EPXDLCO3 + HO_2 → EPXDLCO3H	KAPHO2*rco3_ooh	Rickard and Pascoe (2009)
G44418	TrGAroCN	EPXDLCO3 + NO → C3DIALO2 + CO_2 + NO_2	KAPNO	Rickard and Pascoe (2009)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G44419	TrGAroCN	EPXDLCO3 + NO_2 → EPXDLPAN	k_CH3CO3_NO2	Rickard and Pascoe (2009)
G44420	TrGAroCN	EPXDLCO3 + NO_3 → C3DIALO2 + CO_2 + NO_2	KRO2NO3*1.74	Rickard and Pascoe (2009)
G44421	TrGAroC	EPXDLCO3 → C3DIALO2 + CO_2	k1_RO2RCO3	Rickard and Pascoe (2009)*
G44422	TrGAroC	MALNHYOHCO + OH → CO + CO + CO + CO_2 + HO_2	5.68E-12	Rickard and Pascoe (2009)
G44423	TrGAroCN	MALDIAL + NO_3 → MALDIALCO3 + HNO_3	2*KNO3AL*2.0	Rickard and Pascoe (2009)
G44424	TrGAroC	MALDIAL + O_3 → 1.0675 GLYOX + .125 HCHO + .1125 $HCOCO_2H$ + .0675 H_2O_2 + .82 HO_2 + .57 OH + 1.265 CO + .25 CO_2	2.00E-18	Rickard and Pascoe (2009)*
G44425	TrGAroC	MALDIAL + OH → .83 MALDIALCO3 + .17 MALDIALO2	5.20E-11	Rickard and Pascoe (2009)*
G44426	TrGAroC	MALANHYOOH + OH → MALNHYOHCO + OH	4.66E-11	Rickard and Pascoe (2009)
G44427	TrGAroCN	MALDIALPAN + OH → GLYOX + CO + CO + NO_2	3.70E-11	Rickard and Pascoe (2009)
G44428	TrGAroCN	MALDIALPAN → MALDIALCO3 + NO_2	k_PAN_M	Rickard and Pascoe (2009)
G44429a	TrGAroC	MALANHYO2 + HO_2 → MALANHYOOH	KRO2HO2(4)*(1-rcoch2o2_oh-rchohch2o2_oh)	Rickard and Pascoe (2009), Sander et al. (2018)
G44429b	TrGAroC	MALANHYO2 + HO_2 → HCOCOHCO3 + CO_2 + OH	KRO2HO2(4)*(rcoch2o2_oh+rchohch2o2_oh)	Rickard and Pascoe (2009), Sander et al. (2018)
G44430	TrGAroCN	MALANHYO2 + NO → HCOCOHCO3 + CO_2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G44431	TrGAroCN	MALANHYO2 + NO_3 → HCOCOHCO3 + CO_2 + NO_2	KRO2NO3	Rickard and Pascoe (2009)*
G44432	TrGAroC	MALANHYO2 → HCOCOHCO3 + CO_2	k1_RO2sORO2	Rickard and Pascoe (2009)*
G44433	TrGAroC	EPXDLCO3H + OH → EPXDLCO3	2.62E-11	Rickard and Pascoe (2009)
G44434	TrGAroC	CO2C4DIAL + OH → CO + CO + CO + CO + HO_2	2.45E-11	Rickard and Pascoe (2009)
G44435a	TrGAroCN	NBZFUO2 + HO_2 → NBZFUOOH	KRO2HO2(4)*(1-rcoch2o2_oh)	Rickard and Pascoe (2009), Sander et al. (2018)
G44435b	TrGAroCN	NBZFUO2 + HO_2 → .5 CO14O3CHO + .5 NO_2 + .5 NBZFUONE + .5 HO_2 + OH	KRO2HO2(4)*rcoch2o2_oh	Rickard and Pascoe (2009), Sander et al. (2018)
G44436	TrGAroCN	NBZFUO2 + NO → .5 CO14O3CHO + .5 NO_2 + .5 NBZFUONE + .5 HO_2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G44437	TrGAroCN	NBZFUO2 + NO_3 → .5 CO14O3CHO + .5 NO_2 + .5 NBZFUONE + .5 HO_2 + NO_2	KRO2NO3	Rickard and Pascoe (2009)*
G44438	TrGAroCN	NBZFUO2 → .5 CO14O3CHO + .5 NO_2 + .5 NBZFUONE + .5 HO_2	k1_RO2sORO2	Rickard and Pascoe (2009)*
G44439	TrGAroC	MALDALCO2H + OH → .6 MALANHY + HO_2 + .4 GLYOX + .4 CO + .4 CO_2	3.70E-11	Rickard and Pascoe (2009)*
G44440	TrGAroCN	EPXC4DIAL + NO_3 → EPXDLCO3 + HNO_3	2*KNO3AL*4.0	Rickard and Pascoe (2009)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G44441	TrGAroC	EPXC4DIAL + OH → EPXDLCO3	4.32E-11	Rickard and Pascoe (2009)
G44442a	TrGAroC	MECOACETO2 + HO_2 → MECOACEOOH	KRO2HO2(4)*(1-rcoch2o2_oh)	Rickard and Pascoe (2009), Sander et al. (2018)
G44442b	TrGAroC	MECOACETO2 + HO_2 → $CH_3C(O)OO$ + HCHO + CO_2 + OH	KRO2HO2(4)*rcoch2o2_oh	Rickard and Pascoe (2009), Sander et al. (2018)
G44443	TrGAroCN	MECOACETO2 + NO → $CH_3C(O)OO$ + HCHO + CO_2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G44444	TrGAroCN	MECOACETO2 + NO_3 → $CH_3C(O)OO$ + HCHO + CO_2 + NO_2	KRO2NO3	Rickard and Pascoe (2009)*
G44445	TrGAroC	MECOACETO2 → $CH_3C(O)OO$ + HCHO + CO_2	k1_RO2pORO2	Rickard and Pascoe (2009)*
G44446	TrGAroCN	CO14O3CHO + NO_3 → CO + $HCOCH_2O_2$ + CO_2 + HNO_3	KNO3AL*8.0	Rickard and Pascoe (2009)
G44447	TrGAroC	CO14O3CHO + OH → CO + $HCOCH_2O_2$ + CO_2	3.44E-11	Rickard and Pascoe (2009)
G44448	TrGAroCN	NBZFUONE + OH → BZFUCO + NO_2	1.16E-12	Rickard and Pascoe (2009)
G44449a	TrGAroC	BZFUO2 + HO_2 → BZFUOOH	KRO2HO2(4)*(1-rcoch2o2_oh-rchohch2o2_oh)	Rickard and Pascoe (2009), Sander et al. (2018)
G44449b	TrGAroC	BZFUO2 + HO_2 → CO14O3CHO + HO_2 + OH	KRO2HO2(4)*(rcoch2o2_oh+rchohch2o2_oh)	Rickard and Pascoe (2009), Sander et al. (2018)
G44450	TrGAroCN	BZFUO2 + NO → CO14O3CHO + HO_2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G44451	TrGAroCN	BZFUO2 + NO_3 → CO14O3CHO + HO_2 + NO_2	KRO2NO3	Rickard and Pascoe (2009)*
G44452	TrGAroC	BZFUO2 → CO14O3CHO + HO_2	k1_RO2sORO2	Rickard and Pascoe (2009)*
G44453	TrGAroC	BZFUCO + OH → CO14O3CHO + HO_2	1.78E-11	Rickard and Pascoe (2009)
G44456a	TrGAroC	MALDIALO2 + HO_2 → MALDIALOOH	KRO2HO2(4)*(1-rcoch2o2_oh-rchohch2o2_oh)	Rickard and Pascoe (2009)
G44456b	TrGAroC	MALDIALO2 + HO_2 → GLYOX + GLYOX + HO_2 + OH	KRO2HO2(4)*(rcoch2o2_oh+rchohch2o2_oh)	Rickard and Pascoe (2009)
G44457	TrGAroCN	MALDIALO2 + NO → GLYOX + GLYOX + HO_2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G44458	TrGAroCN	MALDIALO2 + NO_3 → GLYOX + GLYOX + HO_2 + NO_2	KRO2NO3	Rickard and Pascoe (2009)*
G44459	TrGAroC	MALDIALO2 → GLYOX + GLYOX + HO_2	k1_RO2sORO2	Rickard and Pascoe (2009)*
G44460	TrGAroCN	EPXDLPAN + OH → HCOCOCHO + CO + NO_2	2.29E-11	Rickard and Pascoe (2009)
G44461	TrGAroCN	EPXDLPAN → EPXDLCO3 + NO_2	k_PAN_M	Rickard and Pascoe (2009)*
G44462	TrGAroC	MECOACEOOH + OH → MECOACETO2	3.59E-12	Rickard and Pascoe (2009)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G45000	TrGC	C_5H_8 + O_3 → .3508 MACR + .01518 MACO2H + .2440 MVK + .7085 HCHO + .11 CH_2OO + .1275 C_3H_6 + .1575 $CH_3C(O)$ + .0510 CH_3 + .2625 HO_2 + .27 OH + .09482 H_2O_2 + .255 CO_2 + .522 CO + .07182 HCHO + .03618 $HCOCH_2O_2$ + .01782 CO + 0.05408 LCARBON	`1.03E-14*EXP(-1995./temp)`	Atkinson et al. (2006), Sander et al. (2018)
G45001	TrGC	C_5H_8 + OH → .63 LISOPAB + .30 LISOPCD + .07 LISOPEFO2	`2.7E-11*EXP(390./temp)`	Atkinson et al. (2006), Sander et al. (2018)
G45002	TrGCN	C_5H_8 + NO_3 → NISOPO2	`3.0E-12*EXP(-450./temp)`	Atkinson et al. (2006)
G45003a	TrGC	LISOPAB + O_2 → LISOPACO2	`5.530E-13`	Sander et al. (2018)
G45003b	TrGC	LISOPAB + O_2 → ISOPBO2	`3.E-12`	Sander et al. (2018)
G45004a	TrGC	LISOPCD + O_2 → LDISOPACO2	`6.780E-13`	Sander et al. (2018)
G45004b	TrGC	LISOPCD + O_2 → ISOPDO2	`3.E-12`	Sander et al. (2018)
G45005	TrGC	LISOPACO2 → LISOPAB + O_2	`3.1E12*exp(-7900./temp)*.6+ 7.8E13*exp(-8600./temp)*.4`	Sander et al. (2018)
G45006	TrGC	ISOPBO2 → LISOPAB + O_2	`3.7E14*exp(-9570./temp) +4.2E14*exp(-9970./temp)`	Sander et al. (2018)
G45007	TrGC	LDISOPACO2 → LISOPCD + O_2	`5.65E12*exp(-8410./temp) *.42+1.4E14*exp(-9110./temp)*.58`	Sander et al. (2018)
G45008	TrGC	ISOPDO2 → LISOPCD + O_2	`5.0E14*exp(-10120./temp) +8.25E14*exp(-10220/temp)`	Sander et al. (2018)
G45009a	TrGC	LISOPACO2 → C1ODC2O2C4OOH	`K16HSZ14 * 2./3.*(1-fhpal)`	Sander et al. (2018)
G45009b	TrGC	LISOPACO2 → LZCODC23DBCOOH + HO_2	`K16HSZ14 * (2./3.*fhpal + 1./3.)`	Sander et al. (2018)
G45010a	TrGC	LDISOPACO2 → C1OOHC3O2C4OD	`k16HSZ41 * 2./3.*(1-fhpal)`	Sander et al. (2018)
G45010b	TrGC	LDISOPACO2 → LZCODC23DBCOOH + HO_2	`k16HSZ41 * (2./3.*fhpal + 1./3.)`	Sander et al. (2018)
G45011	TrGC	LISOPACO2 → .9 LISOPACO + .1 ISOPAOH	`k1_RO2LISOPACO2`	Rickard and Pascoe (2009), Sander et al. (2018)
G45012	TrGC	LISOPACO2 + HO_2 → LISOPACOOH + 0.024 BLOV + 0.119 BSOV	`KRO2HO2(5)`	Rickard and Pascoe (2009)
G45013a	TrGCN	LISOPACO2 + NO → LISOPACO + NO_2 + 0.003 BLOV + 0.101 BSOV	`KRO2NO*(1.-alpha_AN(6,1,0,0,0, temp,cair))`	Lockwood et al. (2010), Paulot et al. (2009a), Sander et al. (2018)
G45013b	TrGCN	LISOPACO2 + NO → LISOPACNO3	`KRO2NO*alpha_AN(6,1,0,0,0,temp, cair)`	Lockwood et al. (2010), Paulot et al. (2009a), Sander et al. (2018)
G45014	TrGCN	LISOPACO2 + NO_3 → LISOPACO + NO_2	`KRO2NO3`	Rickard and Pascoe (2009)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G45015	TrGC	LDISOPACO2 → .9 LISOPACO + .1 ISOPAOH	k1_RO2LISOPACO2	Rickard and Pascoe (2009), Sander et al. (2018)
G45016	TrGC	LDISOPACO2 + HO_2 → LISOPACOOH + 0.024 BLOV + 0.119 BSOV	KRO2HO2(5)	Rickard and Pascoe (2009)
G45017a	TrGCN	LDISOPACO2 + NO → LISOPACO + NO_2 + 0.003 BLOV + 0.101 BSOV	KRO2NO*(1.-alpha_AN(6,1,0,0,0, temp,cair))	Lockwood et al. (2010), Paulot et al. (2009a), Sander et al. (2018)
G45017b	TrGCN	LDISOPACO2 + NO → LISOPACNO3	KRO2NO*alpha_AN(6,1,0,0,0,temp, cair)	Lockwood et al. (2010), Paulot et al. (2009a), Sander et al. (2018)
G45018	TrGCN	LDISOPACO2 + NO_3 → LISOPACO + NO_2	KRO2NO3	Rickard and Pascoe (2009)
G45019a	TrGC	LISOPACOOH + OH → LISOPACO2	k_roohro	Sander et al. (2018)
G45019b	TrGC	LISOPACOOH + OH → LZCODC23DBCOOH + HO_2	k_s*f_allyl*f_soh	Sander et al. (2018)
G45019c	TrGC	LISOPACOOH + OH → LHC4ACCHO + OH	(k_s*f_sooh*f_allyl+ k_rohro)	Sander et al. (2018)
G45019d	TrGC	LISOPACOOH + OH → LIEPOX + OH	(k_adt+k_ads)*a_ch2oh*a_ch2ooh	Sander et al. (2018)*
G45020	TrGC	ISOPAOH + OH → LHC4ACCHO + HO_2	(k_adt+k_ads)*a_ch2oh*a_ch2oh+k_ s*f_soh*f_allyl+k_rohro	Sander et al. (2018)
G45021	TrGCN	LISOPACNO3 + OH → LISOPACNO3O2	(k_adt+k_ads)*a_ch2ono2*a_ch2oh	Sander et al. (2018)*
G45022	TrGC	ISOPBO2 → .8 MVK + .8 HCHO + .8 HO_2 + .2 ISOPBOH	k1_RO2ISOPBO2	Rickard and Pascoe (2009)
G45023a	TrGC	ISOPBO2 + HO_2 → ISOPBOOH	KRO2HO2(5)*(1.-rchohch2o2_oh)	Sander et al. (2018)
G45023b	TrGC	ISOPBO2 + HO_2 → MVK + HCHO + HO_2 + OH + 0.024 BLOV + 0.119 BSOV	KRO2HO2(5)*rchohch2o2_oh	Sander et al. (2018)
G45024a	TrGCN	ISOPBO2 + NO → MVK + HCHO + HO_2 + NO_2 + 0.003 BLOV + 0.101 BSOV	KRO2NO*(1.-alpha_AN(6,3,0,0,0, temp,cair))	Lockwood et al. (2010), Sander et al. (2018)
G45024b	TrGCN	ISOPBO2 + NO → ISOPBNO3	KRO2NO*alpha_AN(6,3,0,0,0,temp, cair)	Lockwood et al. (2010), Sander et al. (2018)
G45025	TrGCN	ISOPBO2 + NO_3 → MVK + .75 HCHO + .75 HO_2 + .25 CH_3 + NO_2	KRO2NO3	Rickard and Pascoe (2009)
G45026a	TrGC	ISOPBOOH + OH → LIEPOX + OH	(k_ads+k_adp)*a_ch2ooh	Paulot et al. (2009b), Sander et al. (2018)
G45026b	TrGC	ISOPBOOH + OH → ISOPBO2	k_roohro	Sander et al. (2018)
G45026c	TrGC	ISOPBOOH + OH → MGLYOX + $HOCH_2CHO$	k_rohro+k_s*f_alk*f_soh	Sander et al. (2018)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G45027	TrGC	ISOPBOOH + O_3 → .1368 MACROOH + .1368 H_2O_2 + .2280 HO_2 + .4332 CH_3COCH_2OH + .2280 CO_2 + .6384 OH + .2052 CO + .57 HCHO + .43 MACROOH + .06880 HO_2 + .06880 OH + .2709 CO + .1591 CH_2OO	`1.E-17`	Sander et al. (2018)
G45028	TrGC	ISOPBOH + OH → MVK + .75 HCHO + .75 HO_2 + .25 CH_3	`k_s*f_alk*f_soh+(k_adp+k_ads) *a_ch2oh`	Sander et al. (2018)
G45029	TrGCN	ISOPBNO3 + OH → ISOPBDNO3O2	`(k_adt+k_adp)*f_ch2ono2`	Sander et al. (2018)
G45030	TrGC	ISOPDO2 → .8 MACR + .8 HCHO + .8 HO_2 + .1 HCOC5 + .1 ISOPDOH	`k1_RO2ISOPDO2`	Rickard and Pascoe (2009)
G45031a	TrGC	ISOPDO2 + HO_2 → ISOPDOOH	`KRO2HO2(5)*(1.-rchohch2o2_oh)`	Sander et al. (2018)
G45031b	TrGC	ISOPDO2 + HO_2 → MACR + HCHO + HO_2 + OH + 0.024 BLOV + 0.119 BLOV	`KRO2HO2(5)*rchohch2o2_oh`	Sander et al. (2018)
G45032a	TrGCN	ISOPDO2 + NO → MACR + HCHO + HO_2 + NO_2 + 0.003 BLOV + 0.101 BSOV	`KRO2NO*(1.-alpha_AN(6,2,0,0,0, temp,cair))`	Lockwood et al. (2010), Sander et al. (2018)
G45032b	TrGCN	ISOPDO2 + NO → ISOPDNO3	`KRO2NO*alpha_AN(6,2,0,0,0,temp, cair)`	Lockwood et al. (2010), Sander et al. (2018)
G45033	TrGCN	ISOPDO2 + NO_3 → MACR + HCHO + HO_2 + NO_2	`KRO2NO3`	Rickard and Pascoe (2009)
G45034a	TrGC	ISOPDOOH + OH → LIEPOX + OH	`(k_adt+k_adp)*a_ch2ooh`	Paulot et al. (2009b), Sander et al. (2018)
G45034b	TrGC	ISOPDOOH + OH → ISOPDO2	`k_roohro`	Sander et al. (2018)
G45034c	TrGC	ISOPDOOH + OH → HCOC5 + OH	`k_t*f_tooh*f_allyl*f_pch2oh`	Sander et al. (2018)
G45034d	TrGC	ISOPDOOH + OH → CH_3COCH_2OH + GLYOX + OH	`k_s*f_pch2oh*f_soh`	Sander et al. (2018)
G45035	TrGC	ISOPDOOH + O_3 → 1.393 OH + BIACETOH + .67 HCHO + .05280 HO_2 + .2079 CO + .1221 CH_2OO	`1.E-17`	Sander et al. (2018)
G45036	TrGC	ISOPDOH + OH → HCOC5 + HO_2	`2.*k_rohro+(k_t*f_toh*f_allyl+k_ s*f_soh)*f_pch2oh+(k_adt+k_adp) *a_ch2oh`	Sander et al. (2018)
G45037	TrGCN	ISOPDNO3 + OH → ISOPBDNO3O2	`(k_adp+k_ads)*a_ch2ono2`	Sander et al. (2018)*
G45038	TrGCN	NISOPO2 → .8 NC4CHO + .6 HO_2 + .2 LISOPACNO3	`k1_RO2LISOPACO2`	Rickard and Pascoe (2009)
G45039	TrGCN	NISOPO2 + HO_2 → NISOPOOH	`KRO2HO2(5)`	Rickard and Pascoe (2009)
G45040	TrGCN	NISOPO2 + NO → NC4CHO + HO_2 + NO_2	`KRO2NO`	Rickard and Pascoe (2009)*
G45041	TrGCN	NISOPO2 + NO_3 → NC4CHO + HO_2 + NO_2	`KRO2NO3`	Rickard and Pascoe (2009)
G45042	TrGCN	NISOPOOH + OH → NC4CHO + OH	`1.03E-10`	Rickard and Pascoe (2009)
G45043	TrGCN	NC4CHO + OH → LNISO3	`(k_adt+k_ads)*a_cho*a_ch2ono2`	Sander et al. (2018)*

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G45044	TrGCN	NC4CHO + O_3 → .27 NOA + .027 $HCOCO_2H$ + .0162 GLYOX + .0162 H_2O_2 + .1458 HCOCO + .0405 HCOOH + .0405 CO + .8758 OH + .365 MGLYOX + .73 NO_2 + 0.7705 HCHO + .4055 CO_2 + .73 GLYOX	2.40E-17	Sander et al. (2018)
G45045	TrGCN	NC4CHO + NO_3 → LNISO3 + HNO_3	KNO3AL*4.25	Rickard and Pascoe (2009)
G45046	TrGCN	LNISO3 + HO_2 → LNISOOH	0.5*KRO2HO2(5)+0.5*KAPHO2	Rickard and Pascoe (2009)
G45047	TrGCN	LNISO3 + NO → NOA + .5 HOCHCHO + .5 CO + .5 HO_2 + NO_2 + .5 CO_2	0.5*KAPNO+0.5*KRO2NO	Rickard and Pascoe (2009)*
G45048	TrGCN	LNISO3 + NO_3 → NOA + .5 HOCHCHO + .5 CO + .5 HO_2 + NO_2 + .5 CO_2	KRO2NO3*1.37	Rickard and Pascoe (2009)
G45049	TrGCN	LNISOOH + OH → LNISO3	2.65E-11	Rickard and Pascoe (2009)
G45050a	TrGC	LHC4ACCHO + OH → LC578O2	(k_adtertprim+k_ads)*a_cho*a_ch2oh	Sander et al. (2018)
G45050b	TrGC	LHC4ACCHO + OH → LHC4ACCO3	k_t*f_o	Sander et al. (2018)
G45050c	TrGC	LHC4ACCHO + OH → C4MDIAL + HO_2	k_s*f_soh*f_allyl	Sander et al. (2018)
G45051	TrGC	LHC4ACCHO + O_3 → .2225 $CH_3C(O)$ + .89 CO + .0171875 $HOCH_2CO_2H$ + .075625 H_2O_2 + .0171875 $HCOCO_2H$ + .2775 CH_3COCH_2OH + .6675 HO_2 + .2603125 GLYOX + .2225 HCHO + .89 OH + .2603125 $HOCH_2CHO$ + .5 MGLYOX	2.40E-17	Rickard and Pascoe (2009)
G45052	TrGCN	LHC4ACCHO + NO_3 → LHC4ACCO3 + HNO_3	KNO3AL*4.25	Rickard and Pascoe (2009)
G45053	TrGC	LC578O2 → .25 CH_3COCH_2OH + .75 MGLYOX + .25 HOCHCHO + .75 $HOCH_2CHO$ + .75 HO_2	k1_RO2tORO2	Rickard and Pascoe (2009)
G45054a	TrGC	LC578O2 + HO_2 → MGLYOX + $HOCH_2CHO$ + OH	KRO2HO2(5)*rcoch2o2_oh	Rickard and Pascoe (2009)
G45054b	TrGC	LC578O2 + HO_2 → LC578OOH	KRO2HO2(5)*rcoch2o2_ooh	Rickard and Pascoe (2009)
G45055	TrGCN	LC578O2 + NO → .25 CH_3COCH_2OH + .75 MGLYOX + .25 HOCHCHO + .75 $HOCH_2CHO$ + .75 HO_2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G45056	TrGCN	LC578O2 + NO_3 → .25 CH_3COCH_2OH + .75 MGLYOX + .25 HOCHCHO + .75 $HOCH_2CHO$ + .75 HO_2 + NO_2	KRO2NO3	Rickard and Pascoe (2009)
G45057	TrGC	LC578O2 → .25 CH_3COCH_2OH + .75 MGLYOX + .25 $HOCH_2CHO$ + .75 $HOCH_2CHO$ + HO_2 + OH	KHSB	Sander et al. (2018)
G45058a	TrGC	LC578OOH + OH → LC578O2	k_roohro	Sander et al. (2018)
G45058b	TrGC	LC578OOH + OH → C1ODC2OOHC4OD + HO_2	k_t*f_o*f_tch2oh*f_alk+k_t*f_toh*f_pch2oh*f_pch2oh+k_s*f_soh*f_pch2oh	Sander et al. (2018)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G45059a	TrGC	LHC4ACCO3 $\rightarrow$ OH + .5 MACRO2 + .5 LHMVKABO2 + CO_2	k1_RO2RCO3*0.9	Sander et al. (2018)
G45059b	TrGC	LHC4ACCO3 $\rightarrow$ LHC4ACCO2H	k1_RO2RCO3*0.1	Sander et al. (2018)
G45060a	TrGC	LHC4ACCO3 + HO_2 $\rightarrow$ 2 OH + .5 MACRO2 + .5 LHMVKABO2 + CO_2	KAPHO2*rco3_oh	Sander et al. (2018)
G45060b	TrGC	LHC4ACCO3 + HO_2 $\rightarrow$ LHC4ACCO3H	KAPHO2*rco3_ooh	Sander et al. (2018)
G45060c	TrGC	LHC4ACCO3 + HO_2 $\rightarrow$ LHC4ACCO2H + O_3	KAPHO2*rco3_o3	Sander et al. (2018)
G45061	TrGCN	LHC4ACCO3 + NO $\rightarrow$ .5 MACRO2 + .5 LHMVKABO2 + NO_2 + CO_2	KAPNO	Sander et al. (2018)
G45062	TrGCN	LHC4ACCO3 + NO_2 $\rightarrow$ LC5PAN1719	k_CH3CO3_NO2	Rickard and Pascoe (2009)
G45063	TrGCN	LHC4ACCO3 + NO_3 $\rightarrow$ .5 MACRO2 + .5 LHMVKABO2 + NO_2 + CO_2	KRO2NO3*1.74	Sander et al. (2018)
G45064a	TrGC	LHC4ACCO2H + OH $\rightarrow$ OH + .5 MACRO2 + .5 LHMVKABO2 + CO_2	2.52E-11	Sander et al. (2018)
G45064b	TrGC	LHC4ACCO3H + OH $\rightarrow$ LHC4ACCO3	2.88E-11	Rickard and Pascoe (2009)
G45065	TrGCN	LC5PAN1719 $\rightarrow$ LHC4ACCO3 + NO_2	k_PAN_M	Rickard and Pascoe (2009)
G45066	TrGCN	LC5PAN1719 + OH $\rightarrow$ .5 MACROH + .5 HO12CO3C4 + CO + NO_2	2.52E-11	Rickard and Pascoe (2009)
G45067	TrGC	HCOC5 + OH $\rightarrow$ C59O2	3.81E-11	Rickard and Pascoe (2009)
G45068	TrGC	HCOC5 + O_3 $\rightarrow$ BIACETOH + .335 H_2O_2 + .67 HCHO + .2079 CO + .1221 CH_2OO + .05280 OH	7.51E-16*EXP(-1521./temp)	Sander et al. (2018)
G45069	TrGC	C59O2 $\rightarrow$ CH_3COCH_2OH + HOCH2CO	k1_RO2tORO2	Sander et al. (2018)
G45070a	TrGC	C59O2 + HO_2 $\rightarrow$ OH + CH_3COCH_2OH + HOCH2CO	KRO2HO2(5)*rcoch2o2_oh	Sander et al. (2018)
G45070b	TrGC	C59O2 + HO_2 $\rightarrow$ C59OOH	KRO2HO2(5)*rcoch2o2_ooh	Sander et al. (2018)
G45071	TrGCN	C59O2 + NO $\rightarrow$ CH_3COCH_2OH + HOCH2CO + NO_2	KRO2NO	Sander et al. (2018)*
G45072	TrGCN	C59O2 + NO_3 $\rightarrow$ CH_3COCH_2OH + HOCH2CO + NO_2	KRO2NO3	Sander et al. (2018)
G45073	TrGC	C59OOH + OH $\rightarrow$ C59O2	9.7E-12	Rickard and Pascoe (2009)
G45074	TrGC	LIEPOX + OH $\rightarrow$ DB1O2 + H_2O	5.78E-11*EXP(-400./temp) *(1.52/3.+0.98*2./3.)/1.51	Paulot et al. (2009b), Bates et al. (2014), Sander et al. (2018)*
G45075	TrGC	ISOPBO2 $\rightarrow$ MVK + HCHO + OH	KHSB	Sander et al. (2018)
G45076	TrGC	ISOPDO2 $\rightarrow$ MACR + HCHO + OH	KHSD	Sander et al. (2018)
G45077a	TrGC	LZCODC23DBCOOH + OH $\rightarrow$ .6 C1ODC2O2C4OOH + .4 C1OOHC2O2C4OD	k_adt*a_cho*a_ch2ooh	Sander et al. (2018)
G45077b	TrGC	LZCODC23DBCOOH + OH $\rightarrow$ .6 C1ODC3O2C4OOH + .4 C1OOHC3O2C4OD	k_ads*a_cho*a_ch2ooh	Sander et al. (2018)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G45077c	TrGC	LZCODC23DBCOOH + OH → LZCO3HC23DBCOD	k_t*f_o*f_alk+k_roohro	Sander et al. (2018)
G45077d	TrGC	LZCODC23DBCOOH + OH → C4MDIAL + OH	k_s*f_sooh*f_allyl	Sander et al. (2018)
G45078	TrGC	LZCODC23DBCOOH + O_3 → .4672 OH + .2336 $HCOCOCH_2O_2$ + .2336 CO + .2336 $CH_3C(O)$ + .4672 HOOCH2CHO + .1728 MGLYOX + .1901 OH + .0864 GLYOX + .02765 HOOCH2CHO + .02765 H_2O_2 + .02592 CH_3OOH + .02592 CO_2 + .01037 HCOCO + .01555 CH_2OO + .01555 CO + .006908 $HOOCH_2CO_3$ + .2628 OH + .1314 MGLYOX + .1314 OH + .1314 $HCOCOCH_2OOH$ + .2628 GLYOX + .0972 $CH_3COCH_2O_2H$ + .00972 $HCOCO_2H$ + .005832 GLYOX + .005832 H_2O_2 + .05249 OH + .05249 HCOCO + .01458 HCHO + .01458 CO_2 + .01458 HCOOH + .01458 CO	2.4E-17	Sander et al. (2018)
G45079	TrGC	C1OOHC2O2C4OD → .78 $CH_3COCH_2O_2H$ + .78 HOCHCHO + .22 CO2H3CHO + .22 HCHO + .22 OH	k1_RO2tORO2	Sander et al. (2018)
G45080	TrGCN	C1OOHC2O2C4OD + NO → .78 $CH_3COCH_2O_2H$ + .78 HOCHCHO + .22 CO2H3CHO + .22 HCHO + .22 OH + NO_2	KRO2NO	Sander et al. (2018)*
G45081a	TrGC	C1OOHC2O2C4OD + HO_2 → C1OOHC2OOHC4OD	KRO2HO2(5)*rcoch2o2_ooh	Sander et al. (2018)
G45081b	TrGC	C1OOHC2O2C4OD + HO_2 → .78 $CH_3COCH_2O_2H$ + .78 HOCHCHO + .22 CO2H3CHO + .22 HCHO + 1.22 OH	KRO2HO2(5)*rcoch2o2_oh	Sander et al. (2018)
G45082	TrGC	C1OOHC2O2C4OD → $CH_3COCH_2O_2H$ + GLYOX + OH	KHSB	Sander et al. (2018)
G45083	TrGC	C1ODC2O2C4OOH → OH + C1ODC2OOHC4OD	K15HSDHB	Sander et al. (2018)
G45084a	TrGC	C1OOHC2OOHC4OD + OH → C1ODC2OOHC4OD + OH	2.*k_s*f_sooh*f_tch2oh	Sander et al. (2018)
G45084b	TrGC	C1OOHC2OOHC4OD + OH → $CH_3COCH_2O_2H$ + 2 CO + 2 HO_2 + OH	k_t*f_toh*f_pch2oh*f_pch2oh	Sander et al. (2018)
G45084c	TrGC	C1OOHC2OOHC4OD + OH → C1OOHC2O2C4OD	k_roohro	Sander et al. (2018)
G45085	TrGC	C1ODC2OOHC4OD + OH → CO2H3CHO + CO + H_2O + OH	k_t*f_o*f_tch2oh+k_t*f_toh*f_toh*f_cho	Sander et al. (2018)
G45086	TrGC	C1ODC3O2C4OOH → MGLYOX + HOOCH2CHO + HO_2	k1_RO2sORO2	Sander et al. (2018)
G45087	TrGCN	C1ODC3O2C4OOH + NO → MGLYOX + HOOCH2CHO + HO_2 + NO_2	KRO2NO	Sander et al. (2018)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G45088	TrGC	C1ODC3O2C4OOH + HO_2 → .5 $CH_3C(O)$ + .5 CO + .5 MGLYOX + .5 HO_2 + $HOOCH_2CO_3$	KRO2HO2(5)	Sander et al. (2018)
G45089	TrGC	C1ODC3O2C4OOH → MGLYOX + OH + HOOCH2CHO	KHSD	Sander et al. (2018)
G45090	TrGC	C1OOHC3O2C4OD → .625 MGLYOX + 2 CO + 1.625 HO_2 + .375 $CH_3C(O)$ + .375 CO_2 + OH	K15HSDHB	Sander et al. (2018)
G45091	TrGC	LHC4ACCO3 → LZCO3HC23DBCOD + HO_2	K16HS	Sander et al. (2018)
G45092a	TrGC	C4MDIAL + OH → C1ODC2O2C4OD	(k_adt+k_ads)*a_cho*a_cho	Sander et al. (2018)*
G45092b	TrGC	C4MDIAL + OH → LZCO3C23DBCOD	2*k_t*f_o*f_alk	Sander et al. (2018)*
G45093	TrGCN	C4MDIAL + NO_3 → LZCO3C23DBCOD + HNO_3	KNO3AL*4.25*2.	Sander et al. (2018)*
G45094a	TrGC	C1ODC2O2C4OD + HO_2 → OH + MGLYOX + HOCHCHO	KRO2HO2(5)*rcoch2o2_oh	Sander et al. (2018)
G45094b	TrGC	C1ODC2O2C4OD + HO_2 → C1ODC2OOHC4OD	KRO2HO2(5)*rcoch2o2_ooh	Sander et al. (2018)
G45095	TrGCN	C1ODC2O2C4OD + NO → NO_2 + MGLYOX + HOCHCHO	KRO2NO	Sander et al. (2018)*
G45096	TrGC	C1ODC2O2C4OD → MGLYOX + HOCHCHO	k1_RO2tORO2	Sander et al. (2018)
G45097a	TrGC	C1ODC2OOHC4OD + OH → MGLYOX + 2 CO	(2.*k_t*f_o*f_tch2oh*f_alk+k_t*f_toh*f_cho*f_pch2oh)*.5	Sander et al. (2018)
G45097b	TrGC	C1ODC2OOHC4OD + OH → MGLYOX + 2 CO + OH	(2.*k_t*f_o*f_tch2oh*f_alk+k_t*f_toh*f_cho*f_pch2oh)*.5	Sander et al. (2018)
G45098	TrGCN	LISOPACNO3O2 + NO → .21 NOA + .21 $HOCH_2CHO$ + .21 HO_2 + .49 HO12CO3C4 + .49 HCHO + .49 NO_2 + .045 MVKNO3 + .045 HCHO + .255 CH_3COCH_2OH + .255 NO_3CH2CHO + .225 H_2O_2 + NO_2	KRO2NO	Sander et al. (2018)*
G45099	TrGCN	LISOPACNO3O2 → .21 NOA + .21 $HOCH_2CHO$ + .21 HO_2 + .49 HO12CO3C4 + .49 HCHO + .49 NO_2 + .045 MVKNO3 + .045 HCHO + .255 CH_3COCH_2OH + .255 NO_3CH2CHO + .225 H_2O_2	k1_RO2tORO2+KRO2HO2(5)*c(ind_HO2)	Sander et al. (2018)
G45100	TrGCN	ISOPBDNO3O2 + NO → .6 CH_3COCH_2OH + .6 $HOCH_2CHO$ + .26 MACRNO3 + .14 MVKNO3 + .4 HCHO + .4 HO_2 + 1.6 NO_2	KRO2NO	Sander et al. (2018)*
G45101	TrGCN	ISOPBDNO3O2 → .6 CH_3COCH_2OH + .6 $HOCH_2CHO$ + .26 MACRNO3 + .14 MVKNO3 + .4 HCHO + .4 HO_2 + .6 NO_2	k1_RO2sORO2+KRO2HO2(5)*c(ind_HO2)	Sander et al. (2018)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G45102	TrGCN	LISOPACNO3 + O_3 → .8704 OH + .365 HO_2 + .73 MGLYOX + .4325 NO_3CH2CHO + .135 CH_3COCH_2OH + .0675 GLYOX + .4325 NO_2 + .0891 H_2O_2 + .135 NOA + .0675 HOCHCHO + .3866 $HOCH_2CHO$ + .0405 CH_3OH + .0405 CO + .0054 HOCH2CO	`2.8E-17`	Feierabend et al. (2008), Sander et al. (2018)
G45103	TrGC	DB1O2 → DB1O2	`k1_RO2sOR02`	Sander et al. (2018)
G45104a	TrGC	DB1O2 + HO_2 → DB1OOH	`KRO2HO2(5)*(1.-rchohch2o2_oh)`	Sander et al. (2018)*
G45104b	TrGC	DB1O2 + HO_2 → DB1O2 + OH	`KRO2HO2(5)*rchohch2o2_oh`	Sander et al. (2018)
G45105a	TrGCN	DB1O2 + NO → DB1O2 + NO_2	`KRO2NO*(1.-alpha_AN(7,2,0,0,0, temp,cair))`	Sander et al. (2018)
G45105b	TrGCN	DB1O2 + NO → DB1NO3	`KRO2NO*alpha_AN(7,2,0,0,0,temp, cair)`	Sander et al. (2018)
G45106	TrGCN	DB1O2 + NO_3 → DB1O2 + NO_2	`KRO2NO3`	Sander et al. (2018)
G45107	TrGC	DB1O2 → DB1O2 + OH	`1.E4`	Peeters and Nguyen (2012)*
G45108a	TrGC	DB1O2 → DB1O2	`KDEC*0.72`	see note*
G45108b	TrGC	DB1O2 → .5 HVMK + .5 HMAC + HCHO + HO_2	`KDEC*0.28`	see note*
G45109	TrGC	DB1O2 → .48 CH_3COCH_2OH + .52 $HOCH_2CHO$ + .52 MGLYOX + .48 GLYOX + HO_2	`k1_RO2sOR02`	Sander et al. (2018)
G45110a	TrGC	DB1O2 + HO_2 → DB2OOH	`KRO2HO2(5)*(1.-rchohch2o2_oh)`	Sander et al. (2018)
G45110b	TrGC	DB1O2 + HO_2 → .48 CH_3COCH_2OH + .52 $HOCH_2CHO$ + .52 MGLYOX + .48 GLYOX + HO_2 + OH	`KRO2HO2(5)*rchohch2o2_oh`	Sander et al. (2018)
G45111	TrGCN	DB1O2 + NO → .48 CH_3COCH_2OH + .52 $HOCH_2CHO$ + .52 MGLYOX + .48 GLYOX + HO_2 + NO_2	`KRO2NO`	see note*
G45112	TrGCN	DB1O2 + NO_3 → .48 CH_3COCH_2OH + .52 $HOCH_2CHO$ + .52 MGLYOX + .48 GLYOX + HO_2 + NO_2	`KRO2NO3`	Sander et al. (2018)
G45113	TrGC	DB1O2 → .48 MACROOH + .52 LHMVKABOOH + CO + OH	`K14HSAL`	Sander et al. (2018)
G45114a	TrGC	DB1OOH + OH → DB1O2	`k_roohro`	Sander et al. (2018)
G45114b	TrGC	DB1OOH + OH → HCOOH + HO_2 + CH_3COCHO_2CHO	`k_adt`	Sander et al. (2018)*
G45115	TrGC	DB1OOH + HCOOH → C1ODC2OOHC4OD + HCOOH	`4.67E-26*temp**3.286*EXP(4509./ (1.987*temp))`	Sander et al. (2018), da Silva (2010)*
G45116	TrGCN	DB1NO3 + OH → HCOOH + NO_2 + CH_3COCHO_2CHO	`k_adt`	Sander et al. (2018)*
G45117	TrGC	DB2OOH + OH → DB1O2	`k_roohro`	Sander et al. (2018)*

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G45118	TrGC	LISOPACOOH + O_3 → 1.3272 OH + .36986 HO_2 + .0432 H_2O_2 + .08422 CO + .2025 CH_3OOH + .01215 CH_2OO + .3704 HCHO + .00405 CH_3OH + .0405 CO_2 + .1825 HOCH2COCH2O2 + .365 MGLYOX + .3866 HOOCH2CHO + .135 CH_3COCH_2OH + .0675 GLYOX + .00324 HCOCO + .3866 $HOCH_2CHO$ + .135 $CH_3COCH_2O_2H$ + .0675 HOCHCHO + .0054 HOCH2CO	`4.829E-16`	Sander et al. (2018)
G45119a	TrGC	LZCO3HC23DBCOD + OH → .62 CO2H3CHO + .62 OH + .62 CO_2 + .38 MGLYOX + .38 $HCOCO_3H$ + .38 HO_2	`k_adt*a_cho*a_co2h`	Sander et al. (2018)
G45119b	TrGC	LZCO3HC23DBCOD + OH → .62 CH_3COCO_3H + 1.24 CO + 1.24 HO_2 + .38 MGLYOX + .38 HO_2 + .38 CO + .38 HO_2 + .38 OH + .38 CO_2	`k_ads*a_cho*a_co2h`	Sander et al. (2018)
G45120	TrGC	LISOPEFO2 → LISOPEFO	`k1_RO2pORO2`	Sander et al. (2018)
G45121a	TrGCN	LISOPEFO2 + NO → LISOPEFO + NO_2	`KRO2NO*(1.-alpha_AN(6,1,0,0,0,temp,cair))`	Sander et al. (2018)
G45121b	TrGCN	LISOPEFO2 + NO → ISOPDNO3	`KRO2NO*alpha_AN(6,1,0,0,0,temp,cair)`	Sander et al. (2018)*
G45122a	TrGC	LISOPEFO2 + HO_2 → .7143 ISOPDOOH + .2857 ISOPBOOH	`KRO2HO2(5)*(1.-rchohch2o2_oh)`	Sander et al. (2018)
G45122b	TrGC	LISOPEFO2 + HO_2 → LISOPEFO + OH	`KRO2HO2(5)*rchohch2o2_oh`	Sander et al. (2018)
G45123	TrGCN	LISOPEFO2 + NO_3 → LISOPEFO + NO_2	`KRO2NO3`	Sander et al. (2018)
G45124	TrGC	LISOPEFO2 → .7143 MACR + .2857 MVK + HCHO + OH	`0.7143*KHSD+.2857*KHSB`	Sander et al. (2018)
G45125	TrGC	LISOPEFO → .7143 MACR + .2857 MVK + HCHO + HO_2	`KDEC`	Sander et al. (2018)
G45126a	TrGC	LISOPACO → 3METHYLFURAN + HO_2	`KDEC*0.37`	Sander et al. (2018), Paulot et al. (2009a), Francisco-Marquez et al. (2003)
G45126b	TrGC	LISOPACO → .65 LHC4ACCHO + .65 HO_2 + .35 DB1O2	`KDEC*(1.-0.37)`	Sander et al. (2018), Paulot et al. (2009a), Francisco-Marquez et al. (2003)
G45127a	TrGC	LISOPACO → 3METHYLFURAN + HO_2	`KDEC*0.37`	Sander et al. (2018), Paulot et al. (2009a), Francisco-Marquez et al. (2003)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G45127b	TrGC	LISOPACO → .65 LHC4ACCHO + .65 HO_2 + .35 DB1O2	`KDEC*(1.-0.37)`	Sander et al. (2018), Paulot et al. (2009a), Francisco-Marquez et al. (2003)
G45128	TrGC	3METHYLFURAN + OH → L3METHYLFURANO2	`3.2E-11*EXP(310./temp)`	Sander et al. (2018)*
G45129	TrGCN	3METHYLFURAN + NO_3 → L3METHYLFURANO2 + NO_2	`1.9E-11`	Sander et al. (2018), Atkinson et al. (2006)*
G45130	TrGC	L3METHYLFURANO2 → C4MDIAL + HO_2	`k1_RO2sOR02`	Sander et al. (2018)
G45131	TrGCN	L3METHYLFURANO2 + NO → C4MDIAL + HO_2 + NO_2	`KRO2NO`	Sander et al. (2018)*
G45132	TrGC	L3METHYLFURANO2 + HO_2 → C4MDIAL + HO_2	`KRO2HO2(5)`	Sander et al. (2018)*
G45133	TrGC	LZCO3C23DBCOD → .62 EZCH3CO2CHCHO + .38 EZCHOCCH3CHO2 + CO_2	`k1_RO2RCO3`	Sander et al. (2018)
G45134a	TrGC	LZCO3C23DBCOD + HO_2 → .62 EZCH3CO2CHCHO + .38 EZCHOCCH3CHO2 + CO_2 + OH	`KAPHO2*rco3_oh`	Sander et al. (2018)
G45134b	TrGC	LZCO3C23DBCOD + HO_2 → LZCO3HC23DBCOD	`KAPHO2*(rco3_ooh+rco3_o3)`	Sander et al. (2018)*
G45135	TrGCN	LZCO3C23DBCOD + NO → .62 EZCH3CO2CHCHO + .38 EZCHOCCH3CHO2 + CO_2 + NO_2	`KAPNO`	Sander et al. (2018)
G45136	TrGCN	LZCO3C23DBCOD + NO_2 → LZCPANC23DBCOD	`k_CH3CO3_NO2`	Rickard and Pascoe (2009)
G45137	TrGCN	LZCO3C23DBCOD + NO_3 → .62 EZCH3CO2CHCHO + .38 EZCHOCCH3CHO2 + CO_2 + NO_2	`KRO2NO3*1.74`	Sander et al. (2018)
G45138	TrGCN	LZCPANC23DBCOD → LZCO3C23DBCOD + NO_2	`k_PAN_M`	Rickard and Pascoe (2009)
G45139	TrGCN	LZCPANC23DBCOD + OH → .62 EZCH3CO2CHCHO + .38 EZCHOCCH3CHO2 + CO_2 + NO_2	`2.52E-11`	Sander et al. (2018)*
G45200	TrGTerC	C511O2 → $CH_3C(O)$ + HCOCH2CHO	`k1_RO2sOR02`	Rickard and Pascoe (2009)
G45201	TrGTerCN	C511O2 + NO → $CH_3C(O)$ + HCOCH2CHO + NO_2	`KRO2NO`	Rickard and Pascoe (2009)*
G45202a	TrGTerC	C511O2 + HO_2 → C511OOH	`KRO2HO2(5)*rcoch2o2_ooh`	Rickard and Pascoe (2009), Sander et al. (2018)
G45202b	TrGTerC	C511O2 + HO_2 → $CH_3C(O)$ + HCOCH2CHO + OH	`KRO2HO2(5)*rcoch2o2_oh`	Rickard and Pascoe (2009), Sander et al. (2018)
G45203	TrGTerC	C511OOH + OH → C511O2	`7.49E-11`	Rickard and Pascoe (2009)
G45204	TrGTerC	CO23C4CHO + OH → CO23C4CO3	`6.65E-11`	Rickard and Pascoe (2009)
G45205	TrGTerCN	CO23C4CHO + NO_3 → CO23C4CO3 + HNO_3	`KNO3AL*5.5`	Rickard and Pascoe (2009)
G45206	TrGTerC	CO23C4CO3 → $CH_3COCOCH_2O_2$ + CO_2	`k1_RO2RCO3`	Rickard and Pascoe (2009)
G45207	TrGTerCN	CO23C4CO3 + NO → $CH_3COCOCH_2O_2$ + CO_2 + NO_2	`KAPNO`	Rickard and Pascoe (2009)*
G45208	TrGTerCN	CO23C4CO3 + NO_2 → C5PAN9	`k_CH3CO3_NO2`	Rickard and Pascoe (2009)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G45209a	TrGTerC	CO23C4CO3 + HO_2 → CO23C4CO3H	KAPHO2*(rco3_ooh+rco3_o3)	Rickard and Pascoe (2009)
G45209b	TrGTerC	CO23C4CO3 + HO_2 → $CH_3COCOCH_2O_2$ + CO_2 + OH	KAPHO2*rco3_oh	Rickard and Pascoe (2009)
G45210	TrGTerCN	C5PAN9 → CO23C4CO3 + NO_2	k_PAN_M	Rickard and Pascoe (2009)
G45211	TrGTerCN	C5PAN9 + OH → $CH_3COCOCHO$ + CO + NO_2	3.12E-13	Rickard and Pascoe (2009)
G45212	TrGTerC	C512O2 → C513O2	k1_RO2pRO2	Rickard and Pascoe (2009)
G45213	TrGTerC	C512O2 + HO_2 → C512OOH	KRO2HO2(5)	Rickard and Pascoe (2009)
G45214	TrGTerCN	C512O2 + NO → C513O2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G45215	TrGTerC	C512OOH + OH → CO13C4CHO + OH	1.01E-10	Rickard and Pascoe (2009)
G45216	TrGTerC	C513O2 → GLYOX + $HOC_2H_4CO_3$	k1_RO2sORO2	Rickard and Pascoe (2009)
G45217	TrGTerCN	C513O2 + NO → GLYOX + $HOC_2H_4CO_3$ + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G45218a	TrGTerC	C513O2 + HO_2 → C513OOH	KRO2HO2(5)*rcoch2o2_ooh	Rickard and Pascoe (2009), Sander et al. (2018)
G45218b	TrGTerC	C513O2 + HO_2 → GLYOX + $HOC_2H_4CO_3$ + OH	KRO2HO2(5)*rcoch2o2_oh	Rickard and Pascoe (2009), Sander et al. (2018)
G45219	TrGTerC	CO13C4CHO + OH → CHOC3COCO3	1.33E-10	Rickard and Pascoe (2009)
G45220	TrGTerCN	CO13C4CHO + NO_3 → CHOC3COCO3 + HNO_3	2.*KNO3AL*5.5	Rickard and Pascoe (2009)
G45221	TrGTerC	C513OOH + OH → C513CO + OH	9.23E-11	Rickard and Pascoe (2009)
G45222	TrGTerC	CHOC3COCO3 → CHOC3COO2 + CO_2	k1_RO2RCO3	Rickard and Pascoe (2009)
G45223	TrGTerC	CHOC3COCO3 + HO_2 → CHOC3COOOH	KAPHO2	Rickard and Pascoe (2009)
G45224	TrGTerCN	CHOC3COCO3 + NO_2 → CHOC3COPAN	k_CH3CO3_NO2	Rickard and Pascoe (2009)
G45225	TrGTerCN	CHOC3COCO3 + NO → CHOC3COO2 + CO_2 + NO_2	KAPNO	Rickard and Pascoe (2009)*
G45226	TrGTerC	C513CO + OH → $HOC_2H_4CO_3$ + CO + CO	2.64E-11	Rickard and Pascoe (2009)
G45227	TrGTerC	C514O2 + HO_2 → C514OOH	KRO2HO2(5)	Rickard and Pascoe (2009)
G45228a	TrGTerCN	C514O2 + NO → CO13C4CHO + HO_2 + NO_2	KRO2NO*(1.-alpha_AN(7,2,0,1,0, temp,cair))	Rickard and Pascoe (2009), Sander et al. (2018)
G45228b	TrGTerCN	C514O2 + NO → C514NO3	KRO2NO*alpha_AN(7,2,0,1,0,temp, cair)	Rickard and Pascoe (2009), Sander et al. (2018)
G45229	TrGTerCN	C514O2 + NO_3 → CO13C4CHO + HO_2 + NO_2	KRO2NO3	Rickard and Pascoe (2009)
G45230	TrGTerC	C514O2 → CO13C4CHO + HO_2	k1_RO2sRO2	Rickard and Pascoe (2009)
G45231	TrGTerC	C514OOH + OH → CO13C4CHO + OH	1.10E-10	Rickard and Pascoe (2009)
G45232	TrGTerCN	C514NO3 + OH → CO13C4CHO + NO_2	4.33E-11	Rickard and Pascoe (2009)
G45233	TrGTerC	CHOC3COOOH + OH → CHOC3COCO3	7.55E-11	Rickard and Pascoe (2009)
G45234	TrGTerCN	CHOC3COPAN → CHOC3COCO3 + NO_2	k_PAN_M	Rickard and Pascoe (2009)
G45235	TrGTerCN	CHOC3COPAN + OH → C4CODIAL + CO + NO_2	7.19E-11	Rickard and Pascoe (2009)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G45236	TrGTerC	MBO + OH → LMBOABO2	8.1E-12*EXP(610./TEMP)	Rickard and Pascoe (2009), Sander et al. (2018)*
G45237a	TrGTerC	MBO + O_3 → HCHO + .16 CH_3COCH_3 + .16 HO_2 + .16 CO + .16 OH + .84 MBOOO	1.0E-17*0.57	Rickard and Pascoe (2009), Sander et al. (2018)
G45237b	TrGTerC	MBO + O_3 → IBUTALOH + .63 CO + .37 $HOCH_2OOH$ + .16 OH + .16 HO_2	1.0E-17*0.43	Rickard and Pascoe (2009), Sander et al. (2018)
G45238	TrGTerCN	MBO + NO_3 → LNMBOABO2	4.6E-14*EXP(-400./TEMP)	Rickard and Pascoe (2009), Sander et al. (2018)
G45239	TrGTerC	LMBOABO2 + HO_2 → LMBOABOOH	KRO2HO2(5)	Rickard and Pascoe (2009), Sander et al. (2018)
G45240a	TrGTerCN	LMBOABO2 + NO → LMBOABNO3	KRO2NO*(.67*alpha_AN(7,2,0,0,0, temp,cair)+.33*alpha_AN(7,1,0,0, 0,temp,cair))	Rickard and Pascoe (2009), Sander et al. (2018)
G45240b	TrGTerCN	LMBOABO2 + NO → $HOCH_2CHO$ + CH_3COCH_3 + HO_2 + NO_2	KRO2NO*(1.-(.67*alpha_AN(7,2,0, 0,0,temp,cair)+.33*alpha_AN(7,1, 0,0,0,temp,cair)))*.67	Rickard and Pascoe (2009), Sander et al. (2018)
G45240c	TrGTerCN	LMBOABO2 + NO → IBUTALOH + HCHO + HO_2 + NO_2	KRO2NO*(1.-(.67*alpha_AN(7,2,0, 0,0,temp,cair)+.33*alpha_AN(7,1, 0,0,0,temp,cair)))*.33	Rickard and Pascoe (2009), Sander et al. (2018)
G45241a	TrGTerC	LMBOABO2 → $HOCH_2CHO$ + CH_3COCH_3 + HO_2	k1_RO2sORO2*.67	Rickard and Pascoe (2009), Sander et al. (2018)
G45241b	TrGTerC	LMBOABO2 → IBUTALOH + HCHO + HO_2	k1_RO2pORO2*.33	Rickard and Pascoe (2009), Sander et al. (2018)
G45242a	TrGTerC	LMBOABOOH + OH → MBOACO	0.67*2.93E-11+.33*2.05E-12	Rickard and Pascoe (2009), Sander et al. (2018)
G45242b	TrGTerC	LMBOABOOH + OH → LMBOABO2	k_roohro	Rickard and Pascoe (2009), Sander et al. (2018)
G45243	TrGTerCN	LMBOABNO3 + OH → MBOACO + NO_2	0.67*1.75E-12+.33*2.69E-12	Rickard and Pascoe (2009), Sander et al. (2018)
G45244	TrGTerC	MBOACO + OH → MBOCOCO + HO_2	3.79E-12	Rickard and Pascoe (2009)
G45245	TrGTerC	MBOCOCO + OH → CO + IPRHOCO3	1.38E-11	Rickard and Pascoe (2009)
G45246	TrGTerCN	LNMBOABO2 + HO_2 → LNMBOABOOH	KRO2HO2(5)	Rickard and Pascoe (2009), Sander et al. (2018)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G45247	TrGTerCN	LNMBOABO2 + NO → .65 NO_3CH2CHO + .65 CH_3COCH_3 + .65 HO_2 + .35 IBUTALOH + .35 HCHO + .35 NO_2 + NO_2	`KRO2NO`	Rickard and Pascoe (2009), Sander et al. (2018)*
G45248	TrGTerCN	LNMBOABO2 + NO_3 → .65 NO_3CH2CHO + .65 CH_3COCH_3 + .65 HO_2 + .35 IBUTALOH + .35 HCHO + .35 NO_2 + NO_2	`KRO2NO3`	Rickard and Pascoe (2009), Sander et al. (2018)
G45249	TrGTerCN	LNMBOABO2 → .65 NO_3CH2CHO + .65 CH_3COCH_3 + .65 HO_2 + .35 IBUTALOH + .35 HCHO + .35 NO_2	`k1_RO2sORO2`	Rickard and Pascoe (2009), Sander et al. (2018)
G45250a	TrGTerCN	LNMBOABOOH + OH → .65 C4MCONO3OH + .35 NMBOBCO	`0.65*4.89E-12+.35*2.52E-12`	Rickard and Pascoe (2009), Sander et al. (2018)
G45250b	TrGTerCN	LNMBOABOOH + OH → LNMBOABO2	`k_roohro`	Rickard and Pascoe (2009), Sander et al. (2018)
G45251	TrGTerCN	NMBOBCO + OH → NC4OHCO3	`4.26E-12`	Rickard and Pascoe (2009)
G45252a	TrGTerCN	NC4OHCO3 + HO_2 → IBUTALOH + CO_2 + NO_2 + OH	`KAPHO2*rco3_oh`	Rickard and Pascoe (2009), Sander et al. (2018)
G45252b	TrGTerCN	NC4OHCO3 + HO_2 → NC4OHCO3H	`KAPHO2*(rco3_o3+rco3_ooh)`	Rickard and Pascoe (2009), Sander et al. (2018)
G45253	TrGTerCN	NC4OHCO3 + NO → IBUTALOH + CO_2 + NO_2 + NO_2	`KAPNO`	Rickard and Pascoe (2009)
G45254	TrGTerCN	NC4OHCO3 + NO_2 → NC4OHCPAN	`k_CH3CO3_NO2`	Rickard and Pascoe (2009)
G45255	TrGTerCN	NC4OHCO3 + NO_3 → IBUTALOH + CO_2 + NO_2 + NO_2	`KRO2NO3*1.74`	Rickard and Pascoe (2009)
G45256	TrGTerCN	NC4OHCO3 → IBUTALOH + CO_2 + NO_2	`k1_RO2RCO3`	Rickard and Pascoe (2009)
G45257	TrGTerCN	NC4OHCO3H + OH → NC4OHCO3	`4.50E-12`	Rickard and Pascoe (2009)
G45258	TrGTerCN	NC4OHCPAN + OH → IBUTALOH + CO + NO_2 + NO_2	`1.27E-12`	Rickard and Pascoe (2009)
G45259	TrGTerCN	NC4OHCPAN → NC4OHCO3 + NO_2	`K_PAN_M`	Rickard and Pascoe (2009)
G45260	TrGTerCN	C4MCONO3OH + OH → CH_3COCH_3 + HCHO + CO_2 + NO_2	`1.23E-12`	Rickard and Pascoe (2009), Sander et al. (2018)
G45400	TrGAroCN	NC4MDCO2HN + OH → MMALANHY + NO_2	`k_roohro`	Rickard and Pascoe (2009)*
G45401	TrGAroCN	C54CO + NO_3 → 3 CO + $CH_3C(O)OO$ + HNO_3	`KNO3AL*5.5`	Rickard and Pascoe (2009)
G45402	TrGAroC	C54CO + OH → 3 CO + $CH_3C(O)OO$	`1.72E-11`	Rickard and Pascoe (2009)
G45403a	TrGAroCN	NTLFUO2 + HO_2 → NTLFUOOH	`KRO2HO2(5)*(1-rcoch2o2_oh)`	Rickard and Pascoe (2009)
G45403b	TrGAroCN	NTLFUO2 + HO_2 → ACCOMECHO + NO_2 + OH	`KRO2HO2(5)*rcoch2o2_oh`	Rickard and Pascoe (2009)
G45404	TrGAroCN	NTLFUO2 + NO → ACCOMECHO + NO_2 + NO_2	`KRO2NO`	Rickard and Pascoe (2009)*
G45405	TrGAroCN	NTLFUO2 + NO_3 → ACCOMECHO + NO_2 + NO_2	`KRO2NO3`	Rickard and Pascoe (2009)*
G45406	TrGAroCN	NTLFUO2 → ACCOMECHO + NO_2	`k1_RO2tORO2`	Rickard and Pascoe (2009)*
G45407	TrGAroC	C5134CO2OH + OH → C54CO + HO_2	`7.48E-11`	Rickard and Pascoe (2009)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G45408	TrGAroCN	C5COO2NO2 + OH → MGLYOX + CO + CO + NO_2	5.43E-11	Rickard and Pascoe (2009)
G45409	TrGAroCN	C5COO2NO2 → C5CO14O2 + NO_2	k_PAN_M	Rickard and Pascoe (2009)*
G45410	TrGAroC	C5DIALOOH + OH → C5DIALCO + OH	7.52E-11	Rickard and Pascoe (2009)
G45411a	TrGAroC	C4CO2DBCO3 + HO_2 → C4CO2DCO3H	KAPHO2*(rco3_ooh+rco3_o3)	Rickard and Pascoe (2009)
G45411b	TrGAroC	C4CO2DBCO3 + HO_2 → HO_2 + CO + HCOCOCHO + CO_2 + OH	KAPHO2*rco3_oh	Rickard and Pascoe (2009), Sander et al. (2018)
G45412	TrGAroCN	C4CO2DBCO3 + NO → HO_2 + CO + HCOCOCHO + CO_2 + NO_2	KAPNO	Rickard and Pascoe (2009)
G45413	TrGAroCN	C4CO2DBCO3 + NO_2 → C4CO2DBPAN	k_CH3CO3_NO2	Rickard and Pascoe (2009)*
G45414	TrGAroCN	C4CO2DBCO3 + NO_3 → HO_2 + CO + HCOCOCHO + CO_2 + NO_2	KRO2NO3*1.74	Rickard and Pascoe (2009)
G45415	TrGAroC	C4CO2DBCO3 → HO_2 + CO + HCOCOCHO + CO_2	k1_RO2RCO3	Rickard and Pascoe (2009)
G45416	TrGAroC	MMALANHY + OH → MMALANHYO2	1.50E-12	Rickard and Pascoe (2009)
G45421a	TrGAroC	MMALANHYO2 + HO_2 → MMALNHYOOH	KRO2HO2(5)*(1-rcoch2o2_oh-rchohch2o2_oh)	Rickard and Pascoe (2009), Sander et al. (2018)
G45421b	TrGAroC	MMALANHYO2 + HO_2 → CO2H3CO3 + CO_2 + OH	KRO2HO2(5)*(rcoch2o2_oh+rchohch2o2_oh)	Rickard and Pascoe (2009), Sander et al. (2018)
G45422	TrGAroCN	MMALANHYO2 + NO → CO2H3CO3 + CO_2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G45423	TrGAroCN	MMALANHYO2 + NO_3 → CO2H3CO3 + CO_2 + NO_2	KRO2NO3	Rickard and Pascoe (2009)*
G45424	TrGAroC	MMALANHYO2 → CO2H3CO3 + CO_2	k1_RO2tORO2	Rickard and Pascoe (2009)*
G45428	TrGAroCN	C4CO2DBPAN + OH → HCOCOCHO + CO_2 + CO + NO_2	2.74E-11	Rickard and Pascoe (2009)
G45429	TrGAroCN	C4CO2DBPAN → C4CO2DBCO3 + NO_2	k_PAN_M	Rickard and Pascoe (2009)*
G45430a	TrGAroC	C5CO14O2 + HO_2 → .83 MALANHY + .83 CH_3 + .17 MGLYOX + .17 HO_2 + .17 CO + .17 CO_2 + OH	KAPHO2*rco3_oh	Rickard and Pascoe (2009)*
G45430b	TrGAroC	C5CO14O2 + HO_2 → C5CO14OH + O_3	KAPHO2*rco3_o3	Rickard and Pascoe (2009)
G45430c	TrGAroC	C5CO14O2 + HO_2 → C5CO14OOH	KAPHO2*rco3_ooh	Rickard and Pascoe (2009)
G45431	TrGAroCN	C5CO14O2 + NO → .83 MALANHY + .83 CH_3 + .17 MGLYOX + .17 HO_2 + .17 CO + .17 CO_2 + NO_2	KAPNO	Rickard and Pascoe (2009)*
G45432	TrGAroCN	C5CO14O2 + NO_2 → C5COO2NO2	k_CH3CO3_NO2	Rickard and Pascoe (2009)*
G45433	TrGAroCN	C5CO14O2 + NO_3 → .83 MALANHY + .83 CH_3 + .17 MGLYOX + .17 HO_2 + .17 CO + .17 CO_2 + NO_2	KRO2NO3*1.74	Rickard and Pascoe (2009)*
G45434	TrGAroC	C5CO14O2 → .83 MALANHY + .83 CH_3 + .17 MGLYOX + .17 HO_2 + .17 CO + .17 CO_2	k1_RO2RCO3	Rickard and Pascoe (2009)*

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G45436	TrGAroC	C5CO14OH + OH → .83 MALANHY + .83 CH_3 + .17 MGLYOX + .17 HO_2 + .17 CO + .17 CO_2	5.44E-11	Rickard and Pascoe (2009)*
G45441	TrGAroCN	C5DICARB + NO_3 → C5CO14O2 + HNO_3	KNO3AL*2.75	Rickard and Pascoe (2009)
G45442	TrGAroC	C5DICARB + O_3 → .5338 GLYOX + .063 CH_3CHO + .348 $CH_3C(O)OO$ + .918 CO + .57 OH + .473 HO_2 + .0563 CH_3COCO_2H + .5338 MGLYOX + .676 H_2O_2 + .063 HCHO + .0563 $HCOCO_2H$ + .2465 CO_2	2.00E-18	Rickard and Pascoe (2009)
G45443	TrGAroC	C5DICARB + OH → .48 C5CO14O2 + .52 C5DICARBO2	6.2E-11	Rickard and Pascoe (2009)
G45444	TrGAroC	MC3ODBCO2H + OH → .35 GLYOX + .35 CH_3 + .35 CO + .35 CO_2 + .65 MMALANHY + .65 HO_2	4.38E-11	Rickard and Pascoe (2009)*
G45451	TrGAroCN	TLFUONE + NO_3 → NTLFUO2	1.00E-12	Rickard and Pascoe (2009)
G45452	TrGAroC	TLFUONE + O_3 → .5 CO + .5 OH + .5 MECOACETO2 + .3125 C24O3CCO2H + .1875 ACCOMECHO + .1875 H_2O_2	8.00E-19	see note*
G45453	TrGAroC	TLFUONE + OH → TLFUO2	6.90E-11	Rickard and Pascoe (2009)
G45454a	TrGAroC	ACCOMECO3 + HO_2 → ACCOMECO3H	KAPHO2*(rco3_ooh+rco3_o3)	Rickard and Pascoe (2009)
G45454b	TrGAroC	ACCOMECO3 + HO_2 → MECOACETO2 + CO_2 + OH	KAPHO2*rco3_oh	Rickard and Pascoe (2009)
G45455	TrGAroCN	ACCOMECO3 + NO → MECOACETO2 + CO_2 + NO_2	KAPNO	Rickard and Pascoe (2009)
G45456	TrGAroCN	ACCOMECO3 + NO_2 → ACCOMEPAN	k_CH3CO3_NO2	Rickard and Pascoe (2009)*
G45457	TrGAroCN	ACCOMECO3 + NO_3 → MECOACETO2 + CO_2 + NO_2	KRO2NO3*1.74	Rickard and Pascoe (2009)
G45458	TrGAroC	ACCOMECO3 → MECOACETO2 + CO_2	k1_RO2RCO3	Rickard and Pascoe (2009)
G45459	TrGAroC	C4CO2DCO3H + OH → C4CO2DBCO3	3.06E-11	Rickard and Pascoe (2009)
G45464	TrGAroCN	ACCOMECHO + NO_3 → ACCOMECO3 + HNO_3	KNO3AL*5.5	Rickard and Pascoe (2009)
G45465	TrGAroC	ACCOMECHO + OH → ACCOMECO3	7.09E-11	Rickard and Pascoe (2009)
G45466	TrGAroC	MMALNHYOOH + OH → MMALANHYO2	1.69E-11	Rickard and Pascoe (2009)
G45467a	TrGAroC	C5DICAROOH + OH → C5134CO2OH + OH	1.21E-10	Rickard and Pascoe (2009)
G45467b	TrGAroC	C5DICAROOH + OH → C5DICARBO2	k_roohro	Rickard and Pascoe (2009)
G45468	TrGAroC	C24O3CCO2H + OH → MECOACETO2 + CO_2	8.76E-13	Rickard and Pascoe (2009)
G45469	TrGAroCN	NTLFUOOH + OH → NTLFUO2	4.44E-12	Rickard and Pascoe (2009)
G45470	TrGAroCN	ACCOMEPAN + OH → METACETHO + CO + CO + NO_2	1.00E-14	Rickard and Pascoe (2009)
G45471	TrGAroCN	ACCOMEPAN → ACCOMECO3 + NO_2	k_PAN_M	Rickard and Pascoe (2009)
G45476a	TrGAroC	TLFUO2 + HO_2 → TLFUOOH	KRO2HO2(5)*(1-rcoch2o2_oh-rchohch2o2_oh)	Rickard and Pascoe (2009)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G45476b	TrGAroC	TLFUO2 + HO_2 → ACCOMECHO + HO_2 + OH	`KRO2HO2(5)*(rcoch2o2_oh+rchohch2o2_oh)`	Rickard and Pascoe (2009)*
G45477	TrGAroCN	TLFUO2 + NO → ACCOMECHO + HO_2 + NO_2	`KRO2NO`	Rickard and Pascoe (2009)*
G45478	TrGAroCN	TLFUO2 + NO_3 → ACCOMECHO + HO_2 + NO_2	`KRO2NO3`	Rickard and Pascoe (2009)*
G45479	TrGAroC	TLFUO2 → ACCOMECHO + HO_2	`k1_RO2tORO2`	Rickard and Pascoe (2009)*
G45480	TrGAroC	C5CO14OOH + OH → C5CO14O2	`3.59E-12`	Rickard and Pascoe (2009)
G45483	TrGAroC	TLFUOOH + OH → TLFUO2	`2.53E-11`	Rickard and Pascoe (2009)
G45485	TrGAroC	ACCOMECO3H + OH → ACCOMECO3	`3.59E-12`	Rickard and Pascoe (2009)
G45486a	TrGAroC	C5DIALO2 + HO_2 → C5DIALOOH	`KRO2HO2(5)*(1-rcoch2o2_oh)`	Rickard and Pascoe (2009)
G45486b	TrGAroC	C5DIALO2 + HO_2 → MALDIAL + CO + HO_2 + OH	`KRO2HO2(5)*rcoch2o2_oh`	Rickard and Pascoe (2009)*
G45487	TrGAroCN	C5DIALO2 + NO → MALDIAL + CO + HO_2 + NO_2	`KRO2NO`	Rickard and Pascoe (2009)*
G45488	TrGAroCN	C5DIALO2 + NO_3 → MALDIAL + CO + HO_2 + NO_2	`KRO2NO3`	Rickard and Pascoe (2009)*
G45489	TrGAroC	C5DIALO2 → MALDIAL + CO + HO_2	`k1_RO2sORO2`	Rickard and Pascoe (2009)*
G45490a	TrGAroC	C5DICARBO2 + HO_2 → C5DICAROOH	`KRO2HO2(5)*(rco3_ooh+rco3_o3)`	Rickard and Pascoe (2009)
G45491b	TrGAroC	C5DICARBO2 + HO_2 → MGLYOX + GLYOX + HO_2 + OH	`KRO2HO2(5)*rco3_oh`	Rickard and Pascoe (2009)*
G45492	TrGAroCN	C5DICARBO2 + NO → MGLYOX + GLYOX + HO_2 + NO_2	`KRO2NO`	Rickard and Pascoe (2009)*
G45493	TrGAroCN	C5DICARBO2 + NO_3 → MGLYOX + GLYOX + HO_2 + NO_2	`KRO2NO3`	Rickard and Pascoe (2009)*
G45494	TrGAroC	C5DICARBO2 → MGLYOX + GLYOX + HO_2	`k1_RO2sORO2`	Rickard and Pascoe (2009)*
G46200a	TrGTerC	CO235C6O2 + HO_2 → CO235C6OOH	`KRO2HO2(6)*rcoch2o2_ooh`	Rickard and Pascoe (2009), Sander et al. (2018)
G46200b	TrGTerC	CO235C6O2 + HO_2 → CO23C4CO3 + HCHO + OH	`KRO2HO2(6)*rcoch2o2_oh`	Rickard and Pascoe (2009), Sander et al. (2018)
G46201	TrGTerCN	CO235C6O2 + NO → CO23C4CO3 + HCHO + NO_2	`KRO2NO`	Rickard and Pascoe (2009)*
G46202	TrGTerC	CO235C6O2 → CO23C4CO3 + HCHO	`k1_RO2pORO2`	Rickard and Pascoe (2009)
G46203	TrGTerC	CO235C6OOH + OH → CO235C6O2	`1.01E-11`	Rickard and Pascoe (2009)
G46204	TrGTerC	C614O2 → CO23C4CHO + HCHO + HO_2	`k1_RO2sORO2`	Rickard and Pascoe (2009)
G46205a	TrGTerCN	C614O2 + NO → CO23C4CHO + HCHO + HO_2 + NO_2	`KRO2NO*(1.-alpha_AN(9,2,0,1,0,temp,cair))`	Rickard and Pascoe (2009)
G46205b	TrGTerCN	C614O2 + NO → C614NO3	`KRO2NO*alpha_AN(9,2,0,1,0,temp,cair)`	Rickard and Pascoe (2009)
G46206a	TrGTerC	C614O2 + HO_2 → C614OOH	`KRO2HO2(6)*(1.-rchohch2o2_oh)`	Rickard and Pascoe (2009), Sander et al. (2018)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G46206b	TrGTerC	C614O2 + HO_2 → CO23C4CHO + HCHO + HO_2 + OH	KRO2HO2(6)*rchohch2o2_oh	Rickard and Pascoe (2009), Sander et al. (2018)
G46207	TrGTerCN	C614NO3 + OH → C614CO + NO_2	7.11E-12	Rickard and Pascoe (2009)
G46208	TrGTerC	C614OOH + OH → C614CO + OH	8.69E-11	Rickard and Pascoe (2009)
G46209	TrGTerC	C614CO + OH → CO235C5CHO + HO_2	3.22E-12	Rickard and Pascoe (2009)
G46210	TrGTerC	CO235C5CHO + OH → CO23C4CO3 + CO	1.33E-11	Rickard and Pascoe (2009)
G46211	TrGTerCN	CO235C5CHO + NO_3 → CO23C4CO3 + CO + HNO_3	KNO3AL*5.5	Rickard and Pascoe (2009)
G46400	TrGAroC	PHENOOH + OH → PHENO2	1.16E-10	Rickard and Pascoe (2009)
G46401	TrGAroC	C6CO4DB + OH → CO + CO + HO_2 + CO + HCOCOCHO	7.70E-11	Rickard and Pascoe (2009)
G46402	TrGAroC	C5CO2DCO3H + OH → C5CO2DBCO3	3.60E-11	Rickard and Pascoe (2009)
G46403	TrGAroCN	NDNPHENOOH + OH → NDNPHENO2	k_roohro	Rickard and Pascoe (2009)
G46404a	TrGAroC	C615CO2O2 + HO_2 → C615CO2OOH	KRO2HO2(6)*(1.-rcoch2o2_oh)	Rickard and Pascoe (2009)
G46404b	TrGAroC	C615CO2O2 + HO_2 → C5DICARB + CO + HO_2 + OH	KRO2HO2(6)*rcoch2o2_oh	Rickard and Pascoe (2009)*
G46405	TrGAroCN	C615CO2O2 + NO → C5DICARB + CO + HO_2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G46406	TrGAroCN	C615CO2O2 + NO_3 → C5DICARB + CO + HO_2 + NO_2	KRO2NO3	Rickard and Pascoe (2009)*
G46407	TrGAroC	C615CO2O2 → C5DICARB + CO + HO_2	k1_RO2sOR02	Rickard and Pascoe (2009)*
G46408	TrGAroCN	BZEMUCPAN + OH → MALDIAL + CO + CO_2 + NO_2	4.05E-11	Rickard and Pascoe (2009)
G46409	TrGAroCN	BZEMUCPAN → BZEMUCCO3 + NO_2	k_PAN_M	Rickard and Pascoe (2009)
G46410	TrGAroCN	BZBIPERNO3 + OH → BZOBIPEROH + NO_2	7.30E-11	Rickard and Pascoe (2009)
G46411	TrGAroCN	HOC6H4NO2 + NO_3 → NPHEN1O + HNO_3	9.00E-14	Rickard and Pascoe (2009)
G46412	TrGAroCN	HOC6H4NO2 + OH → NPHEN1O	9.00E-13	Rickard and Pascoe (2009)
G46413a	TrGAroCN	NDNPHENO2 + HO_2 → NDNPHENOOH	KRO2HO2(6)*(1-rchohch2o2_oh)	Rickard and Pascoe (2009)
G46413b	TrGAroCN	NDNPHENO2 + HO_2 → NC4DCO2H + HNO_3 + CO + CO + NO_2 + OH	KRO2HO2(6)*rchohch2o2_oh	Rickard and Pascoe (2009)*
G46414	TrGAroCN	NDNPHENO2 + NO → NC4DCO2H + HNO_3 + CO + CO + NO_2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G46415	TrGAroCN	NDNPHENO2 + NO_3 → NC4DCO2H + HNO_3 + CO + CO + NO_2 + NO_2	KRO2NO3	Rickard and Pascoe (2009)*
G46416	TrGAroCN	NDNPHENO2 → NC4DCO2H + HNO_3 + CO + CO + NO_2	k1_RO2ISOPDO2	Rickard and Pascoe (2009)*
G46417	TrGAroC	PBZQCO + OH → C5CO2OHCO3	6.07E-11	Rickard and Pascoe (2009)
G46418	TrGAroCN	CATECHOL + NO_3 → CATEC1O + HNO_3	9.9E-11	Rickard and Pascoe (2009)*
G46419	TrGAroC	CATECHOL + O_3 → MALDALCO2H + $HCOCO_2H$ + HO_2 + OH	9.2E-18	Rickard and Pascoe (2009)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G46420	TrGAroC	CATECHOL + OH → CATEC1O	1.0E-10	Rickard and Pascoe (2009)
G46421	TrGAroC	C5COOHCO3H + OH → C5CO2OHCO3	8.01E-11	Rickard and Pascoe (2009)
G46422	TrGAroCN	NCATECHOL + NO_3 → NNCATECO2	2.60E-12	Rickard and Pascoe (2009)
G46423	TrGAroCN	NCATECHOL + OH → NCATECO2	3.47E-12	Rickard and Pascoe (2009)
G46424a	TrGAroC	C5CO2OHCO3 + HO_2 → C5COOHCO3H	KAPHO2*(rco3_ooh+rco3_o3)	Rickard and Pascoe (2009)
G46424b	TrGAroC	C5CO2OHCO3 + HO_2 → HOCOC4DIAL + HO_2 + CO + CO_2 + OH	KAPHO2*rco3_oh	Rickard and Pascoe (2009)
G46425	TrGAroCN	C5CO2OHCO3 + NO → HOCOC4DIAL + HO_2 + CO + CO_2 + NO_2	KAPNO	Rickard and Pascoe (2009)
G46426	TrGAroCN	C5CO2OHCO3 + NO_2 → C5CO2OHPAN	k_CH3CO3_NO2	Rickard and Pascoe (2009)*
G46427	TrGAroCN	C5CO2OHCO3 + NO_3 → HOCOC4DIAL + HO_2 + CO + CO_2 + NO_2	KRO2NO3*1.74	Rickard and Pascoe (2009)
G46428	TrGAroC	C5CO2OHCO3 → HOCOC4DIAL + HO_2 + CO + CO_2	k1_RO2RCO3	Rickard and Pascoe (2009)
G46429	TrGAroCN	BZEPOXMUC + NO_3 → BZEMUCCO3 + HNO_3	2*KNO3AL*2.75	Rickard and Pascoe (2009)
G46430	TrGAroC	BZEPOXMUC + O_3 → EPXC4DIAL + .125 HCHO + .1125 $HCOCO_2H$ + .0675 GLYOX + .0675 H_2O_2 + .82 HO_2 + .57 OH + 1.265 CO + .25 CO_2	2.00E-18	Rickard and Pascoe (2009)*
G46431	TrGAroC	BZEPOXMUC + OH → .31 BZEMUCCO3 + .69 BZEMUCO2	6.08E-11	Rickard and Pascoe (2009)
G46432a	TrGAroCN	NCATECO2 + HO_2 → NCATECOOH	KRO2HO2(6)*(1-rchohch2o2_oh)	Rickard and Pascoe (2009)
G46432b	TrGAroCN	NCATECO2 + HO_2 → NC4DCO2H + $HCOCO_2H$ + HO_2 + OH	KRO2HO2(6)*rchohch2o2_oh	Rickard and Pascoe (2009)*
G46433	TrGAroCN	NCATECO2 + NO → NC4DCO2H + $HCOCO_2H$ + HO_2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G46434	TrGAroCN	NCATECO2 + NO_3 → NC4DCO2H + $HCOCO_2H$ + HO_2 + NO_2	KRO2NO3	Rickard and Pascoe (2009)*
G46435	TrGAroCN	NCATECO2 → NC4DCO2H + $HCOCO_2H$ + HO_2	k1_RO2ISOPDO2	Rickard and Pascoe (2009)*
G46436	TrGAroCN	NPHEN1OOH + OH → NPHEN1O2	9.00E-13	Rickard and Pascoe (2009)
G46437a	TrGAroCN	NPHENO2 + HO_2 → NPHENOOH	KRO2HO2(6)*(1-rchohch2o2_oh)	Rickard and Pascoe (2009)
G46437b	TrGAroCN	NPHENO2 + HO_2 → MALDALCO2H + GLYOX + NO_2 + OH	KRO2HO2(6)*rchohch2o2_oh	Rickard and Pascoe (2009)*
G46438	TrGAroCN	NPHENO2 + NO → MALDALCO2H + GLYOX + NO_2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G46439	TrGAroCN	NPHENO2 + NO_3 → MALDALCO2H + GLYOX + NO_2 + NO_2	KRO2NO3	Rickard and Pascoe (2009)*

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G46440	TrGAroCN	NPHENO2 $\rightarrow$ MALDALCO2H + GLYOX + NO_2	k1_RO2ISOPDO2	Rickard and Pascoe (2009)*
G46441	TrGAroC	BENZENE + OH $\rightarrow$ .352 BZBIPERO2 + .118 BZEPOXMUC + .118 HO_2 + .53 PHENOL + .53 HO_2	2.3E-12*EXP(-190/TEMP)	Rickard and Pascoe (2009)*
G46442	TrGAroCN	C5CO2OHPAN + OH $\rightarrow$ HOCOC4DIAL + CO + CO + NO_2	7.66E-11	Rickard and Pascoe (2009)
G46443	TrGAroCN	C5CO2OHPAN $\rightarrow$ C5CO2OHCO3 + NO_2	k_PAN_M	Rickard and Pascoe (2009)
G46444	TrGAroCN	CATEC1O + NO_2 $\rightarrow$ NCATECHOL	k_C6H5O_NO2	Rickard and Pascoe (2009), Platz et al. (1998)
G46445	TrGAroC	CATEC1O + O_3 $\rightarrow$ CATEC1O2	k_C6H5O_O3	Rickard and Pascoe (2009), Tao and Li (1999)
G46446	TrGAroC	BZEMUCCO + OH $\rightarrow$ EPXDLCO3 + GLYOX	9.20E-11	Rickard and Pascoe (2009)
G46447a	TrGAroCN	NNCATECO2 + HO_2 $\rightarrow$ NNCATECOOH	KRO2HO2(6)*(1-rchohch2o2_oh)	Rickard and Pascoe (2009)
G46447b	TrGAroCN	NNCATECO2 + HO_2 $\rightarrow$ NC4DCO2H + $HCOCO_2H$ + NO_2 + OH	KRO2HO2(6)*rchohch2o2_oh	Rickard and Pascoe (2009)*
G46448	TrGAroCN	NNCATECO2 + NO $\rightarrow$ NC4DCO2H + $HCOCO_2H$ + NO_2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G46449	TrGAroCN	NNCATECO2 + NO_3 $\rightarrow$ NC4DCO2H + $HCOCO_2H$ + NO_2 + NO_2	KRO2NO3	Rickard and Pascoe (2009)*
G46450	TrGAroCN	NNCATECO2 $\rightarrow$ NC4DCO2H + $HCOCO_2H$ + NO_2	k1_RO2ISOPDO2	Rickard and Pascoe (2009)*
G46451	TrGAroC	BZEMUCCO2H + OH $\rightarrow$ C5DIALO2 + CO_2	4.06E-11	Rickard and Pascoe (2009)
G46452	TrGAroCN	NNCATECOOH + OH $\rightarrow$ NNCATECO2	k_roohro	Rickard and Pascoe (2009)
G46453	TrGAroCN	NPHEN1O + NO_2 $\rightarrow$ DNPHEN	k_C6H5O_NO2	Rickard and Pascoe (2009), Platz et al. (1998)
G46454	TrGAroCN	NPHEN1O + O_3 $\rightarrow$ NPHEN1O2	k_C6H5O_O3	Rickard and Pascoe (2009), Tao and Li (1999)
G46455	TrGAroCN	DNPHEN + NO_3 $\rightarrow$ NDNPHENO2	2.25E-15	Rickard and Pascoe (2009)
G46456	TrGAroCN	DNPHEN + OH $\rightarrow$ DNPHENO2	3.00E-14	Rickard and Pascoe (2009)
G46457	TrGAroCN	PHENOL + NO_3 $\rightarrow$ .742 C6H5O + .742 HNO_3 + .258 NPHENO2	3.8E-12	Rickard and Pascoe (2009)*
G46458	TrGAroC	PHENOL + OH $\rightarrow$ .06 C6H5O + .8 CATECHOL + .8 HO_2 + .14 PHENO2	4.7E-13*EXP(1220/TEMP)	Rickard and Pascoe (2009)*
G46459	TrGAroCN	PBZQONE + NO_3 $\rightarrow$ NBZQO2	3.00E-13	Rickard and Pascoe (2009)
G46460	TrGAroC	PBZQONE + OH $\rightarrow$ PBZQO2	4.6E-12	Rickard and Pascoe (2009)
G46461a	TrGAroC	PHENO2 + HO_2 $\rightarrow$ PHENOOH	KRO2HO2(6)*(1-rchohch2o2_oh)	Rickard and Pascoe (2009)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G46461b	TrGAroC	PHENO2 + HO_2 → .71 MALDALCO2H + .71 GLYOX + .29 PBZQONE + HO_2 + OH	`KRO2HO2(6)*rchohch2o2_oh`	Rickard and Pascoe (2009)*
G46462	TrGAroCN	PHENO2 + NO → .71 MALDALCO2H + .71 GLYOX + .29 PBZQONE + HO_2 + NO_2	`KRO2NO`	Rickard and Pascoe (2009)*
G46463	TrGAroCN	PHENO2 + NO_3 → .71 MALDALCO2H + .71 GLYOX + .29 PBZQONE + HO_2 + NO_2	`KRO2NO3`	Rickard and Pascoe (2009)*
G46464	TrGAroC	PHENO2 → .71 MALDALCO2H + .71 GLYOX + .29 PBZQONE + HO_2	`k1_RO2ISOPDO2`	Rickard and Pascoe (2009)*
G46465	TrGAroC	C615CO2OOH + OH → C6125CO + OH	`9.42E-11`	Rickard and Pascoe (2009)
G46466a	TrGAroC	C5CO2DBCO3 + HO_2 → C5CO2DCO3H	`KAPHO2*(rco3_ooh+rco3_o3)`	Rickard and Pascoe (2009)
G46466b	TrGAroC	C5CO2DBCO3 + HO_2 → $CH_3C(O)$ + HCOCOCHO + CO_2 + OH	`KAPHO2*rco3_oh`	Rickard and Pascoe (2009)
G46467	TrGAroCN	C5CO2DBCO3 + NO → $CH_3C(O)$ + HCOCOCHO + CO_2 + NO_2	`KAPNO`	Rickard and Pascoe (2009)
G46468	TrGAroCN	C5CO2DBCO3 + NO_2 → C5CO2DBPAN	`k_CH3CO3_NO2`	Rickard and Pascoe (2009)*
G46469	TrGAroCN	C5CO2DBCO3 + NO_3 → $CH_3C(O)$ + HCOCOCHO + CO_2 + NO_2	`KRO2NO3*1.74`	Rickard and Pascoe (2009)
G46470	TrGAroC	C5CO2DBCO3 → $CH_3C(O)$ + HCOCOCHO + CO_2	`k1_RO2RCO3`	Rickard and Pascoe (2009)
G46471	TrGAroCN	NPHEN1O2 + HO_2 → NPHEN1OOH	`KRO2HO2(6)`	Rickard and Pascoe (2009)
G46472a	TrGAroCN	NPHEN1O2 + NO → NPHEN1O + NO_2	`KRO2NO`	Rickard and Pascoe (2009)
G46472b	TrGAroCN	NPHEN1O2 + NO_2 → NPHEN1O + NO_3	`k_C6H5O2_NO2`	Jagiella and Zabel (2007)*
G46473	TrGAroCN	NPHEN1O2 + NO_3 → NPHEN1O + NO_2	`KRO2NO3`	Rickard and Pascoe (2009)
G46474	TrGAroCN	NPHEN1O2 → NPHEN1O	`k1_RO2sRO2`	Rickard and Pascoe (2009)
G46475	TrGAroCN	NPHENOOH + OH → NPHENO2	`1.07E-10`	Rickard and Pascoe (2009)
G46476	TrGAroCN	C6H5O + NO_2 → HOC6H4NO2	`k_C6H5O_NO2`	Rickard and Pascoe (2009), Platz et al. (1998)*
G46477	TrGAroC	C6H5O + O_3 → C6H5O2	`k_C6H5O_O3`	Rickard and Pascoe (2009), Tao and Li (1999)
G46478	TrGAroCN	NCATECOOH + OH → NCATECO2	`k_roohro`	Rickard and Pascoe (2009)
G46479	TrGAroC	PBZQOOH + OH → PBZQCO + OH	`1.23E-10`	Rickard and Pascoe (2009)
G46480a	TrGAroC	PBZQO2 + HO_2 → PBZQOOH	`KRO2HO2(6)*(1-rchohch2o2_oh-rcoch2o2_oh)`	Rickard and Pascoe (2009)
G46480b	TrGAroC	PBZQO2 + HO_2 → C5CO2OHCO3 + OH	`KRO2HO2(6)*(rchohch2o2_oh+rcoch2o2_oh)`	Rickard and Pascoe (2009)*
G46481	TrGAroCN	PBZQO2 + NO → C5CO2OHCO3 + NO_2	`KRO2NO`	Rickard and Pascoe (2009)*

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G46482	TrGAroCN	PBZQO2 + NO_3 → C5CO2OHCO3 + NO_2	KRO2NO3	Rickard and Pascoe (2009)*
G46483	TrGAroC	PBZQO2 → C5CO2OHCO3	k1_RO2sORO2	Rickard and Pascoe (2009)*
G46484	TrGAroC	BZOBIPEROH + OH → MALDIALCO3 + GLYOX	8.16E-11	Rickard and Pascoe (2009)
G46485a	TrGAroCN	DNPHENO2 + HO_2 → DNPHENOOH	KRO2HO2(6)*(1-rchohch2o2_oh)	Rickard and Pascoe (2009)
G46485b	TrGAroCN	DNPHENO2 + HO_2 → NC4DCO2H + $HCOCO_2H$ + NO_2 + OH	KRO2HO2(6)*rchohch2o2_oh	Rickard and Pascoe (2009)*
G46486	TrGAroCN	DNPHENO2 + NO → NC4DCO2H + $HCOCO_2H$ + NO_2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G46487	TrGAroCN	DNPHENO2 + NO_3 → NC4DCO2H + $HCOCO_2H$ + NO_2 + NO_2	KRO2NO3	Rickard and Pascoe (2009)*
G46488	TrGAroCN	DNPHENO2 → NC4DCO2H + $HCOCO_2H$ + NO_2	k1_RO2ISOPDO2	Rickard and Pascoe (2009)*
G46489	TrGAroC	BZBIPEROOH + OH → BZOBIPEROH + OH	9.77E-11	Rickard and Pascoe (2009)
G46490a	TrGAroC	BZEMUCO2 + HO_2 → BZEMUCOOH	KRO2HO2(6)	Rickard and Pascoe (2009)
G46490b	TrGAroC	BZEMUCO2 + HO_2 → .5 EPXC4DIAL + .5 GLYOX + .5 HO_2 + .5 C3DIALO2 + .5 C32OH13CO + OH	KRO2HO2(6)	Rickard and Pascoe (2009)*
G46491a	TrGAroCN	BZEMUCO2 + NO → BZEMUCNO3	KRO2NO*alpha_AN(10,2,0,1,0, temp,cair)	Rickard and Pascoe (2009)
G46491b	TrGAroCN	BZEMUCO2 + NO → .5 EPXC4DIAL + .5 GLYOX + .5 HO_2 + .5 C3DIALO2 + .5 C32OH13CO + NO_2	KRO2NO*(1.-alpha_AN(10,2,0,1,0, temp,cair))	Rickard and Pascoe (2009)*
G46492	TrGAroCN	BZEMUCO2 + NO_3 → .5 EPXC4DIAL + .5 GLYOX + .5 HO_2 + .5 C3DIALO2 + .5 C32OH13CO + NO_2	KRO2NO3	Rickard and Pascoe (2009)*
G46493	TrGAroC	BZEMUCO2 → .5 EPXC4DIAL + .5 GLYOX + .5 HO_2 + .5 C3DIALO2 + .5 C32OH13CO	k1_RO2sORO2	Rickard and Pascoe (2009)*
G46494	TrGAroCN	C5CO2DBPAN + OH → HCOCOCHO + CH_3CHO + CO_2 + NO_2	3.28E-11	Rickard and Pascoe (2009)
G46495	TrGAroCN	C5CO2DBPAN → C5CO2DBCO3 + NO_2	k_PAN_M	Rickard and Pascoe (2009)
G46496	TrGAroCN	NBZQOOH + OH → NBZQO2	6.68E-11	Rickard and Pascoe (2009)
G46497	TrGAroC	CATEC1OOH + OH → CATEC1O2	k_roohro	Rickard and Pascoe (2009)
G46498	TrGAroC	C6125CO + OH → C5CO14O2 + CO	6.45E-11	Rickard and Pascoe (2009)
G46499a	TrGAroCN	NBZQO2 + HO_2 → NBZQOOH	KRO2HO2(6)*(1-rcoch2o2_oh)	Rickard and Pascoe (2009)
G46499b	TrGAroCN	NBZQO2 + HO_2 → C6CO4DB + NO_2 + OH	KRO2HO2(6)*rcoch2o2_oh	Rickard and Pascoe (2009)*
G46500	TrGAroCN	NBZQO2 + NO → C6CO4DB + NO_2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G46501	TrGAroCN	NBZQO2 + NO_3 → C6CO4DB + NO_2 + NO_2	KRO2NO3	Rickard and Pascoe (2009)*
G46502	TrGAroCN	NBZQO2 → C6CO4DB + NO_2	k1_RO2sORO2	Rickard and Pascoe (2009)*
G46503	TrGAroCN	DNPHENOOH + OH → DNPHENO2	k_roohro	Rickard and Pascoe (2009)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G46504	TrGAroC	CATEC1O2 + HO_2 → CATEC1OOH	KRO2HO2(6)	Rickard and Pascoe (2009)
G46505a	TrGAroCN	CATEC1O2 + NO → CATEC1O + NO_2	KRO2NO	Rickard and Pascoe (2009)
G46505b	TrGAroCN	CATEC1O2 + NO_2 → CATEC1O + NO_3	K_C6H5O2_NO2	Jagiella and Zabel (2007)*
G46506	TrGAroCN	CATEC1O2 + NO_3 → CATEC1O + NO_2	KRO2NO3	Rickard and Pascoe (2009)
G46507	TrGAroC	CATEC1O2 → CATEC1O	k1_RO2sORO2	Rickard and Pascoe (2009)
G46508	TrGAroC	BZEMUCCO3H + OH → BZEMUCCO3	4.37E-11	Rickard and Pascoe (2009)
G46509	TrGAroC	C6H5OOH + OH → C6H5O2	3.60E-12	Rickard and Pascoe (2009)
G46510	TrGAroC	BZEMUCOOH + OH → BZEMUCCO + OH	1.31E-10	Rickard and Pascoe (2009)
G46511a	TrGAroC	BZEMUCCO3 + HO_2 → BZEMUCCO2H + O_3	KAPHO2*rco3_o3	Rickard and Pascoe (2009)
G46511b	TrGAroC	BZEMUCCO3 + HO_2 → BZEMUCCO3H	KAPHO2*rco3_ooh	Rickard and Pascoe (2009)
G46511c	TrGAroC	BZEMUCCO3 + HO_2 → C5DIALO2 + CO_2 + OH	KAPHO2*rco3_oh	Rickard and Pascoe (2009)
G46512	TrGAroCN	BZEMUCCO3 + NO → C5DIALO2 + CO_2 + NO_2	KAPNO	Rickard and Pascoe (2009)
G46513	TrGAroCN	BZEMUCCO3 + NO_2 → BZEMUCPAN	k_CH3CO3_NO2	Rickard and Pascoe (2009)
G46514	TrGAroCN	BZEMUCCO3 + NO_3 → C5DIALO2 + CO_2 + NO_2	KRO2NO3*1.74	Rickard and Pascoe (2009)
G46515	TrGAroC	BZEMUCCO3 → C5DIALO2 + CO_2	k1_RO2RCO3	Rickard and Pascoe (2009)*
G46516	TrGAroC	C6H5O2 + HO_2 → C6H5OOH	KRO2HO2(6)	Rickard and Pascoe (2009)
G46517a	TrGAroCN	C6H5O2 + NO → C6H5O + NO_2	KRO2NO	Rickard and Pascoe (2009)
G46517b	TrGAroCN	C6H5O2 + NO_2 → C6H5O + NO_3	K_C6H5O2_NO2	Jagiella and Zabel (2007)*
G46518	TrGAroCN	C6H5O2 + NO_3 → C6H5O + NO_2	KRO2NO3	Rickard and Pascoe (2009)
G46519	TrGAroC	C6H5O2 → C6H5O	k1_RO2sRO2	Rickard and Pascoe (2009)
G46521	TrGAroCN	BZEMUCNO3 + OH → BZEMUCCO + NO_2	4.38E-11	Rickard and Pascoe (2009)
G46522a	TrGAroC	BZBIPERO2 + HO_2 → BZBIPEROOH	KRO2HO2(6)*(1.-rbipero2_oh)	Rickard and Pascoe (2009)
G46522b	TrGAroC	BZBIPERO2 + HO_2 → OH + GLYOX + HO_2 + .5 BZFUONE + .5 BZFUONE	KRO2HO2(6)*rbipero2_oh	Rickard and Pascoe (2009), Birdsall et al. (2010)*
G46523a	TrGAroCN	BZBIPERO2 + NO → BZBIPERNO3	KRO2NO*alpha_AN(9,2,0,0,1,temp, cair)	Rickard and Pascoe (2009)
G46523b	TrGAroCN	BZBIPERO2 + NO → NO_2 + GLYOX + HO_2 + .5 BZFUONE + .5 BZFUONE	KRO2NO*(1.-alpha_AN(9,2,0,0,1, temp,cair))	Rickard and Pascoe (2009)*
G46524	TrGAroCN	BZBIPERO2 + NO_3 → NO_2 + GLYOX + HO_2 + .5 BZFUONE + .5 BZFUONE	KRO2NO3	Rickard and Pascoe (2009)*
G46525	TrGAroC	BZBIPERO2 → GLYOX + HO_2 + BZFUONE	k1_RO2sORO2	Rickard and Pascoe (2009)*
G47200	TrGTerCN	CO235C6CHO + NO_3 → CO235C6CO3 + HNO_3	KNO3AL*5.5	Rickard and Pascoe (2009)
G47201	TrGTerC	CO235C6CHO + OH → CO235C6CO3	6.70E-11	Rickard and Pascoe (2009)
G47202a	TrGTerC	CO235C6CO3 + HO_2 → C235C6CO3H	KAPHO2*(rco3_ooh+rco3_o3)	Rickard and Pascoe (2009)
G47202b	TrGTerC	CO235C6CO3 + HO_2 → CO235C6O2 + CO_2 + OH	KAPHO2*rco3_oh	Rickard and Pascoe (2009)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G47203	TrGTerCN	CO235C6CO3 + NO → CO235C6O2 + CO_2 + NO_2	KAPNO	Rickard and Pascoe (2009)
G47204	TrGTerCN	CO235C6CO3 + NO_2 → C7PAN3	k_CH3CO3_NO2	Rickard and Pascoe (2009)
G47205	TrGTerC	CO235C6CO3 → CO235C6O2 + CO_2	k1_RO2RCO3	Rickard and Pascoe (2009)
G47206	TrGTerC	C235C6CO3H + OH → CO235C6CO3	4.75E-12	Rickard and Pascoe (2009)
G47207	TrGTerCN	C7PAN3 + OH → CO235C5CHO + CO + NO_2	8.83E-13	Rickard and Pascoe (2009)
G47208	TrGTerCN	C7PAN3 → CO235C6CO3 + NO_2	k_PAN_M	Rickard and Pascoe (2009)
G47209a	TrGTerC	C716O2 + HO_2 → C716OOH	KRO2HO2(7)*rcoch2o2_ooh	Rickard and Pascoe (2009), Sander et al. (2018)
G47209b	TrGTerC	C716O2 + HO_2 → CO13C4CHO + $CH_3C(O)$ + OH	KRO2HO2(7)*rcoch2o2_oh	Rickard and Pascoe (2009), Sander et al. (2018)
G47210	TrGTerCN	C716O2 + NO → CO13C4CHO + $CH_3C(O)$ + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G47211	TrGTerC	C716O2 → CO13C4CHO + $CH_3C(O)$	k1_RO2sORO2	Rickard and Pascoe (2009)
G47212	TrGTerC	C716OOH + OH → CO235C6CHO + OH	1.20E-10	Rickard and Pascoe (2009)
G47213	TrGTerC	C721O2 + HO_2 → C721OOH	KRO2HO2(7)	Rickard and Pascoe (2009)
G47214	TrGTerCN	C721O2 + NO → C722O2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G47215	TrGTerC	C721O2 → C722O2	k1_RO2pRO2	Rickard and Pascoe (2009)
G47216	TrGTerC	C721OOH + OH → C721O2	1.27E-11	Rickard and Pascoe (2009)
G47217	TrGTerC	C722O2 + HO_2 → C722OOH	KRO2HO2(7)	Rickard and Pascoe (2009)
G47218	TrGTerCN	C722O2 + NO → CH_3COCH_3 + C44O2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G47219	TrGTerC	C722O2 → CH_3COCH_3 + C44O2	k1_RO2tRO2	Rickard and Pascoe (2009)
G47220	TrGTerC	C722OOH + OH → C722O2	3.31E-11	Rickard and Pascoe (2009)
G47221	TrGTerC	ROO6R3O2 → ROO6R5O2	5.68E10*EXP(-8745./TEMP)	Vereecken and Peeters (2012)
G47222	TrGTerCN	ROO6R3O2 + NO → ROO6R3O + NO_2	KRO2NO	Vereecken and Peeters (2012)*
G47223	TrGTerC	ROO6R3O2 + HO_2 → 7 LCARBON	KRO2HO2(7)	Vereecken and Peeters (2012)*
G47224	TrGTerC	ROO6R3O2 → ROO6R3O	k1_RO2sRO2	Vereecken and Peeters (2012)
G47225	TrGTerC	ROO6R3O → 7 LCARBON + HO_2	5.7E10*EXP(-2949./TEMP)	Vereecken and Peeters (2012)*
G47226	TrGTerC	ROO6R5O2 → 7 LCARBON + OH	9.17E10*EXP(-8706./TEMP)	Vereecken and Peeters (2012)*
G47400	TrGAroC	TOLUENE + OH → .07 C6H5CH2O2 + .18 CRESOL + .18 HO_2 + .65 TLBIPERO2 + .10 TLEPOXMUC + .10 HO_2 + .04 AELV	1.8E-12*EXP(340/TEMP)	Rickard and Pascoe (2009)*
G47401	TrGAroC	C6H5CH2O2 + HO_2 → C6H5CH2OOH	1.5E-13*EXP(1310/TEMP)	Rickard and Pascoe (2009)
G47402a	TrGAroCN	C6H5CH2O2 + NO → C6H5CH2NO3	KRO2NO*alpha_AN(7,1,0,0,0,temp, cair)	Rickard and Pascoe (2009)*
G47402b	TrGAroCN	C6H5CH2O2 + NO → BENZAL + HO_2 + NO_2	KRO2NO*(1.-alpha_AN(7,1,0,0,0, temp,cair))	Rickard and Pascoe (2009)*

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G47403	TrGAroCN	C6H5CH2O2 + NO_3 → BENZAL + HO_2 + NO_2	KRO2NO3	Rickard and Pascoe (2009)*
G47404	TrGAroC	C6H5CH2O2 → BENZAL + HO_2	2.*(k_CH3O2*2.4E-14*EXP(1620./TEMP))**0.5*RO2	Rickard and Pascoe (2009)*
G47405	TrGAroCN	CRESOL + NO_3 → .103 CRESO2 + .103 HNO_3 + .506 NCRESO2 + .391 TOL1O + .391 HNO_3	1.4E-11	Rickard and Pascoe (2009)*
G47406	TrGAroC	CRESOL + OH → .2 CRESO2 + .727 MCATECHOL + .727 HO_2 + .073 TOL1O	4.65E-11	Rickard and Pascoe (2009)*
G47407a	TrGAroC	TLBIPERO2 + HO_2 → TLBIPEROOH	KRO2HO2(7)*(1.-rbipero2_oh)	Rickard and Pascoe (2009)
G47407b	TrGAroC	TLBIPERO2 + HO_2 → OH + .6 GLYOX + .4 MGLYOX + HO_2 + .2 C4MDIAL + .2 C5DICARB + .2 TLFUONE + .2 BZFUONE + .2 MALDIAL + 0.780 ALOV	KRO2HO2(7)*rbipero2_oh	Rickard and Pascoe (2009), Birdsall et al. (2010)*
G47408a	TrGAroCN	TLBIPERO2 + NO → NO_2 + .6 GLYOX + .4 MGLYOX + HO_2 + .2 C4MDIAL + .2 C5DICARB + .2 TLFUONE + .2 BZFUONE + .2 MALDIAL + 0.097 ALOV + 0.748 ASOV	KRO2NO*(1.-alpha_AN(11,2,0,0,1,temp,cair))	Rickard and Pascoe (2009)*
G47408b	TrGAroCN	TLBIPERO2 + NO → TLBIPERNO3	KRO2NO*alpha_AN(11,2,0,0,1,temp,cair)	Rickard and Pascoe (2009)*
G47409	TrGAroCN	TLBIPERO2 + NO_3 → NO_2 + .6 GLYOX + .4 MGLYOX + HO_2 + .2 C4MDIAL + .2 C5DICARB + .2 TLFUONE + .2 BZFUONE + .2 MALDIAL	KRO2NO3	Rickard and Pascoe (2009)*
G47410	TrGAroC	TLBIPERO2 → .6 GLYOX + .4 MGLYOX + HO_2 + .2 C4MDIAL + .2 C5DICARB + .2 TLFUONE + .2 BZFUONE + .2 MALDIAL	k1_RO2sORO2	Rickard and Pascoe (2009)*
G47411	TrGAroCN	TLEPOXMUC + NO_3 → TLEMUCCO3 + HNO_3	KNO3AL*2.75	Rickard and Pascoe (2009)
G47412	TrGAroC	TLEPOXMUC + O_3 → EPXC4DIAL + .125 CH_3CHO + .695 $CH_3C(O)$ + .57 CO + .57 OH + .125 HO_2 + .1125 CH_3COCO_2H + .0675 MGLYOX + .0675 H_2O_2 + .25 CO_2	5.00E-18	Rickard and Pascoe (2009)*
G47413	TrGAroC	TLEPOXMUC + OH → .31 TLEMUCCO3 + .69 TLEMUCO2	7.99E-11	Rickard and Pascoe (2009)*
G47414	TrGAroC	C6H5CH2OOH + OH → BENZAL + OH	2.05E-11	Rickard and Pascoe (2009)
G47415	TrGAroCN	C6H5CH2NO3 + OH → BENZAL + NO_2	6.03E-12	Rickard and Pascoe (2009)
G47416	TrGAroCN	BENZAL + NO_3 → C6H5CO3 + HNO_3	2.40E-15	Rickard and Pascoe (2009)
G47417	TrGAroC	BENZAL + OH → C6H5CO3	5.9E-12*EXP(225/TEMP)	Rickard and Pascoe (2009)
G47418a	TrGAroC	CRESO2 + HO_2 → CRESOOH	KRO2HO2(7)*(1-rchohch2o2_oh)	Rickard and Pascoe (2009)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G47418b	TrGAroC	CRESO2 + HO_2 → .68 C5CO14OH + .68 GLYOX + HO_2 + .32 PTLQONE + OH	KRO2HO2(7)*rchohch2o2_oh	Rickard and Pascoe (2009)*
G47419	TrGAroCN	CRESO2 + NO → .68 C5CO14OH + .68 GLYOX + HO_2 + .32 PTLQONE + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G47420	TrGAroCN	CRESO2 + NO_3 → .68 C5CO14OH + .68 GLYOX + HO_2 + .32 PTLQONE + NO_2	KRO2NO3	Rickard and Pascoe (2009)*
G47421	TrGAroC	CRESO2 → .68 C5CO14OH + .68 GLYOX + HO_2 + .32 PTLQONE	k1_RO2ISOPDO2	Rickard and Pascoe (2009)*
G47422a	TrGAroCN	NCRESO2 + HO_2 → NCRESOOH	KRO2HO2(7)*(1-rchohch2o2_oh)	Rickard and Pascoe (2009)
G47422b	TrGAroCN	NCRESO2 + HO_2 → C5CO14OH + GLYOX + NO_2 + OH	KRO2HO2(7)*rchohch2o2_oh	Rickard and Pascoe (2009)*
G47423	TrGAroCN	NCRESO2 + NO → C5CO14OH + GLYOX + NO_2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G47424	TrGAroCN	NCRESO2 + NO_3 → C5CO14OH + GLYOX + NO_2 + NO_2	KRO2NO3	Rickard and Pascoe (2009)*
G47425	TrGAroCN	NCRESO2 → C5CO14OH + GLYOX + NO_2	k1_RO2ISOPDO2	Rickard and Pascoe (2009)*
G47426	TrGAroCN	TOL1O + NO_2 → TOL1OHNO2	k_C6H5O_NO2	Rickard and Pascoe (2009), Platz et al. (1998)*
G47427	TrGAroC	TOL1O + O_3 → OXYL1O2	k_C6H5O_O3	Rickard and Pascoe (2009), Tao and Li (1999)
G47428	TrGAroCN	MCATECHOL + NO_3 → MCATEC1O + HNO_3	1.7E-10*1.0	Rickard and Pascoe (2009)
G47429	TrGAroC	MCATECHOL + O_3 → MC3ODBCO2H + $HCOCO_2H$ + HO_2 + OH	2.8E-17	Rickard and Pascoe (2009)*
G47430	TrGAroC	MCATECHOL + OH → MCATEC1O	2.0E-10*1.0	Rickard and Pascoe (2009)
G47431	TrGAroC	TLBIPEROOH + OH → TLOBIPEROH + OH	9.64E-11	Rickard and Pascoe (2009)
G47432	TrGAroCN	TLBIPERNO3 + OH → TLOBIPEROH + NO_2	7.16E-11	Rickard and Pascoe (2009)
G47433	TrGAroC	TLOBIPEROH + OH → C5CO14O2 + GLYOX	7.99E-11	Rickard and Pascoe (2009)
G47434a	TrGAroC	TLEMUCCO3 + HO_2 → C615CO2O2 + CO_2 + OH	KAPHO2*rco3_oh	Rickard and Pascoe (2009)
G47434b	TrGAroC	TLEMUCCO3 + HO_2 → TLEMUCCO2H + O_3	KAPHO2*rco3_o3	Rickard and Pascoe (2009)
G47434c	TrGAroC	TLEMUCCO3 + HO_2 → TLEMUCCO3H	KAPHO2*rco3_ooh	Rickard and Pascoe (2009)
G47435	TrGAroCN	TLEMUCCO3 + NO → C615CO2O2 + CO_2 + NO_2	KAPNO	Rickard and Pascoe (2009)
G47436	TrGAroCN	TLEMUCCO3 + NO_2 → TLEMUCPAN	k_CH3CO3_NO2	Rickard and Pascoe (2009)*
G47437	TrGAroCN	TLEMUCCO3 + NO_3 → C615CO2O2 + CO_2 + NO_2	KRO2NO3*1.74	Rickard and Pascoe (2009)
G47438	TrGAroC	TLEMUCCO3 → C615CO2O2 + CO_2	k1_RO2RCO3	Rickard and Pascoe (2009)*
G47439a	TrGAroC	TLEMUCO2 + HO_2 → TLEMUCOOH	KRO2HO2(7)*(1-rchohch2o2_oh-rcoch2o2_oh)	Rickard and Pascoe (2009)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G47439b	TrGAroC	TLEMUCO2 + HO_2 → .5 C3DIALO2 + .5 CO2H3CHO + .5 EPXC4DIAL + .5 MGLYOX + .5 HO_2 + OH	`KRO2HO2(7)*(rchohch2o2_oh+rcoch2o2_oh)`	Rickard and Pascoe (2009)*
G47440a	TrGAroCN	TLEMUCO2 + NO → TLEMUCNO3	`KRO2NO*alpha_AN(11,2,1,0,0,temp,cair)`	Rickard and Pascoe (2009)
G47440b	TrGAroCN	TLEMUCO2 + NO → .5 C3DIALO2 + .5 CO2H3CHO + .5 EPXC4DIAL + .5 MGLYOX + .5 HO_2 + NO_2	`KRO2NO*(1.-alpha_AN(11,2,1,0,0,temp,cair))`	Rickard and Pascoe (2009)*
G47441	TrGAroCN	TLEMUCO2 + NO_3 → .5 C3DIALO2 + .5 CO2H3CHO + .5 EPXC4DIAL + .5 MGLYOX + .5 HO_2 + NO_2	`KRO2NO3`	Rickard and Pascoe (2009)*
G47442	TrGAroC	TLEMUCO2 → .5 C3DIALO2 + .5 CO2H3CHO + .5 EPXC4DIAL + .5 MGLYOX + .5 HO_2	`k1_RO2sORO2`	Rickard and Pascoe (2009)*
G47443a	TrGAroC	C6H5CO3 + HO_2 → C6H5CO3H	`1.1E-11*EXP(364./temp)*0.65`	Roth et al. (2010)
G47443b	TrGAroC	C6H5CO3 + HO_2 → C6H5O2 + CO_2 + OH	`1.1E-11*EXP(364./temp)*0.20`	Roth et al. (2010)
G47443c	TrGAroC	C6H5CO3 + HO_2 → PHCOOH + O_3	`1.1E-11*EXP(364./temp)*0.15`	Roth et al. (2010)
G47444	TrGAroCN	C6H5CO3 + NO → C6H5O2 + CO_2 + NO_2	`KAPNO`	Rickard and Pascoe (2009)
G47445	TrGAroCN	C6H5CO3 + NO_2 → PBZN	`k_CH3CO3_NO2`	Rickard and Pascoe (2009)*
G47446	TrGAroCN	C6H5CO3 + NO_3 → C6H5O2 + CO_2 + NO_2	`KRO2NO3*1.74`	Rickard and Pascoe (2009)
G47447	TrGAroC	C6H5CO3 → C6H5O2 + CO_2	`k1_RO2RCO3`	Rickard and Pascoe (2009)*
G47448	TrGAroC	CRESOOH + OH → CRESO2	`1.15E-10`	Rickard and Pascoe (2009)
G47449	TrGAroCN	NCRESOOH + OH → NCRESO2	`1.07E-10`	Rickard and Pascoe (2009)
G47450	TrGAroCN	TOL1OHNO2 + NO_3 → NCRES1O + HNO_3	`3.13E-13*1.0`	Rickard and Pascoe (2009)
G47451	TrGAroCN	TOL1OHNO2 + OH → NCRES1O	`2.8E-12`	Rickard and Pascoe (2009)
G47452	TrGAroC	OXYL1O2 + HO_2 → OXYL1OOH	`KRO2HO2(7)`	Rickard and Pascoe (2009)
G47453	TrGAroCN	OXYL1O2 + NO → TOL1O + NO_2	`KRO2NO`	Rickard and Pascoe (2009)
G47454	TrGAroCN	OXYL1O2 + NO_2 → TOL1O + NO_3	`K_C6H5O2_NO2`	Jagiella and Zabel (2007)*
G47455	TrGAroCN	OXYL1O2 + NO_3 → TOL1O + NO_2	`KRO2NO3`	Rickard and Pascoe (2009)
G47456	TrGAroC	OXYL1O2 → TOL1O	`k1_RO2sRO2`	Rickard and Pascoe (2009)
G47457	TrGAroCN	MCATEC1O + NO_2 → MNCATECH	`k_C6H5O_NO2`	Rickard and Pascoe (2009), Platz et al. (1998)
G47458	TrGAroC	MCATEC1O + O_3 → MCATEC1O2	`k_C6H5O_O3`	Rickard and Pascoe (2009), Tao and Li (1999)
G47459	TrGAroC	TLEMUCCO2H + OH → C615CO2O2 + CO_2	`5.98E-11`	Rickard and Pascoe (2009)
G47460	TrGAroC	TLEMUCCO3H + OH → TLEMUCCO3	`6.29E-11`	Rickard and Pascoe (2009)
G47461	TrGAroCN	TLEMUCPAN + OH → C5DICARB + CO + CO_2 + NO_2	`5.96E-11`	Rickard and Pascoe (2009)
G47462	TrGAroCN	TLEMUCPAN → TLEMUCCO3 + NO_2	`k_PAN_M`	Rickard and Pascoe (2009)
G47463	TrGAroC	TLEMUCOOH + OH → TLEMUCCO + OH	`7.04E-11`	Rickard and Pascoe (2009)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G47464	TrGAroCN	TLEMUCNO3 + OH → TLEMUCCO + NO_2	3.06E-11	Rickard and Pascoe (2009)
G47465	TrGAroC	TLEMUCCO + OH → $CH_3C(O)$ + EPXC4DIAL + CO	4.06E-11	Rickard and Pascoe (2009)
G47466	TrGAroC	C6H5CO3H + OH → C6H5CO3	4.66E-12	Rickard and Pascoe (2009)
G47467	TrGAroC	PHCOOH + OH → C6H5O2 + CO_2	1.10E-12	Rickard and Pascoe (2009)
G47468	TrGAroCN	PBZN + OH → C6H5OOH + CO + NO_2	1.06E-12	Rickard and Pascoe (2009)
G47469	TrGAroCN	PBZN → C6H5CO3 + NO_2	k_PAN_M*0.67	Rickard and Pascoe (2009)
G47470	TrGAroCN	PTLQONE + NO_3 → NPTLQO2	1.00E-12	Rickard and Pascoe (2009)
G47471	TrGAroC	PTLQONE + OH → PTLQO2	2.3E-11	Rickard and Pascoe (2009)
G47472	TrGAroCN	NCRES1O + NO_2 → DNCRES	k_C6H5O_NO2	Rickard and Pascoe (2009), Platz et al. (1998)
G47473	TrGAroCN	NCRES1O + O_3 → NCRES1O2	k_C6H5O_O3	Rickard and Pascoe (2009), Tao and Li (1999)
G47474	TrGAroC	OXYL1OOH + OH → OXYL1O2	4.65E-11	Rickard and Pascoe (2009)
G47475	TrGAroCN	MNCATECH + NO_3 → MNNCATECO2	5.03E-12	Rickard and Pascoe (2009)
G47476	TrGAroCN	MNCATECH + OH → MNCATECO2	6.83E-12	Rickard and Pascoe (2009)
G47477	TrGAroC	MCATEC1O2 + HO_2 → MCATEC1OOH	KRO2HO2(7)	Rickard and Pascoe (2009)
G47478	TrGAroCN	MCATEC1O2 + NO → MCATEC1O + NO_2	KRO2NO	Rickard and Pascoe (2009)
G47479	TrGAroCN	MCATEC1O2 + NO_2 → MCATEC1O + NO_3	K_C6H5O2_NO2	Jagiella and Zabel (2007)*
G47480	TrGAroCN	MCATEC1O2 + NO_3 → MCATEC1O + NO_2	KRO2NO3	Rickard and Pascoe (2009)
G47481	TrGAroC	MCATEC1O2 → MCATEC1O	k1_RO2sORO2	Rickard and Pascoe (2009)
G47482a	TrGAroCN	NPTLQO2 + HO_2 → NPTLQOOH	KRO2HO2(7)*(1-rcoch2o2_oh)	Rickard and Pascoe (2009)
G47482b	TrGAroCN	NPTLQO2 + HO_2 → C7CO4DB + NO_2 + OH	KRO2HO2(7)*rcoch2o2_oh	Rickard and Pascoe (2009)*
G47483	TrGAroCN	NPTLQO2 + NO → C7CO4DB + NO_2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G47484	TrGAroCN	NPTLQO2 + NO_3 → C7CO4DB + NO_2 + NO_2	KRO2NO3	Rickard and Pascoe (2009)*
G47485	TrGAroCN	NPTLQO2 → C7CO4DB + NO_2	k1_RO2sORO2	Rickard and Pascoe (2009)*
G47486a	TrGAroC	PTLQO2 + HO_2 → PTLQOOH	KRO2HO2(7)*(1-rchohch2o2_ oh-rcoch2o2_oh)	Rickard and Pascoe (2009)
G47486b	TrGAroC	PTLQO2 + HO_2 → C6CO2OHCO3 + OH	KRO2HO2(7)*(rchohch2o2_oh+ rcoch2o2_oh)	Rickard and Pascoe (2009)*
G47487	TrGAroCN	PTLQO2 + NO → C6CO2OHCO3 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G47488	TrGAroCN	PTLQO2 + NO_3 → C6CO2OHCO3 + NO_2	KRO2NO3	Rickard and Pascoe (2009)*
G47489	TrGAroC	PTLQO2 → C6CO2OHCO3	k1_RO2sORO2	Rickard and Pascoe (2009)*
G47490	TrGAroCN	DNCRES + NO_3 → NDNCRESO2	7.83E-15	Rickard and Pascoe (2009)
G47491	TrGAroCN	DNCRES + OH → DNCRESO2	5.10E-14	Rickard and Pascoe (2009)
G47492	TrGAroCN	NCRES1O2 + HO_2 → NCRES1OOH	KRO2HO2(7)	Rickard and Pascoe (2009)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G47493	TrGAroCN	NCRES1O2 + NO → NCRES1O + NO_2	KRO2NO	Rickard and Pascoe (2009)
G47494	TrGAroCN	NCRES1O2 + NO_2 → NCRES1O + NO_3	K_C6H5O2_NO2	Jagiella and Zabel (2007)*
G47495	TrGAroCN	NCRES1O2 + NO_3 → NCRES1O + NO_2	KRO2NO3	Rickard and Pascoe (2009)
G47496	TrGAroCN	NCRES1O2 → NCRES1O	k1_RO2sRO2	Rickard and Pascoe (2009)
G47497a	TrGAroCN	MNNCATECO2 + HO_2 → MNNCATCOOH	KRO2HO2(7)*(1-rchohch2o2_oh)	Rickard and Pascoe (2009)
G47497b	TrGAroCN	MNNCATECO2 + HO_2 → NC4MDCO2HN + $HCOCO_2H$ + NO_2 + OH	KRO2HO2(7)*rchohch2o2_oh	Rickard and Pascoe (2009)*
G47498	TrGAroCN	MNNCATECO2 + NO → NC4MDCO2HN + $HCOCO_2H$ + NO_2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G47499	TrGAroCN	MNNCATECO2 + NO_3 → NC4MDCO2HN + $HCOCO_2H$ + NO_2 + NO_2	KRO2NO3	Rickard and Pascoe (2009)*
G47500	TrGAroCN	MNNCATECO2 → NC4MDCO2HN + $HCOCO_2H$ + NO_2	k1_RO2ISOPDO2	Rickard and Pascoe (2009)
G47501a	TrGAroCN	MNCATECO2 + HO_2 → MNCATECOOH	KRO2HO2(7)*(1-rchohch2o2_oh)	Rickard and Pascoe (2009)
G47501b	TrGAroCN	MNCATECO2 + HO_2 → NC4MDCO2HN + $HCOCO_2H$ + HO_2 + OH	KRO2HO2(7)*rchohch2o2_oh	Rickard and Pascoe (2009)*
G47502	TrGAroCN	MNCATECO2 + NO → NC4MDCO2HN + $HCOCO_2H$ + HO_2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G47503	TrGAroCN	MNCATECO2 + NO_3 → NC4MDCO2HN + $HCOCO_2H$ + HO_2 + NO_2	KRO2NO3	Rickard and Pascoe (2009)*
G47504	TrGAroCN	MNCATECO2 → NC4MDCO2HN + $HCOCO_2H$ + HO_2	k1_RO2ISOPDO2	Rickard and Pascoe (2009)*
G47505	TrGAroC	MCATEC1OOH + OH → MCATEC1O2	2.05E-10	Rickard and Pascoe (2009)
G47506	TrGAroCN	NPTLQOOH + OH → NPTLQO2	8.56E-11	Rickard and Pascoe (2009)
G47507	TrGAroC	PTLQOOH + OH → PTLQCO + OH	1.42E-10	Rickard and Pascoe (2009)
G47508	TrGAroC	PTLQCO + OH → C6CO2OHCO3	7.95E-11	Rickard and Pascoe (2009)
G47509a	TrGAroCN	NDNCRESO2 + HO_2 → NDNCRESOOH	KRO2HO2(7)*(1-rchohch2o2_oh)	Rickard and Pascoe (2009)
G47509b	TrGAroCN	NDNCRESO2 + HO_2 → NC4MDCO2HN + HNO_3 + 2 CO + NO_2 + OH	KRO2HO2(7)*rchohch2o2_oh	Rickard and Pascoe (2009)*
G47510	TrGAroCN	NDNCRESO2 + NO → NC4MDCO2HN + HNO_3 + 2 CO + NO_2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G47511	TrGAroCN	NDNCRESO2 + NO_3 → NC4MDCO2HN + HNO_3 + 2 CO + NO_2 + NO_2	KRO2NO3	Rickard and Pascoe (2009)*
G47512	TrGAroCN	NDNCRESO2 → NC4MDCO2HN + HNO_3 + 2 CO + NO_2	k1_RO2ISOPDO2	Rickard and Pascoe (2009)*
G47513a	TrGAroCN	DNCRESO2 + HO_2 → DNCRESOOH	KRO2HO2(7)*(1-rchohch2o2_oh)	Rickard and Pascoe (2009)
G47513b	TrGAroCN	DNCRESO2 + HO_2 → NC4MDCO2HN + $HCOCO_2H$ + NO_2 + OH	KRO2HO2(7)*rchohch2o2_oh	Rickard and Pascoe (2009)*

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G47514	TrGAroCN	DNCRESO2 + NO → NC4MDCO2HN + $HCOCO_2H$ + NO_2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G47515	TrGAroCN	DNCRESO2 + NO_3 → NC4MDCO2HN + $HCOCO_2H$ + NO_2 + NO_2	KRO2NO3	Rickard and Pascoe (2009)*
G47516	TrGAroCN	DNCRESO2 → NC4MDCO2HN + $HCOCO_2H$ + NO_2	k1_RO2ISOPDO2	Rickard and Pascoe (2009)*
G47517	TrGAroCN	NCRES1OOH + OH → NCRES1O2	1.53E-12	Rickard and Pascoe (2009)
G47518	TrGAroCN	MNNCATCOOH + OH → MNNCATECO2	k_roohro	Rickard and Pascoe (2009)
G47519	TrGAroCN	MNCATECOOH + OH → MNCATECO2	k_roohro	Rickard and Pascoe (2009)
G47520	TrGAroC	C7CO4DB + OH → CO + CO + $CH_3C(O)$ + HCOCOCHO	9.58E-11	Rickard and Pascoe (2009)
G47521a	TrGAroC	C6CO2OHCO3 + HO_2 → C5134CO2OH + HO_2 + CO + CO_2 + OH	KAPHO2*rco3_oh	Rickard and Pascoe (2009)
G47521b	TrGAroC	C6CO2OHCO3 + HO_2 → C6COOHCO3H	KAPHO2*(rco3_ooh+rco3_o3)	Rickard and Pascoe (2009)
G47522	TrGAroCN	C6CO2OHCO3 + NO → C5134CO2OH + HO_2 + CO + CO_2 + NO_2	KAPNO	Rickard and Pascoe (2009)
G47523	TrGAroCN	C6CO2OHCO3 + NO_2 → C6CO2OHPAN	k_CH3CO3_NO2	Rickard and Pascoe (2009)
G47524	TrGAroCN	C6CO2OHCO3 + NO_3 → C5134CO2OH + HO_2 + CO + CO_2 + NO_2	KRO2NO3*1.74	Rickard and Pascoe (2009)
G47525	TrGAroC	C6CO2OHCO3 → C5134CO2OH + HO_2 + CO + CO_2	k1_RO2RCO3	Rickard and Pascoe (2009)
G47526	TrGAroCN	NDNCRESOOH + OH → NDNCRESO2	k_roohro	Rickard and Pascoe (2009)
G47527	TrGAroCN	DNCRESOOH + OH → DNCRESO2	k_roohro	Rickard and Pascoe (2009)
G47528	TrGAroC	C6COOHCO3H + OH → C6CO2OHCO3	9.29E-11	Rickard and Pascoe (2009)
G47529	TrGAroCN	C6CO2OHPAN + OH → C5134CO2OH + CO + CO + NO_2	8.96E-11	Rickard and Pascoe (2009)
G47530	TrGAroCN	C6CO2OHPAN → C6CO2OHCO3 + NO_2	k_PAN_M	Rickard and Pascoe (2009)
G48200	TrGTerC	C85O2 → C86O2	k1_RO2tRO2	Rickard and Pascoe (2009)
G48201	TrGTerC	C85O2 + HO_2 → C85OOH	KRO2HO2(8)	Rickard and Pascoe (2009)
G48202	TrGTerCN	C85O2 + NO → C86O2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G48203	TrGTerC	C85OOH + OH → C85O2	1.29E-11	Rickard and Pascoe (2009)
G48204	TrGTerC	C86O2 → C511O2 + CH_3COCH_3	k1_RO2tRO2	Rickard and Pascoe (2009)
G48205	TrGTerCN	C86O2 + NO → C511O2 + CH_3COCH_3 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G48206	TrGTerC	C86O2 + HO_2 → C86OOH	KRO2HO2(8)	Rickard and Pascoe (2009)
G48207	TrGTerC	C86OOH + OH → C86O2	3.45E-11	Rickard and Pascoe (2009)
G48208	TrGTerC	C811O2 → C812O2	k1_RO2pRO2	Rickard and Pascoe (2009)
G48209	TrGTerC	C811O2 + HO_2 → 8 LCARBON	KRO2HO2(8)	Rickard and Pascoe (2009)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G48210	TrGTerCN	C811O2 + NO → C812O2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G48211	TrGTerC	C812O2 → C813O2	k1_RO2tORO2	Rickard and Pascoe (2009)
G48212	TrGTerCN	C812O2 + NO → C813O2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G48213	TrGTerC	C812O2 + HO_2 → C812OOH	KRO2HO2(8)	Rickard and Pascoe (2009)
G48214	TrGTerC	C812OOH + OH → C812O2	1.09E-11	Rickard and Pascoe (2009)
G48215	TrGTerC	C813O2 → CH_3COCH_3 + C512O2	k1_RO2tRO2	Rickard and Pascoe (2009)
G48216	TrGTerCN	C813O2 + NO → CH_3COCH_3 + C512O2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G48217	TrGTerC	C813O2 + HO_2 → C813OOH	KRO2HO2(8)	Rickard and Pascoe (2009)
G48218	TrGTerC	C813OOH + OH → C813O2	1.86E-11	Rickard and Pascoe (2009)
G48219	TrGTerCN	C721CHO + NO_3 → C721CO3 + HNO_3	KNO3AL*8.5	Rickard and Pascoe (2009)
G48220	TrGTerC	C721CHO + OH → C721CO3	2.63E-11	Rickard and Pascoe (2009)
G48221a	TrGTerC	C721CO3 + HO_2 → C721CO3H	KAPHO2*rco3_ooh	Rickard and Pascoe (2009)
G48221b	TrGTerC	C721CO3 + HO_2 → C721O2 + CO_2 + OH	KAPHO2*rco3_oh	Rickard and Pascoe (2009)
G48221c	TrGTerC	C721CO3 + HO_2 → NORPINIC + O_3	KAPHO2*rco3_o3	Rickard and Pascoe (2009)
G48222	TrGTerCN	C721CO3 + NO → C721O2 + CO_2 + NO_2	KAPNO	Rickard and Pascoe (2009)*
G48223	TrGTerCN	C721CO3 + NO_2 → C721PAN	k_CH3CO3_NO2	Rickard and Pascoe (2009)
G48224	TrGTerCN	C721CO3 + NO_3 → C721O2 + CO_2 + NO_2	KRO2NO3*1.74	Rickard and Pascoe (2009)
G48225	TrGTerC	C721CO3 → C721O2 + CO_2	k1_RO2RCO3*0.9	Sander et al. (2018)
G48226	TrGTerC	C721CO3 → NORPINIC	k1_RO2RCO3*0.1	Sander et al. (2018)
G48227	TrGTerC	C721CO3H + OH → C721CO3	9.65E-12	Rickard and Pascoe (2009)
G48228	TrGTerC	NORPINIC + OH → C721O2 + CO_2	6.57E-12	Rickard and Pascoe (2009)
G48229	TrGTerCN	C721PAN + OH → C721OOH + CO + NO_2	2.96E-12	Rickard and Pascoe (2009)
G48230	TrGTerCN	C721PAN → C721CO3 + NO_2	k_PAN_M	Rickard and Pascoe (2009)
G48231	TrGTerC	C8BC + OH → C8BCO2	3.04E-12	Rickard and Pascoe (2009)
G48232	TrGTerC	C8BCO2 + HO_2 → C8BCOOH	KRO2HO2(8)	Rickard and Pascoe (2009)
G48233a	TrGTerCN	C8BCO2 + NO → C89O2 + NO_2	KRO2NO*(1.-alpha_AN(8,2,0,0,0, temp,cair))	Rickard and Pascoe (2009)
G48233b	TrGTerCN	C8BCO2 + NO → C8BCNO3	KRO2NO*alpha_AN(8,2,0,0,0,temp, cair)	Rickard and Pascoe (2009)
G48234	TrGTerC	C8BCO2 → C89O2	k1_RO2sRO2	Rickard and Pascoe (2009)
G48235	TrGTerC	C8BCOOH + OH → C8BCCO + OH	1.62E-11	Rickard and Pascoe (2009)
G48236	TrGTerCN	C8BCNO3 + OH → C8BCCO + NO_2	1.84E-12	Rickard and Pascoe (2009)
G48237	TrGTerC	C8BCCO + OH → C89O2	3.94E-12	Rickard and Pascoe (2009)
G48238	TrGTerC	C89O2 + HO_2 → C89OOH	KRO2HO2(8)	Rickard and Pascoe (2009)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G48239a	TrGTerCN	C89O2 + NO → C810O2 + NO_2	KRO2NO*(1.-alpha_AN(7,2,0,0,0, temp,cair))	Rickard and Pascoe (2009)
G48239b	TrGTerCN	C89O2 + NO → C89NO3	KRO2NO*alpha_AN(7,2,0,0,0,temp, cair)	Rickard and Pascoe (2009)
G48240	TrGTerCN	C89O2 + NO_3 → C810O2 + NO_2	KRO2NO3	Rickard and Pascoe (2009)
G48241	TrGTerC	C89O2 → C810O2	k1_RO2tRO2	Rickard and Pascoe (2009)
G48242	TrGTerC	C89OOH + OH → C89O2	3.61E-11	Rickard and Pascoe (2009)
G48243	TrGTerCN	C89NO3 + OH → CH_3COCH_3 + CO13C4CHO + NO_2	2.56E-11	Rickard and Pascoe (2009)
G48244	TrGTerC	C810O2 + HO_2 → C810OOH	KRO2HO2(8)	Rickard and Pascoe (2009)
G48245a	TrGTerCN	C810O2 + NO → CH_3COCH_3 + C514O2 + NO_2	KRO2NO*(1.-alpha_AN(10,3,0,0,0, temp,cair))	Rickard and Pascoe (2009)
G48245b	TrGTerCN	C810O2 + NO → C810NO3	KRO2NO*alpha_AN(10,3,0,0,0, temp,cair)	Rickard and Pascoe (2009)
G48246	TrGTerCN	C810O2 + NO_3 → CH_3COCH_3 + C514O2 + NO_2	KRO2NO3	Rickard and Pascoe (2009)
G48247	TrGTerC	C810O2 → CH_3COCH_3 + C514O2	k1_RO2tRO2	Rickard and Pascoe (2009)
G48248	TrGTerC	C810OOH + OH → C810O2	8.35E-11	Rickard and Pascoe (2009)
G48249	TrGTerCN	C810NO3 + OH → CH_3COCH_3 + CO13C4CHO + NO_2	4.96E-11	Rickard and Pascoe (2009)
G48400a	TrGAroC	LXYL + OH → TLEPOXMUC + HO_2 + LCARBON + 0.063 ALOV + 0.424 ASOV	0.401E-11	Rickard and Pascoe (2009)*
G48400b	TrGAroC	LXYL + OH → C6H5CH2O2 + LCARBON	0.101E-11	Rickard and Pascoe (2009)*
G48400c	TrGAroC	LXYL + OH → CRESOL + LCARBON	0.261E-11	Rickard and Pascoe (2009)*
G48400d	TrGAroC	LXYL + OH → TLBIPERO2 + HO_2 + LCARBON	0.932E-11	Rickard and Pascoe (2009)*
G48401	TrGAroCN	LXYL + NO_3 → C6H5CH2O2 + HNO_3 + LCARBON	3.9E-16	Rickard and Pascoe (2009)*
G48402	TrGAroC	EBENZ + OH → .10 TLEPOXMUC + .07 C6H5CH2O2 + .18 CRESOL + .65 TLBIPERO2 + .28 HO_2 + LCARBON	7.00E-12	Rickard and Pascoe (2009)*
G48403	TrGAroCN	EBENZ + NO_3 → C6H5CH2O2 + HNO_3 + LCARBON	1.20E-16	Rickard and Pascoe (2009)*
G48404	TrGAroCN	STYRENE + NO_3 → NSTYRENO2	1.50E-12	Rickard and Pascoe (2009)
G48405	TrGAroC	STYRENE + O_3 → .545 HCHO + .1 BENZENE + .28 C6H5O2 + .56 CO + .36 OH + .28 HO_2 + .075 PHCOOH + .545 BENZAL + .09 H_2O_2 + .075 HCOOH + .2 CO_2	1.70E-17	Rickard and Pascoe (2009)*
G48406	TrGAroC	STYRENE + OH → STYRENO2	5.80E-11	Rickard and Pascoe (2009)
G48407	TrGAroCN	NSTYRENO2 + HO_2 → NSTYRENOOH	KRO2HO2(8)	Rickard and Pascoe (2009)
G48408	TrGAroCN	NSTYRENO2 + NO → NO_2 + NO_2 + HCHO + BENZAL	KRO2NO	Rickard and Pascoe (2009)*
G48409	TrGAroCN	NSTYRENO2 + NO_3 → NO_2 + NO_2 + HCHO + BENZAL	KRO2NO3	Rickard and Pascoe (2009)*

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G48410	TrGAroCN	NSTYRENO2 → NO_2 + HCHO + BENZAL	k1_RO2sRO2	Rickard and Pascoe (2009)*
G48411	TrGAroCN	NSTYRENOOH + OH → NSTYRENO2	6.16E-11	Rickard and Pascoe (2009)
G48412a	TrGAroC	STYRENO2 + HO_2 → STYRENOOH	KRO2HO2(8)*(1-rchohch2o2_oh)	Rickard and Pascoe (2009)
G48412b	TrGAroC	STYRENO2 + HO_2 → HO_2 + OH + HCHO + BENZAL	KRO2HO2(8)*rchohch2o2_oh	Rickard and Pascoe (2009)*
G48413	TrGAroCN	STYRENO2 + NO → NO_2 + HO_2 + HCHO + BENZAL	KRO2NO	Rickard and Pascoe (2009)*
G48414	TrGAroCN	STYRENO2 + NO_3 → NO_2 + HO_2 + HCHO + BENZAL	KRO2NO3	Rickard and Pascoe (2009)*
G48415	TrGAroC	STYRENO2 → HO_2 + HCHO + BENZAL	k1_RO2sRO2	Rickard and Pascoe (2009)*
G48416	TrGAroC	STYRENOOH + OH → STYRENO2	6.16E-11	Rickard and Pascoe (2009)
G49200	TrGTerC	C96O2 → C97O2	k1_RO2pRO2	Rickard and Pascoe (2009)
G49201	TrGTerC	C96O2 + HO_2 → C96OOH	KRO2HO2(9)	Rickard and Pascoe (2009)
G49202a	TrGTerCN	C96O2 + NO → C97O2 + NO_2	KRO2NO*(1.-alpha_AN(10,1,0,0,0, temp,cair))	Rickard and Pascoe (2009)
G49202b	TrGTerCN	C96O2 + NO → C96NO3	KRO2NO*alpha_AN(10,1,0,0,0, temp,cair)	Rickard and Pascoe (2009)
G49203	TrGTerCN	C96NO3 + OH → NORPINAL + NO_2	2.88E-12	Rickard and Pascoe (2009)
G49204a	TrGTerC	C96OOH + OH → C96O2	k_roohro	Rickard and Pascoe (2009)
G49205b	TrGTerC	C96OOH + OH → NORPINAL + OH	1.30E-11	Rickard and Pascoe (2009)
G49206	TrGTerC	C97O2 → C98O2	k1_RO2tRO2	Rickard and Pascoe (2009)
G49207	TrGTerCN	C97O2 + NO → C98O2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G49208a	TrGTerC	C97O2 + HO_2 → C97OOH	KRO2HO2(9)*rcoch2o2_ooh	Rickard and Pascoe (2009), Sander et al. (2018)
G49208b	TrGTerC	C97O2 + HO_2 → C98O2 + OH	KRO2HO2(9)*rcoch2o2_oh	Rickard and Pascoe (2009), Sander et al. (2018)
G49209	TrGTerC	C97OOH + OH → C97O2	1.05E-11	Rickard and Pascoe (2009)
G49210	TrGTerC	C98O2 → C614O2 + CH_3COCH_3	k1_RO2tRO2	Rickard and Pascoe (2009)
G49211a	TrGTerCN	C98O2 + NO → C614O2 + CH_3COCH_3 + NO_2	KRO2NO*(1.-alpha_AN(12,3,0,0,0, temp,cair))	Rickard and Pascoe (2009)
G49211b	TrGTerCN	C98O2 + NO → 9 LCARBON + LNITROGEN	KRO2NO*alpha_AN(12,3,0,0,0, temp,cair)	Rickard and Pascoe (2009)
G49212	TrGTerC	C98O2 + HO_2 → C98OOH	KRO2HO2(9)	Rickard and Pascoe (2009)
G49213	TrGTerC	C98OOH + OH → C98O2	2.05E-11	Rickard and Pascoe (2009)
G49214	TrGTerC	NORPINAL + OH → C85CO3	2.64E-11	Rickard and Pascoe (2009)
G49215	TrGTerCN	NORPINAL + NO_3 → C85CO3 + HNO_3	KNO3AL*8.5	Rickard and Pascoe (2009)
G49216	TrGTerC	C85CO3 → C85O2 + CO_2	k1_RO2RCO3	Rickard and Pascoe (2009)
G49217	TrGTerCN	C85CO3 + NO → C85O2 + CO_2 + NO_2	KAPNO	Rickard and Pascoe (2009)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G49218	TrGTerCN	C85CO3 + NO_2 → C9PAN2	k_CH3CO3_NO2	Rickard and Pascoe (2009)
G49219a	TrGTerC	C85CO3 + HO_2 → C85CO3H	KAPHO2*(rco3_ooh+rco3_o3)	Rickard and Pascoe (2009)
G49219b	TrGTerC	C85CO3 + HO_2 → C85O2 + CO_2 + OH	KAPHO2*rco3_oh	Rickard and Pascoe (2009)
G49220	TrGTerCN	C9PAN2 → C85CO3 + NO_2	k_PAN_M	Rickard and Pascoe (2009)
G49221	TrGTerCN	C9PAN2 + OH → C85OOH + CO + NO_2	6.60E-12	Rickard and Pascoe (2009)
G49222	TrGTerC	C85CO3H + OH → C85CO3	1.02E-11	Rickard and Pascoe (2009)
G49223a	TrGTerC	C89CO3 → .8 C811CO3 + .2 C89O2 + .2 CO_2	k1_RO2RCO3*0.9	Sander et al. (2018)
G49223b	TrGTerC	C89CO3 → C89CO2H	k1_RO2RCO3*0.1	Sander et al. (2018)
G49224a	TrGTerC	C89CO3 + HO_2 → C89CO3H	KAPHO2*rco3_ooh	Rickard and Pascoe (2009)
G49224b	TrGTerC	C89CO3 + HO_2 → C89CO2H + O_3	KAPHO2*rco3_o3	Rickard and Pascoe (2009)
G49224c	TrGTerC	C89CO3 + HO_2 → .80 C811CO3 + .20 C89O2 + .2 CO_2 + OH	KAPHO2*rco3_oh	Rickard and Pascoe (2009)
G49225	TrGTerCN	C89CO3 + NO_2 → C89PAN	k_CH3CO3_NO2	Rickard and Pascoe (2009)
G49226	TrGTerCN	C89CO3 + NO → .8 C811CO3 + .2 C89O2 + .2 CO_2 + NO_2	KAPNO	Rickard and Pascoe (2009)
G49227	TrGTerC	C89CO2H + OH → .8 C811CO3 + .2 C89O2 + .2 CO_2	2.69E-11	Rickard and Pascoe (2009)
G49228	TrGTerC	C89CO3H + OH → C89CO3	3.00E-11	Rickard and Pascoe (2009)
G49229	TrGTerCN	C89PAN → C89CO3 + NO_2	k_PAN_M	Rickard and Pascoe (2009)
G49230	TrGTerCN	C89PAN + OH → CH_3COCH_3 + CO13C4CHO + CO + NO_2	2.52E-11	Rickard and Pascoe (2009)
G49231a	TrGTerC	C811CO3 → C811O2 + CO_2	k1_RO2RCO3*0.9	Sander et al. (2018)
G49231b	TrGTerC	C811CO3 → PINIC	k1_RO2RCO3*0.1	Sander et al. (2018)
G49232a	TrGTerC	C811CO3 + HO_2 → C811CO3H	KAPHO2*rco3_ooh	Rickard and Pascoe (2009)
G49232b	TrGTerC	C811CO3 + HO_2 → PINIC + O_3	KAPHO2*rco3_o3	Rickard and Pascoe (2009)
G49232c	TrGTerC	C811CO3 + HO_2 → C811O2 + CO_2 + OH	KAPHO2*rco3_oh	Rickard and Pascoe (2009)
G49233	TrGTerCN	C811CO3 + NO → C811O2 + CO_2 + NO_2	KAPNO	Rickard and Pascoe (2009)
G49234	TrGTerCN	C811CO3 + NO_2 → C811PAN	k_CH3CO3_NO2	Rickard and Pascoe (2009)
G49235	TrGTerC	PINIC + OH → C811O2 + CO_2	7.29E-12	Rickard and Pascoe (2009)
G49236	TrGTerC	NOPINONE + OH → NOPINDO2	1.55E-11	Capouet et al. (2008), Rickard and Pascoe (2009)
G49237a	TrGTerC	NOPINDO2 + HO_2 → NOPINDOOH	KRO2HO2(9)*rcoch2o2_ooh	Rickard and Pascoe (2009), Sander et al. (2018)
G49237b	TrGTerC	NOPINDO2 + HO_2 → C89CO3 + OH	KRO2HO2(9)*rcoch2o2_oh	Rickard and Pascoe (2009), Sander et al. (2018)
G49238	TrGTerCN	NOPINDO2 + NO → C89CO3 + NO_2	KRO2NO	Rickard and Pascoe (2009)*

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G49239	TrGTerC	NOPINDO2 → C89CO3	k1_RO2pORO2	Rickard and Pascoe (2009)
G49240	TrGTerC	NOPINDOOH → NOPINDCO	2.63E-11	Rickard and Pascoe (2009)
G49241	TrGTerC	NOPINDCO + OH → C89CO3	3.07E-12	Rickard and Pascoe (2009)
G49242	TrGTerC	NOPINOO → NOPINONE + H_2O_2	6.00E-18*c(ind_H2O)	Rickard and Pascoe (2009)
G49243	TrGTerC	NOPINOO + CO → NOPINONE + CO_2	1.2E-15	Rickard and Pascoe (2009)
G49244	TrGTerCN	NOPINOO + NO → NOPINONE + NO_2	1.E-14	Rickard and Pascoe (2009)
G49245	TrGTerCN	NOPINOO + NO_2 → NOPINONE + NO_3	1.E-15	Rickard and Pascoe (2009)
G49246	TrGTerC	NORPINENOL + OH → HCOOH + OH + C86O2	k_CH2CHOH_OH_HCOOH	Sander et al. (2018), So et al. (2014)*
G49247	TrGTerC	NORPINENOL + HCOOH → NORPINAL + HCOOH	k_CH2CHOH_HCOOH	Sander et al. (2018), da Silva (2010)*
G49248	TrGTerC	NORPINAL + HCOOH → NORPINENOL + HCOOH	k_ALD_HCOOH	Sander et al. (2018), da Silva (2010)*
G49249	TrGTerC	C811CO3H + OH → C811CO3	1.04E-11	Rickard and Pascoe (2009)
G49250	TrGTerCN	C811PAN → C811CO3 + NO_2	k_PAN_M	Rickard and Pascoe (2009)
G49251	TrGTerCN	C811PAN + OH → C721CHO + CO + NO_2	6.77E-12	Rickard and Pascoe (2009)
G49400a	TrGAroC	LTMB + OH → TLBIPERO2 + HO_2 + 2 LCARBON + .04 ALOV	2.917E-11	Rickard and Pascoe (2009)*
G49400b	TrGAroC	LTMB + OH → C6H5CH2O2 + 2 LCARBON	0.189E-11	Rickard and Pascoe (2009)*
G49400c	TrGAroC	LTMB + OH → CRESOL + 2 LCARBON	0.141E-11	Rickard and Pascoe (2009)*
G49400d	TrGAroC	LTMB + OH → TLBIPERO2 + HO_2 + 2 LCARBON	2.917E-11	Rickard and Pascoe (2009)*
G49401	TrGAroCN	LTMB + NO_3 → C6H5CH2O2 + HNO_3 + 2 LCARBON	1.52E-15	Rickard and Pascoe (2009)*
G40200	TrGTerC	APINENE + OH → .75 LAPINABO2 + .15 MENTHEN6ONE + .15 HO_2 + .10 ROO6R1O2	1.2E-11*EXP(440./TEMP)	Atkinson et al. (2006)*
G40201a	TrGTerCN	LAPINABO2 + NO → PINAL + HO_2 + NO_2 + 0.052 BLOV + 0.184 BSOV	KRO2NO*(1.-(.65*alpha_AN(11,3,0,0,0,temp,cair)+.35*alpha_AN(11,2,0,0,0,temp,cair)))	Rickard and Pascoe (2009), Sander et al. (2018)
G40201b	TrGTerCN	LAPINABO2 + NO → LAPINABNO3	KRO2NO*(.65*alpha_AN(11,3,0,0,0,temp,cair)+.35*alpha_AN(11,2,0,0,0,temp,cair))	Rickard and Pascoe (2009), Sander et al. (2018)
G40202a	TrGTerC	LAPINABO2 + HO_2 → LAPINABOOH	KRO2HO2(10)*(1.-rchohch2o2_oh)	Rickard and Pascoe (2009), Sander et al. (2018)
G40202b	TrGTerC	LAPINABO2 + HO_2 → PINAL + HO_2 + OH	(1-ya_soan1-ya_soan2)*KRO2HO2(10)*rchohch2o2_oh	Rickard and Pascoe (2009), Sander et al. (2018)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G40203	TrGTerC	LAPINABO2 → PINAL + HO_2	`RO2*(0.65*k1_RO2tORO2+.35*k1_RO2sORO2)`	Rickard and Pascoe (2009)*
G40204	TrGTerC	LAPINABOOH + OH → .35 LAPINABO2 + .65 C96CO3	`2.77E-11`	Rickard and Pascoe (2009)*
G40205	TrGTerCN	LAPINABNO3 + OH → .35 PINAL + .65 C96CO3 + NO_2	`4.29E-12`	Rickard and Pascoe (2009)*
G40206	TrGTerC	MENTHEN6ONE + OH → OHMENTHEN6ONEO2	`6.46E-11`	Vereecken et al. (2007)*
G40207	TrGTerCN	OHMENTHEN6ONEO2 + NO → 2OHMENTHEN6ONE + HO_2 + NO_2	`KRO2NO`	Vereecken et al. (2007)*
G40208	TrGTerC	OHMENTHEN6ONEO2 + HO_2 → 2OHMENTHEN6ONE	`KRO2HO2(10)`	Vereecken et al. (2007)
G40209	TrGTerC	OHMENTHEN6ONEO2 → 2OHMENTHEN6ONE + HO_2	`k1_RO2tORO2`	Vereecken et al. (2007)
G40210	TrGTerC	2OHMENTHEN6ONE + OH → 10 LCARBON	`1E-11`	Vereecken et al. (2007)
G40211	TrGTerC	PINAL + OH → .772 C96CO3 + .228 PINALO2	`5.2E-12*EXP(600./TEMP)`	Wallington et al. (2018)*
G40212	TrGTerCN	PINAL + NO_3 → C96CO3 + HNO_3	`2.0E-14`	Wallington et al. (2018)*
G40213a	TrGTerC	C96CO3 → C96O2 + CO_2	`k1_RO2RCO3*0.9`	Rickard and Pascoe (2009)
G40213b	TrGTerC	C96CO3 → PINONIC	`k1_RO2RCO3*0.1`	Rickard and Pascoe (2009)
G40214a	TrGTerC	C96CO3 + HO_2 → PERPINONIC	`KAPHO2*rco3_ooh`	Rickard and Pascoe (2009)
G40214b	TrGTerC	C96CO3 + HO_2 → PINONIC + O_3	`KAPHO2*rco3_o3`	Rickard and Pascoe (2009)
G40214c	TrGTerC	C96CO3 + HO_2 → C96O2 + OH + CO_2	`KAPHO2*rco3_oh`	Rickard and Pascoe (2009)
G40215	TrGTerCN	C96CO3 + NO_2 → C10PAN2	`k_CH3CO3_NO2`	Rickard and Pascoe (2009)
G40216	TrGTerCN	C96CO3 + NO → C96O2 + NO_2 + CO_2	`KAPNO`	Rickard and Pascoe (2009)
G40217	TrGTerCN	C96CO3 + NO_3 → C96O2 + NO_2 + CO_2	`KRO2NO3*1.74`	Rickard and Pascoe (2009)
G40218	TrGTerCN	C10PAN2 → C96CO3 + NO_2	`k_PAN_M`	Rickard and Pascoe (2009)
G40219	TrGTerCN	C10PAN2 + OH → NORPINAL + CO + NO_2	`3.66E-12`	Rickard and Pascoe (2009)
G40220	TrGTerC	PINONIC + OH → C96O2 + CO_2	`6.65E-12`	Rickard and Pascoe (2009)
G40221	TrGTerC	PERPINONIC + OH → C96CO3	`9.73E-12`	Rickard and Pascoe (2009)
G40222	TrGTerC	PINALO2 + HO_2 → PINALOOH	`KRO2HO2(10)`	Rickard and Pascoe (2009)
G40223a	TrGTerCN	PINALO2 + NO → C106O2 + NO_2	`KRO2NO*(1.-alpha_AN(12,3,0,1,0,temp,cair))`	Rickard and Pascoe (2009), Sander et al. (2018)
G40223b	TrGTerCN	PINALO2 + NO → PINALNO3	`KRO2NO*alpha_AN(12,3,0,1,0,temp,cair)`	Rickard and Pascoe (2009), Sander et al. (2018)
G40224	TrGTerC	PINALO2 → C106O2	`k1_RO2tRO2`	Rickard and Pascoe (2009)
G40225	TrGTerC	PINALOOH + OH → PINALO2	`2.75E-11`	Rickard and Pascoe (2009)
G40226	TrGTerCN	PINALNO3 + OH → CO235C6CHO + CH_3COCH_3 + NO_2	`2.25E-11`	Rickard and Pascoe (2009)
G40227	TrGTerC	C106O2 + HO_2 → C106OOH	`KRO2HO2(10)`	Rickard and Pascoe (2009)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G40228a	TrGTerCN	C106O2 + NO → C716O2 + CH_3COCH_3 + NO_2	KRO2NO*0.875*(1.-alpha_AN(13,3,0,0,0,temp,cair))	Rickard and Pascoe (2009), Sander et al. (2018)
G40228b	TrGTerCN	C106O2 + NO → C106NO3	KRO2NO*0.875*alpha_AN(13,3,0,0,0,temp,cair)	Rickard and Pascoe (2009), Sander et al. (2018)
G40229	TrGTerC	C106O2 → C716O2 + CH_3COCH_3	k1_RO2tRO2	Rickard and Pascoe (2009)
G40230	TrGTerC	C106OOH + OH → C106O2	8.01E-11	Rickard and Pascoe (2009)
G40231	TrGTerCN	C106NO3 + OH → CO235C6CHO + CH_3COCH_3 + NO_2	7.03E-11	Rickard and Pascoe (2009)
G40232	TrGTerC	APINENE + O_3 → .09 APINBOO + .08 PINONIC + .77 OH + .33 NORPINAL + .33 CO + .33 HO_2 + .06 APINAOO + .44 C109O2 + 0.07 BELV + 0.13 BLOV	8.05E-16*EXP(-640./TEMP)	Wallington et al. (2018), Ehn et al. (2014)*
G40233	TrGTerC	APINAOO → PINAL + H_2O_2	1.00E-17*c(ind_H2O)	Rickard and Pascoe (2009)
G40234	TrGTerC	APINAOO + CO → PINAL + CO_2	1.20E-15	Rickard and Pascoe (2009)
G40235	TrGTerCN	APINAOO + NO → PINAL + NO_2 + 0.052 BLOV + 0.184 BSOV	1.00E-14	Rickard and Pascoe (2009)
G40236	TrGTerCN	APINAOO + NO_2 → PINAL + NO_3	1.00E-15	Rickard and Pascoe (2009)
G40237a	TrGTerC	APINBOO → PINONIC	1.00E-17*c(ind_H2O)*(0.08+0.15)	Rickard and Pascoe (2009)
G40237b	TrGTerC	APINBOO → PINAL + H_2O_2	1.00E-17*c(ind_H2O)*0.77	Rickard and Pascoe (2009)
G40238	TrGTerC	APINBOO + CO → PINAL + CO_2	1.20E-15	Rickard and Pascoe (2009)
G40239	TrGTerCN	APINBOO + NO → PINAL + NO_2 + 0.052 BLOV + 0.184 BSOV	1.00E-14	Rickard and Pascoe (2009)
G40240	TrGTerCN	APINBOO + NO_2 → PINAL + NO_3	1.00E-15	Rickard and Pascoe (2009)
G40241	TrGTerC	C109O2 → C89CO3 + HCHO	k1_RO2pORO2	Rickard and Pascoe (2009)
G40242	TrGTerCN	C109O2 + NO → C89CO3 + HCHO + NO_2 + 0.052 BLOV + 0.184 BSOV	KRO2NO	Rickard and Pascoe (2009)*
G40243a	TrGTerC	C109O2 + HO_2 → C109OOH	KRO2HO2(10)*rcoch2o2_ooh	Rickard and Pascoe (2009), Sander et al. (2018)
G40243b	TrGTerC	C109O2 + HO_2 → C89CO3 + HCHO + OH	(1-ya_soan1-ya_soan2)*KRO2HO2(10)*rcoch2o2_oh	Rickard and Pascoe (2009), Sander et al. (2018)
G40244	TrGTerC	C109OOH + OH → C109CO + OH	5.47E-11	Rickard and Pascoe (2009)
G40245	TrGTerC	C109CO + OH → C89CO3 + CO	5.47E-11	Rickard and Pascoe (2009)
G40246	TrGTerCN	APINENE + NO_3 → LNAPINABO2	1.2E-12*EXP(490./temp)	Wallington et al. (2018)*
G40247	TrGTerCN	LNAPINABO2 → PINAL + NO_2	(0.65*k1_RO2tRO2 + 0.35*k1_RO2sRO2)	Rickard and Pascoe (2009)
G40248	TrGTerCN	LNAPINABO2 + NO → PINAL + NO_2 + NO_2	KRO2NO	Rickard and Pascoe (2009)*
G40249	TrGTerCN	LNAPINABO2 + HO_2 → LNAPINABOOH	KRO2HO2(10)	Rickard and Pascoe (2009)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G40250	TrGTerCN	LNAPINABO2 + NO_3 → PINAL + NO_2 + NO_2	`KRO2NO3`	Rickard and Pascoe (2009)
G40251	TrGTerCN	LNAPINABOOH + OH → LNAPINABO2	`(.65*6.87E-12+.35*1.23E-11)`	Rickard and Pascoe (2009)
G40252a	TrGTerC	BPINENE + OH → BPINAO2	`1.47E-11*EXP(467./TEMP) *(0.8326*0.3+0.068)/(0.8326+0.068)`	Gill and Hites (2002)*
G40252b	TrGTerC	BPINENE + OH → ROO6R1O2	`1.47E-11*EXP(467./TEMP) *0.8326*0.7/(0.8326+0.068)`	Gill and Hites (2002)*
G40253a	TrGTerC	BPINAO2 + HO_2 → BPINAOOH	`KRO2HO2(10)*rcoch2o2_ooh`	Rickard and Pascoe (2009), Sander et al. (2018)
G40253b	TrGTerC	BPINAO2 + HO_2 → NOPINONE + HCHO + HO_2 + OH	`(1-ya_soan1-ya_soan2) *KRO2HO2(10)*rcoch2o2_oh`	Rickard and Pascoe (2009), Sander et al. (2018)
G40254a	TrGTerCN	BPINAO2 + NO → NOPINONE + HCHO + HO_2 + NO_2 + 0.052 BLOV + 0.184 BSOV	`KRO2NO*(1.-alpha_AN(11,3,0,0,0, temp,cair))`	Rickard and Pascoe (2009), Sander et al. (2018)
G40254b	TrGTerCN	BPINAO2 + NO → BPINANO3	`KRO2NO*alpha_AN(11,3,0,0,0, temp,cair)`	Rickard and Pascoe (2009), Sander et al. (2018)
G40255	TrGTerC	BPINAO2 → NOPINONE + HCHO + HO_2	`k1_RO2tORO2`	Rickard and Pascoe (2009)
G40256	TrGTerC	BPINAOOH + OH → BPINAO2	`1.33E-11`	Rickard and Pascoe (2009)
G40257	TrGTerCN	BPINANO3 + OH → NOPINONE + HCHO + NO_2	`4.70E-12`	Rickard and Pascoe (2009)
G40258a	TrGTerCN	ROO6R1O2 + NO → ROO6R3O2 + CH_3COCH_3 + NO_2 + 0.052 BLOV + 0.184 BSOV	`KRO2NO*(1.-alpha_AN(13,3,0,0,0, temp,cair))`	Vereecken and Peeters (2012)
G40258b	TrGTerCN	ROO6R1O2 + NO → ROO6R1NO3	`KRO2NO*alpha_AN(13,3,0,0,0, temp,cair)`	Vereecken and Peeters (2012)
G40260	TrGTerC	ROO6R1O2 → ROO6R3O2 + CH_3COCH_3	`k1_RO2tORO2`	Vereecken and Peeters (2012)
G40261a	TrGTerCN	RO6R1O2 + NO → RO6R3O2 + NO_2	`KRO2NO*(1.-alpha_AN(12,3,0,0,0, temp,cair))`	Vereecken and Peeters (2012)
G40261b	TrGTerCN	RO6R1O2 + NO → RO6R1NO3	`KRO2NO*alpha_AN(12,3,0,0,0, temp,cair)`	Vereecken and Peeters (2012)
G40262	TrGTerC	RO6R1O2 + HO_2 → 10 LCARBON	`KRO2HO2(10)`	Vereecken and Peeters (2012)*
G40263	TrGTerC	RO6R1O2 → RO6R3O2	`k1_RO2sORO2`	Vereecken and Peeters (2012)
G40264a	TrGTerCN	RO6R3O2 + NO → 9 LCARBON + HCHO + HO_2 + NO_2	`KRO2NO*(1.-alpha_AN(12,3,0,0,0, temp,cair))`	Vereecken and Peeters (2012)
G40264b	TrGTerCN	RO6R3O2 + NO → 10 LCARBON + LNITROGEN	`KRO2NO*alpha_AN(12,3,0,0,0, temp,cair)`	Vereecken and Peeters (2012)
G40265	TrGTerC	RO6R3O2 + HO_2 → 10 LCARBON	`KRO2HO2(10)`	Vereecken and Peeters (2012)
G40266	TrGTerC	RO6R3O2 → 9 LCARBON + HCHO + HO_2	`k1_RO2sRO2`	Vereecken and Peeters (2012)*

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G40267a	TrGTerC	BPINENE + O_3 → NOPINONE + .63 CO + .37 CH_2OO + .16 OH + .16 HO_2 + 0.07 BELV + 0.13 BLOV	`1.35E-15*EXP(-1270./TEMP) *.051/(1-.027)`	Wallington et al. (2018)*
G40267b	TrGTerC	BPINENE + O_3 → NOPINOO + CO_2	`1.35E-15*EXP(-1270./TEMP) *.368/(1-.027)`	Nguyen et al. (2009), Wallington et al. (2018)
G40267c	TrGTerC	BPINENE + O_3 → NOPINDO2 + CO_2 + OH	`1.35E-15*EXP(-1270./TEMP) *.283/(1-.027)`	Nguyen et al. (2009), Wallington et al. (2018)
G40267d	TrGTerC	BPINENE + O_3 → C8BC + 2 CO_2	`1.35E-15*EXP(-1270./TEMP) *(.104+.167)/(1-.027)`	Nguyen et al. (2009), Wallington et al. (2018)
G40268	TrGTerCN	BPINENE + NO_3 → LNBPINABO2	`2.51E-12`	Wallington et al. (2018)*
G40269	TrGTerCN	LNBPINABO2 + HO_2 → LNBPINABOOH	`KRO2HO2(10)`	Rickard and Pascoe (2009)
G40270	TrGTerCN	LNBPINABO2 + NO → NOPINONE + HCHO + NO_2 + NO_2	`KRO2NO`	Rickard and Pascoe (2009)*
G40271	TrGTerCN	LNBPINABO2 + NO_3 → NOPINONE + HCHO + NO_2 + NO_2	`KRO2NO3`	Rickard and Pascoe (2009)
G40272a	TrGTerCN	LNBPINABO2 → NOPINONE + HCHO + NO_2	`k1_RO2tRO2*0.7`	Rickard and Pascoe (2009)
G40272b	TrGTerCN	LNBPINABO2 → BPINANO3	`k1_RO2tRO2*0.3`	Rickard and Pascoe (2009)
G40273	TrGTerCN	LNBPINABOOH + OH → LNBPINABO2	`9.58E-12`	Rickard and Pascoe (2009)
G40274	TrGTerCN	ROO6R1NO3 + OH → ROO6R3O2 + CH_3COCH_3 + NO_2	`9.16E-13`	Vereecken and Peeters (2012), Gill and Hites (2002)*
G40275	TrGTerCN	RO6R1NO3 + OH → 9 LCARBON + HCHO + HO_2 + NO_2	`9.16E-13`	Vereecken and Peeters (2012), Gill and Hites (2002)
G40276	TrGTerC	PINEOL + OH → HCOOH + OH + NORPINAL	`k_CH2CHOH_OH_HCOOH`	Sander et al. (2018), So et al. (2014)*
G40277	TrGTerC	PINEOL + HCOOH → PINAL + HCOOH	`k_CH2CHOH_HCOOH`	Sander et al. (2018), da Silva (2010)*
G40278	TrGTerC	PINAL + HCOOH → PINEOL + HCOOH	`k_ALD_HCOOH`	Sander et al. (2018), da Silva (2010)*
G40279a	TrGC	CARENE + OH → LAPINABO2	`8.8E-11*(.50+.25)`	Atkinson and Arey (2003)
G40279b	TrGC	CARENE + OH → MENTHEN6ONE + HO_2	`8.8E-11*.25*.60`	Atkinson and Arey (2003)
G40279c	TrGC	CARENE + OH → ROO6R1O2	`8.8E-11*.25*.40`	Atkinson and Arey (2003)
G40280a	TrGC	CARENE + O_3 → APINBOO	`3.7E-17*.50*.18`	Atkinson and Arey (2003)
G40280b	TrGC	CARENE + O_3 → PINONIC	`3.7E-17*.50*.16`	Atkinson and Arey (2003)
G40280c	TrGC	CARENE + O_3 → OH + NORPINAL + CO + HO_2	`3.7E-17*.50*.66`	Atkinson and Arey (2003)
G40280d	TrGC	CARENE + O_3 → APINAOO	`3.7E-17*.50*.12`	Atkinson and Arey (2003)
G40280e	TrGC	CARENE + O_3 → OH + C109O2	`3.7E-17*.50*(.22+.66)`	Atkinson and Arey (2003)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G40281	TrGCN	CARENE + NO_3 → LNAPINABO2	9.1E-12	Atkinson and Arey (2003)
G40282a	TrGTerC	SABINENE + OH → BPINAO2	1.47E-11*EXP(467./TEMP) *(0.8326*0.3+0.068)/(0.8326+0.068)	Gill and Hites (2002)*
G40282b	TrGTerC	SABINENE + OH → ROO6R1O2	1.47E-11*EXP(467./TEMP) *0.8326*0.7/(0.8326+0.068)	Vereecken and Peeters (2012), Gill and Hites (2002)*
G40283a	TrGTerC	SABINENE + O_3 → NOPINONE + .63 CO + .37 $HOCH_2OOH$ + .16 OH + .16 HO_2	1.35E-15*EXP(-1270./TEMP) *.051/(1-.027)	Wallington et al. (2018)*
G40283b	TrGTerC	SABINENE + O_3 → NOPINOO + CO_2	1.35E-15*EXP(-1270./TEMP) *.368/(1-.027)	Nguyen et al. (2009), Wallington et al. (2018)
G40283c	TrGTerC	SABINENE + O_3 → NOPINDO2 + CO_2 + OH	1.35E-15*EXP(-1270./TEMP) *.283/(1-.027)	Nguyen et al. (2009), Wallington et al. (2018)
G40283d	TrGTerC	SABINENE + O_3 → C8BC + 2 CO_2	1.35E-15*EXP(-1270./TEMP) *(.104+.167)/(1-.027)	Nguyen et al. (2009), Wallington et al. (2018)
G40284	TrGTerCN	SABINENE + NO_3 → LNBPINABO2	2.51E-12	Wallington et al. (2018)*
G40285a	TrGTerC	CAMPHENE + OH → BPINAO2	1.47E-11*EXP(467./TEMP) *(0.8326*0.3+0.068)/(0.8326+0.068)	Gill and Hites (2002)*
G40285b	TrGTerC	CAMPHENE + OH → ROO6R1O2	1.47E-11*EXP(467./TEMP) *0.8326*0.7/(0.8326+0.068)	Vereecken and Peeters (2012), Gill and Hites (2002)*
G40286a	TrGTerC	CAMPHENE + O_3 → NOPINONE + .63 CO + .37 $HOCH_2OOH$ + .16 OH + .16 HO_2	1.35E-15*EXP(-1270./TEMP) *.051/(1-.027)	Wallington et al. (2018)*
G40286b	TrGTerC	CAMPHENE + O_3 → NOPINOO + CO_2	1.35E-15*EXP(-1270./TEMP) *.368/(1-.027)	Nguyen et al. (2009), Wallington et al. (2018)
G40286c	TrGTerC	CAMPHENE + O_3 → NOPINDO2 + CO_2 + OH	1.35E-15*EXP(-1270./TEMP) *.283/(1-.027)	Nguyen et al. (2009), Wallington et al. (2018)
G40286d	TrGTerC	CAMPHENE + O_3 → C8BC + 2 CO_2	1.35E-15*EXP(-1270./TEMP) *(.104+.167)/(1-.027)	Nguyen et al. (2009), Wallington et al. (2018)
G40287	TrGTerCN	CAMPHENE + NO_3 → LNBPINABO2	2.51E-12	Wallington et al. (2018)*
G40400	TrGAroC	LHAROM + OH → .14 TLEPOXMUC + .03 C6H5CH2O2 + .04 CRESOL + .79 TLBIPERO2 + .18 HO_2 + 4 LCARBON	5.67E-11	Rickard and Pascoe (2009)*
G40401	TrGAroCN	LHAROM + NO_3 → C6H5CH2O2 + HNO_3 + 4 LCARBON	2.60E-15	Rickard and Pascoe (2009)*
G6100	UpStTrGCl	Cl + O_3 → ClO + O_2	2.8E-11*EXP(-250./temp)	Atkinson et al. (2007)
G6102a	StTrGCl	ClO + ClO → Cl_2 + O_2	1.0E-12*EXP(-1590./temp)	Atkinson et al. (2007)
G6102b	StTrGCl	ClO + ClO → 2 Cl + O_2	3.0E-11*EXP(-2450./temp)	Atkinson et al. (2007)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G6102c	StTrGCl	$ClO + ClO \rightarrow Cl + OClO$	3.5E-13*EXP(-1370./temp)	Atkinson et al. (2007)
G6102d	StTrGCl	$ClO + ClO \rightarrow Cl_2O_2$	k_ClO_ClO	Burkholder et al. (2015)
G6103	StTrGCl	$Cl_2O_2 \rightarrow ClO + ClO$	k_ClO_ClO/(2.16E-27*EXP(8537./temp))	Burkholder et al. (2015)*
G6202	StTrGCl	$Cl + H_2O_2 \rightarrow HCl + HO_2$	1.1E-11*EXP(-980./temp)	Atkinson et al. (2007)
G6204	StTrGCl	$ClO + HO_2 \rightarrow HOCl + O_2$	2.2E-12*EXP(340./temp)	Atkinson et al. (2007)*
G6205	StTrGCl	$HCl + OH \rightarrow Cl + H_2O$	1.7E-12*EXP(-230./temp)	Atkinson et al. (2007)
G6300	UpStTrGClN	$ClO + NO \rightarrow NO_2 + Cl$	6.2E-12*EXP(295./temp)	Atkinson et al. (2007)
G6301	StTrGClN	$ClO + NO_2 \rightarrow ClNO_3$	k_3rd_iupac(temp,cair,1.6E-31,3.4,7.E-11,0.,0.4)	Atkinson et al. (2007)
G6302	TrGClN	$ClNO_3 \rightarrow ClO + NO_2$	6.918E-7*EXP(-10909./temp)*cair	Anderson and Fahey (1990)
G6304	StTrGClN	$ClNO_3 + Cl \rightarrow Cl_2 + NO_3$	6.2E-12*EXP(145./temp)	Atkinson et al. (2007)
G6400	StTrGCl	$Cl + CH_4 \rightarrow HCl + CH_3$	6.6E-12*EXP(-1240./temp)	Atkinson et al. (2006)
G6401	StTrGCl	$Cl + HCHO \rightarrow HCl + CO + HO_2$	8.1E-11*EXP(-34./temp)	Atkinson et al. (2006)
G6402	StTrGCl	$Cl + CH_3OOH \rightarrow HCHO + HCl + OH$	5.9E-11	Atkinson et al. (2006)*
G6403	StTrGCl	$ClO + CH_3O_2 \rightarrow HO_2 + Cl + HCHO$	1.8E-12*EXP(-600./temp)	Burkholder et al. (2015)
G6408	StTrGCCl	$CH_3CCl_3 + OH \rightarrow 2\ LCARBON + H_2O + 3\ Cl$	1.64E-12*EXP(-1520./temp)	Burkholder et al. (2015)
G6409	TrGCCl	$Cl + C_2H_4 \rightarrow HOCH_2CH_2O_2 + HCl$	k_3rd_iupac(temp,cair,1.85E-29,3.3,6.0E-10,0.0,0.4)	Atkinson et al. (2006)*
G6410	TrGCCl	$Cl + CH_3CHO \rightarrow HCl + CH_3C(O)$	8.0e-11	Atkinson et al. (2006)
G6411	TrGCCl	$C_2H_2 + Cl \rightarrow LCARBON + CH_3 + HCl$	k_3rd_iupac(temp,cair,6.1e-30,3.0,2.0e-10,0.,0.6)	Atkinson et al. (2006)
G6412	TrGCCl	$C_2H_6 + Cl \rightarrow C_2H_5O_2 + HCl$	8.3E-11*EXP(-100./temp)	Atkinson et al. (2006)
G6413	StTrGClN	$Cl + CH_3ONO_2 \rightarrow HCl + HCHO + NO_2$	1.3E-11*EXP(-1200./temp)	Burkholder et al. (2015)
G6414	StTrGClN	$Cl + CH_3ONO \rightarrow HCl + HCHO + NO$	2.1E-12	Sokolov et al. (1999)
G6415	StTrGCl	$Cl + CH_3O_2 \rightarrow .5\ ClO + .5\ CH_3O + .5\ HCl + .5\ CH_2OO$	1.6E-10	Burkholder et al. (2015)
G6416	TrGCClN	$Cl + CH_3CN \rightarrow NCCH_2O_2 + HCl$	1.6E-11*EXP(-2104./temp)	Tyndall et al. (1996), Tyndall et al. (2001b), Sander et al. (2018)
G8100	TrGI	$I + O_3 \rightarrow IO + O_2$	2.1E-11*EXP(-830./temp)	Atkinson et al. (2007)
G8102	TrGI	$OIO + OIO \rightarrow I(part)$	5.E-11	von Glasow et al. (2002)*
G8103	TrGI	$IO + IO \rightarrow .38\ OIO + 1.62\ I + .62\ O_2$	5.4E-11*EXP(180./temp)	Atkinson et al. (2007)*
G8200	TrGI	$I + HO_2 \rightarrow HI + O_2$	1.5E-11*EXP(-1090./temp)	Atkinson et al. (2007)
G8201	TrGI	$IO + HO_2 \rightarrow HOI + O_2$	1.4E-11*EXP(540./temp)	Atkinson et al. (2007)
G8202	TrGI	$HI + OH \rightarrow I + H_2O$	1.6E-11*EXP(440./temp)	Atkinson et al. (2007)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G8203	TrGI	$OIO + OH \rightarrow HIO_3$	2.2E-10*EXP(243./temp)	Plane et al. (2006)
G8204	TrGI	$I_2 + OH \rightarrow HOI + I$	2.1E-10	Atkinson et al. (2007)
G8205	TrGI	$HOI + OH \rightarrow IO + H_2O$	5.0E-12	Riffault et al. (2005)
G8300	TrGIN	$I + NO_2 \rightarrow INO_2$	k_I_NO2	Atkinson et al. (2007)*
G8301	TrGIN	$I + NO_3 \rightarrow IO + NO_2$	1.E-10	Dillon et al. (2008)
G8302	TrGIN	$IO + NO \rightarrow I + NO_2$	7.15E-12*EXP(300./temp)	Atkinson et al. (2007)
G8303	TrGIN	$IO + NO_2 \rightarrow INO_3$	k_3rd_iupac(temp,cair,7.7E-31, 5.,1.6E-11,0.,0.4)	Atkinson et al. (2007)
G8304	TrGIN	$OIO + NO \rightarrow NO_2 + IO$	1.1E-12*EXP(542./temp)	Atkinson et al. (2007)
G8305	TrGIN	$INO_2 \rightarrow I + NO_2$	k_I_NO2/(3.7E-7*EXP(9568./temp) *1.E6*R_gas*temp/(atm2Pa*N_A))	van den Bergh and Troe (1976), Atkinson et al. (2007)*
G8306	TrGIN	$INO_3 \rightarrow IO + NO_2$	2.1e15*EXP(-13670./temp)	Kaltsoyannis and Plane (2008)
G8307	TrGIN	$I_2 + NO_3 \rightarrow I + INO_3$	1.5E-12	Atkinson et al. (2007)
G8308	TrGIN	$IO + NO_3 \rightarrow OIO + NO_2$	9.E-12	Dillon et al. (2008)
G8309	TrGIN	$I + INO_3 \rightarrow I_2 + NO_3$	9.1E-11*EXP(-146./temp)	Kaltsoyannis and Plane (2008)
G8400	TrGCI	$CH_3CHICH_3 + OH \rightarrow 2\ LCARBON + CH_3O_2 + I$	1.22E-12	Carl and Crowley (2001)
G8401	TrGI	$CH_3O_2 + IO \rightarrow .4\ I + .6\ OIO + HCHO + HO_2$	2.E-12	Dillon et al. (2006b), Bale et al. (2005)*
G8402	TrGIN	$CH_3I + NO_3 \rightarrow HNO_3 + HCHO + IO$	3.4E-17	Wayne et al. (1991)*
G8600	TrGClI	$IO + ClO \rightarrow .2\ ICl + .25\ Cl + .55\ OClO + .8\ I + .45\ O_2$	4.7E-12*EXP(280./temp)	Atkinson et al. (2007)
G9200	StTrGS	$SO_2 + OH \rightarrow HSO_3$	k_3rd(temp,cair,3.E-31,-3.3, 1.5E-12,0.,0.6)	Burkholder et al. (2015)
G9400a	TrGCS	$DMS + OH \rightarrow CH_3SCH_2 + H_2O$	1.13E-11*EXP(-253./temp)	Hynes and Wine (1996)
G9400b	TrGCS	$DMS + OH \rightarrow DMSOH + HO_2$	k_DMS_OH	Hynes and Wine (1996)
G9401	TrGCNS	$DMS + NO_3 \rightarrow CH_3SCH_2 + HNO_3$	1.9E-13*EXP(520./temp)	Karl et al. (2007)
G9402a	TrGCS	$DMSO + OH \rightarrow DMSOHO$	k_DMSO_OH_B1	Karl et al. (2007)
G9402b	TrGCS	$DMSO + OH \rightarrow CH_3SOCH_2 + H_2O$	k_DMSO_OH_B2	Karl et al. (2007)
G9403	TrGS	$CH_3SO_2 \rightarrow SO_2 + CH_3O_2$	1.0E13*EXP((-16500.*fch3so2/ 1.98635)/temp)	Karl et al. (2007)
G9404	TrGS	$CH_3SO_2 + O_3 \rightarrow CH_3SO_3 + O_2$	k_CH3SO2_O3	Karl et al. (2007)
G9405	TrGS	$CH_3SO_3 + HO_2 \rightarrow CH_3SO_3H + O_2$	k_CH3SO3_HO2	Karl et al. (2007)
G9408	StTrGS	$CH_2OO + SO_2 \rightarrow H_2SO_4 + HCHO$	k_CH2OO_SO2	Welz et al. (2012), Stone et al. (2014)*
G9409	TrGTerCS	$NOPINOO + SO_2 \rightarrow NOPINONE + H_2SO_4$	7.E-14	Rickard and Pascoe (2009)
G9410	TrGTerCS	$APINAOO + SO_2 \rightarrow PINAL + H_2SO_4$	7.00E-14	Rickard and Pascoe (2009)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G9411	TrGTerCS	APINBOO + SO_2 → PINAL + H_2SO_4	7.00E-14	Rickard and Pascoe (2009)
G9412	TrGTerCS	MBOOO + SO_2 → IBUTALOH + H_2SO_4	7.00E-14	Rickard and Pascoe (2009)
G9600	TrGCClS	DMS + Cl → CH_3SO_2 + HCl + HCHO	3.3E-10	Atkinson et al. (2004)
G9800	TrGCIS	DMS + IO → DMSO + I	3.2E-13*EXP(-925./temp)	Dillon et al. (2006a)
G3126MK	StTrGN	NO + NO_2 → N_2O_3	7.9E-12	Atkinson et al. (2004)
G3127MK	StTrGN	NO_2 + NO_2 → N_2O_4	1.0E-12	Atkinson et al. (2004)
G3128MK	StTrGN	N_2O_3 → NO + NO_2	3.6E8	Atkinson et al. (2004)
G3129MK	StTrGN	N_2O_4 → NO_2 + NO_2	4.4E6	Atkinson et al. (2004)
G4170aMK	TrGAmiN	MMA + OH → .3 CH2NH + .3 HO_2 + .1 MMAO2 + .6 CH3NH	(1-ya_soan1-ya_soan2-ya_soan5) *k_MMA_OH*xgcamin	Karl (2012)
G4170bMK	TrGAmiN	MMA + OH → BSOV + LCARBON + LNITROGEN	ya_soan1*k_MMA_OH	Karl (2012)
G4170cMK	TrGAmiN	MMA + OH → BLOV + LCARBON + LNITROGEN	ya_soan2*k_MMA_OH	Karl (2012)
G4170dMK	TrGAmiN	MMA + OH → BELV + LCARBON + LNITROGEN	ya_soan5*k_MMA_OH	Karl (2012)
G4171MK	TrGAmiN	CH3NH + NO_2 → .5 MMNNO2 + .5 CH2NH + .5 HONO	10*1.1E-12*xgcnno	Karl (2012)
G4172MK	TrGAmiN	CH3NH + O_2 → CH2NH + HO_2	2.4E-17	Nielsen et al. (2011)
G4173MK	TrGAmiN	MMAO2 + NO → H2NCHO + HO_2 + NO_2	0.5*7.7E-12	Nielsen et al. (2011)
G4174MK	TrGAmiN	H2NCHO + OH → HNCO + H_2O	k_H2NCHO_OH	Karl (2012)
G4175MK	TrGAmiN	MMA + CH2NH → CH2NCH3 + NH_3	4.0E-17*wall	Karl (2012)
G4200aMK	TrGAmiCN	MEA + OH → 0.05 H2NCH2CHO + 0.8 MEABO2 + 0.15 MEAN + 0.05 HO_2	(1-ya_soan1-ya_soan2-ya_soan5) *k_MEA_OH*fkoh_mea*xgcamin	Karl et al. (2012a)
G4200bMK	TrGAmiCN	MEA + OH → BSOV + 2 LCARBON + LNITROGEN	ya_soan1*k_MEA_OH*fkoh_mea	Karl et al. (2012a)
G4200cMK	TrGAmiCN	MEA + OH → BLOV + 2 LCARBON + LNITROGEN	ya_soan2*k_MEA_OH*fkoh_mea	Karl et al. (2012a)
G4200dMK	TrGAmiCN	MEA + OH → BELV + 2 LCARBON + LNITROGEN	ya_soan5*k_MEA_OH*fkoh_mea	Karl et al. (2012a)
G4201MK	TrGAmiCN	H2NCH2CHO + OH → 0.8 H2NCH2CO3 + 0.2 H2NCHO2CHO + H_2O	k_AAC_OH	Karl et al. (2012a)
G4202MK	TrGAmiCN	H2NCH2CO3 + NO → MMAO2 + CO_2 + NO_2	KAPNO	Karl et al. (2012a)
G4203MK	TrGAmiCN	H2NCHO2CHO + NO → H2NCOCHO + HO_2 + NO_2	KRO2NO*2	Karl et al. (2012a)
G4204MK	TrGAmiCN	H2NCOCHO + OH → H2NCOCO3 + H_2O	k_OXA_OH	Karl et al. (2012a)
G4205MK	TrGAmiCN	H2NCOCO3 + NO → H2NCHO + CO_2 + NO_2	KAPNO	Karl et al. (2012a)
G4206MK	TrGAmiCN	MEABO2 + NO → MEABO + NO_2	KRO2NO	Karl et al. (2012a)
G4207MK	TrGAmiCN	MEABO + O_2 → H2NCOCH2OH + HO_2	k_RO_O2*0.25	Karl et al. (2012a)
G4208MK	TrGAmiCN	MEABO → H2NCHO + HCHO	k_DEC*0.2	Karl et al. (2012a)
G4209MK	TrGAmiCN	H2NCOCH2OH + OH → H2NCOCHO + HO_2	k_HAC_OH	Karl et al. (2012a)
G4210MK	TrGAmiCN	MEAN + NO_2 → 0.5 MEANNO2 + 0.5 HNCHCH2OH + 0.5 HONO	1.4E-13 *xgcnno	Karl et al. (2012a)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G4211MK	TrGAmiCN	MEAN + O_2 → HNCHCH2OH + HO_2	1.2E-19	Karl et al. (2012a)
G4212MK	TrGAmiCN	MEAN + NO → MEANNO	8.5E-14 *xgcnno	Karl et al. (2012a)
G4213MK	TrGAmiCN	MEANNO2 + OH → MEANHA + HO_2	k_NITRAMA_OH	Karl et al. (2012a)
G4214MK	TrGAmiCN	HNCHCH2OH + OH → H2NCOCH2OH + HO_2	3.00E-13	Karl et al. (2012a)
G4220aMK	TrGAmiCN	DMA + OH → .4 CH3NCH3 + .5 DMAO2 + .1 CH2NCH3 + .1 HO_2	(1-ya_soan1-ya_soan2-ya_soan5) *k_DMA_OH*xgcamin	Karl (2012)
G4220bMK	TrGAmiCN	DMA + OH → BSOV + 2 LCARBON + LNITROGEN	ya_soan1*k_DMA_OH	Karl (2012)
G4220cMK	TrGAmiCN	DMA + OH → BLOV + 2 LCARBON + LNITROGEN	ya_soan2*k_DMA_OH	Karl (2012)
G4220dMK	TrGAmiCN	DMA + OH → BELV + 2 LCARBON + LNITROGEN	ya_soan5*k_DMA_OH	Karl (2012)
G4221MK	TrGAmiCN	CH3NCH3 + NO → NDMA	2.39E-13 *xgcnno	Nielsen et al. (2011)
G4222MK	TrGAmiCN	CH3NCH3 + O_2 → CH2NCH3 + HO_2	9.54E-20	Nielsen et al. (2011)
G4223MK	TrGAmiCN	CH3NCH3 + NO_2 → .6 CH2NCH3 + .7 HONO + .3 DMNNO2 + .1 CH3NO + .1 CH_3O_2	3.18E-13 *xgcnno	Lazarou et al. (1994)
G4224MK	TrGAmiCN	DMNNO2 + OH → CH2NCH3 + NO_2 + H_2O	4.5E-12	Karl (2012)
G4225MK	TrGAmiCN	DMAO2 + NO → CH3NHCHO + 0.5 NO_2 + 0.5 HO_2 + 0.5 HONO	8.5E-12	Karl (2012)
G4226MK	TrGAmiCN	NDMA + OH → CH2NCH3 + NO + H_2O	3.0E-12	Karl (2012)
G4230aMK	TrGAmiCN	CH2NCH3 + OH → CH3NHCHO + HO_2	(1-ya_soan1-ya_soan2-ya_soan5) *k_MMI_OH	Karl (2012)
G4230bMK	TrGAmiCN	CH2NCH3 + OH → BSOV + 2 LCARBON + LNITROGEN	ya_soan1*k_MMI_OH	Karl (2012)
G4230cMK	TrGAmiCN	CH2NCH3 + OH → BLOV + 2 LCARBON + LNITROGEN	ya_soan2*k_MMI_OH	Karl (2012)
G4230dMK	TrGAmiCN	CH2NCH3 + OH → BELV + 2 LCARBON + LNITROGEN	ya_soan5*k_MMI_OH	Karl (2012)
G4300aMK	TrGAmiCN	TMA + OH → TMAO2 + H_2O	(1-ya_soan1-ya_soan2-ya_soan5) *k_TMA_OH*xgcamin	Karl (2012)
G4300bMK	TrGAmiCN	TMA + OH → BSOV + 3 LCARBON + LNITROGEN	ya_soan1*k_TMA_OH	Karl (2012)
G4300cMK	TrGAmiCN	TMA + OH → BLOV + 3 LCARBON + LNITROGEN	ya_soan2*k_TMA_OH	Karl (2012)
G4300dMK	TrGAmiCN	TMA + OH → BELV + 3 LCARBON + LNITROGEN	ya_soan5*k_TMA_OH	Karl (2012)
G4301MK	TrGAmiCN	TMAO2 + NO → .3 DMNCHO + .3 HONO + .7 TMAO + .7 NO_2 + .7 HO_2	8.5E-12	Karl (2012)
G4302MK	TrGAmiCN	TMAO + O_2 → DMNCHO + HO_2	2.4E-15*5	Karl (2012)
G4303MK	TrGAmiCN	TMAO → CH3NCH3 + HCHO	4.0E+05	Karl (2012)
G4304MK	TrGAmiCN	DMNCHO + OH → DMNCHOO2	1.4E-11*0.25	Karl (2012)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G4305MK	TrGAmiCN	DMNCHOO2 + NO → TMADF + NO_2 + HO_2	8.5E-12	Karl (2012)
G4306MK	TrGAmiCN	HOETNHCHO + OH → HOCH2CONHCHO + HO_2	k_HAC_OH	Karl (2012)
G4310MK	TrGAmiCN	DMCNH2 + O_2 → 0.8 DMCOONH2 + 0.1 CH2CNH2CH3 + 0.1 DMCNH + HO_2	k_RO_O2*0.25	Nielsen and Schade (2012)
G4311MK	TrGAmiCN	DMCOONH2 + NO → H2NCOCH3 + HCHO + NO_2	KAPNO	Nielsen and Schade (2012)
G4312MK	TrGAmiCN	CH3CNH2MOH + NO → 0.5 H2NCOCH3 + 0.5 H2NCOCH2OH + CO_2 + NO_2	KAPNO	Nielsen and Schade (2012)
G4313MK	TrGAmiCN	CH3CNH2MOH + O_2 → 0.8 HNCCH3MOH + 0.1 H2NCCHOHCH3 + 0.1 H2NCCH2MOH + HO_2	k_RO_O2*0.25	Nielsen and Schade (2012)
G4320MK	TrGC	CH3CHOCH3 + O_2 → CH_3COCH_3 + HO_2	k_RO_O2	Rickard and Pascoe (2009)
G4400MK	TrGAmiCN	DEA + OH → .4 HOETNETOH + .5 DEAO2 + .1 HOETNHCH2CHO + .1 HO_2	k_DEA_OH*xgcamin	Karl (2012)
G4401MK	TrGAmiCN	HOETNETOH + NO → NDELA	2.39E-13 *xgcnno	Nielsen et al. (2011)
G4402MK	TrGAmiCN	HOETNETOH + O_2 → HOCH2CHNETOH + HO_2	9.54E-20	Nielsen et al. (2011)
G4403MK	TrGAmiCN	HOETNETOH + NO_2 → .6 HOCH2CHNETOH + .7 HONO + .3 DEANNO2 + .1 HOCH2CH2NO + .1 $HOCH_2CH_2O_2$	3.18E-13 *xgcnno	Lazarou et al. (1994)
G4404MK	TrGAmiCN	DEANNO2 + OH → HOCH2CHNETOH + NO_2 + H_2O	1.74E-11	Karl (2012)
G4405MK	TrGAmiCN	DEAO2 + NO → HOETNHCHO + HCHO + 0.5 NO_2 + 0.5 HO_2 + 0.5 HONO	8.5E-12	Karl (2012)
G4406MK	TrGAmiCN	NDELA + OH → HOCH2CHNETOH + NO + H_2O	1.61E-11	Karl (2012)
G4407MK	TrGAmiCN	HOCH2CHNETOH + OH → HOCH2CONETOH + HO_2	k_MMI_OH	Karl (2012)
G4410aMK	TrGAmiCN	AMP + OH → 0.30 AMPN + 0.65 DMCNH2CHO + 0.5 HO_2 + 0.05 NH_3 + 0.05 MACR	(1-ya_soan1-ya_soan2-ya_soan5) *k_AMP_OH*fkoh_mea	Harris and J. N. Pitts (1983)
G4410bMK	TrGAmiCN	AMP + OH → BSOV + 4 LCARBON + LNITROGEN	ya_soan1*k_AMP_OH*fkoh_mea	Harris and J. N. Pitts (1983)
G4410cMK	TrGAmiCN	AMP + OH → BLOV + 4 LCARBON + LNITROGEN	ya_soan2*k_AMP_OH*fkoh_mea	Harris and J. N. Pitts (1983)
G4410dMK	TrGAmiCN	AMP + OH → BELV + 4 LCARBON + LNITROGEN	ya_soan5*k_AMP_OH*fkoh_mea	Harris and J. N. Pitts (1983)
G4411MK	TrGAmiCN	AMPN + NO → NAMP	0.26*3.18E-13	Nielsen and Schade (2012)
G4412MK	TrGAmiCN	AMPN + NO_2 → AMPNNO2	3.18E-13	Lazarou et al. (1994)
G4413MK	TrGAmiCN	AMPN + O_3 → AMPOX	1.7E-13	Nielsen and Schade (2012)
G4414MK	TrGAmiCN	DMCNH2CHO + OH → DMCNH2CO3 + H_2O	k_AAC_OH	Nielsen and Schade (2012)
G4415MK	TrGAmiCN	DMCNH2CO3 + NO → DMCNH2 + CO_2 + NO_2	KAPNO	Nielsen and Schade (2012)
G4416MK	TrGAmiCN	DMCNH2CO3 + NO_2 → AMPAN	k_CH3CO3_NO2	Sander et al. (2006)
G4417MK	TrGAmiCN	AMPAN → DMCNH2CO3 + NO_2	k_PAN_M	Sander et al. (2006)
G4418MK	TrGAmiCN	AMPO + O_2 → DMOCNH2MOH + HO_2	k_RO_O2*0.25	Nielsen and Schade (2012)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G4419MK	TrGAmiCN	AMPO → CH3CNH2MOH + HCHO	k_DEC*0.2	Nielsen and Schade (2012)
G4420MK	TrGAmiCN	AMPNNO2 + OH → AMPNA + HO_2	k_NITRAMA_OH	Nielsen and Schade (2012)
G4421MK	TrGAmiCN	AMP + NO_3 → 0.5 AMPN + 0.4 DMCNH2CHO + 0.1 AMPO + 0.5 HO_2 + HNO_3	k_AMP_NO3	Carter (2008)
G4500MK	TrGAmiCN	DEANCH2O2 + NO → DEANCHO + NO_2 + HO_2	8.5E-12	Karl (2012)
G4600MK	TrGAmiCN	TEA + OH → TEAO2 + H_2O	k_TEA_OH*xgcamin	Karl (2012)
G4601MK	TrGAmiCN	TEAO2 + NO → .3 DEANCH2CHO + .3 HONO + .7 TEAO + .7 NO_2 + .7 HO_2	8.5E-12	Karl (2012)
G4602MK	TrGAmiCN	TEAO + O_2 → DEANCOCH2OH + HO_2	2.4E-15*5	Karl (2012)
G4603MK	TrGAmiCN	TEAO → HOETNETOH + CH_3CHO	4.0E+05	Karl (2012)
G4604MK	TrGAmiCN	TEAO + O_2 → DEANCH2O2 + HCHO	2.4E-15	Karl (2012)
G4605MK	TrGAmiCN	DEANCH2CHO + OH → DEANCH2COO2	1.4E-11	Karl (2012)
G4606MK	TrGAmiCN	DEANCH2COO2 + NO → DEANCHO + HCHO + NO_2 + HO_2	8.5E-12	Karl (2012)
G4610MK	TrGC	TME + O_3 → CH_3COCH_3 + $CH_3COCH_2O_2$ + OH	3.03E-15*EXP(-294./temp)	Rickard and Pascoe (2009)
G4611MK	TrGC	TME + OH → TMEO2	1.1E-10	Rickard and Pascoe (2009)
G4612MK	TrGCN	TMEO2 + NO → 2.0 CH_3COCH_3 + NO_2 + HO_2	KRO2NO	Rickard and Pascoe (2009)
G4620MK	TrGC	CHEX + OH → CHEXO2	2.88E-17*temp*temp*exp(309./temp)	Rickard and Pascoe (2009)
G4621MK	TrGC	CHEXO2 → 0.6 CHEXO + 0.2 CHEXOL + 0.2 CHEXONE	9.20E-14*RO2	Rickard and Pascoe (2009)
G4622MK	TrGC	CHEXO2 + HO_2 → CHEXOOH	0.770*KRO2HO2(6)	Rickard and Pascoe (2009)
G4623MK	TrGCN	CHEXO2 + NO → CHEXO + NO_2	KRO2NO	Rickard and Pascoe (2009)
G4624MK	TrGC	CHEXO + O_2 → CHEXONE + HO_2	k_RO_O2	Rickard and Pascoe (2009)
G40202c	TrGTerC	LAPINABO2 + HO_2 → BSOV + 10 LCARBON	ya_soan1*KRO2HO2(10)*rchohch2o2_oh	Rickard and Pascoe (2009), Sander et al. (2018)
G40202d	TrGTerC	LAPINABO2 + HO_2 → BLOV + 10 LCARBON	ya_soan2*KRO2HO2(10)*rchohch2o2_oh	Rickard and Pascoe (2009), Sander et al. (2018)
G40243c	TrGTerC	C109O2 + HO_2 → BSOV + 10 LCARBON	ya_soan1*KRO2HO2(10)*rcoch2o2_oh	Rickard and Pascoe (2009), Sander et al. (2018)
G40243d	TrGTerC	C109O2 + HO_2 → BLOV + 10 LCARBON	ya_soan2*KRO2HO2(10)*rcoch2o2_oh	Rickard and Pascoe (2009), Sander et al. (2018)
G40253c	TrGTerC	BPINAO2 + HO_2 → BSOV + 10 LCARBON	ya_soan1*KRO2HO2(10)*rcoch2o2_oh	Rickard and Pascoe (2009), Sander et al. (2018)
G40253d	TrGTerC	BPINAO2 + HO_2 → BLOV + 10 LCARBON	ya_soan2*KRO2HO2(10)*rcoch2o2_oh	Rickard and Pascoe (2009), Sander et al. (2018)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G40259a	TrGTerC	ROO6R1O2 + HO_2 → 10 LCARBON	(1-ya_soan1-ya_soan2) *KRO2HO2(10)	Vereecken and Peeters (2012)*
G40259b	TrGTerC	ROO6R1O2 + HO_2 → BSOV + 10 LCARBON	ya_soan1*KRO2HO2(10)	Rickard and Pascoe (2009), Sander et al. (2018)
G40259ac	TrGTerC	ROO6R1O2 + HO_2 → BLOV + 10 LCARBON	ya_soan2*KRO2HO2(10)	Rickard and Pascoe (2009), Sander et al. (2018)
G40600	TrGTer	BSOV + OH → BLOV	4.0E-11	Tsimpidi et al. (2010)
G40601	TrGTer	BLOV + OH → BELV	4.0E-11	Tsimpidi et al. (2010)
G40602	TrGTer	BSOV → BELV	9.6E-06	Carlton et al. (2010)
G40603	TrGTer	BLOV → BELV	9.6E-06	Carlton et al. (2010)
G40700	TrGTer	ASOV + OH → ALOV	4.0E-11	Tsimpidi et al. (2010)
G40701	TrGTer	ALOV + OH → AELV	4.0E-11	Tsimpidi et al. (2010)
G40702	TrGTer	ASOV → AELV	9.6E-06	Carlton et al. (2010)
G40703	TrGTer	ALOV → AELV	9.6E-06	Carlton et al. (2010)
G40704	TrGTer	PIOV + OH → PSOV	2.0E-11	Lambe et al. (2009)
G40705	TrGTer	PSOV + OH → PELV	2.0E-11	Lambe et al. (2009)
G40706	TrGTer	PELV → PSOV	5.0E-04	Lim and Ziemann (2009)
G9201MK	StTrGS	$HSO_3 + O_2 \rightarrow HO_2 + SO_3$	1.3E-12*exp(-330/temp)	Burkholder et al. (2015)
G9202MK	StTrGS	$SO_3 + H_2O \rightarrow H_2SO_4$	2.4E-15	Reiner and Arnold (1993)
G9420aMK	TrGCS	DMSOH + O_2 → DMSO + HO_2	5.0E-13	Karl et al. (2007)
G9420bMK	TrGCS	DMSOH + O_2 → DMSOHOO	3.0E-14	Karl et al. (2007)
G9421MK	TrGCS	DMSOH → $CH_3SOH + CH_3O_2$	5.0E+5	Karl et al. (2007)
G9422MK	TrGCNS	DMSOHOO + NO → DMSOHO + NO_2	5.0E-12	Karl et al. (2007)
G9423MK	TrGCS	DMSOHO + O_2 → $DMSO_2$ + HO_2	1.0E-13	Karl et al. (2007)
G9424MK	TrGCS	DMSOHO → $CH_3SOOH + CH_3O_2$	6.0E+5	Karl et al. (2007)
G9425MK	TrGCS	$CH_3SOCH_2 + O_2 \rightarrow CH_3SOCH_2O_2$	0.5E-13	Karl et al. (2007)
G9426MK	TrGCNS	$CH_3SOCH_2O_2$ + NO → CH_3SO + HCHO + NO_2	1.0E-11	Karl et al. (2007)
G9427MK	TrGCS	$CH_3SCH_2 + O_2 \rightarrow CH_3SCH_2OO$	5.7E-12	Karl et al. (2007)
G9428MK	TrGCNS	CH_3SCH_2OO + NO → HCHO + CH_3S + NO_2	1.9E-11	Karl et al. (2007)
G9429MK	TrGCS	$CH_3SCH_2OO + HO_2 \rightarrow CH_3SCH_2OOH + O_2$	1.5E-12	Karl et al. (2007)
G9430MK	TrGCS	$CH_3SCH_2OO + CH_3O_2 \rightarrow CH_3S$ + 2 HCHO + HO_2	1.8E-13	Karl et al. (2007)
G9431MK	TrGS	$CH_3S + O_3 \rightarrow CH_3SO + O_2$	1.36E-12*EXP(374./temp)	Karl et al. (2007)
G9432MK	TrGNS	$CH_3S + NO_2 \rightarrow CH_3SO$ + NO	2.8E-11*EXP(240./temp)	Karl et al. (2007)
G9433aMK	TrGS	$CH_3S + O_2 \rightarrow CH_3SOO$	1.5E-14	Karl et al. (2007)
G9433bMK	TrGS	$CH_3S + O_2 \rightarrow CH_3O_2 + SO_2$	3.0E-18	Karl et al. (2007)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G9434MK	TrGS	$CH_3S + CH_3O_2 \rightarrow CH_3SO + HCHO + HO_2$	6.1E-11	Karl et al. (2007)
G9435MK	TrGS	$CH_3SO + O_3 \rightarrow 0.5\ CH_3SO_2 + 0.5\ O_2 + 0.5\ CH_3O_2 + 0.5\ SO_2$	3.2E-13	Karl et al. (2007)
G9436aMK	TrGNS	$CH_3SO + NO_2 \rightarrow CH_3SO_2 + NO$	1.2E-11	Karl et al. (2007)
G9436bMK	TrGNS	$CH_3SO + NO_2 \rightarrow CH_3O_2 + SO_2 + NO$	3.5E-12	Karl et al. (2007)
G9437MK	TrGS	$CH_3SO + O_2 \rightarrow CH_3SOO_2$	7.7E-18	Karl et al. (2007)
G9438MK	TrGS	$CH_3SO + CH_3O_2 \rightarrow CH_3SO_2 + HCHO + HO_2$	3.0E-12	Karl et al. (2007)
G9439MK	TrGS	$CH_3SOH + OH \rightarrow CH_3SO + H_2O$	1.10E-10	Karl et al. (2007)
G9440MK	TrGNS	$CH_3SOH + NO_3 \rightarrow CH_3SO + HNO_3$	3.4E-12	Karl et al. (2007)
G9441MK	TrGS	$CH_3SOH + HO_2 \rightarrow CH_3SO + H_2O_2$	8.5E-13	Karl et al. (2007)
G9442MK	TrGS	$CH_3SOH + CH_3O_2 \rightarrow CH_3SO + CH_3OOH$	8.5E-13	Karl et al. (2007)
G9443MK	TrGS	$CH_3SOO + O_3 \rightarrow CH_3SO + 2O_2$	8.0E-13	Karl et al. (2007)
G9444MK	TrGNS	$CH_3SOO + NO \rightarrow CH_3SO + NO_2$	1.1E-11	Karl et al. (2007)
G9445MK	TrGS	$CH_3SOO + O_2 \rightarrow CH_3O_2 + SO_2$	2.0E-17	Karl et al. (2007)
G9446aMK	TrGS	$CH_3SOO \rightarrow CH_3S + O_2$	1.0E+5	Karl et al. (2007)
G9446bMK	TrGS	$CH_3SOO \rightarrow CH_3SO_2$	8.0	Karl et al. (2007)
G9447MK	TrGS	$CH_3SOO + CH_3O_2 \rightarrow CH_3SO + HCHO + HO_2$	5.5E-12	Karl et al. (2007)
G9448MK	TrGS	$CH_3SOOH + OH \rightarrow CH_3SO_2 + H_2O$	k_MSIA_OH	Karl et al. (2007)
G9449MK	TrGNS	$CH_3SOOH + NO_3 \rightarrow CH_3SO_2 + HNO_3$	1.0E-13	Karl et al. (2007)
G9450MK	TrGS	$CH_3SOOH + HO_2 \rightarrow CH_3SOO + H_2O_2$	1.0E-15	Karl et al. (2007)
G9451MK	TrGS	$CH_3SOOH + CH_3O_2 \rightarrow CH_3SO_2 + CH_3OOH$	1.0E-15	Karl et al. (2007)
G9452MK	TrGS	$CH_3SO_2 + HO_2 \rightarrow CH_3SO_3 + OH$	2.5E-13	Karl et al. (2007)
G9453MK	TrGNS	$CH_3SO_2 + NO_2 \rightarrow CH_3SO_3 + NO$	2.2E-11	Karl et al. (2007)
G9454MK	TrGS	$CH_3SO_2 + CH_3O_2 \rightarrow CH_3SO_3 + HCHO + HO_2$	2.5E-13	Karl et al. (2007)
G9455MK	TrGNS	$CH_3SO_3 + NO_2 \rightarrow CH_3SOO_2NO_2$	3.0E-15	Karl et al. (2007)
G9456MK	TrGS	$CH_3SO_3 \rightarrow SO_3 + CH_3O_2$	k_CH3SO3_dec	Karl et al. (2007)
G9457MK	TrGNS	$CH_3SOO_2 + NO \rightarrow CH_3SO_2 + NO_2$	1.0E-11	Karl et al. (2007)
G9458MK	TrGS	$CH_3SOO_2 + HO_2 \rightarrow CH_3SOO_2H + O_2$	3.0E-12	Karl et al. (2007)
G9459MK	TrGS	$CH_3SOO_2 \rightarrow CH_3SO + O_2$	170.	Karl et al. (2007)
G9460MK	TrGS	$CH_3SOO_2 + CH_3O_2 \rightarrow CH_3SO_2 + HCHO + HO_2$	5.5E-12	Karl et al. (2007)
G9461MK	TrGNS	$CH_3SO_2O_2 + NO \rightarrow CH_3SO_3 + NO_2$	1.0E-11	Karl et al. (2007)
G9462MK	TrGS	$CH_3SO_2O_2 + HO_2 \rightarrow CH_3SO_4H$	2.0E-12	Karl et al. (2007)
G9463MK	TrGS	$CH_3SO_2O_2 \rightarrow CH_3SO_2 + O_2$	170.	Karl et al. (2007)
G9464MK	TrGS	$CH_3SO_2O_2 + CH_3O_2 \rightarrow CH_3SO_3 + HCHO + HO_2$	5.5E-12	Karl et al. (2007)
G9465MK	TrGNS	$CH_3SO_2O_2 + NO_2 \rightarrow CH_3SO_2O_2NO_2$	5.89E-12	Karl et al. (2007)

Table 1: Gas phase reactions (... continued)

#	labels	reaction	rate coefficient	reference
G9466MK	TrGNS	$CH_3SOO_2NO_2 \rightarrow CH_3SO_3 + NO_2$	1.15E-2	Karl et al. (2007)
G9467MK	TrGNS	$CH_3SO_2O_2NO_2 \rightarrow CH_3SO_2O_2 + NO_2$	1.15E-2	Karl et al. (2007)
G9468MK	TrGCS	$DMSO_2 + OH \rightarrow DMSO_2OO$	k_DMSO2_OH	Karl et al. (2007)
G9469MK	TrGCNS	$DMSO_2OO + NO \rightarrow DMSO_2O + NO_2$	5.0E-12	Karl et al. (2007)
G9470MK	TrGCS	$DMSO_2OO + HO_2 \rightarrow DMSO_2OOH + O_2$	1.5E-12	Karl et al. (2007)
G9471MK	TrGCS	$DMSO_2O \rightarrow CH_3SO_2 + HCHO$	1.0E+1	Karl et al. (2007)
G9472MK	TrGS	$CH_3SO_3H + H_2O \rightarrow MSA(H_2O)$	6.58E-10	Dawson et al. (2012)
G9473MK	TrGS	$MSA(H_2O) \rightarrow CH_3SO_3H + H_2O$	5.53E08*EXP(-669.71/temp)	Dawson et al. (2012)
G9900MK	TrGAmiCNS	$DMA + MSA(H_2O) \rightarrow MSA(DMA)(H_2O)$	1.32E-09	Dawson et al. (2012)
G9901MK	TrGAmiCNS	$MSA(DMA)(H_2O) \rightarrow DMA + MSA(H_2O)$	30.9 *EXP(-4118.74/temp)	Dawson et al. (2012)
G9902MK	TrGAmiCNS	$MSA(DMA)(H_2O) \rightarrow MSA(DMA) + H_2O$	9.96E08*EXP(-736.69/temp)	Dawson et al. (2012)
G9903MK	TrGAmiCNS	$MSA(DMA) + H_2O \rightarrow MSA(DMA)(H_2O)$	1.66E-09	Dawson et al. (2012)
G9910MK	TrGAmiCNS	$TMA + MSA(H_2O) \rightarrow MSA(TMA)(H_2O)$	1.44E-09	Dawson et al. (2012)
G9911MK	TrGAmiCNS	$MSA(TMA)(H_2O) \rightarrow TMA + MSA(H_2O)$	5.29 *EXP(-4487.08/temp)	Dawson et al. (2012)
G9912MK	TrGAmiCNS	$MSA(TMA)(H_2O) \rightarrow MSA(TMA) + H_2O$	7.91E06*EXP(-1674.28/temp)	Dawson et al. (2012)
G9913MK	TrGAmiCNS	$MSA(TMA) + H_2O \rightarrow MSA(TMA)(H_2O)$	1.49E-09	Dawson et al. (2012)

General notes

Three-body reactions

Rate coefficients for three-body reactions are defined via the function `k_3rd`(T, M, k_0^{300}, n, k_{inf}^{300}, m, f_{c}). In the code, the temperature T is called `temp` and the concentration of "air molecules" M is called `cair`. Using the auxiliary variables $k_0(T)$, $k_{\mathrm{inf}}(T)$, and k_{ratio}, `k_3rd` is defined as:

$$k_0(T) = k_0^{300} \times \left(\frac{300\mathrm{K}}{T}\right)^n \quad (1)$$

$$k_{\mathrm{inf}}(T) = k_{\mathrm{inf}}^{300} \times \left(\frac{300\mathrm{K}}{T}\right)^m \quad (2)$$

$$k_{\mathrm{ratio}} = \frac{k_0(T)M}{k_{\mathrm{inf}}(T)} \quad (3)$$

$$\texttt{k\_3rd} = \frac{k_0(T)M}{1+k_{\mathrm{ratio}}} \times f_{\mathrm{c}}^{\left(\frac{1}{1+(\log_{10}(k_{\mathrm{ratio}}))^2}\right)} \quad (4)$$

A similar function, called `k_3rd_iupac` here, is used by Wallington et al. (2018) for three-body reactions. It has the same function parameters as `k_3rd` and it is defined as:

$$k_0(T) = k_0^{300} \times \left(\frac{300\mathrm{K}}{T}\right)^n \quad (5)$$

$$k_{\mathrm{inf}}(T) = k_{\mathrm{inf}}^{300} \times \left(\frac{300\mathrm{K}}{T}\right)^m \quad (6)$$

$$k_{\mathrm{ratio}} = \frac{k_0(T)M}{k_{\mathrm{inf}}(T)} \quad (7)$$

$$N = 0.75 - 1.27 \times \log_{10}(f_{\mathrm{c}}) \quad (8)$$

$$\texttt{k\_3rd\_iupac} = \frac{k_0(T)M}{1+k_{\mathrm{ratio}}} \times f_{\mathrm{c}}^{\left(\frac{1}{1+(\log_{10}(k_{\mathrm{ratio}})/N)^2}\right)} \quad (9)$$

Structure-Activity Relationships (SAR)

Some unmeasured rate coefficients are estimated with structure-activity relationships, using the following parameters and substituent factors:

k for H-abstraction by OH in $\mathrm{cm^{-3}s^{-1}}$	
`k_p`	$4.49 \times 10^{-18} \times (T/\mathrm{K})^2 \exp(-320\,\mathrm{K}/T)$
`k_s`	$4.50 \times 10^{-18} \times (T/\mathrm{K})^2 \exp(253\,\mathrm{K}/T)$
`k_t`	$2.12 \times 10^{-18} \times (T/\mathrm{K})^2 \exp(696\,\mathrm{K}/T)$
`k_rohro`	$2.1 \times 10^{-18} \times (T/\mathrm{K})^2 \exp(-85\,\mathrm{K}/T)$
`k_co2h`	$0.7 \times k_{\mathrm{CH_3CO_2H+OH}}$
`k_roohro`	$0.6 \times k_{\mathrm{CH_3OOH+OH}}$
`f_alk`	1.23
`f_soh`	3.44
`f_toh`	2.68
`f_sooh`	8.
`f_tooh`	8.
`f_ono2`	0.04
`f_ch2ono2`	0.20
`f_cpan`	0.25
`f_allyl`	3.6
`f_cho`	0.55
`f_co2h`	1.67
`f_co`	0.73
`f_o`	8.15
`f_pch2oh`	1.29
`f_tch2oh`	0.53

k for OH-addition to double bonds in $\mathrm{cm^{-3}s^{-1}}$	
`k_adp`	$4.5 \times 10^{-12} \times (T/300\,\mathrm{K})^{-0.85}$
`k_ads`	$1/4 \times (1.1 \times 10^{-11} \times \exp(485\,\mathrm{K}/T)+$ $1.0 \times 10^{-11} \times \exp(553\,\mathrm{K}/T))$
`k_adt`	$1.922 \times 10^{-11} \times \exp(450\,\mathrm{K}/T) - k_{\mathrm{ads}}$
`k_adsecprim`	3.0×10^{-11}
`k_adtertprim`	5.7×10^{-11}
`a_pan`	0.56
`a_cho`	0.31
`a_coch3`	0.76
`a_ch2oh`	1.7
`a_ch2ooh`	1.7
`a_coh`	2.2
`a_cooh`	2.2
`a_co2h`	0.25
`a_ch2ono2`	0.64

$\mathrm{RO_2}$ self and cross reactions

The self and cross reactions of organic peroxy radicals are treated according to the permutation reaction formalism as implemented in the MCM (Rickard and Pascoe, 2009), as decribed by Jenkin et al. (1997). Every organic peroxy radical reacts in a pseudo-first-order reaction with a rate constant that is expressed as $k^{\mathrm{1st}} = 2 \times \sqrt{k_{\mathrm{self}} \times \texttt{k\_CH3O2}} \times [\mathrm{RO_2}]$ where k_{self} = second-order rate coefficient of the self reaction of the organic peroxy radical, `k_CH3O2` = second-order rate coefficient of the self reaction of $\mathrm{CH_3O_2}$, and $[\mathrm{RO_2}]$ = sum of the concentrations of all organic peroxy radicals.

Specific notes

G2110: The rate coefficient is: k_HO2_HO2 = (3.0E-13*EXP(460./temp)+2.1E-33*EXP(920./temp) *cair)*(1.+1.4E-21*EXP(2200./temp)*C(ind_H2O)).

G2117: Converted to Kc [molec-1 cm3]= Kp*R*T/NA , where R is 82.05736 [cm3 atm K-1 mol-1].

G2118: Assuming fast equilibrium.

G3109: The rate coefficient is: k_NO3_NO2 = k_3rd(temp,cair,2.4E-30,3.0,1.6E-12,-0.1,0.6).

G3110: The rate coefficient is defined as backward reaction divided by equilibrium constant.

G3203: The rate coefficient is: k_NO2_HO2 = k_3rd(temp,cair,1.9E-31,3.4,4.0E-12,0.3,0.6).

G3206: The rate coefficient is: k_HNO3_OH = 1.32E-14 * EXP(527/temp) + 1 / (1 / (7.39E-32 * EXP(453/temp)*cair) + 1 / (9.73E-17 * EXP(1910/temp)))

G3207: The rate coefficient is defined as backward reaction divided by equilibrium constant.

G4104b: Methyl nitrate yield according to Banic et al. (2003) but reduced by a factor of 10 according to the upper limit derived from measurements by Munger et al. (1999).

G4109: Same temperature dependence as for CH_3CHO+NO_3 assumed.

G4115: The rate coefficient is defined as backward reaction divided by equilibrium constant.

G4116: Same value as for PAN + OH.

G4126: Same as for G4104 but scaled to match the recommeded value at 298K.

G4127: Same as for CH3O2 + NO3 in G4105.

G4130a: SAR for H-abstraction by OH.

G4130b: SAR for H-abstraction by OH.

G4132: SAR for H-abstraction by OH.

G4133: Lower limit of the rate constant. Products uncertain but CH_3OH can be excluded because of a likely high energy barrier (L. Vereecken, pers. comm.). CH_2OO production cannot be excluded.

G4134: Estimate based on the decomposition lifetime of 3 s (Olzmann et al., 1997) and a 20 kcal/mol energy barrier (Vereecken and Francisco, 2012).

G4135: Rate constant for $CH_2OO + NO_2$ (G4138) multiplied by the factor from Ouyang et al. (2013).

G4136: Average of two measurements.

G4137: Upper limit.

G4138: Average of 7.E-12 and 1.5E-12.

G4141: $HOOCH_2OCHO$ forms and then decomposes to formic anhydride (Gruzdev et al., 1993) which hydrolyses in the humid atmosphere (Conn et al., 1942).

G4142: High-pressure limit.

G4143: Generic estimate for reaction with alcohols.

G4144: Generic estimate for reaction with RO_2.

G4148: Same value as for NO_2+CH_3O_2.

G4149: Barnes et al. (1985) estimated a decomposition rate equal to that of $CH_3O_2NO_2$.

G4150: Value for $CH_3O_2NO_2$ + OH, H-abstraction enhanced by the HO-group by f_soh.

G4154: Products assumed to be $CH_3O_2 + O_2$ (could also be HCHO + O_2 + OH).

G4160b: Half of the H-yield is attributed to fast secondary chemistry.

G4160c: The NH + CO channel is also significant but neglected here.

G4161: No studies below 450 K and only the major channel is considered.

G4164: Upper limit. Dominant pathway under atmospheric conditions.

G42001: The product distribution is from Rickard and Pascoe (2009), after substitution of the energized Criegee intermediate, CH_2OO, by its decomposition products and reaction of the stabilized CI with the water dimer.

G42010: Only major channel considered as the end products are essentially the same.

G42013: The rate coefficient is: k_CH3CO3_NO2 = k_3rd(temp,cair,9.7E-29,5.6,9.3E-12,1.5,0.6).

G42018: The rate coefficient is the same as for the CH_3 channel in G4107 (CH_3OOH+OH).

G42021: The rate coefficient is k_PAN_M = k_CH3CO3_NO2/9.0E-29*EXP(-14000./temp), i.e. the rate coefficient is defined as backward reaction divided by equilibrium constant.

G42022a: Quantum yields and products are from Glowacki et al. (2012).

G42022b: Quantum yields and products are from Glowacki et al. (2012).

G42024a: Rate constant is the high-pressure limit as recommended by Atkinson et al. (2006).

G42024b: Rate constant is the high-pressure limit as recommended by Atkinson et al. (2006).

G42047: Orlando et al. (1998) estimated that about 25% of the $HOCH_2CH_2O$ in this reaction is produced with sufficient excess energy that it decomposes promptly. The decomposition products are 2 HCHO + HO_2.

G42051a: Same as for the CH_3O_2 channel in G4107: CH_3OOH+OH.

G42058b: The aldehydic H is assumed to be like the analogous H of $HOCH_2CHO$.

G42074a: Factor of 3 to match the estimate of k = 1.E-11 molec/cm3/s by Paulot et al. (2009a).

G42074b: Factor of 3 to match the estimate of k = 1.E-11 molec/cm3/s by Paulot et al. (2009a).

G42075: $NO_3CH_2CO_2H$ and $NO_3CH_2CO_3H$ neglected.

G42078: $NO_3CH_2CO_2H$ neglected.

G42082: Same rate constant as for PAN + OH.

G42083a: Rate constant is the high-pressure limit as recommended by Atkinson et al. (2006).

G42083b: Rate constant is the high-pressure limit as recommended by Atkinson et al. (2006).

G42085a: Uncertainties on the kinetics at pressures $<$ 0.1 bar.

G42085b: Channel proposed by Hynes and Wine 1991, OH + HCHO + HOCN, could not be confirmed by Tyndall et al. (2001b). There is no alternative mechanism at the moment. Products assumed to be OH + CH3CO3 + NO

G42086b: Assuming HCN is from channel 2h, HCO + H + HCN. HCO is replaced by H + CO.

G42086c: Assuming exothermic channels 2b and 2d are equally important.

G42087: HCOCN is produced but replaced here by its likely oxidation products (HCN + CO_2) as studied by Tyndall et al. (2001b). The rate constant for a typical RO_2 + NO reaction is used.

G42088: $NCCH_2OOH$ is produced but replaced here by its likely oxidation products (HCN + CO_2) as studied by Tyndall et al. (2001b). The rate constant for a typical RO_2 + HO_2 reaction is used.

G42089a: The minor channel with k=5.2E-12 is combined with the major one producing HCOOH.

G42090: Theoretical keto-enol tautomerization catalyzed by formic acid (Grenfell et al., 2006).

G42091: Theoretical keto-enol tautomerization catalyzed by formic acid (Grenfell et al., 2006).

G43001a: Branching ratios according to Rickard et al. (1999).

G43001b: Branching ratios according to Rickard et al. (1999).

G43004: The value for the generic RO_2 + HO_2 reaction from Atkinson (1997) is used here.

G43008: The value for the generic RO_2 + HO_2 reaction from Atkinson (1997) is used here.

G43011: Strong positive deviation of k below 240 K compared to the expression recommended by JPL (Burkholder et al., 2015).

G43015a: The same value as for G4107 (CH_3OOH + OH) is used, multiplied by the branching ratio of the CH_3O_2 channel.

G43028: Alkyl nitrate formation neglected. (also not considered in MCM).

G43037: Alkyl nitrate formation neglected. (also not considered in MCM).

G43040a: Rate coefficient estimated with SAR (Taraborrelli, 2010).

G43040b: Rate coefficient estimated with SAR (Taraborrelli, 2010).

G43044: Alkyl nitrate formation neglected.

G43045c: Rate coefficient assumed to equal to the one of hydroxyacetone (ACETOL) for this channel.

G43048: Using the high-pressure limit.

G43049: The pressure fall-off between 1000 and 100 mbar is only 3% (Kirchner et al., 1999).

G43050: Value for $CH_3O_2NO_2$ + OH, H-abstraction enhanced by the CH_3CO-group by f_co.

G43051c: Products approximated with C_2H_5CHO + HO_2.

G43052: Only major H-abstraction channel considered.

G43059: Products approximated with the major end-product CH_3CHO.

G43060b: Products approximated with the major end-product CH_3CHO.

G43061: Products approximated with the likely end-product CH_3CHO.

G43065: As for $HCOCO_3$.

G43070a: Branching ratios estimated with SAR for H-abstraction rate constants by OH.

G43070b: Branching ratios estimated with SAR for H-abstraction rate constants by OH.

G43071a: Only this channel considered as the intermediate radical is likely more stable than $CHCH(OH)_2$.

G43072: Theoretical keto-enol tautomerization catalyzed by formic acid (Grenfell et al., 2006).

G43073: Theoretical keto-enol tautomerization catalyzed by formic acid (Grenfell et al., 2006).

G43074: HCOCOCHO would be produced but undergoes fast photolysis (faster than MGLYOX) and is substituted with its products.

G43223: Products simplified

G43419: KDEC C3DIALO $\rightarrow$ GLYOX + CO + HO2

G43420: KDEC C3DIALO $\rightarrow$ GLYOX + CO + HO2

G43421: Permutation reaction (minor channels removed).

G44000: The $LC_4H_9O_2$ composition ($nC_4H_9O_2$:$sC_4H_9O_2$ ratio) is assumed to be equal to the ratio of the production rates at 298K: k_p/(k_p+k_s) = 0.1273 and k_s/(k_p+k_s) = 0.8727.

G44001b: $sC_4H_9O_2$ products are substituted with 0.636 MEK + HO_2 and 0.364 CH_3CHO + $C_2H_5O_2$ at 1 bar and 298 K.

G44003c: The alkyl nitrate yield is the weighted average yield for the two isomers forming from $nC_4H_9O_2$ and $sC_4H_9O_2$.

G44010b: H-abstraction from primary C and substitution of the resulting peroxy radical with its products from the reaction with NO.

G44011: H-abstraction from primary C and substitution of the resulting peroxy radical with its products from the reaction with NO.

G44015b: Products assumed to be only from H-abstraction from a secondary C bearing the -OOH group.

G44016: Products assumed to be only from H-abstraction from a secondary C bearing the $-ONO_2$ group.

G44018: LHMVKABO2 is 0.12 HMVKAO2 + 0.88 HMVKBO2.

G44019: LMEKO2 represents 0.62 MEKBO2 + 0.38 MEKAO2.

G44021a: The products of MEKAO are substituted with HCHO + CO_2 + $HOCH_2CH_2O_2$.

G44023a: Products from H-abstraction from the tertiary carbon bearing the ONO_2 group.

G44023b: Products from H-abstraction from the secondary carbon bearing the ONO_2 group.

G44025: Same value as for PAN.

G44026: Products as in G4415. Only the main channels for each isomer are considered. Weighted average for the isomers.

G44035: Rate constant replaced with the one of beta hydroxy RO_2.

G44046b: Using value for secondary nitrate (88% of total).

G44061a: Using value for secondary nitrate (88% of total).

G44061b: Using value for secondary nitrate (88% of total).

G44062a: Simplified products.

G44062b: Simplified products.

G44066: Alkyl nitrate formation neglected.

G44070: Alkyl nitrate formation neglected.

G44076: Alkyl nitrate formation neglected.

G44078: Other channel neglected.

G44081: Alkyl nitrate formation neglected.

G44082: Other channel neglected.

G44085: k for CH_3CHCO from Hatakeyama et al. (1985) adjusted.

G44086: Simplified product distribution.

G44089: The nitrated RO_2 is replaced by its products upon reaction with NO.

G44096: Both LBUT1ENO2 isomers mostly C_2H_5CHO.

G44097a: Branching ratios according to Rickard et al. (1999). CH_3CHO_2CHO is replaced with its major products CH_3CHO + CO + HO_2.

G44097b: Branching ratios according to Rickard et al. (1999).

G44098: The nitrated RO_2 is replaced by its products upon reaction with NO.

G44103b: MEKCOH replaced by its major oxidation products.

G44104: Carbonyl nitrate replaced by its major oxidation products.

G44106: CH3CHOOA products as from C_3H_6 + O_3 reaction.

G44107: The nitrated RO_2 is replaced by its products upon reaction with NO.

G44110: The nitrated RO_2 is replaced by its products upon reaction with NO.

G44124b: Skipping intermediate steps mostly leading to acetone.

G44126: Skipping intermediate steps mostly leading to acetone.

G44127: Only this channel considered as the intermediate radical is likely more stable than $CHCH(OH)_2$.

G44128: Theoretical keto-enol tautomerization catalyzed by formic acid (Grenfell et al., 2006).

G44129: Theoretical keto-enol tautomerization catalyzed by formic acid (Grenfell et al., 2006).

G44130: Only this channel considered as the intermediate radical is likely more stable than $CHCH(OH)_2$.

G44131: Theoretical keto-enol tautomerization catalyzed by formic acid (Grenfell et al., 2006).

G44132: Theoretical keto-enol tautomerization catalyzed by formic acid (Grenfell et al., 2006).

G44133: Only this channel considered as the intermediate radical is likely more stable than $CHCH(OH)_2$.

G44134: Theoretical keto-enol tautomerization catalyzed by formic acid (Grenfell et al., 2006).

G44135: Theoretical keto-enol tautomerization catalyzed by formic acid (Grenfell et al., 2006).

G44136: Only this channel considered as the intermediate radical is likely more stable than $CHCH(OH)_2$.

G44137: Theoretical keto-enol tautomerization catalyzed by formic acid (Grenfell et al., 2006).

G44138: Theoretical keto-enol tautomerization catalyzed by formic acid (Grenfell et al., 2006).

G44139: Simplified oxidation.

G44140: Simplified oxidation.

G44141: Simplified oxidation.

G44142: Simplified oxidation.

G44202: Alkyl nitrate formation neglected.

G44203a: Rate coefficient estimated with SAR (Taraborrelli, 2010).

G44205: Alkyl nitrate formation neglected.

G44210: Alkyl nitrate formation neglected.

G44221: Same k as for MGLYOX + OH (Tyndall et al., 1995).

G44402: KDEC NC4DCO2→ MALANHY + NO2

G44406c: KDEC MALDIALCO2 → 0.6 MALANHY + HO2 + 0.4 GLYOX + 0.4 CO + 0.4 CO2

G44407: KDEC MALDIALCO2 → 0.6 MALANHY + HO2 + 0.4 GLYOX + 0.4 CO + 0.4 CO2

G44409: KDEC MALDIALCO2 → 0.6 MALANHY + HO2 + 0.4 GLYOX + 0.4 CO + 0.4 CO2

G44410: KDEC MALDIALCO2 → 0.6 MALANHY + HO2 + 0.4 GLYOX + 0.4 CO + 0.4 CO2

G44412: KDEC BZFUONOOA → 0.5 BZFUONOO + 0.5 CO + 0.5 CO2 + 0.5 HCOCH2O2 + 0.5 OH and BZFUONOO → 0.625 CO14O3CO2H + 0.375 CO14O3CHO + 0.375 H2O2

G44421: Only major channel.

G44424: KDEC: GLYOOA → 0.125 HCHO + 0.18 GLYOO + 0.82 HO2 + 0.57 OH + 1.265 CO + 0.25 CO2 and H2O substitution GLYOO → 0.625 HCOCO2H + 0.375 GLYOX + 0.375 H2O2

G44425: Merged equations.

G44430: KDEC MALANHYO → HCOCOHCO3

G44431: KDEC MALANHYO → HCOCOHCO3

G44432: Only major channel. KDEC MALANHYO → HCOCOHCO3

G44436: KDEC NBZFUO → 0.5 CO14O3CHO + 0.5 NO2 + 0.5 NBZFUONE + 0.5 HO2

G44437: KDEC NBZFUO → 0.5 CO14O3CHO + 0.5 NO2 + 0.5 NBZFUONE + 0.5 HO2

G44438: KDEC NBZFUO → 0.5 CO14O3CHO + 0.5 NO2 + 0.5 NBZFUONE + 0.5 HO2 and RO2 Only major channel.

G44439: KDEC MALDIALCO2 → 0.6 MALANHY + HO2 + 0.4 GLYOX + 0.4 CO + 0.4 CO2

G44443: KDEC MECOACETO → CH3CO3 + HCHO

G44444: KDEC MECOACETO → CH3CO3 + HCHO

G44445: KDEC MECOACETO → CH3CO3 + HCHO

G44450: KDEC BZFUO → CO14O3CHO + HO2

G44451: KDEC BZFUO → CO14O3CHO + HO2

G44452: KDEC BZFUO → CO14O3CHO + HO2. Only major channel.

G44457: KDEC MALDIALO → GLYOX + GLYOX + HO2

G44458: KDEC MALDIALO → GLYOX + GLYOX + HO2

G44459: KDEC MALDIALO → GLYOX + GLYOX + HO2. Only major channel.

G44461: KBPAN → k_PAN_M

G45019d: Delta-1 and delta-2 LIEPOX are not considered and replaced by beta-LIEPOX formed by ISOPBOOH and ISOPDOOH.

G45021: SAR estimate within uncertainty range of the experimentally determined rate constant by Solberg et al. (1997), 1.1E-11.

G45037: SAR estimate within uncertainty range of the experimentally determined rate constant by Solberg et al. (1997), 4.2E-11.

G45040: Alkyl nitrate formation neglected.

G45043: Old MCM rate constant 4.16E-11.

G45047: Alkyl nitrate formation neglected.

G45055: Alkyl nitrate formation neglected.

G45071: Alkyl nitrate formation neglected.

G45074: Formic acid production consistent with results of Bates et al. (2014). Here, the high yields of formic acid and hydroxycarbonyls at low NO from oxidation of cis-beta-LIEPOX (the most abundant isomer) are approximated with the production of DB1O which undergo both the Dibble double H-transfer to DB2O2 and HOCH2 elimination yielding HVMK and HMAC (keto-vinyl alcohol potentially arising from decomposition of the alkoxy radical resulting from the ring opening after H-abstraction). The rate constant is from Paulot et al. (2009b) and adjusted based on Bates et al. (2014) that determined the single rate constants for the cis- and trans- beta isomer.

G45080: Alkyl nitrate formation neglected.

G45092a: C4MDIAL = CM4DIAL in MCM only from aromatics.

G45092b: Only one acyl peroxy radical considered.

G45093: Two aldehydic sites reacting with NO_3 but only one isomer product considered.

G45095: Alkyl nitrate formation neglected.

G45098: Alkyl nitrate formation neglected.

G45100: Alkyl nitrate formation neglected.

G45104a: DB1OOH is a hydroperoxide bearing a vinyl alcohol moiety that upon reaction with OH yields HCOOH (Davis et al., 1998).

G45107: OH production here is to take into account the hydroperoxidic function formed by the shift of the enolic hydrogen and not present in DB2O2. This approximation leads to spurious HO_2 production.

G45108a: Consistent with the results of Bates et al. (2014).

G45108b: Consistent with the results of Bates et al. (2014). Assuming that the enol alkoxy radical partly decomposes yielding a substitute vinyl alcohol.

G45111: Alkyl nitrate formation neglected.

G45114b: Here, formic acid is mechanistically produced by the OH-addition to the vinyl alcohol which, upon RO_2-to-RO conversion (skipped here), yields the HOCHOH fragment which in turn reacts with O_2 forming $HCOOH + HO_2$. Along $CH_3COCHOOHCHO$ should be produced but not in the mechanism. Only CH_3COCHO_2CHO. The rate constant is consistent with predictions by Ganzeveld et al. (2006) for ENOL. OH-addition to the OH-bearing carbon is considered the dominant channel as it is already for the ENOL (Ganzeveld et al., 2006).

G45115: Theoretical keto-enol tautomerization catalyzed by formic acid (Grenfell et al., 2006). The product should be C1ODC3OOHC4OD but it is neglected in the mechanism.

G45116: As for DB1OOH + OH.

G45117: Additional sinks for DB2OOH are neglected.

G45121b: Nitrate assumed to be major isomer that is mostly similar to products of ISOPDO2-chemistry.

G45128: Rate constant by Liljegren and Stevens (2013). A lumped RO_2 that upon conversion to RO yields 100% 2-methyl-butenedial (C4MDIAL) although Aschmann et al. (2014) quantified a 38% yield of the Z/E mixture.

G45129: As for 3METHYLFURAN + OH but with additional NO_2 production for mass conservation.

G45131: Alkyl nitrate formation neglected.

G45132: Hydroperoxide formation neglected.

G45134b: ZCO2HC23DBCOD formation is neglected. However, it is produced in MCM and in aromatic-related reactions under the name of MC3ODBCO2H.

G45139: LZCPANC23DBCOD is assumed to react like LC5PAN1719.

G45201: Alkyl nitrate formation neglected.

G45207: Alkyl nitrate formation neglected.

G45214: Alkyl nitrate formation neglected.

G45217: Alkyl nitrate formation neglected.

G45225: Alkyl nitrate formation neglected.

G45236: LMBOABO2 = 0.67 MBOAO2 + 0.33 MBOBO2

G45247: Alkyl nitrate formation neglected.

G45400: KDEC NC4MDCO2 → MMALANHY + NO2

G45404: KDEC NTLFUO → ACCOMECHO + NO2

G45405: KDEC NTLFUO → ACCOMECHO + NO2

G45406: KDEC NTLFUO → ACCOMECHO

G45409: KBPAN → k_PAN_M(renaming)

G45413: KFPAN → k_CH3CO3_NO2 (renaming)

G45422: KDEC MMALANHYO→CO2H3CO3

G45423: KDEC MMALANHYO→CO2H3CO3

G45424: KDEC MMALANHYO→CO2H3CO3 and Only major channel.

G45429: KBPAN → k_PAN_M (renamed)

G45430a: KDEC C5CO14CO2 → 0.83 MALANHY + 0.83 CH3 + 0.17 MGLYOX + 0.17 HO2 + 0.17 CO + 0.17 CO2

G45431: KDEC C5CO14CO2 → 0.83 MALANHY + 0.83 CH3 + 0.17 MGLYOX + 0.17 HO2 + 0.17 CO + 0.17 CO2

G45432: KFPAN →k_CH3CO3_NO2 (renaming)

G45433: KDEC C5CO14CO2 → 0.83 MALANHY + 0.83 CH3 + 0.17 MGLYOX + 0.17 HO2 + 0.17 CO + 0.17 CO2

G45434: KDEC C5CO14CO2 → 0.83 MALANHY + 0.83 CH3 + 0.17 MGLYOX + 0.17 HO2 + 0.17 CO + 0.17 CO2 and only major channel.

G45436: KDEC C5CO14CO2 → 0.83 MALANHY + 0.83 CH3 + 0.17 MGLYOX + 0.17 HO2 + 0.17 CO + 0.17 CO2

G45444: KDEC MC3CODBCO2 → 0.35 GLYOX + 0.35 CH3 + 0.35 CO + 0.35 CO2 + 0.65 MMALANHY + 0.65 HO2

G45452: KDEC TLFUONOOA → 0.5 CO + 0.5 OH + 0.5 MECOACETO2 + 0.5 TLFUONOO and H2O subs TLFUONOO → 0.625 C24O3CCO2H + 0.375 ACCOMECHO + 0.375 H2O2

G45456: KFPAN →k_CH3CO3_NO2 (renaming)

G45476b: KDEC NTLFUO → ACCOMECHO + NO2 and reactions with KRO2HO2.

G45477: KDEC NTLFUO → ACCOMECHO + NO2

G45478: KDEC NTLFUO → ACCOMECHO + NO2

G45479: KDEC NTLFUO → ACCOMECHO + NO2

G45486b: KDEC C5DIALO →MALDIAL + CO + HO2 and reactions with KRO2HO2.

G45487: KDEC C5DIALO →MALDIAL

G45488: KDEC C5DIALO →MALDIAL

G45489: KDEC C5DIALO →MALDIAL

G45491b: Reactions with KRO2HO2.

G45492: MGLYOX + GLYOX + HO2 from KDEC substitution

G45493: MGLYOX + GLYOX + HO2 from KDEC substitution

G45494: Permutation reaction (minor channels removed).

G46201: Alkyl nitrate formation neglected.

G46404b: Reactions with KRO2HO2 and KDEC C615CO2O → C5DICARB + CO + HO2.

G46405: KDEC C615CO2O → C5DICARB + CO + HO2

G46406: KDEC C615CO2O → C5DICARB + CO + HO2

G46407: Only major channel.

G46413b: Reactions with KRO2HO2 and KDEC NDNPHENO → NC4DCO2H + HNO3 + CO + CO + NO2.

G46414: KDEC NDNPHENO → NC4DCO2H + HNO3 + CO + CO + NO2

G46415: KDEC NDNPHENO → NC4DCO2H + HNO3 + CO + CO + NO2

G46416: KDEC NDNPHENO → NC4DCO2H + HNO3 + CO + CO + NO2

G46418: KDEC CATECOOA → MALDALCO2H + HCOCO2H + HO2 + OH

G46426: KFPAN →k_CH3CO3_NO2

G46430: KDEC GLYOOA → .125 HCHO + .18 GLYOO + .82 HO2 + .57 OH + 1.265 CO

G46432b: Reactions with KRO2HO2 and KDEC NCATECO → NC4DCO2H + HCOCO2H + HO2

G46433: KDEC NCATECO → NC4DCO2H + HCOCO2H + HO2

G46434: KDEC NCATECO → NC4DCO2H + HCOCO2H + HO2

G46435: KDEC NCATECO → NC4DCO2H + HCOCO2H + HO2

G46437b: Reactions with KRO2HO2 and KDEC NPHENO → MALDALCO2H + GLYOX + NO2

G46438: KDEC NPHENO → MALDALCO2H + GLYOX + NO2

G46439: KDEC NPHENO → MALDALCO2H + GLYOX + NO2

G46440: KDEC NPHENO → MALDALCO2H + GLYOX + NO2

G46441: Merged equations.

G46447b: reactions with KRO2HO2 and KDEC NNCATECO → NC4DCO2H + HCOCO2H + NO2

G46448: KDEC NNCATECO → NC4DCO2H + HCOCO2H + NO2

G46449: KDEC NNCATECO → NC4DCO2H + HCOCO2H + NO2

G46450: KDEC NNCATECO → NC4DCO2H + HCOCO2H + NO2

G46457: Merged equations.

G46458: Merged equations.

G46461b: Reactions with KRO2HO2 and KDEC PHENO → 0.71 MALDALCO2H + 0.71 GLYOX + 0.29 PBZQONE + HO2

G46462: KDEC PHENO → 0.71 MALDALCO2H + 0.71 GLYOX + 0.29 PBZQONE + HO2

G46463: KDEC PHENO → 0.71 MALDALCO2H + 0.71 GLYOX + 0.29 PBZQONE + HO2

G46464: KDEC PHENO → 0.71 MALDALCO2H + 0.71 GLYOX + 0.29 PBZQONE + HO2 and Only major channel.

G46468: KFPAN →k_CH3CO3_NO2

G46472b: new channel

G46476: HOC6H4NO2 is a nitro-phenol

G46480b: Reactions with KRO2HO2 and KDEC PBZQO →C5CO2OHCO3

G46481: KDEC PBZQO →C5CO2OHCO3

G46482: KDEC PBZQO →C5CO2OHCO3

G46483: KDEC PBZQO →C5CO2OHCO3 and Only major channel.

G46485b: Reactions with KRO2HO2 and KDEC DNPHENO → NC4DCO2H + HCOCO2H + NO2

G46486: KDEC DNPHENO → NC4DCO2H + HCOCO2H + NO2

G46487: KDEC DNPHENO → NC4DCO2H + HCOCO2H + NO2

G46488: KDEC DNPHENO → NC4DCO2H + HCOCO2H + NO2

G46490b: Reactions with KRO2HO2 and KDEC BZEMUCO → 0.5 EPXC4DIAL + 0.5 GLYOX + 0.5 HO2 + 0.5 C3DIALO2 + 0.5 C32OH13CO.

G46491b: KDEC BZEMUCO → 0.5 EPXC4DIAL + 0.5 GLYOX + 0.5 HO2 + 0.5 C3DIALO2 + 0.5 C32OH13CO.

G46492: KDEC BZEMUCO → 0.5 EPXC4DIAL + 0.5 GLYOX + 0.5 HO2 + 0.5 C3DIALO2 + 0.5 C32OH13CO

G46493: KDEC BZEMUCO → 0.5 EPXC4DIAL + 0.5 GLYOX + 0.5 HO2 + 0.5 C3DIALO2 + 0.5 C32OH13CO and Only major channel.

G46499b: Reactions with KRO2HO2 and KDEC NBZQO → C6CO4DB + NO2.

G46500: KDEC NBZQO → C6CO4DB + NO2

G46501: KDEC NBZQO → C6CO4DB + NO2

G46502: KDEC NBZQO → C6CO4DB + NO2

G46505b: New channel.

G46515: Only major channel.

G46517b: New channel.

G46522b: In analogy to TLBIPERO2 from toluene (Birdsall et al., 2010).

G46523b: KDEC BZBIPERO → GLYOX + HO2 + 0.5 BZFUONE + 0.5 BZFUONE

G46524: KDEC BZBIPERO → GLYOX + HO2 + 0.5 BZFUONE + 0.5 BZFUONE

G46525: KDEC BZBIPERO → GLYOX + HO2 + 0.5 BZFUONE + 0.5 BZFUONE and Only major channel.

G47210: Alkyl nitrate formation neglected.

G47214: Alkyl nitrate formation neglected.

G47218: Alkyl nitrate formation neglected.

G47222: Alkyl nitrate formation neglected.

G47223: ROO6R3OOH produced but no sink for it.

G47225: ROO6R4P produced but no sink for it.

G47226: ROO6R5P produced but no sink for it

G47400: Merged.

G47402a: KROPRIM*O2 fast reaction C6H5CH2O = BENZAL + HO2.

G47402b: KROPRIM*O2 fast reaction C6H5CH2O = BENZAL + HO2.

G47403: KROPRIM*O2 fast reaction C6H5CH2O = BENZAL + HO2.

G47404: KROPRIM*O2 fast reaction C6H5CH2O = BENZAL + HO2. C6H5CH2OH replaced by its oxidation product BENZAL.

G47405: Merged.

G47406: Merged.

G47407b: According to Birdsall et al. (2010), the branching ratio rbipero2_oh is set to 0.4 in order to take into account the OH-recycling and summed yield of butendial and methylbutendial.

G47408a: KDEC TLBIPERO → 0.6 GLYOX + 0.4 MGLYOX + HO2 + 0.2 C4MDIAL + 0.2 C5DICARB + 0.2 TLFUONE + 0.2 BZFUONE + 0.2 MALDIAL

G47408b: KDEC TLBIPERO → 0.6 GLYOX + 0.4 MGLYOX + HO2 + 0.2 ZCODC23DB COD + 0.2 C5DICARB + 0.2 TLFUONE + 0.2 BZFUONE + 0.2 MALDIAL

G47409: KDEC TLBIPERO → 0.6 GLYOX + 0.4 MGLYOX + HO2 + 0.2 ZCODC23DB COD + 0.2 C5DICARB + 0.2 TLFUONE + 0.2 BZFUONE + 0.2 MALDIAL

G47410: Only major channel and KDEC TLBIPERO → 0.6 GLYOX + 0.4 MGLYOX + HO2 + 0.2 ZCODC23DB COD + 0.2 C5DICARB + 0.2 TLFUONE + 0.2 BZFUONE + 0.2 MALDIAL

G47412: KDEC MGLOOB → 0.125 CH3CHO + 0.695 CH3CO + 0.57 CO + 0.57 OH + 0.125 HO2 + 0.18 MGLOO + 0.25 CO2

G47413: Merged.

G47418b: Reactions with KRO2HO2 and KDEC CRESO → 0.68 C5CO14OH + 0.68 GLYOX + HO2 + 0.32 PTLQONE.

G47419: KDEC CRESO → 0.68 C5CO14OH + 0.68 GLYOX + HO2 + 0.32 PTLQONE

G47420: KDEC CRESO → 0.68 C5CO14OH + 0.68 GLYOX + HO2 + 0.32 PTLQONE

G47421: KDEC CRESO → 0.68 C5CO14OH + 0.68 GLYOX + HO2 + 0.32 PTLQONE and Only major channel.

G47422b: Reactions with KRO2HO2 and KDEC NCRESO → C5CO14OH + GLYOX + NO2

G47423: KDEC NCRESO → C5CO14OH + GLYOX + NO2

G47424: KDEC NCRESO → C5CO14OH + GLYOX + NO2

G47425: KDEC NCRESO → C5CO14OH + GLYOX + NO2 and Only major channel.

G47426: TOL1OHNO2 is a nitro-phenol

G47429: KDEC MCATECOOA → MC3ODBCO2H + HCOCO2H + HO2 + OH

G47436: KFPAN →k_CH3CO3_NO2

G47438: Only major channel.

G47439b: Reactions with KRO2HO2 and KDEC TLEMUCO→ 0.5 C3DIALO2 + 0.5 CO2H3CHO + 0.5 EPXC4DIAL + 0.5 MGLYOX + 0.5 HO2

G47440b: KDEC TLEMUCO→ 0.5 C3DIALO2 + 0.5 CO2H3CHO + 0.5 EPXC4DIAL + 0.5 MGLYOX + 0.5 HO2

G47441: KDEC TLEMUCO→ 0.5 C3DIALO2 + 0.5 CO2H3CHO + 0.5 EPXC4DIAL + 0.5 MGLYOX + 0.5 HO2

G47442: KDEC TLEMUCO→ 0.5 C3DIALO2 + 0.5 CO2H3CHO + 0.5 EPXC4DIAL + 0.5 MGLYOX + 0.5 HO2 and Only major channel.

G47445: KFPAN →k_CH3CO3_NO2

G47447: Only major channel.

G47454: New channel.

G47479: New channel.

G47482b: Reactions with KRO2HO2 and KDEC NPTLQO → C7CO4DB + NO2

G47483: KDEC NPTLQO → C7CO4DB + NO2

G47484: KDEC NPTLQO → C7CO4DB + NO2

G47485: KDEC NPTLQO → C7CO4DB + NO2

G47486b: Reactions with KRO2HO2 and KDEC PTLQO → C6CO2OHCO3

G47487: KDEC PTLQO → C6CO2OHCO3

G47488: KDEC PTLQO → C6CO2OHCO3

G47489: Only major channel. KDEC PTLQO → C6CO2OHCO3.

G47494: New channel.

G47497b: Reactions with KRO2HO2 and KDEC MNNCATECO → NC4MDCO2H + HCOCO2H + NO2

G47498: KDEC MNNCATECO → NC4MDCO2H + HCOCO2H + NO2

G47499: KDEC MNNCATECO → NC4MDCO2H + HCOCO2H + NO2

G47501b: Reactions with KRO2HO2 and KDEC MNCATECO → NC4MDCO2H + HCOCO2H + HO2

G47502: KDEC MNCATECO → NC4MDCO2H + HCOCO2H + HO2

G47503: KDEC MNCATECO → NC4MDCO2H + HCOCO2H + HO2

G47504: KDEC MNCATECO → NC4MDCO2H + HCOCO2H + HO2

G47509b: Reactions with KRO2HO2 and KDEC NDNCRESO → NC4MDCO2H + HNO3 + CO + CO + NO2

G47510: KDEC NDNCRESO → NC4MDCO2H + HNO3 + CO + CO + NO2

G47511: KDEC NDNCRESO → NC4MDCO2H + HNO3 + CO + CO + NO2

G47512: KDEC NDNCRESO → NC4MDCO2H + HNO3 + CO + CO + NO2

G47513b: Reactions with KRO2HO2 and KDEC DNCRESO → NC4MDCO2H + HCOCO2H + NO2

G47514: KDEC DNCRESO → NC4MDCO2H + HCOCO2H + NO2

G47515: KDEC DNCRESO → NC4MDCO2H + HCOCO2H + NO2

G47516: KDEC DNCRESO → NC4MDCO2H + HCOCO2H + NO2

G48202: Alkyl nitrate formation neglected.

G48205: Alkyl nitrate formation neglected.

G48210: Alkyl nitrate formation neglected.

G48212: Alkyl nitrate formation neglected.

G48216: Alkyl nitrate formation neglected.

G48222: Alkyl nitrate formation neglected.

G48400a: Same products as for toluene. Assuming a 1:1:1 proportion in xylenes emissions the analogous toluene product is produced with a rate constant equal to (1.36E-11*0.24 + 2.31E-11*0.29 + 1.43E-11*0.155)/3, where k and coefficients are for the single isomers ortho, meta and para from MCM.

G48400b: Same products as for toluene. Assuming a 1:1:1 proportion in xylenes emissions the analogous toluene product is produced with a rate constant equal to (1.36E-11*0.05 + 2.31E-11*0.04 + 1.43E-11*0.10)/3, where k and coefficients are for the single isomers ortho, meta and para from MCM.

G48400c: Same products as for toluene. Assuming a 1:1:1 proportion in xylenes emissions the analogous toluene product is produced with a rate constant equal to (1.36E-11*0.16 + 2.31E-11*0.17 + 1.43E-11*0.12)/3, where k and coefficients are for the single isomers ortho, meta and para from MCM.

G48400d: Same products as for toluene. Assuming a 1:1:1 proportion in xylenes emissions the analogous toluene product is produced with a rate constant equal to (1.36E-11*0.55 + 2.31E-11*0.50 + 1.43E-11*0.625)/3, where k and coefficients are for the single isomers ortho, meta and para from MCM.

G48401: Same products as for toluene. The rate constant is the average of m, p, o k=(4.10E-16+2.60E-16+5.00E-16)/3 = 3.9E-16.

G48402: merged under same rate constant

G48403: Same products as for toluene

G48405: KDEC CH2OOB → 0.24 CH2OO + 0.40 CO + 0.36 HO2 + 0.36 CO + 0.36 OH and H2O + PHCHOO → 0.625 PHCOOH + 0.375 BENZAL + 0.375 H2O2 + 0.2 CO2

G48408: KDEC NSTYRENEO → NO2 + HCHO + BENZAL

G48409: KDEC NSTYRENEO → NO2 + HCHO + BENZAL

G48410: KDEC NSTYRENEO → NO2 + HCHO + BENZAL

G48412b: KDEC STYRENO → HO2 + HCHO + BENZAL and reactions with KRO2HO2.

G48413: KDEC STYRENO → HO2 + HCHO + BENZAL

G48414: KDEC STYRENO → HO2 + HCHO + BENZAL

G48415: KDEC STYRENO → HO2 + HCHO + BENZAL

G49207: Alkyl nitrate formation neglected.

G49238: Alkyl nitrate formation neglected.

G49246: Only this channel considered as the intermediate radical is likely more stable than $CHCH(OH)_2$.Instead of the (lacking) carbonyl a product of further degradation is assumed.

G49247: Theoretical keto-enol tautomerization catalyzed by formic acid (Grenfell et al., 2006).

G49248: Theoretical keto-enol tautomerization catalyzed by formic acid (Grenfell et al., 2006).

G49400a: Same products as for toluene. Assuming a 1:1:1 proportion in xylenes emissions the analogous toluene product is produced with a rate constant equal to (3.27E-11*0.70 + 3.25E-11*0.61 + 5.67E-11*0.79)/3, where k and coefficients are for the single isomers 1,2,3-, 1,3,4- and 1,3,5- from MCM.

G49400b: Same products as for toluene. Assuming a 1:1:1 proportion in xylenes emissions the analogous toluene product is produced with a rate constant equal to (3.27E-11*0.06 + 3.25E-11*0.06 + 5.67E-11*0.03)/3, where k and coefficients are for the single isomers 1,2,3-, 1,3,4- and 1,3,5- from MCM.

G49400c: Same products as for toluene. Assuming a 1:1:1 proportion in xylenes emissions the analogous toluene product is produced with a rate constant equal to (3.27E-11*0.03 + 3.25E-11*0.03 + 5.67E-11*.04)/3, where k and coefficients are for the single isomers 1,2,3-, 1,3,4- and 1,3,5- from MCM.

G49400d: Same products as for toluene. Assuming a 1:1:1 proportion in xylenes emissions the analogous toluene product is produced with a rate constant equal to (3.27E-11*0.70 + 3.25E-11*0.61 + 5.67E-11*0.79)/3, where k and coefficients are for the single isomers 1,2,3-, 1,3,4- and 1,3,5- from MCM.

G49401: Same products as for toluene. The rate constant is the average of m, p, o k=(1.90+1.80+0.88)E-15/3=1.52E-15.

G40200: Products from Vereecken et al. (2007). LAPINABO2 = 0.65 APINAO2 + 0.35 APINBO2

G40203: Weighted average for isomers A and B, `k = 0.33*9.20E-14+0.67*8.80E-13`.

G40204: Weighted average for isomers A and B, `k = 0.35*1.83E-11+0.65*3.28E-11`.

G40205: Weighted average for isomers A and B, `k = 0.35*5.50E-12+0.65*3.64E-12`.

G40206: SAR-estimated rate constant, `(kads+kadt)*acoch3 = 6.46E-11` where `kads = 3.0E-11`, `kadt = 5.5E-11`, `acoch3 = 0.76`

G40207: Alkyl nitrate formation neglected.

G40211: Products from Rickard and Pascoe (2009).

G40212: Products from Rickard and Pascoe (2009).

G40232: Products from Capouet et al. (2008). BELV from Ehn et al. (2014). BLOV guess

G40242: Alkyl nitrate formation neglected.

G40246: Products from Rickard and Pascoe (2009).

G40248: Alkyl nitrate formation neglected.

G40252a: Products from Vereecken and Peeters (2012).

G40252b: Products from Vereecken and Peeters (2012).

G40262: RO6R1OOH is produced but no sink for it.

G40266: Rate constant modified according to MCM protocol.

G40267a: Products from Nguyen et al. (2009). BELV from Ehn et al. (2014). BLOV guess

G40268: Products from Rickard and Pascoe (2009).

G40270: Alkyl nitrate neglected.

G40274: As for RO6R1NO3 in G4085.

G40276: Only this channel considered as the intermediate radical is likely more stable than $CHCH(OH)_2$.

G40277: Theoretical keto-enol tautomerization catalyzed by formic acid (Grenfell et al., 2006).

G40278: Theoretical keto-enol tautomerization catalyzed by formic acid (Grenfell et al., 2006).

G40282a: Products from Vereecken and Peeters (2012).

G40282b: Products from Vereecken and Peeters (2012).

G40283a: Products from Nguyen et al. (2009).

G40284: Products from Rickard and Pascoe (2009).

G40285a: Products from Vereecken and Peeters (2012).

G40285b: Products from Vereecken and Peeters (2012).

G40286a: Products from Nguyen et al. (2009).

G40287: Products from Rickard and Pascoe (2009).

G40400: DIET35TOL(from MCM) as representative of higher aromatics

G40401: Same products as for toluene.

G6103: The rate coefficient is defined as backward reaction divided by equilibrium constant.

G6204: At low temperatures, there may be a minor reaction channel leading to O_3+HCl. See Finkbeiner et al. (1995) for details. It is neglected here.

G6402: The initial products are probably HCl and CH_2OOH (Atkinson et al., 2006). It is assumed that CH_2OOH dissociates into HCHO and OH.

G6409: It is assumed that the reaction liberates all Cl atoms in the form of HCl.

G8102: Consistent with O'Dowd and Hoffmann (2005), it is assumed that the reaction produces new particles.

G8103: The yield of 38 % OIO is from Atkinson et al. (2007). It is assumed here that the remaining 62 % produce 2 I + O_2.

G8300: The rate coefficient is: `k_I_NO2 = k_3rd_iupac(temp,cair,3.E-31,1.,6.6E-11,0.,0.63)`.

G8305: The rate coefficient is defined as backward reaction (Atkinson et al., 2007) divided by equilibrium constant (van den Bergh and Troe, 1976).

G8401: The rate coefficient is from Dillon et al. (2006b), the yield of I atoms is a lower limit given on page 2170 of Bale et al. (2005).

G8402: The products are from Nakano et al. (2005).

G9408: Average of 3.9E-11 and 3.42E-11.

G40259a: ROO6R1OOH is produced but no sink for it.

Table 2: Photolysis reactions

#	labels	reaction	rate coefficient	reference
J (gas)				
J1000a	UpStTrGJ	$O_2 + h\nu \rightarrow O(^3P) + O(^3P)$	jx(ip_O2)	Sander et al. (2014)
J1001a	UpStTrGJ	$O_3 + h\nu \rightarrow O(^1D) + O_2$	jx(ip_O1D)	Sander et al. (2014)
J1001b	UpStTrGJ	$O_3 + h\nu \rightarrow O(^3P) + O_2$	jx(ip_O3P)	Sander et al. (2014)
J2101	UpStTrGJ	$H_2O_2 + h\nu \rightarrow 2\ OH$	jx(ip_H2O2)	Sander et al. (2014)
J3101	UpStTrGJN	$NO_2 + h\nu \rightarrow NO + O(^3P)$	jx(ip_NO2)	Sander et al. (2014)
J3103a	UpStTrGJN	$NO_3 + h\nu \rightarrow NO_2 + O(^3P)$	jx(ip_NO2O)	Sander et al. (2014)
J3103b	UpStTrGJN	$NO_3 + h\nu \rightarrow NO + O_2$	jx(ip_NOO2)	Sander et al. (2014)
J3104	StTrGJN	$N_2O_5 + h\nu \rightarrow NO_2 + NO_3$	jx(ip_N2O5)	Sander et al. (2014)
J3200	TrGJN	$HONO + h\nu \rightarrow NO + OH$	jx(ip_HONO)	Sander et al. (2014)
J3201	StTrGJN	$HNO_3 + h\nu \rightarrow NO_2 + OH$	jx(ip_HNO3)	Sander et al. (2014)
J3202	StTrGJN	$HNO_4 + h\nu \rightarrow .667\ NO_2 + .667\ HO_2 + .333\ NO_3 + .333\ OH$	jx(ip_HNO4)	Sander et al. (2014)
J41000	StTrGJ	$CH_3OOH + h\nu \rightarrow CH_3O + OH$	jx(ip_CH3OOH)	Sander et al. (2014)
J41001a	StTrGJ	$HCHO + h\nu \rightarrow H_2 + CO$	jx(ip_COH2)	Sander et al. (2014)
J41001b	StTrGJ	$HCHO + h\nu \rightarrow H + CO + HO_2$	jx(ip_CHOH)	Sander et al. (2014)
J41004	StTrGJN	$CH_3ONO + h\nu \rightarrow CH_3O + NO$	jx(ip_CH3ONO)	Sander et al. (2014)
J41005	StTrGJN	$CH_3ONO_2 + h\nu \rightarrow CH_3O + NO_2$	jx(ip_CH3NO3)	Sander et al. (2014)
J41006	StTrGJN	$CH_3O_2NO_2 + h\nu \rightarrow .667\ NO_2 + .667\ CH_3O_2 + .333\ NO_3 + .333\ CH_3O$	jx(ip_CH3O2NO2)	Sander et al. (2014)*
J41007	StTrGJ	$HOCH_2OOH + h\nu \rightarrow HCOOH + OH + HO_2$	jx(ip_CH3OOH)	Sander et al. (2014)
J41008	StTrGJ	$CH_3O_2 + h\nu \rightarrow HCHO + OH$	jx(ip_CH3O2)	Sander et al. (2014)
J41009	StTrGJ	$HCOOH + h\nu \rightarrow CO + HO_2 + OH$	jx(ip_HCOOH)	Sander et al. (2014)
J41010	StTrGJN	$HOCH_2O_2NO_2 + h\nu \rightarrow .667\ NO_2 + .667\ HOCH_2O_2 + .333\ NO_3 + .333\ HCOOH + .333\ HO_2$	jx(ip_CH3O2NO2)	Sander et al. (2014)
J42000	TrGJC	$C_2H_5OOH + h\nu \rightarrow CH_3CHO + HO_2 + OH$	jx(ip_CH3OOH)	von Kuhlmann (2001)
J42001a	TrGJC	$CH_3CHO + h\nu \rightarrow CH_3 + HO_2 + CO$	jx(ip_CH3CHO)	Sander et al. (2014)
J42001b	TrGJC	$CH_3CHO + h\nu \rightarrow CH_2CHOH$	jx(ip_CH3CHO2VINY)	Clubb et al. (2012)
J42002	TrGJC	$CH_3C(O)OOH + h\nu \rightarrow CH_3 + OH + CO_2$	jx(ip_CH3CO3H)	Sander et al. (2014)
J42004	TrGJCN	$PAN + h\nu \rightarrow .7\ CH_3C(O) + .7\ NO_2 + .3\ CH_3 + .3\ CO_2 + .3\ NO_3$	jx(ip_PAN)	Sander et al. (2014)*
J42005a	TrGJC	$HOCH_2CHO + h\nu \rightarrow HCHO + 2\ HO_2 + CO$	jx(ip_HOCH2CHO)*0.83	Sander et al. (2014)*
J42005b	TrGJC	$HOCH_2CHO + h\nu \rightarrow OH + HCOCH_2O_2$	jx(ip_HOCH2CHO)*0.07	Sander et al. (2014)*
J42005c	TrGJC	$HOCH_2CHO + h\nu \rightarrow CH_3OH + CO$	jx(ip_HOCH2CHO)*0.10	Sander et al. (2014)*
J42006	TrGJC	$HOCH_2CO_3H + h\nu \rightarrow HCHO + HO_2 + OH + CO_2$	jx(ip_CH3OOH)	Rickard and Pascoe (2009)

Table 2: Photolysis reactions (... continued)

#	labels	reaction	rate coefficient	reference
J42007	TrGJCN	PHAN + hν $\rightarrow$ .7 HOCH2CO + .7 NO_2 + .3 HCHO + .3 HO_2 + .3 CO_2 + .3 NO_3	jx(ip_PAN)	see note*
J42008	TrGJC	GLYOX + hν $\rightarrow$ 2 CO + 2 HO_2	jx(ip_GLYOX)	Sander et al. (2014)
J42009	TrGJC	$HCOCO_2H$ + hν $\rightarrow$ 2 HO_2 + CO + CO_2	jx(ip_MGLYOX)	Rickard and Pascoe (2009)
J42010	TrGJC	$HCOCO_3H$ + hν $\rightarrow$ HO_2 + CO + OH + CO_2	jx(ip_CH3OOH)+jx(ip_HOCH2CHO)	Rickard and Pascoe (2009)
J42011	TrGJC	HYETHO2H + hν $\rightarrow$ $HOCH_2CH_2O$ + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J42012	TrGJCN	ETHOHNO3 + hν $\rightarrow$ HO_2 + 2 HCHO + NO_2	j_IC3H7NO3	Rickard and Pascoe (2009)
J42013	TrGJC	HOOCH2CO3H + hν $\rightarrow$ OH + HCHO + CO_2 + OH	2*jx(ip_CH3OOH)	Sander et al. (2018)
J42014	TrGC	HOOCH2CO2H + hν $\rightarrow$ OH + HCHO + HO_2 + CO_2	jx(ip_CH3OOH)	Sander et al. (2018)
J42015	TrGC	CH2CO + hν $\rightarrow$ .4 CO_2 + .8 H + .34 CO + .34 OH + .34 HO_2 + .16 HCHO + .16 $O(^3P)$ + .1 HCOOH + CO	j_ketene* 0.36	Sander et al. (2018)
J42016	TrGC	CH3CHOHOOH + hν $\rightarrow$ CH_3 + HCOOH + OH	jx(ip_CH3OOH)	Sander et al. (2018)
J42017	TrGJCN	NO_3CH2CHO + hν $\rightarrow$ HO_2 + CO + HCHO + NO_2	(jx(ip_C2H5NO3)+jx(ip_CH3CHO)) *(jx(ip_NOA)+1E-10)/(0.59*j_IC3H7NO3+jx(ip_CH3COCH3)+1E-10)	Sander et al. (2018)*
J42018	TrGJC	HOOCH2CHO + hν $\rightarrow$ OH + HCHO + CO + HO_2	jx(ip_CH3OOH)+jx(ip_HOCH2CHO)	Sander et al. (2018)
J42019	TrGJCN	$C_2H_5ONO_2$ + hν $\rightarrow$ CH_3CHO + HO_2 + NO_2	jx(ip_C2H5NO3)	Sander et al. (2018)
J42020	TrGJCN	NO_3CH2CHO + hν $\rightarrow$ .7 NO_3CH2CO_3 + .7 NO_2 + .3 HCHO + .3 NO_2 + .3 CO_2 + .3 NO_3	jx(ip_PAN)	Sander et al. (2018)*
J42021	StTrGJCN	$C_2H_5O_2NO_2$ + hν $\rightarrow$ .667 NO_2 + .667 $C_2H_5O_2$ + .333 NO_3 + .333 CH_3CHO + .333 HO_2	jx(ip_CH3O2NO2)	Sander et al. (2018)*
J43000	TrGJC	iC_3H_7OOH + hν $\rightarrow$ CH_3COCH_3 + HO_2 + OH	jx(ip_CH3OOH)	von Kuhlmann (2001)
J43001	TrGJC	CH_3COCH_3 + hν $\rightarrow$ $CH_3C(O)$ + CH_3	jx(ip_CH3COCH3)	Sander et al. (2014)
J43002	TrGJC	CH_3COCH_2OH + hν $\rightarrow$ .5 $CH_3C(O)$ + .5 HCHO + .5 HO_2 + .5 HOCH2CO + .5 CH_3	j_ACETOL	Sander et al. (2014)*
J43003	TrGJC	MGLYOX + hν $\rightarrow$ $CH_3C(O)$ + CO + HO_2	jx(ip_MGLYOX)	Sander et al. (2014)
J43004	TrGJC	$CH_3COCH_2O_2H$ + hν $\rightarrow$ $CH_3C(O)$ + HCHO + OH	jx(ip_CH3OOH)+j_ACETOL	Rickard and Pascoe (2009)
J43005	TrGJC	HOCH2COCH2OOH + hν $\rightarrow$ HOCH2CO + HCHO + OH	jx(ip_CH3OOH)+j_ACETOL	Sander et al. (2018)
J43006	TrGJCN	$iC_3H_7ONO_2$ + hν $\rightarrow$ CH_3COCH_3 + NO_2 + HO_2	j_IC3H7NO3	von Kuhlmann et al. (2003)*
J43007	TrGJCN	NOA + hν $\rightarrow$ $CH_3C(O)$ + HCHO + NO_2	jx(ip_NOA)	Barnes et al. (1993)
J43009	TrGJC	HYPROPO2H + hν $\rightarrow$ CH_3CHO + HCHO + HO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J43010	TrGJCN	PR2O2HNO3 + hν $\rightarrow$ NOA + HO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J43011	TrGJC	HOCH2COCHO + hν $\rightarrow$ HOCH2CO + CO + HO_2	jx(ip_MGLYOX)	Rickard and Pascoe (2009)
J43012	TrGJC	$HCOCOCH_2OOH$ + hν $\rightarrow$ HCOCO + HCHO + OH	jx(ip_CH3OOH)+j_ACETOL	Sander et al. (2018)

Table 2: Photolysis reactions (... continued)

#	labels	reaction	rate coefficient	reference
J43013	TrGJC	$HCOCOCH_2OOH + h\nu \rightarrow HOOCH_2CO_3 + CO + HO_2$	jx(ip_MGLYOX)	Sander et al. (2018)
J43014	TrGJTerC	$HCOCH2CHO + h\nu \rightarrow HCOCH_2O_2 + HO_2 + CO$	jx(ip_HOCH2CHO)*2.	Rickard and Pascoe (2009)
J43015	TrGJTerC	$HCOCH2CO2H + h\nu \rightarrow HCOCH_2O_2 + CO_2 + HO_2$	jx(ip_HOCH2CHO)	Rickard and Pascoe (2009)
J43016	TrGJTerC	$HOC2H4CO3H + h\nu \rightarrow HOCH_2CH_2O_2 + CO_2 + OH$	jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J43017	TrGJC	$HCOCOCHO + h\nu \rightarrow HCOCO + HO_2 + CO$	2.*jx(ip_MGLYOX)	Sander et al. (2018)
J43018	TrGJC	$CH_3COCO_2H + h\nu \rightarrow .32\ CH_3CHO + .16\ CH_2CHOH + .54\ CO_2 + .38\ CH_3C(O) + .38\ HO_2 + .38\ CO_2 + .07\ CH_3COOH + .07\ CO + .05\ CH_3C(O) + .05\ CO + .05\ OH$	jx(IP_CH3COCO2H)	Sander et al. (2018)*
J43019	TrGC	$CH_3COCO_3H + h\nu \rightarrow CH_3C(O) + OH + CO_2$	jx(IP_MGLYOX)+jx(ip_CH3OOH)	Sander et al. (2018)
J43020	TrGC	$CH3CHCO + h\nu \rightarrow C_2H_4 + CO$	j_ketene*0.36*2.	Sander et al. (2018)
J43021	TrGCN	$PROPOLNO3 + h\nu \rightarrow HOCH_2CHO + HCHO + HO_2 + NO_2$	j_IC3H7NO3	Sander et al. (2018)
J43022	TrGCN	$CH_3COCH_2OONO_2 + h\nu \rightarrow CH_3C(O) + HCHO + NO_3$	jx(ip_CH3O2NO2)+jx(ip_CH3COCH3)	Sander et al. (2018)
J43023	TrGJC	$C_3H_7OOH + h\nu \rightarrow C_2H_5CHO + HO_2 + OH$	jx(ip_CH3OOH)	von Kuhlmann (2001)
J43024	TrGJCN	$C_3H_7ONO_2 + h\nu \rightarrow C_2H_5CHO + NO_2 + HO_2$	0.59*j_IC3H7NO3	see note*
J43025a	TrGJC	$C_2H_5CHO + h\nu \rightarrow C_2H_5O_2 + HO_2 + CO$	jx(ip_C2H5CHO2HCO)	see note*
J43025b	TrGJC	$C_2H_5CHO + h\nu \rightarrow CH_2CHCH_2OH$	jx(ip_C2H5CHO2ENOL)	Andrews et al. (2012), Sander et al. (2018)*
J43026	TrGJCN	$PPN + h\nu \rightarrow .7\ C_2H_5CO_3 + .7\ NO_2 + .3\ C_2H_5O_2 + .3\ CO_2 + .3\ NO_3$	jx(ip_PAN)	Sander et al. (2014)
J43027	TrGJC	$C_2H_5CO_3H + h\nu \rightarrow C_2H_5O_2 + CO_2 + OH$	jx(ip_CH3OOH)	von Kuhlmann (2001)
J43028a	TrGJC	$HCOCOCH_2OOH + h\nu \rightarrow HOOCH_2CO_3 + CO + HO_2$	jx(ip_MGLYOX)	Sander et al. (2018)
J43028b	TrGJC	$HCOCOCH_2OOH + h\nu \rightarrow HCOCO + HCHO + OH$	jx(ip_HOCH2CHO)+jx(ip_CH3OOH)	Sander et al. (2018)
J43200	TrGJTerC	$HCOCH2CO3H + h\nu \rightarrow HCOCH_2O_2 + CO_2 + OH$	jx(ip_HOCH2CHO)+jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J43400	TrGJAroC	$C3DIALOOH + h\nu \rightarrow GLYOX + CO + HO_2 + OH$	jx(ip_HOCH2CHO)*2+jx(ip_CH3OOH)	Rickard and Pascoe (2009)*
J43401	TrGJAroC	$C32OH13CO + h\nu \rightarrow GLYOX + HO_2 + HO_2 + CO$	jx(ip_HOCH2CHO)*2	Rickard and Pascoe (2009)
J43402	TrGJAroC	$HCOCOHCO3H + h\nu \rightarrow GLYOX + HO_2 + CO_2 + OH$	jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J44000a	TrGJC	$LC_4H_9OOH + h\nu \rightarrow OH + C_3H_7CHO + HO_2$	jx(ip_CH3OOH)*(k_p/(k_p+k_s))	Rickard and Pascoe (2009), Sander et al. (2018)
J44000b	TrGJC	$LC_4H_9OOH + h\nu \rightarrow OH + .636\ MEK + .636\ HO_2 + .364\ CH_3CHO + .364\ C_2H_5O_2$	jx(ip_CH3OOH)*(k_s/(k_p+k_s))	Rickard and Pascoe (2009), Sander et al. (2018)
J44001	TrGJC	$MVK + h\nu \rightarrow .5\ C_3H_6 + .5\ CH_3C(O) + .5\ HCHO + CO + .5\ HO_2$	jx(ip_MVK)	Sander et al. (2014)

Table 2: Photolysis reactions (... continued)

#	labels	reaction	rate coefficient	reference
J44002	TrGJC	MEK + hν → $CH_3C(O)$ + $C_2H_5O_2$	0.42*jx(ip_CHOH)	von Kuhlmann et al. (2003)
J44003	TrGJC	LMEKOOH + hν → .62 $CH_3C(O)$ + .62 CH_3CHO + .38 HCHO + .38 CO_2 + .38 $HOCH_2CH_2O_2$ + OH	jx(ip_CH3OOH)+0.42*jx(ip_CHOH)	Sander et al. (2018)
J44004	TrGJC	BIACET + hν → 2 $CH_3C(O)$	2.15*jx(ip_MGLYOX)	see note*
J44005a	TrGJCN	LC4H9NO3 + hν → NO_2 + C_3H_7CHO + HO_2	j_IC3H7NO3*(k_p/(k_p+k_s))	see note*
J44005b	TrGJCN	LC4H9NO3 + hν → NO_2 + MEK + HO_2	j_IC3H7NO3*(k_s/(k_p+k_s))	see note*
J44006	TrGJCN	MPAN + hν → .7 MACO3 + .7 NO_2 + .3 MACO2 + .3 NO_3	jx(ip_PAN)	see note*
J44007a	TrGJC	CO2H3CO3H + hν → MGLYOX + HO_2 + OH + CO_2	jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J44007b	TrGJC	CO2H3CO3H + hν → $CH_3C(O)$ + HO_2 + $HCOCO_3H$	j_ACETOL	Rickard and Pascoe (2009)
J44008	TrGJC	MACR + hν → .5 MACO3 + .5 $CH_3C(O)$ + .5 HCHO + .5 CO + HO_2	jx(ip_MACR)	Sander et al. (2014)
J44009	TrGJC	MACROOH + hν → MACRO + OH	jx(ip_CH3OOH)+2.77*jx(ip_HOCH2CHO)	Sander et al. (2018)*
J44010	TrGJC	MACROH + hν → CH_3COCH_2OH + CO + HO_2 + HO_2	2.77*jx(ip_HOCH2CHO)	see note*
J44011	TrGJC	MACO3H + hν → MACO2 + OH	jx(ip_CH3OOH)	Sander et al. (2018)
J44012	TrGJC	LHMVKABOOH + hν → .12 MGLYOX + .12 HO_2 + .88 $CH_3C(O)$ + .88 $HOCH_2CHO$ + .12 HCHO + OH	jx(ip_CH3OOH)+j_ACETOL	Sander et al. (2018)
J44013	TrGJC	CO2H3CHO + hν → MGLYOX + CO + HO_2 + HO_2	jx(ip_HOCH2CHO)+j_ACETOL	Sander et al. (2018)
J44014	TrGJC	HO12CO3C4 + hν → $CH_3C(O)$ + $HOCH_2CHO$ + HO_2	j_ACETOL	Rickard and Pascoe (2009)
J44015	TrGJC	BIACETOH + hν → $CH_3C(O)$ + HOCH2CO	2.15*jx(ip_MGLYOX)	see note*
J44016	TrGC	$HCOCCH_3CO$ + hν → .5 OH + .5 CH_3CHO + CO + .5 CH3CHCO + .5 CO	j_KETENE	Sander et al. (2018)
J44017a	TrGC	$CH_3COCHCO$ + hν → .0192 CH_3COCO_2H + .1848 H_2O_2 + .2208 MGLYOX + .36 OH + .36 CO + .56 $CH_3C(O)$ + .2 CH_3CHO + .2 CO_2 + .2 HCHO + .2 HO_2 + CO	j_KETENE*0.5	Sander et al. (2018),Rickard and Pascoe (2009)*
J44017b	TrGC	$CH_3COCHCO$ + hν → CH3CHCO + CO	j_KETENE*0.5	Sander et al. (2018)
J44018a	TrGJC	$CH_3COCOCHO$ + hν → $CH_3C(O)$ + 2 CO + HO_2	jx(ip_MGLYOX)	Sander et al. (2018)
J44018b	TrGJC	$CH_3COCOCHO$ + hν → HCOCO + $CH_3C(O)$	2.15*jx(ip_MGLYOX)	Sander et al. (2018)
J44019	TrGJC	CH3COCOCO2H + hν → $CH_3C(O)$ + CO + CO_2 + HO_2	3.15*jx(ip_MGLYOX)	Sander et al. (2018)
J44020a	TrGJTerC	$CH_3COCOCH_2OOH$ + hν → $CH_3C(O)$ + OH + HCHO + CO	jx(ip_CH3OOH)+j_ACETOL	Rickard and Pascoe (2009)
J44020b	TrGJTerC	$CH_3COCOCH_2OOH$ + hν → $CH_3C(O)$ + HCOCO	2.15*jx(ip_MGLYOX)	Rickard and Pascoe (2009)
J44021	TrGJTerC	C44OOH + hν → HCOCH2CHO + CO_2 + HO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J44022	TrGJTerC	C413COOOH + hν → HCOCH2CO3 + HCHO + OH	jx(ip_CH3OOH)+jx(ip_HOCH2CHO)+j_ACETOL	Rickard and Pascoe (2009)

Table 2: Photolysis reactions (... continued)

#	labels	reaction	rate coefficient	reference
J44023a	TrGJTerC	C4CODIAL + hν $\rightarrow$ $HCOCOCH_2O_2$ + HO_2 + CO	`jx(ip_HOCH2CHO)`	Rickard and Pascoe (2009)
J44023b	TrGJTerC	C4CODIAL + hν $\rightarrow$ HCOCH2CO3 + HO_2 + CO	`jx(ip_MGLYOX)`	Rickard and Pascoe (2009)
J44024	TrGJTerC	C312COCO3H + hν $\rightarrow$ $HCOCOCH_2O_2$ + CO_2 + OH	`jx(ip_CH3OOH)+jx(ip_MGLYOX)`	Rickard and Pascoe (2009)
J44025	TrGJCN	LMEKNO3 + hν $\rightarrow$ .62 $CH_3C(O)$ + .62 CH_3CHO + .38 HCHO + .38 CO_2 + .38 $HOCH_2CH_2O_2$ + NO_2	`jx(ip_MEKNO3)`	Barnes et al. (1993), Sander et al. (2018)*
J44026	TrGJCN	MVKNO3 + hν $\rightarrow$ $CH_3C(O)$ + $HOCH_2CHO$ + NO_2	`jx(ip_MEKNO3)`	Barnes et al. (1993), Sander et al. (2018)*
J44027	TrGJCN	MACRNO3 + hν $\rightarrow$ CH_3COCH_2OH + CO + HO_2 + NO_2	`(2.84*j_IC3H7NO3+jx(ip_CH3CHO)) *(jx(ip_MEKNO3)+1E-10)/(j_ IC3H7NO3+0.42*jx(ip_CHOH)+1E-10)`	Müller et al. (2014), Sander et al. (2018)*
J44028	TrGJCN	TC4H9NO3 + hν $\rightarrow$ CH_3COCH_3 + CH_3 + NO_2	`2.84*j_IC3H7NO3`	Sander et al. (2018)
J44029	TrGJC	TC_4H_9OOH + hν $\rightarrow$ CH_3COCH_3 + CH_3 + OH	`jx(ip_CH3OOH)`	Sander et al. (2018)
J44030	TrGJCN	IBUTOLBNO3 + hν $\rightarrow$ CH_3COCH_3 + HCHO + HO_2 + NO_2	`2.84*j_IC3H7NO3`	Sander et al. (2018)
J44031	TrGJC	IBUTOLBOOH + hν $\rightarrow$ CH_3COCH_3 + HCHO + HO_2 + OH	`jx(ip_CH3OOH)`	Sander et al. (2018)
J44032	TrGJC	LBUT1ENOOH + hν $\rightarrow$ C_2H_5CHO + HCHO + HO_2 + OH	`jx(ip_CH3OOH)`	Sander et al. (2018)
J44033	TrGJCN	LBUT1ENNO3 + hν $\rightarrow$ C_2H_5CHO + HCHO + HO_2 + NO_2	`j_IC3H7NO3`	Sander et al. (2018)
J44034	TrGJC	BUT2OLOOH + hν $\rightarrow$ 2 CH_3CHO + HO_2 + OH	`jx(ip_CH3OOH)`	Sander et al. (2018)
J44035	TrGJCN	BUT2OLNO3 + hν $\rightarrow$ 2 CH_3CHO + HO_2 + NO_2	`j_IC3H7NO3`	Sander et al. (2018)
J44036	TrGJC	BUT2OLO + hν $\rightarrow$ $CH_3C(O)$ + HOCH2CO	`j_ACETOL`	Sander et al. (2018)
J44037a	TrGJC	C_3H_7CHO + hν $\rightarrow$ $C_3H_7O_2$ + CO + HO_2	`jx(ip_C3H7CHO2HCO)`	Sander et al. (2018)
J44037b	TrGJC	C_3H_7CHO + hν $\rightarrow$ C_2H_4 + CH_2CHOH	`jx(ip_C3H7CHO2VINY)`	Sander et al. (2018)*
J44038	TrGJC	IPRCHO + hν $\rightarrow$ $iC_3H_7O_2$ + CO + HO_2	`jx(ip_IPRCHO2HCO)`	Sander et al. (2018)
J44039	TrGJCN	IC4H9NO3 + hν $\rightarrow$ IPRCHO + NO_2	`j_IC3H7NO3`	Sander et al. (2018)
J44040	TrGJC	IC_4H_9OOH + hν $\rightarrow$ IPRCHO + HO_2 + OH	`jx(ip_CH3OOH)`	Sander et al. (2018)
J44041	TrGJC	PERIBUACID + hν $\rightarrow$ $iC_3H_7O_2$ + CO_2 + OH	`jx(ip_CH3OOH)`	Sander et al. (2018)
J44042	TrGJCN	PIPN + hν $\rightarrow$ .7 IPRCO3 + .7 NO_2 + .3 $iC_3H_7O_2$ + .3 CO_2 + .3 NO_3	`jx(ip_PAN)`	Sander et al. (2018), Sander et al. (2014)
J44043	TrGJC	HVMK + hν $\rightarrow$ MGLYOX + CO + 2 OH	`jx(ip_PeDIONE24)`	Sander et al. (2018), Nakanishi et al. (1977), Messaadia et al. (2015), Yoon et al. (1999)*
J44044	TrGJC	HMAC + hν $\rightarrow$ $HCOCCH_3CO$ + 2 OH	`jx(ip_PeDIONE24)`	Sander et al. (2018), Nakanishi et al. (1977), Messaadia et al. (2015), Yoon et al. (1999)*

Table 2: Photolysis reactions (... continued)

#	labels	reaction	rate coefficient	reference
J44045a	TrGJC	CO2C3CHO + hν → $CH_3COCH_2O_2$ + HO_2 + CO	jx(ip_C2H5CHO2HCO)	Rickard and Pascoe (2009)
J44045b	TrGJC	CO2C3CHO + hν → HVMK	jx(ip_C2H5CHO2ENOL)	Andrews et al. (2012), Sander et al. (2018)
J44046a	TrGJC	IBUTDIAL + hν → CH_3CHO + CO + HO_2 + CO_2 + H_2O	jx(ip_C2H5CHO2HCO)*2.	see note*
J44046b	TrGJC	IBUTDIAL + hν → HMAC	jx(ip_C2H5CHO2ENOL)*2.	Andrews et al. (2012), Sander et al. (2018)
J44200	TrGJTerC	IBUTALOH + hν → CH_3COCH_3 + HO_2 + HO_2 + CO	j_ACETOL	Rickard and Pascoe (2009)
J44201	TrGJTerC	IPRHOCO3H + hν → CH_3COCH_3 + HO_2 + CO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J44400a	TrGJAroC	MALDIALOOH + hν → C32OH13CO + CO + OH + HO_2	jx(ip_HOCH2CHO)*2	Rickard and Pascoe (2009)
J44400b	TrGJAroC	MALDIALOOH + hν → GLYOX + GLYOX + HO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)*
J44401	TrGJAroC	BZFUOOH + hν → CO14O3CHO + HO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)*
J44402	TrGJAroC	HOCOC4DIAL + hν → HCOCOHCO3 + HO_2 + CO	jx(ip_MGLYOX)+jx(ip_HOCH2CHO)	Rickard and Pascoe (2009)
J44403	TrGJAroCN	NBZFUOOH + hν → .5 CO14O3CHO + .5 NO_2 + .5 NBZFUONE + .5 HO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)*
J44404a	TrGJAroC	MALDALCO3H + hν → $HCOCO_3H$ + HO_2 + CO + HO_2 + CO	jx(ip_MACR)	Rickard and Pascoe (2009)
J44404b	TrGJAroC	MALDALCO3H + hν → .6 MALANHY + HO_2 + .4 GLYOX + .4 CO + .4 CO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)*
J44405	TrGJAroC	EPXDLCO2H + hν → C3DIALO2 + CO_2 + HO_2	2.77*jx(ip_HOCH2CHO)	Rickard and Pascoe (2009)
J44406	TrGJAroC	MALDIAL + hν → .4 BZFUONE + .6 MALDIALCO3 + .6 HO_2	jx(ip_NO2)*0.14	Rickard and Pascoe (2009)
J44407	TrGJAroC	MALANHYOOH + hν → HCOCOHCO3 + CO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)*
J44408	TrGJAroC	EPXDLCO3H + hν → C3DIALO2 + OH + CO_2	jx(ip_CH3OOH)+2.77*jx(ip_HOCH2CHO)	Rickard and Pascoe (2009)
J44409	TrGJAroC	CO2C4DIAL + hν → CO + CO + HO_2 + HO_2 + CO + CO	jx(ip_MGLYOX)*2	Rickard and Pascoe (2009)
J44410	TrGJAroC	MALDALCO2H + hν → $HCOCO_2H$ + HO_2 + CO + HO_2 + CO	jx(ip_MACR)	Rickard and Pascoe (2009)
J44411	TrGJAroC	EPXC4DIAL + hν → C3DIALO2 + CO + HO_2	2.77*jx(ip_HOCH2CHO)*2	Rickard and Pascoe (2009)
J44412	TrGJAroC	CO14O3CHO + hν → HO_2 + CO + $HCOCH_2O_2$ + CO_2	jx(ip_MGLYOX)	Rickard and Pascoe (2009)
J44414	TrGJAroC	MECOACEOOH + hν → $CH_3C(O)$ + HCHO + CO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)*
J45002	TrGJC	LISOPACOOH + hν → LISOPACO + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J45003	TrGJCN	LISOPACNO3 + hν → LISOPACO + NO_2	0.59*j_IC3H7NO3	see note*
J45004	TrGJC	ISOPBOOH + hν → MVK + HCHO + HO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J45005	TrGJCN	ISOPBNO3 + hν → MVK + HCHO + HO_2 + NO_2	2.84*j_IC3H7NO3	see note*

Table 2: Photolysis reactions (... continued)

#	labels	reaction	rate coefficient	reference
J45006	TrGJC	ISOPDOOH + hν $\rightarrow$ MACR + HCHO + HO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J45007	TrGJCN	ISOPDNO3 + hν $\rightarrow$ MACR + HCHO + HO_2 + NO_2	j_IC3H7NO3	see note*
J45008	TrGJCN	NISOPOOH + hν $\rightarrow$ NC4CHO + HO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J45009	TrGJCN	NC4CHO + hν $\rightarrow$ LHC4ACCO3 + NO_2	(.59*j_IC3H7NO3+jx(ip_MACR)) *(jx(ip_MEKNO3)+1E-10)/(j_IC3H7NO3+0.42*jx(ip_CHOH)+1E-10)	Müller et al. (2014), Sander et al. (2018)*
J45010	TrGJCN	LNISOOH + hν $\rightarrow$ NOA + OH + .5 HOCHCHO + .5 CO + .5 HO_2 + .5 CO_2	jx(ip_CH3OOH)	Taraborrelli et al. (2009), Sander et al. (2018)
J45011	TrGJC	LHC4ACCHO + hν $\rightarrow$ .5 LHC4ACCO3 + .5 HO_2 + .5 CO + .5 OH + .25 MACRO2 + .25 LHMVKABO2	jx(ip_MACR)	Sander et al. (2018)
J45012	TrGJC	LC578OOH + hν $\rightarrow$ .25 CH_3COCH_2OH + .75 MGLYOX + .25 HOCHCHO + .75 $HOCH_2CHO$ + .75 HO_2 + OH	jx(ip_CH3OOH)+ 2.77*jx(ip_HOCH2CHO)	Sander et al. (2018)
J45013	TrGJC	LHC4ACCO3H + hν $\rightarrow$ OH + .5 MACRO2 + .5 LHMVKABO2 + OH + CO_2	j_HPALD	Sander et al. (2018)
J45014	TrGJCN	LC5PAN1719 + hν $\rightarrow$ .7 LHC4ACCO3 + .7 NO_2 + .15 MACRO2 + .15 LHMVKABO2 + .3 CO_2 + .3 NO_3	jx(ip_PAN)	Sander et al. (2018)
J45015	TrGJC	HCOC5 + hν $\rightarrow$ .65 CH_3 + .65 CO + .65 HCHO + .35 OH + .35 $CH_3COCH_2O_2$ + HOCH2CO	0.5*jx(ip_MVK)	Sander et al. (2018)*
J45016	TrGJC	C59OOH + hν $\rightarrow$ CH_3COCH_2OH + HOCH2CO + OH	j_ACETOL+jx(ip_CH3OOH)	Sander et al. (2018)
J45017	TrGJTerC	C511OOH + hν $\rightarrow$ $CH_3C(O)$ + HCOCH2CHO + OH	jx(ip_CH3OOH)+jx(ip_HOCH2CHO)	Rickard and Pascoe (2009)
J45018a	TrGJTerC	CO23C4CHO + hν $\rightarrow$ $CH_3COCOCH_2O_2$ + HO_2 + CO	jx(ip_HOCH2CHO)	Rickard and Pascoe (2009)
J45018b	TrGJTerC	CO23C4CHO + hν $\rightarrow$ $CH_3C(O)$ + HCOCH2CO3	2.15*jx(ip_MGLYOX)	Rickard and Pascoe (2009)
J45019	TrGJTerC	CO23C4CO3H + hν $\rightarrow$ $CH_3COCOCH_2O_2$ + CO_2 + OH	jx(ip_CH3OOH)+jx(ip_HOCH2CHO)	Rickard and Pascoe (2009)
J45020	TrGJTerC	C512OOH + hν $\rightarrow$ C513O2 + OH	jx(ip_CH3OOH)+jx(ip_HOCH2CHO)	Rickard and Pascoe (2009)
J45021	TrGJTerC	CO13C4CHO + hν $\rightarrow$ CHOC3COO2 + CO + HO_2	jx(ip_HOCH2CHO)*2.	Rickard and Pascoe (2009)
J45022	TrGJTerC	C513OOH + hν $\rightarrow$ GLYOX + $HOC_2H_4CO_3$ + OH	jx(ip_CH3OOH)+jx(ip_HOCH2CHO)	Rickard and Pascoe (2009)
J45023	TrGJTerC	C513CO + hν $\rightarrow$ $HOC_2H_4CO_3$ + HO_2 + CO + CO	jx(ip_MGLYOX)+2.15*jx(ip_MGLYOX)	Rickard and Pascoe (2009)
J45024	TrGJTerC	C514OOH + hν $\rightarrow$ CO13C4CHO + HO_2 + OH	jx(ip_CH3OOH)+jx(ip_HOCH2CHO)*2.	Rickard and Pascoe (2009)
J45025	TrGJTerCN	C514NO3 + hν $\rightarrow$ CO13C4CHO + HO_2 + NO_2	j_IC3H7NO3+jx(ip_HOCH2CHO)*2.	Rickard and Pascoe (2009)
J45026a	TrGJC	LZCODC23DBCOOH + hν $\rightarrow$ OH + CO + HVMK + OH	j_HPALD*0.6*0.5	Sander et al. (2018), Jenkin et al. (2015), Peeters et al. (2014)
J45026b	TrGJC	LZCODC23DBCOOH + hν $\rightarrow$ OH + CO + $CH_3C(O)$ + $HOCH_2CHO$	j_HPALD*0.6*0.5	Sander et al. (2018), Jenkin et al. (2015), Peeters et al. (2014)

Table 2: Photolysis reactions (... continued)

#	labels	reaction	rate coefficient	reference
J45026c	TrGJC	LZCODC23DBCOOH + hν $\rightarrow$ OH + CO + HMAC + OH	j_HPALD*0.4*0.5	Sander et al. (2018), Jenkin et al. (2015), Peeters et al. (2014)
J45026d	TrGJC	LZCODC23DBCOOH + hν $\rightarrow$ OH + CO + CO + CH_3COCH_2OH + HO_2	j_HPALD*0.4*0.5	Sander et al. (2018), Jenkin et al. (2015), Peeters et al. (2014)
J45027	TrGJC	LZCO3HC23DBCOD + hν $\rightarrow$ .62 EZCH3CO2CHCHO + .38 EZCHOCCH3CHO2 + OH + CO_2	j_HPALD	Sander et al. (2018)
J45028a	TrGJC	C1OOHC2OOHC4OD + hν $\rightarrow$ $CH_3COCH_2O_2H$ + OH + 2 CO + HO_2	2.77*jx(IP_HOCH2CHO)	Sander et al. (2018)
J45028b	TrGJC	C1OOHC2OOHC4OD + hν $\rightarrow$ .5 $CH_3COCH_2O_2H$ + .5 HOCHCHO + .5 CO2H3CHO + .5 HCHO + 1.5 OH	2.*jx(IP_CH3OOH)	Sander et al. (2018)
J45029	TrGC	DB1OOH + hν $\rightarrow$ DB1O2 + OH	jx(IP_CH3OOH)	Sander et al. (2018)
J45030	TrGC	DB2OOH + hν $\rightarrow$ .48 CH_3COCH_2OH + .52 $HOCH_2CHO$ + .52 MGLYOX + .48 GLYOX + HO_2 + OH	jx(ip_CH3OOH)	Sander et al. (2018)
J45031a	TrGJC	C1ODC2OOHC4OD + hν $\rightarrow$ MGLYOX + HOCHCHO + OH	jx(ip_CH3OOH)	Sander et al. (2018)
J45031b	TrGJC	C1ODC2OOHC4OD + hν $\rightarrow$ CO2H3CHO + CO + HO_2 + OH	2.*2.77*jx(IP_HOCH2CHO)	Sander et al. (2018)
J45032	TrGJC	C4MDIAL + hν $\rightarrow$ .5 $CH_3COCHCO$ + .5 $HCOCCH_3CO$ + CO + HO_2 + OH	jx(ip_NO2)*0.1*0.5	Sander et al. (2018)*
J45033	TrGCN	DB1NO3 + hν $\rightarrow$ DB1O2 + NO_2	j_IC3H7NO3	Sander et al. (2018)
J45034	TrGJTerC	CHOC3COOOH + hν $\rightarrow$ CHOC3COO2 + CO_2 + OH	jx(ip_CH3OOH)+jx(ip_HOCH2CHO) +j_ACETOL	Rickard and Pascoe (2009)
J45200a	TrGJTerC	LMBOABOOH + hν $\rightarrow$ $HOCH_2CHO$ + CH_3COCH_3 + HO_2 + OH	jx(ip_CH3OOH)*.67	Rickard and Pascoe (2009), Sander et al. (2018)
J45200b	TrGJTerC	LMBOABOOH + hν $\rightarrow$ IBUTALOH + HCHO + HO_2 + OH	jx(ip_CH3OOH)*.33	Rickard and Pascoe (2009), Sander et al. (2018)
J45201	TrGJTerC	MBOACO + hν $\rightarrow$ HCHO + HO_2 + IPRHOCO3	j_ACETOL	Rickard and Pascoe (2009)
J45202	TrGJTerC	MBOCOCO + hν $\rightarrow$ CO + HO_2 + IPRHOCO3	jx(ip_MGLYOX)	Rickard and Pascoe (2009)
J45203a	TrGJTerCN	LNMBOABOOH + hν $\rightarrow$ NO_3CH2CHO + CH_3COCH_3 + HO_2 + OH	jx(ip_CH3OOH)*.65	Rickard and Pascoe (2009), Sander et al. (2018)

Table 2: Photolysis reactions (... continued)

#	labels	reaction	rate coefficient	reference
J45203b	TrGJTerCN	LNMBOABOOH + hν $\rightarrow$ IBUTALOH + HCHO + NO_2 + OH	jx(ip_CH3OOH)*.35	Rickard and Pascoe (2009), Sander et al. (2018)
J45204	TrGJTerCN	NC4OHCO3H + hν $\rightarrow$ IBUTALOH + CO_2 + NO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J45400	TrGJAroC	C54CO + hν $\rightarrow$ HO_2 + CO + CO + CO + $CH_3C(O)$	jx(ip_MGLYOX)+2.15*jx(ip_MGLYOX)*2	Rickard and Pascoe (2009)
J45401	TrGJAroC	C5134CO2OH + hν $\rightarrow$ $CH_3COCOCHO$ + HO_2 + CO + HO_2	jx(ip_HOCH2CHO)+2.15*jx(ip_MGLYOX)	Rickard and Pascoe (2009)
J45402	TrGJAroC	C5DIALOOH + hν $\rightarrow$ MALDIAL + CO + HO_2 + OH	jx(ip_CH3OOH)+jx(ip_MACR)	Rickard and Pascoe (2009)*
J45406	TrGJAroC	C5CO14OH + hν $\rightarrow$ $CH_3C(O)$ + $HCOCO_2H$ + HO_2 + CO	jx(ip_MVK)	Rickard and Pascoe (2009)
J45407	TrGJAroC	C5DICARB + hν $\rightarrow$ .6 C5CO14O2 + .6 HO_2 + .4 TLFUONE	jx(ip_NO2)*0.2	Rickard and Pascoe (2009)*
J45408	TrGJAroC	MC3ODBCO2H + hν $\rightarrow$ CH_3COCO_2H + HO_2 + CO + HO_2 + CO	jx(ip_MACR)	Rickard and Pascoe (2009)
J45409	TrGJAroC	ACCOMECHO + hν $\rightarrow$ MECOACETO2 + HO_2 + CO	jx(ip_HOCH2CHO)	Rickard and Pascoe (2009)
J45410	TrGJAroC	MMALNHYOOH + hν $\rightarrow$ CO2H3CO3 + CO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)*
J45411	TrGJAroC	C5DICAROOH + hν $\rightarrow$ MGLYOX + GLYOX + HO_2 + OH	jx(ip_CH3OOH)+jx(ip_HOCH2CHO)+j_ACETOL	Rickard and Pascoe (2009)*
J45412	TrGJAroCN	NTLFUOOH + hν $\rightarrow$ ACCOMECHO + NO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)*
J45414	TrGJAroC	C5CO14OOH + hν $\rightarrow$ .83 MALANHY + .83 CH_3 + .17 MGLYOX + .17 HO_2 + .17 CO + .17 CO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)*
J45415	TrGJAroC	TLFUOOH + hν $\rightarrow$ ACCOMECHO + HO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)*
J45417	TrGJAroC	ACCOMECO3H + hν $\rightarrow$ MECOACETO2 + CO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J45418	TrGJAroC	C5DIALCO + hν $\rightarrow$ MALDIALCO3 + CO + HO_2	jx(ip_MGLYOX)+jx(ip_MACR)	Rickard and Pascoe (2009)
J46200	TrGJTerCN	C614NO3 + hν $\rightarrow$ CO23C4CHO + HCHO + HO_2 + NO_2	2.15*jx(ip_MGLYOX)	Rickard and Pascoe (2009)
J46201	TrGJTerC	C614OOH + hν $\rightarrow$ CO23C4CHO + HCHO + HO_2 + OH	jx(ip_CH3OOH)+2.15*jx(ip_MGLYOX)	Rickard and Pascoe (2009)
J46202	TrGJTerC	CO235C5CHO + hν $\rightarrow$ CO23C4CO3 + CO + HO_2	jx(ip_MGLYOX)	Rickard and Pascoe (2009)
J46203	TrGJTerC	CO235C6OOH + hν $\rightarrow$ CO23C4CO3 + HCHO + OH	jx(ip_CH3OOH)+2.15*jx(ip_MGLYOX)	Rickard and Pascoe (2009)
J46400	TrGJAroC	PHENOOH + hν $\rightarrow$ .71 MALDALCO2H + .71 GLYOX + .29 PBZQONE + HO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)*
J46401	TrGJAroC	C6CO4DB + hν $\rightarrow$ C4CO2DBCO3 + HO_2 + CO	jx(ip_MGLYOX)*2	Rickard and Pascoe (2009)

Table 2: Photolysis reactions (... continued)

#	labels	reaction	rate coefficient	reference
J46402	TrGJAroC	C5CO2DCO3H + hν $\rightarrow$ $CH_3C(O)$ + HCOCOCHO + CO_2 + OH	jx(ip_CH3OOH)+jx(ip_MGLYOX)	Rickard and Pascoe (2009)
J46403	TrGJAroCN	NDNPHENOOH + hν $\rightarrow$ NC4DCO2H + HNO_3 + CO + CO + NO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)*
J46404	TrGJAroCN	BZBIPERNO3 + hν $\rightarrow$ GLYOX + HO_2 + .5 BZFUONE + .5 BZFUONE + NO_2	j_IC3H7NO3	Rickard and Pascoe (2009)*
J46405	TrGJAroCN	HOC6H4NO2 + hν $\rightarrow$ HONO + CPDKETENE	jx(ip_HOC6H4NO2)	Chen et al. (2011)*
J46406	TrGJAroC	CPDKETENE + hν $\rightarrow$ CO_2 + CO + 2 HO_2 + MALDIAL	j_KETENE	see note*
J46407	TrGJAroC	C5COOHCO3H + hν $\rightarrow$ HOCOC4DIAL + HO_2 + CO + CO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J46408	TrGJAroC	BZEPOXMUC + hν $\rightarrow$ .5 C5DIALO2 + 1.5 HO_2 + 1.5 CO + .5 MALDIAL	4.E3*jx(ip_MVK)*0.1	Rickard and Pascoe (2009)
J46409	TrGJAroCN	NPHEN1OOH + hν $\rightarrow$ NPHEN1O + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J46410	TrGJAroC	BZEMUCCO + hν $\rightarrow$ HCOCOHCO3 + C3DIALO2	jx(ip_HOCH2CHO)*2+j_ACETOL	Rickard and Pascoe (2009)
J46411	TrGJAroC	BZEMUCCO2H + hν $\rightarrow$ C5DIALO2 + CO_2 + HO_2	jx(ip_MACR)	Rickard and Pascoe (2009)
J46412	TrGJAroCN	NNCATECOOH + hν $\rightarrow$ NC4DCO2H + $HCOCO_2H$ + NO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)*
J46413	TrGJAroC	C615CO2OOH + hν $\rightarrow$ C5DICARB + CO + HO_2 + OH	jx(ip_MVK)+jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J46414	TrGJAroCN	NPHENOOH + hν $\rightarrow$ MALDALCO2H + GLYOX + OH + NO_2	j_IC3H7NO3 + jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J46415	TrGJAroCN	NCATECOOH + hν $\rightarrow$ NC4DCO2H + $HCOCO_2H$ + HO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)*
J46416	TrGJAroC	PBZQOOH + hν $\rightarrow$ C5CO2OHCO3 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)*
J46417	TrGJAroC	BZOBIPEROH + hν $\rightarrow$ MALDIALCO3 + GLYOX + HO_2	j_ACETOL	Rickard and Pascoe (2009)
J46418	TrGJAroC	BZBIPEROOH + hν $\rightarrow$ GLYOX + HO_2 + .5 BZFUONE + .5 BZFUONE + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)*
J46419	TrGJAroCN	NBZQOOH + hν $\rightarrow$ C6CO4DB + NO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)*
J46420	TrGJAroC	CATEC1OOH + hν $\rightarrow$ CATEC1O + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J46421	TrGJAroC	C6125CO + hν $\rightarrow$ C5CO14O2 + CO + HO_2	jx(ip_MGLYOX)+jx(ip_MVK)	Rickard and Pascoe (2009)
J46422	TrGJAroCN	DNPHENOOH + hν $\rightarrow$ NC4DCO2H + $HCOCO_2H$ + NO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)*
J46423	TrGJAroC	BZEMUCCO3H + hν $\rightarrow$ C5DIALO2 + CO_2 + OH	jx(ip_CH3OOH)+jx(ip_MACR)	Rickard and Pascoe (2009)
J46424	TrGJAroC	C6H5OOH + hν $\rightarrow$ C6H5O + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J46425	TrGJAroC	BZEMUCOOH + hν $\rightarrow$ .5 EPXC4DIAL + .5 GLYOX + .5 HO_2 + .5 C3DIALO2 + .5 C32OH13CO + OH	jx(ip_CH3OOH)+jx(ip_HOCH2CHO)*2	Rickard and Pascoe (2009)*

Table 2: Photolysis reactions (... continued)

#	labels	reaction	rate coefficient	reference
J46427	TrGJAroCN	BZEMUCNO3 + hν → EPXC4DIAL + NO_2 + GLYOX + HO_2	2.77*jx(ip_HOCH2CHO)	Rickard and Pascoe (2009)
J46428	TrGJAroCN	DNPHEN + hν → HONO + NCPDKETENE	jx(ip_HOC6H4NO2)	Sander et al. (2018)
J46429	TrGJAroCN	NCPDKETENE + hν → CO_2 + CO + 2 HO_2 + NC4DCO2H	j_KETENE	see note*
J47200	TrGJTerC	CO235C6CHO + hν → CHOC3COCO3 + $CH_3C(O)$	2.15*jx(ip_MGLYOX)	Rickard and Pascoe (2009)
J47201	TrGJTerC	C235C6CO3H + hν → CO235C6O2 + CO_2 + OH	jx(ip_CH3OOH)+2.15*jx(ip_MGLYOX)	Rickard and Pascoe (2009)
J47202	TrGJTerC	C716OOH + hν → CO13C4CHO + $CH_3C(O)$ + OH	jx(ip_CH3OOH)+jx(ip_HOCH2CHO)	Rickard and Pascoe (2009)
J47203	TrGJTerC	C721OOH + hν → C722O2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J47204	TrGJTerC	C722OOH + hν → CH_3COCH_3 + C44O2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J47400	TrGJAroC	TLEPOXMUC + hν → .5 C615CO2O2 + HO_2 + CO + .5 EPXC4DIAL + .5 $CH_3C(O)$	4.E3*jx(ip_MVK)*0.1	Rickard and Pascoe (2009)
J47401	TrGJAroC	C6H5CH2OOH + hν → BENZAL + HO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)*
J47402	TrGJAroCN	C6H5CH2NO3 + hν → BENZAL + HO_2 + NO_2	0.59*j_IC3H7NO3	Rickard and Pascoe (2009)*
J47403	TrGJAroC	BENZAL + hν → HO_2 + CO + C6H5O2	jx(ip_BENZAL)	Wallington et al. (2018)
J47404	TrGJAroC	TLBIPEROOH + hν → .6 GLYOX + .4 MGLYOX + HO_2 + .2 C4MDIAL + .2 C5DICARB + .2 TLFUONE + .2 BZFUONE + .2 MALDIAL + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)*
J47405	TrGJAroCN	TLBIPERNO3 + hν → .6 GLYOX + .4 MGLYOX + HO_2 + .2 C4MDIAL + .2 C5DICARB + .2 TLFUONE + .2 BZFUONE + .2 MALDIAL + NO_2	j_IC3H7NO3	Rickard and Pascoe (2009)*
J47406	TrGJAroC	TLOBIPEROH + hν → C5CO14O2 + GLYOX + HO_2	j_ACETOL	Rickard and Pascoe (2009)
J47407	TrGJAroC	CRESOOH + hν → .68 C5CO14OH + .68 GLYOX + HO_2 + .32 PTLQONE + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)*
J47408a	TrGJAroCN	NCRESOOH + hν → .68 C5CO14OH + .68 GLYOX + HO_2 + .32 PTLQONE + OH + NO_2	j_IC3H7NO3	Rickard and Pascoe (2009)*
J47408b	TrGJAroCN	NCRESOOH + hν → C5CO14OH + GLYOX + NO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)*
J47409	TrGJAroCN	TOL1OHNO2 + hν → HONO + MCPDKETENE	jx(ip_HOPh3Me2NO2)	see note*
J47410	TrGJAroC	TLEMUCCO2H + hν → C615CO2O2 + CO_2 + HO_2	jx(ip_MACR)	Rickard and Pascoe (2009)
J47411	TrGJAroC	TLEMUCCO3H + hν → C615CO2O2 + CO_2 + OH	jx(ip_CH3OOH)+jx(ip_MACR)	Rickard and Pascoe (2009)
J47412	TrGJAroC	TLEMUCOOH + hν → .5 C3DIALO2 + .5 CO2H3CHO + .5 EPXC4DIAL + .5 MGLYOX + .5 HO_2 + OH	jx(ip_CH3OOH)+2.77*jx(ip_HOCH2CHO)+j_ACETOL	Rickard and Pascoe (2009)*
J47413	TrGJAroCN	TLEMUCNO3 + hν → EPXC4DIAL + NO_2 + $CH_3C(O)$ + CO + HO_2	2.77*jx(ip_HOCH2CHO)+j_ACETOL	Rickard and Pascoe (2009)

Table 2: Photolysis reactions (... continued)

#	labels	reaction	rate coefficient	reference
J47414	TrGJAroC	TLEMUCCO + hν $\rightarrow$ $CH_3C(O)$ + EPXC4DIAL + CO + HO_2	`2.77*jx(ip_HOCH2CHO)+2.15*jx(ip_MGLYOX)`	Rickard and Pascoe (2009)
J47415	TrGJAroC	C6H5CO3H + hν $\rightarrow$ C6H5O2 + CO_2 + OH	`jx(ip_CH3OOH)`	Rickard and Pascoe (2009)
J47416	TrGJAroC	OXYL1OOH + hν $\rightarrow$ TOL1O + OH	`jx(ip_CH3OOH)`	Rickard and Pascoe (2009)
J47417	TrGJAroCN	MNCATECH + hν $\rightarrow$ HONO + MCPDKETENE	`jx(ip_HOPh3Me2NO2)`	see note*
J47418	TrGJAroC	MCPDKETENE + hν $\rightarrow$ CO_2 + CO + 2 HO_2 + C4MDIAL	`j_KETENE`	see note*
J47419	TrGJAroCN	DNCRES + hν $\rightarrow$ HONO + MNCPDKETENE	`jx(ip_HOPh3Me2NO2)`	see note*
J47420	TrGJAroCN	MNCPDKETENE + hν $\rightarrow$ CO_2 + CO + 2 HO_2 + NC4MDCO2HN	`j_KETENE`	see note*
J47421	TrGJAroC	MCATEC1OOH + hν $\rightarrow$ MCATEC1O + OH	`jx(ip_CH3OOH)`	Rickard and Pascoe (2009)
J47422	TrGJAroCN	NPTLQOOH + hν $\rightarrow$ C7CO4DB + NO_2 + OH	`jx(ip_CH3OOH)`	Rickard and Pascoe (2009)*
J47423	TrGJAroC	PTLQOOH + hν $\rightarrow$ C6CO2OHCO3 + OH	`jx(ip_CH3OOH)`	Rickard and Pascoe (2009)*
J47424	TrGJAroCN	NCRES1OOH + hν $\rightarrow$ NCRES1O + OH	`jx(ip_CH3OOH)`	Rickard and Pascoe (2009)
J47425	TrGJAroCN	MNNCATCOOH + hν $\rightarrow$ NC4MDCO2HN + $HCOCO_2H$ + NO_2 + OH	`jx(ip_CH3OOH)`	Rickard and Pascoe (2009)*
J47426	TrGJAroCN	MNCATECOOH + hν $\rightarrow$ NC4MDCO2HN + $HCOCO_2H$ + HO_2 + OH	`jx(ip_CH3OOH)`	Rickard and Pascoe (2009)*
J47427	TrGJAroC	C7CO4DB + hν $\rightarrow$ C5CO2DBCO3 + HO_2 + CO	`jx(ip_MGLYOX)*2`	Rickard and Pascoe (2009)
J47428	TrGJAroCN	NDNCRESOOH + hν $\rightarrow$ NC4MDCO2HN + HNO_3 + CO + CO + NO_2 + OH	`jx(ip_CH3OOH)`	Rickard and Pascoe (2009)*
J47429	TrGJAroCN	DNCRESOOH + hν $\rightarrow$ NC4MDCO2HN + $HCOCO_2H$ + NO_2 + OH	`jx(ip_CH3OOH)`	Rickard and Pascoe (2009)*
J47430	TrGJAroC	C6COOHCO3H + hν $\rightarrow$ C5134CO2OH + HO_2 + CO + CO_2 + OH	`jx(ip_CH3OOH)`	Rickard and Pascoe (2009)
J48200	TrGJTerC	C86OOH + hν $\rightarrow$ C511O2 + CH_3COCH_3 + OH	`jx(ip_CH3OOH)+ jx(ip_HOCH2CHO)`	Rickard and Pascoe (2009)
J48201	TrGJTerC	C812OOH + hν $\rightarrow$ C813O2 + OH	`jx(ip_CH3OOH)`	Rickard and Pascoe (2009)
J48202	TrGJTerC	C813OOH + hν $\rightarrow$ CH_3COCH_3 + C512O2 + OH	`jx(ip_CH3OOH)+jx(ip_MGLYOX)`	Rickard and Pascoe (2009)
J48203	TrGJTerC	C721CHO + hν $\rightarrow$ C721O2 + CO + HO_2	`jx(ip_HOCH2CHO)`	Rickard and Pascoe (2009)
J48204	TrGJTerC	C721CO3H + hν $\rightarrow$ C721O2 + CO_2 + OH	`jx(ip_CH3OOH)`	Rickard and Pascoe (2009)
J48205	TrGJTerC	C8BCOOH + hν $\rightarrow$ C89O2 + OH	`jx(ip_CH3OOH)`	Rickard and Pascoe (2009)
J48206	TrGJTerC	C89OOH + hν $\rightarrow$ C810O2 + OH	`jx(ip_CH3OOH)+jx(ip_HOCH2CHO)`	Rickard and Pascoe (2009)
J48207	TrGJTerCN	C89NO3 + hν $\rightarrow$ C810O2 + NO_2	`jx(ip_CH3OOH)+jx(ip_HOCH2CHO)`	Rickard and Pascoe (2009)
J48208	TrGJTerC	C810OOH + hν $\rightarrow$ CH_3COCH_3 + C514O2 + OH	`jx(ip_CH3OOH)+jx(ip_HOCH2CHO)`	Rickard and Pascoe (2009)

Table 2: Photolysis reactions (... continued)

#	labels	reaction	rate coefficient	reference
J48209	TrGJTerCN	C810NO3 + hν $\rightarrow$ CH_3COCH_3 + C514O2 + NO_2	2.84*j_IC3H7NO3+jx(ip_HOCH2CHO)	Rickard and Pascoe (2009)
J48210	TrGJTerCN	C8BCNO3 + hν $\rightarrow$ C89O2 + NO_2	j_IC3H7NO3	Rickard and Pascoe (2009)
J48211	TrGJTerC	C85OOH + hν $\rightarrow$ C86O2 + OH	jx(ip_CH3OOH)+j_ACETOL	Rickard and Pascoe (2009)
J48400	TrGJAroC	STYRENOOH + hν $\rightarrow$ HO_2 + HCHO + BENZAL + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)*
J49200	TrGJTerC	C96OOH + hν $\rightarrow$ C97O2 + OH	jx(ip_CH3OOH)+j_ACETOL	Rickard and Pascoe (2009)
J49201	TrGJTerC	C97OOH + hν $\rightarrow$ C98O2 + OH	jx(ip_CH3OOH)+j_ACETOL	Rickard and Pascoe (2009)
J49202	TrGJTerC	C98OOH + hν $\rightarrow$ C614O2 + CH_3COCH_3 + OH	(jx(ip_CH3OOH)+2.15*jx(ip_MGLYOX))	Rickard and Pascoe (2009)
J49203a	TrGJTerC	NORPINAL + hν $\rightarrow$ C85O2 + CO + HO_2	jx(ip_PINAL2HCO)	Rickard and Pascoe (2009), Sander et al. (2018)
J49203b	TrGJTerC	NORPINAL + hν $\rightarrow$ NORPINENOL	jx(ip_PINAL2ENOL)	Sander et al. (2018), Andrews et al. (2012)
J49204	TrGJTerC	C85CO3H + hν $\rightarrow$ C85O2 + CO_2 + OH	jx(ip_CH3OOH)+j_ACETOL	Rickard and Pascoe (2009)
J49205	TrGJTerC	C89CO2H + hν $\rightarrow$ .8 C811CO3 + .2 C89O2 + .2 CO_2 + HO_2	jx(ip_HOCH2CHO)	Rickard and Pascoe (2009)
J49206	TrGJTerC	C89CO3H + hν $\rightarrow$ .8 C811CO3 + .2 C89O2 + .2 CO_2 + OH	jx(ip_CH3OOH)+jx(ip_HOCH2CHO)	Rickard and Pascoe (2009)
J49207	TrGJTerC	C811CO3H + hν $\rightarrow$ C811O2 + CO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J49208	TrGJTerC	NOPINDOOH + hν $\rightarrow$ C89CO3 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J40200	TrGJTerC	LAPINABOOH + hν $\rightarrow$ PINAL + HO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J40201	TrGJTerC	MENTHEN6ONE + hν $\rightarrow$ RO6R1O2 + OH	jx(ip_CH3OOH)	Vereecken et al. (2007)
J40202	TrGJTerC	2OHMENTHEN6ONE + hν $\rightarrow$ 10 LCARBON + OH	jx(ip_CH3OOH)	Vereecken et al. (2007)
J40203a	TrGJTerC	PINAL + hν $\rightarrow$ C96O2 + CO + HO_2	jx(ip_PINAL2HCO)	Rickard and Pascoe (2009)
J40203b	TrGJTerC	PINAL + hν $\rightarrow$ PINEOL	jx(ip_PINAL2ENOL)	Sander et al. (2018), Andrews et al. (2012)*
J40204	TrGJTerC	PERPINONIC + hν $\rightarrow$ C96O2 + CO_2 + OH	jx(ip_CH3OOH)+j_ACETOL	Rickard and Pascoe (2009)
J40205	TrGJTerC	PINALOOH + hν $\rightarrow$ C106O2 + OH	jx(ip_CH3OOH)+jx(ip_HOCH2CHO)	Rickard and Pascoe (2009)
J40206	TrGJTerCN	PINALNO3 + hν $\rightarrow$ C106O2 + NO_2	j_IC3H7NO3+jx(ip_HOCH2CHO)	Rickard and Pascoe (2009)
J40207	TrGJTerC	C106OOH + hν $\rightarrow$ C716O2 + CH_3COCH_3 + OH	jx(ip_CH3OOH)+jx(ip_HOCH2CHO)	Rickard and Pascoe (2009)
J40208	TrGJTerCN	C106NO3 + hν $\rightarrow$ C716O2 + CH_3COCH_3 + NO_2	j_IC3H7NO3+ jx(ip_HOCH2CHO)	Rickard and Pascoe (2009)
J40209	TrGJTerC	C109OOH + hν $\rightarrow$ C89CO3 + HCHO + OH	jx(ip_CH3OOH)+jx(ip_HOCH2CHO)	Rickard and Pascoe (2009)
J40210	TrGJTerC	C109CO + hν $\rightarrow$ C89CO3 + CO + HO_2	jx(ip_MGLYOX)+jx(ip_HOCH2CHO)	Rickard and Pascoe (2009)
J40211	TrGJTerCN	LNAPINABOOH + hν $\rightarrow$ PINAL + NO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J40212	TrGJTerC	BPINAOOH + hν $\rightarrow$ NOPINONE + HCHO + HO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)
J40213	TrGJTerCN	LNBPINABOOH + hν $\rightarrow$ NOPINONE + HCHO + NO_2 + OH	jx(ip_CH3OOH)	Rickard and Pascoe (2009)

Table 2: Photolysis reactions (... continued)

#	labels	reaction	rate coefficient	reference
J40214	TrGJTerCN	ROO6R1NO3 + hν $\rightarrow$ ROO6R3O2 + CH_3COCH_3 + NO_2	2.84*j_IC3H7NO3+jx(ip_CH3OOH)	Sander et al. (2018)
J40215	TrGJTerCN	RO6R1NO3 + hν $\rightarrow$ 9 LCARBON + HCHO + HO_2 + NO_2	2.84*j_IC3H7NO3	Sander et al. (2018)
J6000	StTrGJCl	Cl_2 + hν $\rightarrow$ Cl + Cl	jx(ip_Cl2)	Sander et al. (2014)
J6100	StTrGJCl	Cl_2O_2 + hν $\rightarrow$ 2 Cl	jx(ip_Cl2O2)	Sander et al. (2014)
J6101	StTrGJCl	OClO + hν $\rightarrow$ ClO + O(^{3}P)	jx(ip_OClO)	Sander et al. (2014)
J6201	StTrGJCl	HOCl + hν $\rightarrow$ OH + Cl	jx(ip_HOCl)	Sander et al. (2014)
J6300	TrGJClN	$ClNO_2$ + hν $\rightarrow$ Cl + NO_2	jx(ip_ClNO2)	Sander et al. (2014)
J6301a	StTrGJClN	$ClNO_3$ + hν $\rightarrow$ Cl + NO_3	jx(ip_ClNO3)	Sander et al. (2014)
J6301b	StTrGJClN	$ClNO_3$ + hν $\rightarrow$ ClO + NO_2	jx(ip_ClONO2)	Sander et al. (2014)
J8000	TrGJI	I_2 + hν $\rightarrow$ I + I	jx(ip_I2)	Sander et al. (2014)
J8100	TrGJI	IO + hν $\rightarrow$ I + O(^{3}P)	jx(ip_IO)	Sander et al. (2014)
J8200	TrGJI	HOI + hν $\rightarrow$ I + OH	jx(ip_HOI)	Sander et al. (2014)
J8300	TrGJIN	INO_2 + hν $\rightarrow$ I + NO_2	jx(ip_INO2)	Sander et al. (2014)
J8301	TrGJIN	INO_3 + hν $\rightarrow$ I + NO_3	jx(ip_INO3)	Sander et al. (2014)
J8400	TrGJI	CH_2I_2 + hν $\rightarrow$ 2 I + 2 HO_2 + CO	jx(ip_CH2I2)	Sander et al. (2014)
J8401	TrGJI	CH_3I + hν $\rightarrow$ I + CH_3	jx(ip_CH3I)	Sander et al. (2014)
J8402	TrGJCI	CH_3CHICH_3 + hν $\rightarrow$ 2 LCARBON + I + CH_3	jx(ip_C3H7I)	Sander et al. (2014)
J8403	TrGJClI	CH_2ClI + hν $\rightarrow$ I + Cl + 2 HO_2 + CO	jx(ip_CH2ClI)	Sander et al. (2014)
J8600	TrGJClI	ICl + hν $\rightarrow$ I + Cl	jx(ip_ICl)	Sander et al. (2014)
J4200MK	TrGAJCN	H2NCH2CHO + hν $\rightarrow$ MMAO2 + HCHO	jx(ip_CH3CHO)	Karl et al. (2012a)
J4201MK	TrGAJCN	H2NCOCHO + hν $\rightarrow$ H2NCHO + HCHO	jx(ip_CH3CHO)	Karl et al. (2012a)
J4202MK	TrGAJCN	MEANNO + hν $\rightarrow$ MEAN + NO	0.33*jx(ip_NO2)	Karl et al. (2012a)
J4203MK	TrGAJCN	NDMA + hν $\rightarrow$ CH3NCH3 + NO	0.25*jx(ip_NO2)	Karl et al. (2012a)
J4300MK	TrGJCN	IPN + hν $\rightarrow$ CH3CHOCH3 + NO	0.08*jx(ip_NO2)	Karl (2012)
J4400MK	TrGAJCN	NDELA + hν $\rightarrow$ HOETNETOH + NO	0.33*jx(ip_NO2)	Karl et al. (2012a)
J4401MK	TrGAJCN	NAMP + hν $\rightarrow$ AMPN + NO	0.34*jx(ip_NO2)	Nielsen and Schade (2012)
J4402MK	TrGAJCN	DMOCNH2MOH + hν $\rightarrow$ CH3CNH2MOH + HCHO	jx(ip_CH3CHO)	Karl (2012)
J4403MK	TrGAJCN	DMCNH2CHO + hν $\rightarrow$ DMCNH2 + HCHO	jx(ip_CH3CHO)	Karl (2012)
J4600MK	TrGJC	CHEXONE + hν $\rightarrow$ C_2H_4 + C_3H_6 + CO	0.42*jx(ip_CHOH)	von Kuhlmann et al. (2003)
PH (aqueous)				
PH11000_a01	TrAa01JFe	$FeOH^{2+}$(aq) + hν $\rightarrow$ Fe^{2+}(aq) + OH(aq)	xaer(01)*4.51E-3*0.312	Herrmann et al. (2000)
PH11001_a01	TrAa01JFe	$Fe(OH)_2^+$(aq) + hν $\rightarrow$ Fe^{2+}(aq) + OH(aq) + OH^-(aq)	xaer(01)*5.77E-3*0.255	Herrmann et al. (2000)
PH11003_a01	TrAa01JFeS	$FeSO_4^+$(aq) + hν $\rightarrow$ Fe^{2+}(aq) + SO_4^-(aq)	xaer(01)*6.43E-3*7.9E-3	Herrmann et al. (2000)

Table 2: Photolysis reactions (... continued)

#	labels	reaction	rate coefficient	reference
PH2100_a01	TrAa01J	$H_2O_2(aq) + h\nu \rightarrow 2\ OH(aq)$	xaer(01)*jx(ip_H2O2) *7.11E-1	Ervens et al. (2003)
PH3200_a01	TrAa01JN	$NO_3^-(aq) + h\nu \rightarrow NO_2(aq) + OH(aq) + OH^-(aq)$	xaer(01)*jx(ip_NO2O) *1.91E-6	Ervens et al. (2003)
PH3201_a01	TrAa01JN	$NO_2^-(aq) + h\nu \rightarrow NO(aq) + OH^-(aq)$	xaer(01)*jx(ip_NO2) *2.50E-3	Ervens et al. (2003)
PH4101_a01	TrAa01AmiJN	$MMNNO2(aq) + h\nu \rightarrow NO_2(aq) + CH_3NH_2^+(aq) + OH^-(aq)$	xaer(01)*jx(ip_ NO2) *5.3E-3 *xaqcnno	Karl et al. (2012b)
PH4102_a01	TrAa01AmiJCN	$MEANNO2(aq) + h\nu \rightarrow NO_2(aq) + HOCH_2CH_2NH_2^+(aq) + OH^-(aq)$	xaer(01)*jx(ip_ NO2) *5.3E-3 *xaqcnno	Karl et al. (2012b)
PH4201_a01	TrAa01AmiJCN	$NDMA(aq) + h\nu \rightarrow NO(aq) + (CH_3)_2NH^+(aq) + OH^-(aq)$	xaer(01)*jx(ip_ NO2) *1.30E-1 *xaqcnno	Karl et al. (2012b)
PH4202_a01	TrAa01AmiJCN	$DMNNO2(aq) + h\nu \rightarrow NO_2(aq) + (CH_3)_2NH^+(aq) + OH^-(aq)$	xaer(01)*jx(ip_ NO2) *5.3E-3 *xaqcnno	Karl et al. (2012b)
PH4203_a01	TrAa01AmiJCN	$MEANNO(aq) + h\nu \rightarrow NO(aq) + HOCH_2CH_2NH_2^+(aq) + OH^-(aq)$	xaer(01)*jx(ip_ NO2) *1.30E-1 *xaqcnno	Karl et al. (2012b)
PH4401_a01	TrAa01AmiJCN	$NDELA(aq) + h\nu \rightarrow NO(aq) + (HOET)_2NH^+(aq) + OH^-(aq)$	xaer(01)*jx(ip_ NO2) *5.3E-2 *xaqcnno	Karl et al. (2012b)
PH4402_a01	TrAa01AmiJCN	$DEANNO2(aq) + h\nu \rightarrow NO_2(aq) + (HOET)_2NH^+(aq) + OH^-(aq)$	xaer(01)*jx(ip_ NO2) *5.3E-3 *xaqcnno	Karl et al. (2012b)

General notes

j-values are calculated with an external module (e.g., JVAL) and then supplied to the MECCA chemistry.

Values that originate from the Master Chemical Mechanism (MCM) by Rickard and Pascoe (2009) are translated according in the following way:
j(11) → jx(ip_COH2)
j(12) → jx(ip_CHOH)
j(15) → jx(ip_HOCH2CHO)
j(18) → jx(ip_MACR)
j(22) → jx(ip_ACETOL)
j(23)+j(24) → jx(ip_MVK)
j(31)+j(32)+j(33) → jx(ip_GLYOX)
j(34) → jx(ip_MGLYOX)
j(41) → jx(ip_CH3OOH)
j(53) → j(isopropyl nitrate)
j(54) → j(isopropyl nitrate)
j(55) → j(isopropyl nitrate)
j(56)+j(57) → jx(ip_NOA)

Specific notes

J41006: product distribution as for HNO4

J42004: Quantum yields from Burkholder et al. (2015).

J42005a: Quantum yields from Burkholder et al. (2015).

J42005b: Quantum yields from Burkholder et al. (2015).

J42005c: Quantum yields from Burkholder et al. (2015).

J42007: It is assumed that J(PHAN) is the same as J(PAN).

J42017: Enhancement of j according to Müller et al. (2014).

J42020: It is assumed that j(NO_3CH2CHO) is the same as j(PAN).

J42021: In analogy to what is assumed for $CH_3O_2NO_2$ photolysis as in (Sander et al., 2014).

J43002: Following von Kuhlmann et al. (2003), we use j(CH_3COCH_2OH) = 0.11*jx(ip_CHOH). As an additional factor, the quantum yield of 0.65 is taken from Orlando et al. (1999a).

J43006: Following von Kuhlmann et al. (2003), we use J($iC_3H_7ONO_2$) = 3.7*jx(ip_PAN).

J43018: One third of the acetaldehyde channel is considered to be CH2CHOH according to Hjorth (2002) EUPHORE Report.

J43024: Assuming J($C_3H_7ONO_2$) = 0.59 × J($iC_3H_7ONO_2$), consistent with the photolysis rate coefficients used in the MCM (Rickard and Pascoe, 2009).

J43025a: Photolysis frequencies very similar to the ones of CH_3CHO.

J43025b: Photolysis frequencies very similar to the ones of CH_3CHO.

J43400: KDEC C3DIALO → GLYOX + CO + HO2

J44004: It is assumed that J(BIACET) is 2.15 times larger than J(MGLYOX), consistent with the photolysis rate coefficients used in the MCM (Rickard and Pascoe, 2009).

J44005a: It is assumed that J(LC4H9NO3) is the same as J($iC_3H_7ONO_2$).

J44005b: It is assumed that J(LC4H9NO3) is the same as J($iC_3H_7ONO_2$).

J44006: It is assumed that J(MPAN) is the same as J(PAN).

J44009: It is assumed that J(MACROOH) is 2.77 times larger than J($HOCH_2CHO$), consistent with the photolysis rate coefficients used in the MCM (Rickard and Pascoe, 2009).

J44010: It is assumed that J(MACROH) is 2.77 times larger than J($HOCH_2CHO$), consistent with the photolysis rate coefficients used in the MCM (Rickard and Pascoe, 2009).

J44015: It is assumed that J(BIACETOH) is 2.15 times larger than J(MGLYOX), consistent with the photolysis rate coefficients used in the MCM (Rickard and Pascoe, 2009).

J44017a: CO-channel yielding CH_3COCH which upon reaction with O_2 produces an excited Criegee Intermediate assumed to be similar to MGLOOA in MCM. MGLOOA is produced also in other reactions and is substituted by its decomposition products. Furthermore, the stabilized Criegge Intermediate is assumed to solely react with water.

J44025: J values only for the secondary nitrate.

J44026: Like for LMEKNO3 photolysis

J44027: 2.84*J_IC3H7NO3 like for other tertiary alkyl nitrates (see J4505). Enhancement of J according to Müller et al. (2014).

J44037b: Channel which produces just vinyl alcohol and not a larger enol via keto-enol phototautomerization.

J44043: The resulting vinyl peroxy radical is assumed to mostly form with HO_2 a labile hydroperoxide (see ketene formation). The products are further simplified.

J44044: 1,5-H-shift for the resulting vinyl peroxy radical assumed to be dominant.

J44046a: Simplified oxidation.

J44400b: KDEC MALDIALO → GLYOX + GLYOX + HO2

J44401: KDEC BZFUO → CO14O3CHO + HO2

J44403: KDEC NBZFUO → 0.5 CO14O3CHO + 0.5 NO2 + 0.5 NBZFUONE + 0.5 HO2

J44404b: KDEC MALDIALCO2 → 0.6 MALANHY + HO2 + 0.4 GLYOX + 0.4 CO

J44407: KDEC MALANHYO → HCOCOHCO3

J44414: KDEC MECOACETO → CH3CO3 + HCHO

J45003: It is assumed that J(LISOPACNO3) = 0.59 × J($iC_3H_7ONO_2$), consistent with the photolysis rate coefficients used in the MCM (Rickard and Pascoe, 2009).

J45005: It is assumed that J(ISOPBNO3) = 2.84 × J($iC_3H_7ONO_2$), consistent with the photolysis rate coefficients used in the MCM (Rickard and Pascoe, 2009).

J45007: It is assumed that J(ISOPDNO3) is the same as J($iC_3H_7ONO_2$).

J45009: 0.59*J_IC3H7NO3 like for other primary alkyl nitrates (see J4503). Enhancement of J according to Müller et al. (2014).

J45015: Consistent with the MCM (Rickard and Pascoe, 2009), we assume that J(HCOC5) is half as large as J(MVK). With exeption of HOCH2CO the products of MACO2 decomposition without CO_2.

J45032: approximation with 4-oxo-pentenal photolysis combining results of Thüner et al(2004) and Xiang et al(2007)

J45402: KDEC C5DIALO → MALDIAL + CO + HO2

J45407: KDEC TLFUONE → 0.6 C5CO14O2 + 0.6 HO2 + 0.4 TLFUONE

J45410: KDEC MMALANHYO → CO2H3CO3

J45411: KDEC C5DICARBO → MGLYOX + GLYOX + HO2

J45412: KDEC NTLFUO → ACCOMECHO + NO2

J45414: KDEC C5CO14CO2 → 0.83 MALANHY + 0.83 CH3 + .17 MGLYOX + .17 HO2 + .17 CO + .17 CO2

J45415: KDEC TLFUO → ACCOMECHO + HO2

J46400: KDEC PHENO → 0.71 MALDALCO2H + 0.71 GLYOX + 0.29 PBZQONE + HO2

J46403: KDEC NDNPHENO → NC4DCO2H + HNO3 + CO + CO + NO2

J46404: KDEC BZBIPERO → GLYOX + HO2 + 0.5 BZFUONE + 0.5 BZFUONE

J46405: new channel created for nitrophenol decomposition

J46406: new channel created for nitrophenol decomposition

J46412: KDEC NNCATECO → NC4DCO2H + HCOCO2H + NO2

J46415: KDEC NCATECO → NC4DCO2H + HCOCO2H + HO2

J46416: KDEC PBZQO → C5CO2OHCO3

J46418: KDEC BZBIPERO → GLYOX + HO2 + 0.5 BZFUONE + 0.5 BZFUONE

J46419: KDEC NBZQO → C6CO4DB + NO2

J46422: KDEC DNPHENO → NC4DCO2H + HCOCO2H + NO2

J46425: KDEC BZEMUCO → 0.5 EPXC4DIAL + .5 GLYOX + .5 HO2 + .5 C3DIALO2 + .5 C32OH13CO

J46429: new channel

J47401: KROPRIM*O2 fast reaction C6H5CH2O = BENZAL + HO2

J47402: KROPRIM*O2 fast reaction C6H5CH2O = BENZAL + HO2

J47404: KDEC TLBIPERO → 0.6 GLYOX + 0.4 MGLYOX + HO2 + 0.2 C4MDIAL + 0.2 C5DICARB + 0.2 TLFUONE + 0.2 BZFUONE + 0.2 MALDIAL

J47405: KDEC TLBIPERO → 0.6 GLYOX + 0.4 MGLYOX + HO2 + 0.2 C4MDIAL + 0.2 C5DICARB + 0.2 TLFUONE + 0.2 BZFUONE + 0.2 MALDIAL

J47407: KDEC CRESO → 0.68 C5CO14OH + 0.68 GLYOX + HO2 + 0.32 PTLQONE

J47408a: KDEC CRESO → 0.68 C5CO14OH + 0.68 GLYOX + HO2 + 0.32 PTLQONE

J47408b: KDEC NCRESO → C5CO14OH + GLYOX + NO2

J47409: Using J for 3-methyl-2-nitrophenol.

J47412: KDEC TLEMUCO → 0.5 C3DIALO2 + 0.5 CO2H3CHO + 0.5 EPXC4DIAL + 0.5 MGLYOX + 0.5 HO2

J47417: Using J for 3-methyl-2-nitrophenol.

J47418: new channel

J47419: Using J for 3-methyl-2-nitrophenol.

J47420: new channel

J47422: KDEC NPTLQO → C7CO4DB + NO2

J47423: KDEC PTLQO → C6CO2OHCO3

J47425: KDEC MNNCATECO → NC4MDCO2H + HCOCO2H + NO2

J47426: KDEC MNCATECO → NC4MDCO2H + HCOCO2H + HO2

J47428: KDEC NDNCRESO → NC4MDCO2H + HNO3 + CO + CO + NO2

J47429: KDEC DNCRESO → NC4MDCO2H + HCOCO2H + NO2

J48400: KDEC STYRENO → HO2 + HCHO + BENZAL

J40203b: Substituted vinyl alcohol in analogy to CH_3CHO photolysis.

Table 3: Reversible (Henry's law) equilibria and irreversible ("heterogenous") uptake

#	labels	reaction	rate coefficient	reference
H1000f_a01	TrAa01Sc	$O_2 \rightarrow O_2(aq)$	k_exf(01,ind_O2)	see general notes*
H1000b_a01	TrAa01Sc	$O_2(aq) \rightarrow O_2$	k_exb(01,ind_O2)	see general notes*
H1001f_a01	TrAa01MblScScm	$O_3 \rightarrow O_3(aq)$	k_exf(01,ind_O3)	see general notes*
H1001b_a01	TrAa01MblScScm	$O_3(aq) \rightarrow O_3$	k_exb(01,ind_O3)	see general notes*
H2100f_a01	TrAa01Sc	$OH \rightarrow OH(aq)$	k_exf(01,ind_OH)	see general notes*
H2100b_a01	TrAa01Sc	$OH(aq) \rightarrow OH$	k_exb(01,ind_OH)	see general notes*
H2101f_a01	TrAa01Sc	$HO_2 \rightarrow HO_2(aq)$	k_exf(01,ind_HO2)	see general notes*
H2101b_a01	TrAa01Sc	$HO_2(aq) \rightarrow HO_2$	k_exb(01,ind_HO2)	see general notes*
H2102f_a01	TrAa01MblScScm	$H_2O_2 \rightarrow H_2O_2(aq)$	k_exf(01,ind_H2O2)	see general notes*
H2102b_a01	TrAa01MblScScm	$H_2O_2(aq) \rightarrow H_2O_2$	k_exb(01,ind_H2O2)	see general notes*
H3101f_a01	TrAa01ScN	$NO_2 \rightarrow NO_2(aq)$	k_exf(01,ind_NO2)	see general notes*
H3101b_a01	TrAa01ScN	$NO_2(aq) \rightarrow NO_2$	k_exb(01,ind_NO2)	see general notes*
H3102f_a01	TrAa01ScN	$NO_3 \rightarrow NO_3(aq)$	k_exf(01,ind_NO3)	see general notes*
H3102b_a01	TrAa01ScN	$NO_3(aq) \rightarrow NO_3$	k_exb(01,ind_NO3)	see general notes*
H3200f_a01	TrAa01MblScScmN	$NH_3 \rightarrow NH_3(aq)$	k_exf(01,ind_NH3)	see general notes*
H3200b_a01	TrAa01MblScScmN	$NH_3(aq) \rightarrow NH_3$	k_exb(01,ind_NH3)	see general notes*
H3201_a01	TrAa01MblScScmN	$N_2O_5 \rightarrow HNO_3(aq) + HNO_3(aq)$	k_exf_N2O5(01)*C(ind_H2O_a01)	Behnke et al. (1994), Behnke et al. (1997)
H3202f_a01	TrAa01ScN	$HONO \rightarrow HONO(aq)$	k_exf(01,ind_HONO)	see general notes*
H3202b_a01	TrAa01ScN	$HONO(aq) \rightarrow HONO$	k_exb(01,ind_HONO)	see general notes*
H3203f_a01	TrAa01MblScScmN	$HNO_3 \rightarrow HNO_3(aq)$	k_exf(01,ind_HNO3)	see general notes*
H3203b_a01	TrAa01MblScScmN	$HNO_3(aq) \rightarrow HNO_3$	k_exb(01,ind_HNO3)	see general notes*
H3204f_a01	TrAa01ScN	$HNO_4 \rightarrow HNO_4(aq)$	k_exf(01,ind_HNO4)	see general notes*
H3204b_a01	TrAa01ScN	$HNO_4(aq) \rightarrow HNO_4$	k_exb(01,ind_HNO4)	see general notes*
H4100f_a01	TrAa01MblScScm	$CO_2 \rightarrow CO_2(aq)$	k_exf(01,ind_CO2)	see general notes*
H4100b_a01	TrAa01MblScScm	$CO_2(aq) \rightarrow CO_2$	k_exb(01,ind_CO2)	see general notes*
H4101f_a01	TrAa01ScScm	$HCHO \rightarrow HCHO(aq)$	k_exf(01,ind_HCHO)	see general notes*
H4101b_a01	TrAa01ScScm	$HCHO(aq) \rightarrow HCHO$	k_exb(01,ind_HCHO)	see general notes*
H4102f_a01	TrAa01Sc	$CH_3O_2 \rightarrow CH_3OO(aq)$	k_exf(01,ind_CH3O2)	see general notes*
H4102b_a01	TrAa01Sc	$CH_3OO(aq) \rightarrow CH_3O_2$	k_exb(01,ind_CH3O2)	see general notes*
H4103f_a01	TrAa01ScScm	$HCOOH \rightarrow HCOOH(aq)$	k_exf(01,ind_HCOOH)	see general notes*
H4103b_a01	TrAa01ScScm	$HCOOH(aq) \rightarrow HCOOH$	k_exb(01,ind_HCOOH)	see general notes*
H4104f_a01	TrAa01ScScm	$CH_3OOH \rightarrow CH_3OOH(aq)$	k_exf(01,ind_CH3OOH)	see general notes*
H4104b_a01	TrAa01ScScm	$CH_3OOH(aq) \rightarrow CH_3OOH$	k_exb(01,ind_CH3OOH)	see general notes*

Table 3: Reversible (Henry's law) equilibria and irreversible ("heterogenous") uptake

#	labels	reaction	rate coefficient	reference
`H6000f_a01`	TrAa01MblScCl	$Cl_2 \rightarrow Cl_2(aq)$	`k_exf(01,ind_Cl2)`	see general notes*
`H6000b_a01`	TrAa01MblScCl	$Cl_2(aq) \rightarrow Cl_2$	`k_exb(01,ind_Cl2)`	see general notes*
`H6200f_a01`	TrAa01MblScScmCl	$HCl \rightarrow HCl(aq)$	`k_exf(01,ind_HCl)`	see general notes*
`H6200b_a01`	TrAa01MblScScmCl	$HCl(aq) \rightarrow HCl$	`k_exb(01,ind_HCl)`	see general notes*
`H6201f_a01`	TrAa01MblScCl	$HOCl \rightarrow HOCl(aq)$	`k_exf(01,ind_HOCl)`	see general notes*
`H6201b_a01`	TrAa01MblScCl	$HOCl(aq) \rightarrow HOCl$	`k_exb(01,ind_HOCl)`	see general notes*
`H6300_a01`	TrAa01MblClN	$N_2O_5 + Cl^-(aq) \rightarrow ClNO_2 + NO_3^-(aq)$	`k_exf_N2O5(01) * 5.E2`	Behnke et al. (1994), Behnke et al. (1997)
`H6301_a01`	TrAa01MblClN	$ClNO_3 \rightarrow HOCl(aq) + HNO_3(aq)$	`k_exf_ClNO3(01) * C(ind_H2O_a01)`	see general notes*
`H6302_a01`	TrAa01MblClN	$ClNO_3 + Cl^-(aq) \rightarrow Cl_2(aq) + NO_3^-(aq)$	`k_exf_ClNO3(01) * 5.E2`	see general notes*
`H8000f_a01`	TrAa01ScI	$I_2 \rightarrow I_2(aq)$	`k_exf(01,ind_I2)`	see general notes*
`H8000b_a01`	TrAa01ScI	$I_2(aq) \rightarrow I_2$	`k_exb(01,ind_I2)`	see general notes*
`H8100f_a01`	TrAa01MblScI	$IO \rightarrow IO(aq)$	`k_exf(01,ind_IO)`	see general notes*
`H8100b_a01`	TrAa01MblScI	$IO(aq) \rightarrow IO$	`k_exb(01,ind_IO)`	see general notes*
`H8101_a01`	TrAa01I	$OIO \rightarrow HOI(aq) + HO_2(aq)$	`k_exf(01,ind_OIO)`	see general notes*
`H8102_a01`	TrAa01I	$I_2O_2 \rightarrow HOI(aq) + H^+(aq) + IO_2^-(aq)$	`k_exf(01,ind_I2O2)`	see general notes*
`H8200f_a01`	TrAa01MblScI	$HOI \rightarrow HOI(aq)$	`k_exf(01,ind_HOI)`	see general notes*
`H8200b_a01`	TrAa01MblScI	$HOI(aq) \rightarrow HOI$	`k_exb(01,ind_HOI)`	see general notes*
`H8201_a01`	TrAa01MblScI	$HI \rightarrow H^+(aq) + I^-(aq)$	$k_{mt}(HI) \cdot lwc$	see general notes*
`H8202_a01`	TrAa01ScI	$HIO_3 \rightarrow IO_3^-(aq) + H^+(aq)$	$k_{mt}(HIO_3) \cdot lwc$	see general notes*
`H8300_a01`	TrAa01IN	$INO_2 \rightarrow HOI(aq) + HONO(aq)$	`k_exf(01,ind_INO2)`	see general notes*
`H8301_a01`	TrAa01MblIN	$INO_3 \rightarrow HOI(aq) + HNO_3(aq)$	`k_exf(01,ind_INO3)`	see general notes*
`H8600f_a01`	TrAa01MblScClI	$ICl \rightarrow ICl(aq)$	`k_exf(01,ind_ICl)`	see general notes*
`H8600b_a01`	TrAa01MblScClI	$ICl(aq) \rightarrow ICl$	`k_exb(01,ind_ICl)`	see general notes*
`H9100f_a01`	TrAa01MblScScmS	$SO_2 \rightarrow SO_2(aq)$	`xnom7so2*k_exf(01,ind_SO2)`	see general notes*
`H9100b_a01`	TrAa01MblScScmS	$SO_2(aq) \rightarrow SO_2$	`xnom7so2*k_exb(01,ind_SO2)`	see general notes*
`H9200_a01`	TrAa01MblScScmS	$H_2SO_4 \rightarrow H_2SO_4(aq)$	`xnom7sulf*k_exf(01,ind_H2SO4)`	see general notes*
`H9400f_a01`	TrAa01CS	$DMSO \rightarrow DMSO(aq)$	`xnom7dmso*k_exf(01,ind_DMSO)`	see general notes*
`H9400b_a01`	TrAa01CS	$DMSO(aq) \rightarrow DMSO$	`xnom7dmso*k_exb(01,ind_DMSO)`	see general notes*
`H9401_a01`	TrAa01MblS	$CH_3SO_3H \rightarrow CH_3SO_3^-(aq) + H^+(aq)$	`xnom7dmso*k_exf(01,ind_CH3SO3H)`	see general notes*
`H9402f_a01`	TrAa01CS	$DMS \rightarrow DMS(aq)$	`xnom7dmso*k_exf(01,ind_DMS)`	see general notes*
`H9402b_a01`	TrAa01CS	$DMS(aq) \rightarrow DMS$	`xnom7dmso*k_exb(01,ind_DMS)`	see general notes*
`H3103f_a01MK`	TrAa01ScN	$N_2O_3 \rightarrow N_2O_3(aq)$	`xnom7nox*k_exf(01,ind_N2O3)`	Karl et al. (2012b)
`H3103b_a01MK`	TrAa01ScN	$N_2O_3(aq) \rightarrow N_2O_3$	`xnom7nox*k_exb(01,ind_N2O3)`	Karl et al. (2012b)

Table 3: Reversible (Henry's law) equilibria and irreversible ("heterogenous") uptake

#	labels	reaction	rate coefficient	reference
H3104f_a01MK	TrAa01ScN	$N_2O_4 \rightarrow N_2O_4(aq)$	xnom7nox*k_exf(01,ind_N2O4)	Karl et al. (2012b)
H3104b_a01MK	TrAa01ScN	$N_2O_4(aq) \rightarrow N_2O_4$	xnom7nox*k_exb(01,ind_N2O4)	Karl et al. (2012b)
H4106f_a01MK	TrAa01AmiScN	MMA $\rightarrow$ MMA(aq)	xnom7amin*k_exf(01,ind_MMA)	Karl et al. (2012b)
H4106b_a01MK	TrAa01AmiScN	MMA(aq) $\rightarrow$ MMA	xnom7amin*k_exb(01,ind_MMA)	Karl et al. (2012b)
H4107f_a01MK	TrAa01AmiScN	MMNNO2 $\rightarrow$ MMNNO2(aq)	xnom7nno*k_exf(01,ind_MMNNO2)	Karl et al. (2012b)
H4107b_a01MK	TrAa01AmiScN	MMNNO2(aq) $\rightarrow$ MMNNO2	xnom7nno*k_exb(01,ind_MMNNO2)	Karl et al. (2012b)
H4108f_a01MK	TrAa01AmiScN	HNCO $\rightarrow$ HNCO(aq)	xnom7amin*k_exf(01,ind_HNCO)	Karl et al. (2012b)
H4108b_a01MK	TrAa01AmiScN	HNCO(aq) $\rightarrow$ HNCO	xnom7amin*k_exb(01,ind_HNCO)	Karl et al. (2012b)
H4109f_a01MK	TrAa01AmiScN	H2NCHO $\rightarrow$ H2NCHO(aq)	xnom7amin*k_exf(01,ind_H2NCHO)	Karl et al. (2012b)
H4109b_a01MK	TrAa01AmiScN	H2NCHO(aq) $\rightarrow$ H2NCHO	xnom7amin*k_exb(01,ind_H2NCHO)	Karl et al. (2012b)
H4203f_a01MK	TrAa01ScC	$HOCH_2CHO \rightarrow HOCH_2CHO(aq)$	xnom7co2*k_exf(01,ind_HOCH2CHO)	Ervens et al. (2004)
H4203b_a01MK	TrAa01ScC	$HOCH_2CHO(aq) \rightarrow HOCH_2CHO$	xnom7co2*k_exb(01,ind_HOCH2CHO)	Ervens et al. (2004)
H4204f_a01MK	TrAa01ScC	$HOCH_2CO_2H \rightarrow HOCH_2CO_2H(aq)$	xnom7co2*k_exf(01,ind_HOCH2CO2H)	Ervens et al. (2004)
H4204b_a01MK	TrAa01ScC	$HOCH_2CO_2H(aq) \rightarrow HOCH_2CO_2H$	xnom7co2*k_exb(01,ind_HOCH2CO2H)	Ervens et al. (2004)
H4205f_a01MK	TrAa01ScC	$HCOCO_2H \rightarrow HCOCO_2H(aq)$	xnom7co2*k_exf(01,ind_HCOCO2H)	Ervens et al. (2004)
H4205b_a01MK	TrAa01ScC	$HCOCO_2H(aq) \rightarrow HCOCO_2H$	xnom7co2*k_exb(01,ind_HCOCO2H)	Ervens et al. (2004)
H4206f_a01MK	TrAa01ScC	GLYOX $\rightarrow$ GLYOX(aq)	xnom7co2*k_exf(01,ind_GLYOX)	Ervens et al. (2004)
H4206b_a01MK	TrAa01ScC	GLYOX(aq) $\rightarrow$ GLYOX	xnom7co2*k_exb(01,ind_GLYOX)	Ervens et al. (2004)
H4207f_a01MK	TrAa01AmiScCN	DMA $\rightarrow$ DMA(aq)	xnom7amin*k_exf(01,ind_DMA)	Karl et al. (2012b)
H4207b_a01MK	TrAa01AmiScCN	DMA(aq) $\rightarrow$ DMA	xnom7amin*k_exb(01,ind_DMA)	Karl et al. (2012b)
H4208f_a01MK	TrAa01AmiScCN	MEA $\rightarrow$ MEA(aq)	xnom7amin*k_exf(01,ind_MEA)	Karl et al. (2012b)
H4208b_a01MK	TrAa01AmiScCN	MEA(aq) $\rightarrow$ MEA	xnom7amin*k_exb(01,ind_MEA)	Karl et al. (2012b)
H4209f_a01MK	TrAa01AmiScCN	MEANNO $\rightarrow$ MEANNO(aq)	xnom7nno*k_exf(01,ind_MEANNO)	Karl et al. (2012b)
H4209b_a01MK	TrAa01AmiScCN	MEANNO(aq) $\rightarrow$ MEANNO	xnom7nno*k_exb(01,ind_MEANNO)	Karl et al. (2012b)
H4210f_a01MK	TrAa01AmiScCN	MEANNO2 $\rightarrow$ MEANNO2(aq)	xnom7nno*k_exf(01,ind_MEANNO2)	Karl et al. (2012b)
H4210b_a01MK	TrAa01AmiScCN	MEANNO2(aq) $\rightarrow$ MEANNO2	xnom7nno*k_exb(01,ind_MEANNO2)	Karl et al. (2012b)
H4211f_a01MK	TrAa01AmiScCN	NDMA $\rightarrow$ NDMA(aq)	xnom7nno*k_exf(01,ind_NDMA)	Karl et al. (2012b)
H4211b_a01MK	TrAa01AmiScCN	NDMA(aq) $\rightarrow$ NDMA	xnom7nno*k_exb(01,ind_NDMA)	Karl et al. (2012b)
H4212f_a01MK	TrAa01AmiScCN	DMNNO2 $\rightarrow$ DMNNO2(aq)	xnom7nno*k_exf(01,ind_DMNNO2)	Karl et al. (2012b)
H4212b_a01MK	TrAa01AmiScCN	DMNNO2(aq) $\rightarrow$ DMNNO2	xnom7nno*k_exb(01,ind_DMNNO2)	Karl et al. (2012b)
H4213f_a01MK	TrAa01AmiScCN	H2NCOCH2OH $\rightarrow$ H2NCOCH2OH(aq)	xnom7amin*k_exf(01,ind_ H2NCOCH2OH)	Karl et al. (2012b)
H4213b_a01MK	TrAa01AmiScCN	H2NCOCH2OH(aq) $\rightarrow$ H2NCOCH2OH	xnom7amin*k_exb(01,ind_ H2NCOCH2OH)	Karl et al. (2012b)
H4214f_a01MK	TrAa01AmiScCN	CH3NHCHO $\rightarrow$ $CH_3NHCHO(aq)$	xnom7amin*k_exf(01,ind_CH3NHCHO)	Karl et al. (2012b)

Table 3: Reversible (Henry's law) equilibria and irreversible ("heterogenous") uptake

#	labels	reaction	rate coefficient	reference
H4214b_a01MK	TrAa01AmiScCN	$CH_3NHCHO(aq) \rightarrow CH3NHCHO$	xnom7amin*k_exb(01,ind_CH3NHCHO)	Karl et al. (2012b)
H4301f_a01MK	TrAa01ScC	MGLYOX $\rightarrow$ MGLYOX(aq)	xnom7co2*k_exf(01,ind_MGLYOX)	Ervens et al. (2004)
H4301b_a01MK	TrAa01ScC	MGLYOX(aq) $\rightarrow$ MGLYOX	xnom7co2*k_exb(01,ind_MGLYOX)	Ervens et al. (2004)
H4302f_a01MK	TrAa01ScC	MGLYOAC $\rightarrow$ MGLYOAC(aq)	xnom7co2*k_exf(01,ind_MGLYOAC)	Ervens et al. (2004)
H4302b_a01MK	TrAa01ScC	MGLYOAC(aq) $\rightarrow$ MGLYOAC	xnom7co2*k_exb(01,ind_MGLYOAC)	Ervens et al. (2004)
H4303f_a01MK	TrAa01AmiScCN	TMA $\rightarrow$ TMA(aq)	xnom7amin*k_exf(01,ind_TMA)	Karl et al. (2012b)
H4303b_a01MK	TrAa01AmiScCN	TMA(aq) $\rightarrow$ TMA	xnom7amin*k_exb(01,ind_TMA)	Karl et al. (2012b)
H4304f_a01MK	TrAa01AmiScCN	DMNCHO $\rightarrow$ DMNCHO(aq)	xnom7amin*k_exf(01,ind_DMNCHO)	Karl et al. (2012b)
H4304b_a01MK	TrAa01AmiScCN	DMNCHO(aq) $\rightarrow$ DMNCHO	xnom7amin*k_exb(01,ind_DMNCHO)	Karl et al. (2012b)
H4400f_a01MK	TrAa01AmiScCN	DEA $\rightarrow$ DEA(aq)	xnom7amin*k_exf(01,ind_DEA)	Karl et al. (2012b)
H4400b_a01MK	TrAa01AmiScCN	DEA(aq) $\rightarrow$ DEA	xnom7amin*k_exb(01,ind_DEA)	Karl et al. (2012b)
H4401f_a01MK	TrAa01AmiScCN	NDELA $\rightarrow$ NDELA(aq)	xnom7nno*k_exf(01,ind_NDELA)	Karl et al. (2012b)
H4401b_a01MK	TrAa01AmiScCN	NDELA(aq) $\rightarrow$ NDELA	xnom7nno*k_exb(01,ind_NDELA)	Karl et al. (2012b)
H4402f_a01MK	TrAa01AmiScCN	DEANNO2 $\rightarrow$ DEANNO2(aq)	xnom7nno*k_exf(01,ind_DEANNO2)	Karl et al. (2012b)
H4402b_a01MK	TrAa01AmiScCN	DEANNO2(aq) $\rightarrow$ DEANNO2	xnom7nno*k_exb(01,ind_DEANNO2)	Karl et al. (2012b)
H4403f_a01MK	TrAa01Sc	BSOV $\rightarrow$ SUCCAC(aq)	xnom7co2*k_exf(01,ind_BSOV)	Ervens et al. (2004)
H4403b_a01MK	TrAa01Sc	SUCCAC(aq) $\rightarrow$ BSOV	xnom7co2*k_exb(01,ind_BSOV)	Ervens et al. (2004)
H4600f_a01MK	TrAa01AmiScCN	TEA $\rightarrow$ TEA(aq)	xnom7amin*k_exf(01,ind_TEA)	Karl et al. (2012b)
H4600b_a01MK	TrAa01AmiScCN	TEA(aq) $\rightarrow$ TEA	xnom7amin*k_exb(01,ind_TEA)	Karl et al. (2012b)
H4601f_a01MK	TrAa01Sc	BLOV $\rightarrow$ ADIPAC(aq)	xnom7co2*k_exf(01,ind_BLOV)	Ervens et al. (2004)
H4601b_a01MK	TrAa01Sc	ADIPAC(aq) $\rightarrow$ BLOV	xnom7co2*k_exb(01,ind_BLOV)	Ervens et al. (2004)
H9403f_a01MK	TrAa01S	$CH_3SOOH \rightarrow$ MSIA(aq)	xnom7dmso*k_exf(01,ind_CH3SOOH)	see general notes*
H9403b_a01MK	TrAa01S	MSIA(aq) $\rightarrow CH_3SOOH$	xnom7dmso*k_exb(01,ind_CH3SOOH)	see general notes*

General notes

The forward (`k_exf`) and backward (`k_exb`) rate coefficients are calculated in subroutine `mecca_aero_calc_k_ex` in the file `messy_mecca_aero.f90` using accommodation coefficients and Henry's law constants from chemprop (see `chemprop.pdf`).

For uptake of X (X = N_2O_5, $ClNO_3$, or $BrNO_3$) and subsequent reaction with H_2O, Cl^-, and Br^- in H3201, H6300, H6301, H6302, H7300, H7301, H7302, H7601, and H7602, we define:

$$k_{\rm exf}(\mathrm{X})=\frac{k_{\rm mt}(\mathrm{X}) \times \mathrm{LWC}}{[\mathrm{H_2O}] + 5 \times 10^2[\mathrm{Cl^-}] + 3 \times 10^5[\mathrm{Br^-}]}$$

Here, $k_{\rm mt}$ = mass transfer coefficient, and LWC = liquid water content of the aerosol. The total uptake rate of X is only determined by $k_{\rm mt}$. The factors only affect the branching between hydrolysis and the halide reactions. The factor 5×10^2 was chosen such that the chloride reaction dominates over hydrolysis at about $[Cl^-] > 0.1$ M (see Fig. 3 in Behnke et al. (1997)), i.e. when the ratio $[H_2O]/[Cl^-]$ is less than 5×10^2. The ratio $5\times10^2/3\times10^5$ was chosen such that the reactions with chloride and bromide are roughly equal for sea water composition (Behnke et al., 1994). These ratios were measured for uptake of N_2O_5. Here, they are also used for $ClNO_3$ and $BrNO_3$.

Table 4: Heterogeneous reactions

#	labels	reaction	rate coefficient	reference

General notes

Heterogeneous reaction rates are calculated with an external module (e.g., MECCA_KHET) and then supplied to the MECCA chemistry (see `www.messy-interface.org` for details)

Table 5: Acid-base and other equilibria

#	labels	reaction	$K_0[M^{m-n}]$	$-\Delta H/R[K]$	reference
EQ20_a01	TrAa01Sc	$HO_2 \rightleftharpoons O_2^- + H^+$	1.6E-5		Weinstein-Lloyd and Schwartz (1991)
EQ21_a01	TrAa01MblScScm	$H_2O \rightleftharpoons H^+ + OH^-$	1.0E-16	-6716	Chameides (1984)
EQ30_a01	TrAa01MblScScmN	$NH_4^+ \rightleftharpoons H^+ + NH_3$	5.88E-10	-2391	Chameides (1984)
EQ31_a01	TrAa01ScN	$HONO \rightleftharpoons H^+ + NO_2^-$	5.1E-4	-1260	Schwartz and White (1981)
EQ32_a01	TrAa01MblScScmN	$HNO_3 \rightleftharpoons H^+ + NO_3^-$	15	8700	Davis and de Bruin (1964)
EQ33_a01	TrAa01ScN	$HNO_4 \rightleftharpoons NO_4^- + H^+$	1.E-5		Warneck (1999)
EQ40_a01	TrAa01MblScScm	$CO_2 \rightleftharpoons H^+ + HCO_3^-$	4.3E-7	-913	Chameides (1984)*
EQ41_a01	TrAa01ScScm	$HCOOH \rightleftharpoons H^+ + HCOO^-$	1.8E-4		Weast (1980)
EQ60_a01	TrAa01Cl	$Cl_2^- \rightleftharpoons Cl + Cl^-$	7.3E-6		Yu (2004)
EQ61_a01	TrAa01MblScScmCl	$HCl \rightleftharpoons H^+ + Cl^-$	1.7E6	6896	Marsh and McElroy (1985)
EQ62_a01	TrAa01ScCl	$HOCl \rightleftharpoons H^+ + ClO^-$	3.2E-8		Lax (1969)
EQ80_a01	TrAa01MblScClI	$ICl + Cl^- \rightleftharpoons ICl_2^-$	7.7E1		Wang et al. (1989)
EQ90_a01	TrAa01MblScScmS	$SO_2 \rightleftharpoons H^+ + HSO_3^-$	1.7E-2	2090	Chameides (1984)
EQ91_a01	TrAa01MblScScmS	$HSO_3^- \rightleftharpoons H^+ + SO_3^{2-}$	6.0E-8	1120	Chameides (1984)
EQ92_a01	TrAa01MblScScmS	$HSO_4^- \rightleftharpoons H^+ + SO_4^{2-}$	1.2E-2	2720	Seinfeld and Pandis (1998)
EQ93_a01	TrAa01MblScScmS	$H_2SO_4 \rightleftharpoons H^+ + HSO_4^-$	1.0E3		Seinfeld and Pandis (1998)
EQ110_a01	TrAa01Fe	$Fe^{3+} \rightleftharpoons FeOH^{2+} + H^+$	2.34E-3		de Laat and Le (2006)*
EQ111_a01	TrAa01Fe	$FeOH^{2+} \rightleftharpoons Fe(OH)_2^+ + H^+$	2E-4		de Laat and Le (2006)*
EQ112_a01	TrAa01Fe	$Fe^{3+} + H_2O_2 \rightleftharpoons FeHO_2^{2+} + H^+$	3.1E-3		de Laat and Le (2006)
EQ113_a01	TrAa01Fe	$FeOH^{2+} + H_2O_2 \rightleftharpoons Fe(OH)(HO_2)^+ + H^+$	2E-4		de Laat and Le (2006)
EQ114_a01	TrAa01ClFe	$Fe^{3+} + Cl^- \rightleftharpoons FeCl^{2+}$	6.61		de Laat and Le (2006)*
EQ115_a01	TrAa01ClFe	$FeCl^{2+} + Cl^- \rightleftharpoons FeCl_2^+$	1.6		de Laat and Le (2006)*
EQ116_a01	TrAa01FeS	$Fe^{3+} + SO_4^{2-} \rightleftharpoons FeSO_4^+$	120		Brand and van Eldik (1995)*
EQ117_a01	TrAa01FeS	$FeOH^{2+} + HSO_3^- \rightleftharpoons FeSO_3^+$	8.25E2		Warneck (2018)*
EQ118_a01	TrAa01FeS	$Fe^{2+} + SO_3^- \rightleftharpoons FeSO_3^+$	1.6E7		Warneck (2018)
EQ43_a01MK	TrAa01ScC	$HCOCO_2H \rightleftharpoons H^+ + HCOCOO^-$	6.6E-4		Ervens et al. (2004)
EQ44_a01MK	TrAa01ScC	$MGLYOAC \rightleftharpoons H^+ + CH_3COCOO^-$	4.07E-3		Ervens et al. (2004)
EQ45_a01MK	TrAa01Sc	$OXALAC \rightleftharpoons H^+ + HC_2O_4^-$	6.4E-2		Ervens et al. (2004)
EQ46_a01MK	TrAa01Sc	$HC_2O_4^- \rightleftharpoons H^+ + C_2O_4^{2-}$	5.25E-5		Ervens et al. (2004)
EQ47_a01MK	TrAa01Sc	$SUCCAC \rightleftharpoons H^+ + CH_2CH_2HC_2O_4^-$	6.4E-2		Ervens et al. (2004)
EQ48_a01MK	TrAa01Sc	$CH_2CH_2HC_2O_4^- \rightleftharpoons H^+ + CH_2CH_2C_2O_4^{2-}$	5.25E-5		Ervens et al. (2004)
EQ120_a01MK	TrAa01AmiScN	$MMA^+ \rightleftharpoons H^+ + MMA$	2.19E-11		Ge et al. (2011)
EQ121_a01MK	TrAa01AmiScCN	$DMA^+ \rightleftharpoons H^+ + DMA$	1.86E-11		Ge et al. (2011)

Table 5: Acid-base and other equilibria

#	labels	reaction	$K_0[M^{m-n}]$	$-\Delta H/R[K]$	reference
EQ122_a01MK	TrAa01AmiScCN	$TMA^+ \rightleftharpoons H^+ + TMA$	1.74E-10		Das and von Sonntag (1986)
EQ123_a01MK	TrAa01AmiScCN	$MEA^+ \rightleftharpoons H^+ + MEA$	3.98E-10		Kishore et al. (2004)
EQ124_a01MK	TrAa01AmiScCN	$DEA^+ \rightleftharpoons H^+ + DEA$	1.17E-9		Kishore et al. (2004)
EQ125_a01MK	TrAa01AmiScCN	$TEA^+ \rightleftharpoons H^+ + TEA$	1.66E-8		Kishore et al. (2004)
EQ126_a01MK	TrAa01AmiScCN	$(CH_3)_2NH^+CH_2 \rightleftharpoons H^+ + (CH_3)_2NCH_2$	4.00E3		Das and von Sonntag (1986)
EQ127_a01MK	TrAa01AmiScCN	$(CH_3)_3N^+ \rightleftharpoons H^+ + (CH_3)_2NCH_2$	5.71E7		Das and von Sonntag (1986)
EQ128_a01MK	TrAa01AmiScCN	$CH_3NH_2^+CH_2 \rightleftharpoons H^+ + CH_3NHCH_2$	4.00E3		Das and von Sonntag (1986)
EQ129_a01MK	TrAa01AmiScCN	$(CH_3)_2NH^+ \rightleftharpoons H^+ + CH_3NHCH_2$	5.71E7		Das and von Sonntag (1986)
EQ130_a01MK	TrAa01AmiScN	$CH_2NH_3^+ \rightleftharpoons H^+ + CH_2NH_2$	4.00E3		Das and von Sonntag (1986)
EQ131_a01MK	TrAa01AmiScN	$CH_3NH_2^+ \rightleftharpoons H^+ + CH_2NH_2$	5.71E7		Das and von Sonntag (1986)
EQ132_a01MK	TrAa01AmiScCN	$HOCHCH_2NH_3^+ \rightleftharpoons H^+ + NH_2CH_2CHOH$	4.00E3		Das and von Sonntag (1986)
EQ133_a01MK	TrAa01AmiScCN	$HOCH_2CH_2NH_2^+ \rightleftharpoons H^+ + NH_2CH_2CHOH$	5.71E7		Das and von Sonntag (1986)
EQ134_a01MK	TrAa01AmiScCN	$HOETNH_2CH_2CHOH^+ \rightleftharpoons H^+ + HOETNHCH_2CHOH$	4.00E3		Das and von Sonntag (1986)
EQ135_a01MK	TrAa01AmiScCN	$(HOET)_2NH^+ \rightleftharpoons H^+ + HOETNHCH_2CHOH$	5.71E7		Das and von Sonntag (1986)
EQ136_a01MK	TrAa01AmiScCN	$(HOET)_2NH^+CH_2CHOH \rightleftharpoons H^+ + DENCH_2CHOH$	4.00E3		Das and von Sonntag (1986)
EQ137_a01MK	TrAa01AmiScCN	$(HOET)_3N^+ \rightleftharpoons H^+ + DENCH_2CHOH$	5.71E7		Das and von Sonntag (1986)
EQ138_a01MK	TrAa01AmiScN	$HNCO \rightleftharpoons H^+ + NCO^-$	2.0E-4		Roberts et al. (2011)

Specific notes

EQ40_a01: For $pK_a(CO_2)$, see also Dickson and Millero (1987).

EQ110_a01: See also K values listed in Tab. 2.5 of Brand and van Eldik (1995).

EQ111_a01: Equilibrium calculated from K_1 and K_2 in Tab. 1 of de Laat and Le (2006). Rate constant for back reaction assumed. See also K values listed in Tab. 2.5 of Brand and van Eldik (1995).

EQ114_a01: See also K values listed in Tab. 2.5 of Brand and van Eldik (1995).

EQ115_a01: Equilibrium calculated from K_{29} and K_{30} in Tab. 2 of de Laat and Le (2006). Rate constant for forward reaction assumed. See also K values listed in Tab. 2.5 of Brand and van Eldik (1995).

EQ116_a01: Equilibrium at $I = 1$ M. Rate constant for back reaction assumed.

EQ117_a01: Rate of equilibration assumed.

Table 6: Aqueous phase reactions

#	labels	reaction	k_0 $[M^{1-n}s^{-1}]$	$-E_a/R[K]$	reference
A1000_a01	TrAa01Sc	$O_3 + O_2^- \rightarrow OH + OH^-$	1.5E9		Sehested et al. (1983)
A2100_a01	TrAa01Sc	$OH + O_2^- \rightarrow OH^-$	1.0E10		Sehested et al. (1968)
A2101_a01	TrAa01Sc	$OH + OH \rightarrow H_2O_2$	5.5E9		Buxton et al. (1988)
A2102_a01	TrAa01Sc	$HO_2 + O_2^- \rightarrow H_2O_2 + OH^-$	1.0E8	-900	Christensen and Sehested (1988)
A2103_a01	TrAa01Sc	$HO_2 + OH \rightarrow H_2O$	7.1E9		Sehested et al. (1968)
A2104_a01	TrAa01Sc	$HO_2 + HO_2 \rightarrow H_2O_2$	9.7E5	-2500	Christensen and Sehested (1988)
A2105_a01	TrAa01Sc	$H_2O_2 + OH \rightarrow HO_2$	2.7E7	-1684	Christensen et al. (1982)
A3100_a01	TrAa01ScN	$NO_2^- + O_3 \rightarrow NO_3^-$	5.0E5	-6950	Damschen and Martin (1983)
A3101_a01	TrAa01ScN	$NO_2 + NO_2 \rightarrow HNO_3 + HONO$	1.0E8		Lee and Schwartz (1981)
A3102_a01	TrAa01ScN	$NO_4^- \rightarrow NO_2^-$	8.0E1		Warneck (1999)
A3200_a01	TrAa01ScN	$NO_2 + HO_2 \rightarrow HNO_4$	1.8E9		Warneck (1999)
A3201_a01	TrAa01ScN	$NO_2^- + OH \rightarrow NO_2 + OH^-$	1.0E10		Wingenter et al. (1999)
A3202_a01	TrAa01ScN	$NO_3 + OH^- \rightarrow NO_3^- + OH$	8.2E7	-2700	Exner et al. (1992)
A3203_a01	TrAa01ScN	$HONO + OH \rightarrow NO_2$	1.0E10		Barker et al. (1970)
A3204_a01	TrAa01ScN	$HONO + H_2O_2 + H^+ \rightarrow HNO_3 + H^+$	4.6E3	-6800	Damschen and Martin (1983)
A4100_a01	TrAa01Sc	$CO_3^- + O_2^- \rightarrow HCO_3^- + OH^-$	6.5E8		Ross et al. (1992)
A4101_a01	TrAa01Sc	$CO_3^- + H_2O_2 \rightarrow HCO_3^- + HO_2$	4.3E5		Ross et al. (1992)
A4102_a01	TrAa01Sc	$HCOO^- + CO_3^- \rightarrow 2\ HCO_3^- + HO_2$	1.5E5		Ross et al. (1992)
A4103_a01	TrAa01Sc	$HCOO^- + OH \rightarrow OH^- + HO_2 + CO_2$	3.1E9	-1240	Chin and Wine (1994)
A4104_a01	TrAa01Sc	$HCO_3^- + OH \rightarrow CO_3^-$	8.5E6		Ross et al. (1992)
A4105_a01	TrAa01Sc	$HCHO + OH \rightarrow HCOOH + HO_2$	7.7E8	-1020	Chin and Wine (1994)
A4106_a01	TrAa01Sc	$HCOOH + OH \rightarrow HO_2 + CO_2$	1.1E8	-991	Chin and Wine (1994)
A4107_a01	TrAa01Sc	$CH_3OO + O_2^- \rightarrow CH_3OOH + OH^-$	5.0E7		Jacob (1986)
A4108_a01	TrAa01Sc	$CH_3OO + HO_2 \rightarrow CH_3OOH$	4.3E5		Jacob (1986)
A4109_a01	TrAa01Sc	$CH_3OH + OH \rightarrow HCHO + HO_2$	9.7E8		Buxton et al. (1988)
A4110a_a01	TrAa01Sc	$CH_3OOH + OH \rightarrow CH_3OO$	2.7E7	-1715	Jacob (1986)
A4110b_a01	TrAa01Sc	$CH_3OOH + OH \rightarrow HCHO + OH$	1.1E7	-1715	Jacob (1986)
A6000_a01	TrAa01Cl	$Cl + Cl \rightarrow Cl_2$	8.8E7		Wu et al. (1980)
A6001_a01	TrAa01Cl	$Cl_2^- + Cl_2^- \rightarrow Cl_2 + 2\ Cl^-$	3.5E9		Yu (2004)
A6100_a01	TrAa01Cl	$Cl^- + O_3 \rightarrow ClO^-$	3.0E-3		Hoigné et al. (1985)
A6101_a01	TrAa01Cl	$Cl_2 + O_2^- \rightarrow Cl_2^-$	1.0E9		Bjergbakke et al. (1981)
A6102_a01	TrAa01Cl	$Cl_2^- + O_2^- \rightarrow 2\ Cl^-$	1.0E9		Jacobi (1996)*
A6200_a01	TrAa01Cl	$Cl \rightarrow H^+ + ClOH^-$	1.8E5		Yu (2004)
A6201_a01	TrAa01Cl	$Cl + H_2O_2 \rightarrow HO_2 + Cl^- + H^+$	2.7E7	-1684	Christensen et al. (1982)

Table 6: Aqueous phase reactions (...continued)

#	labels	reaction	k_0 $[M^{1-n}s^{-1}]$	$-E_a/R[K]$	reference
A6202_a01	TrAa01Cl	$Cl^- + OH \rightarrow ClOH^-$	4.2E9		Yu (2004)
A6203_a01	TrAa01Cl	$Cl_2 + HO_2 \rightarrow Cl_2^- + H^+$	1.0E9		Bjergbakke et al. (1981)
A6204_a01	TrAa01MblCl	$Cl_2 \rightarrow Cl^- + HOCl + H^+$	21.8	-8012	Wang and Margerum (1994)
A6205_a01	TrAa01Cl	$Cl_2^- + HO_2 \rightarrow 2\ Cl^- + H^+$	1.3E10		Jacobi (1996)
A6206_a01	TrAa01Cl	$HOCl + O_2^- \rightarrow Cl + OH^-$	7.5E6		Long and Bielski (1980)
A6207_a01	TrAa01Cl	$HOCl + HO_2 \rightarrow Cl$	7.5E6		Long and Bielski (1980)
A6208_a01	TrAa01MblCl	$HOCl + Cl^- + H^+ \rightarrow Cl_2$	2.2E4	-3508	Wang and Margerum (1994)
A6209_a01	TrAa01Cl	$ClOH^- \rightarrow Cl^- + OH$	6.0E9		Yu (2004)
A6210_a01	TrAa01Cl	$ClOH^- + H^+ \rightarrow Cl$	2.4E10		Yu (2004)
A6300_a01	TrAa01ClN	$Cl + NO_3^- \rightarrow NO_3 + Cl^-$	1.0E8		Buxton et al. (1999b)
A6301_a01	TrAa01ClN	$Cl^- + NO_3 \rightarrow NO_3^- + Cl$	3.4E8		Buxton et al. (1999b)*
A6302_a01	TrAa01ClN	$Cl_2^- + NO_2^- \rightarrow 2\ Cl^- + NO_2$	6.0E7		Jacobi et al. (1996)
A6400_a01	TrAa01Cl	$Cl_2^- + CH_3OOH \rightarrow 2\ Cl^- + H^+ + CH_3OO$	5.0E4		Jacobi et al. (1996)
A8100_a01	TrAa01MblI	$I^- + O_3 \rightarrow HOI + OH^-$	4.2E9	-9311	Magi et al. (1997)
A8101_a01	TrAa01MblI	$IO + IO \rightarrow HOI + IO_2^- + H^+$	1.5E9		Buxton et al. (1986)
A8200_a01	TrAa01MblI	$IO_2^- + H_2O_2 \rightarrow IO_3^-$	6.0E1		Furrow (1987)
A8201_a01	TrAa01I	$HOI + IO_2^- \rightarrow IO_3^- + I^- + H^+$	6.0E2		Chinake and Simoyi (1996)
A8202_a01	TrAa01MblI	$HOI + I^- + H^+ \rightarrow I_2$	4.4E12		Eigen and Kustin (1962)
A8203_a01	TrAa01MblI	$IO_2^- + I^- + H^+ \rightarrow 2\ HOI + OH^-$	2.0E10		Edblom et al. (1987)
A8600_a01	TrAa01MblClI	$ICl \rightarrow HOI + Cl^- + H^+$	2.4E6		Wang et al. (1989)
A8601_a01	TrAa01MblClI	$I^- + HOCl + H^+ \rightarrow ICl$	3.5E11		Nagy et al. (1988)
A8602_a01	TrAa01ClI	$IO_2^- + HOCl \rightarrow IO_3^- + Cl^- + H^+$	1.5E3		Lengyel et al. (1996)
A8603_a01	TrAa01MblClI	$HOI + Cl^- + H^+ \rightarrow ICl$	2.9E10		Wang et al. (1989)
A8604_a01	TrAa01ClI	$HOI + Cl_2 \rightarrow IO_2^- + 2\ Cl^- + 3H^+$	1.0E6		Lengyel et al. (1996)
A8605_a01	TrAa01ClI	$HOI + HOCl \rightarrow IO_2^- + Cl^- + 2\ H^+$	5.0E5		Citri and Epstein (1988)
A8606_a01	TrAa01ClI	$ICl + I^- \rightarrow I_2 + Cl^-$	1.1E9		Margerum et al. (1986)
A9100_a01	TrAa01ScS	$SO_3^- + O_2 \rightarrow SO_5^-$	1.5E9		Huie and Neta (1987)
A9101_a01	TrAa01MblScScmS	$SO_3^{2-} + O_3 \rightarrow SO_4^{2-}$	1.5E9	-5300	Hoffmann (1986)
A9102_a01	TrAa01ScS	$SO_4^- + O_2^- \rightarrow SO_4^{2-}$	3.5E9		Jiang et al. (1992)
A9103_a01	TrAa01ScS	$SO_4^- + SO_3^{2-} \rightarrow SO_3^- + SO_4^{2-}$	4.6E8		Huie and Neta (1987)
A9104_a01	TrAa01ScS	$SO_5^- + O_2^- \rightarrow HSO_5^- + OH^-$	2.3E8		Buxton et al. (1996)
A9105_a01	TrAa01S	$SO_5^- + SO_3^{2-} \rightarrow .72\ SO_4^- + .72\ SO_4^{2-} + .28\ SO_3^- + .28\ HSO_5^- + .28\ OH^-$	1.3E7		Huie and Neta (1987), Deister and Warneck (1990)*
A9106_a01	TrAa01S	$SO_5^- + SO_5^- \rightarrow O_2 + SO_4^{2-} + LSULFUR$	1.0E8		Ross et al. (1992)*

Table 6: Aqueous phase reactions (...continued)

#	labels	reaction	k_0 $[M^{1-n}s^{-1}]$	$-E_a/R[K]$	reference
A9200_a01	TrAa01ScS	SO_3^{2-} + OH → SO_3^- + OH^-	5.5E9		Buxton et al. (1988)
A9201_a01	TrAa01ScS	SO_4^- + OH → HSO_5^-	1.0E9		Jiang et al. (1992)
A9202_a01	TrAa01ScS	SO_4^- + HO_2 → SO_4^{2-} + H^+	3.5E9		Jiang et al. (1992)
A9203_a01	TrAa01ScS	SO_4^- + H_2O → SO_4^{2-} + H^+ + OH	1.1E1	-1110	Herrmann et al. (1995)
A9204_a01	TrAa01ScS	SO_4^- + H_2O_2 → SO_4^{2-} + H^+ + HO_2	1.2E7		Wine et al. (1989)
A9205_a01	TrAa01ScS	HSO_3^- + O_2^- → SO_4^{2-} + OH	3.0E3		see note*
A9206_a01	TrAa01MblScScmS	HSO_3^- + O_3 → SO_4^{2-} + H^+	3.7E5	-5500	Hoffmann (1986)
A9207_a01	TrAa01ScS	HSO_3^- + OH → SO_3^-	4.5E9		Buxton et al. (1988)
A9208_a01	TrAa01ScS	HSO_3^- + HO_2 → SO_4^{2-} + OH + H^+	3.0E3		see note*
A9209_a01	TrAa01MblScScmS	HSO_3^- + H_2O_2 → SO_4^{2-} + H^+	5.2E6	-3650	Martin and Damschen (1981)
A9210_a01	TrAa01ScS	HSO_3^- + SO_4^- → SO_3^- + SO_4^{2-} + H^+	8.0E8		Huie and Neta (1987)
A9211_a01	TrAa01S	HSO_3^- + SO_5^- → .75 SO_4^- + .75 SO_4^{2-} + .75 H^+ + .25 SO_3^- + .25 HSO_5^-	1.0E5		Huie and Neta (1987)
A9212_a01	TrAa01ScS	HSO_3^- + HSO_5^- + H^+ → 2 HSO_4^- + H^+	7.1E6		Betterton and Hoffmann (1988)
A9301_a01	TrAa01ScNS	SO_4^- + NO_3^- → SO_4^{2-} + NO_3	5.0E4		Exner et al. (1992)
A9302_a01	TrAa01ScNS	SO_4^{2-} + NO_3 → NO_3^- + SO_4^-	1.0E5		Løgager et al. (1993)
A9304_a01	TrAa01ScNS	HSO_3^- + NO_3 → SO_3^- + NO_3^- + H^+	1.4E9	-2000	Exner et al. (1992)
A9305_a01	TrAa01ScNS	HSO_3^- + HNO_4 → HSO_4^- + NO_3^- + H^+	3.1E5		Warneck (1999)
A9400_a01	TrAa01ScS	SO_3^{2-} + HCHO → $CH_2OHSO_3^-$ + OH^-	1.4E4		Boyce and Hoffmann (1984)*
A9401_a01	TrAa01ScS	SO_3^{2-} + CH_3OOH + H^+ → SO_4^{2-} + H^+ + CH_3OH	1.6E7	-3800	Lind et al. (1987)
A9402_a01	TrAa01ScS	HSO_3^- + HCHO → $CH_2OHSO_3^-$	4.3E-1		Boyce and Hoffmann (1984)*
A9403_a01	TrAa01ScS	HSO_3^- + CH_3OOH + H^+ → HSO_4^- + H^+ + CH_3OH	1.6E7	-3800	Lind et al. (1987)
A9404_a01	TrAa01ScS	$CH_2OHSO_3^-$ + OH^- → SO_3^{2-} + HCHO	3.6E3		Seinfeld and Pandis (1998)
A9600_a01	TrAa01ClS	SO_3^{2-} + Cl_2^- → SO_3^- + 2 Cl^-	6.2E7		Jacobi et al. (1996)
A9601_a01	TrAa01MblClS	SO_3^{2-} + HOCl → Cl^- + HSO_4^-	7.6E8		Fogelman et al. (1989)
A9602_a01	TrAa01ClS	SO_4^- + Cl^- → SO_4^{2-} + Cl	2.5E8		Buxton et al. (1999a)
A9603_a01	TrAa01ClS	SO_4^{2-} + Cl → SO_4^- + Cl^-	2.1E8		Buxton et al. (1999a)
A9604_a01	TrAa01ClS	HSO_3^- + Cl_2^- → SO_3^- + 2 Cl^- + H^+	4.7E8	-1082	Shoute et al. (1991)
A9605_a01	TrAa01MblClS	HSO_3^- + HOCl → Cl^- + HSO_4^- + H^+	7.6E8		see note*
A9606_a01	TrAa01ClS	HSO_5^- + Cl^- → HOCl + SO_4^{2-}	1.8E-3	-7352	Fortnum et al. (1960)
A9800_a01	TrAa01IS	HSO_3^- + I_2 → 2 I^- + HSO_4^- + 2 H^+	1.7E9		Yiin and Margerum (1990)
A11101_a01	TrAa01Fe	Fe^{2+} + O_2^- → Fe^{3+} + HO2− + OH^-	1E7		de Laat and Le (2006)
A11102_a01	TrAa01Fe	Fe^{3+} + O_2^- → O_2 + Fe^{2+}	5E7		de Laat and Le (2006)
A11103_a01	TrAa01Fe	Fe^{2+} + O_3 → FeO^{2+} + O_2	8.2E5		Løgager et al. (1992)

Table 6: Aqueous phase reactions (...continued)

#	labels	reaction	k_0 $[M^{1-n}s^{-1}]$	$-E_a/R[K]$	reference
A11201a_a01	TrAa01Fe	$Fe^{2+} + OH \rightarrow Fe^{3+} + OH^-$	2.7E8		de Laat and Le (2006)
A11201b_a01	TrAa01Fe	$FeOH^+ + OH \rightarrow Fe^{3+} + 2\ OH^-$	2.7E8		de Laat and Le (2006)
A11202a_a01	TrAa01Fe	$Fe^{2+} + H_2O_2 \rightarrow Fe^{3+} + OH + OH^-$	5.5E1		de Laat and Le (2006)
A11202b_a01	TrAa01Fe	$FeOH^+ + H_2O_2 \rightarrow Fe^{3+} + OH + 2\ OH^-$	5.9E6		de Laat and Le (2006)
A11203_a01	TrAa01Fe	$FeHO_2^{2+} \rightarrow Fe^{2+} + HO_2$	2.3E-3		de Laat and Le (2006)
A11204_a01	TrAa01Fe	$Fe(OH)(HO_2)^+ \rightarrow Fe^{2+} + HO_2 + OH^-$	2.3E-3		de Laat and Le (2006)
A11206_a01	TrAa01Fe	$Fe^{2+} + HO_2 \rightarrow Fe^{3+} + HO2^-$	1.2E6		de Laat and Le (2006)
A11208a_a01	TrAa01Fe	$FeOH^{2+} + O_2^- \rightarrow Fe^{2+} + O_2 + OH^-$	1.5E8		Rush and Bielski (1985)
A11208b_a01	TrAa01Fe	$Fe(OH)_2^+ + O_2^- \rightarrow Fe^{2+} + O_2 + 2\ OH^-$	1.5E8		Rush and Bielski (1985)
A11209_a01	TrAa01Fe	$Fe^{2+} + O_2^- \rightarrow Fe^{3+} + H_2O_2 + 2\ OH^-$	1.0E7		Rush and Bielski (1985)
A11210_a01	TrAa01Fe	$Fe^{2+} + OH \rightarrow FeOH^{2+}$	4.3E8		Christensen and Sehested (1981)
A11211_a01	TrAa01Fe	$FeO^{2+} + H_2O_2 \rightarrow Fe^{3+} + HO_2 + OH^-$	9.5E3		Løgager et al. (1992)
A11212_a01	TrAa01Fe	$FeO^{2+} \rightarrow Fe^{3+} + OH + OH^-$	1.3E-2		Løgager et al. (1992)
A11213_a01	TrAa01Fe	$FeO^{2+} + HO_2 \rightarrow Fe^{3+} + O_2 + OH^-$	2.0E6		Løgager et al. (1992)
A11214_a01	TrAa01Fe	$FeO^{2+} + OH \rightarrow Fe^{3+} + HO2^-$	1.0E7		Løgager et al. (1992)
A11215_a01	TrAa01Fe	$FeO^{2+} + Fe^{2+} \rightarrow 2\ Fe^{3+} + 2\ OH^-$	1.4E5		Løgager et al. (1992)
A11216_a01	TrAa01Fe	$FeO^{2+} + Fe^{2+} \rightarrow Fe(OH)_2Fe^{4+}$	1.8E4		Jacobsen et al. (1997)
A11217_a01	TrAa01Fe	$Fe(OH)_2Fe^{4+} + H^+ \rightarrow 2\ Fe^{3+} + OH^-$	2.0		Jacobsen et al. (1997)
A11218_a01	TrAa01Fe	$Fe(OH)_2Fe^{4+} \rightarrow 2\ Fe^{3+} + 2\ OH^-$	0.49		Jacobsen et al. (1997)
A11301_a01	TrAa01FeN	$FeO^{2+} + HONO \rightarrow Fe^{3+} + NO_2 + OH^-$	1.1E4		Jacobsen et al. (1998)
A11302_a01	TrAa01FeN	$Fe^{2+} + NO_3 \rightarrow Fe^{3+} + NO_3^-$	8.0E6		Herrmann et al. (2000)*
A11601_a01	TrAa01ClFe	$Fe^{2+} + Cl \rightarrow Fe^{3+} + Cl^-$	5.9E9		Jayson et al. (1973)
A11602a_a01	TrAa01ClFe	$Fe^{2+} + Cl_2^- \rightarrow Fe^{3+} + 2\ Cl^-$	1E7		Thornton and Laurence (1973)
A11602b_a01	TrAa01ClFe	$Fe^{2+} + Cl_2^- \rightarrow FeCl^{2+} + Cl^-$	4E6		Thornton and Laurence (1973)
A11603a_a01	TrAa01ClFe	$FeCl^+ + HO_2 \rightarrow Fe^{3+} + Cl^- + HO2^-$	1.2E6		de Laat and Le (2006)
A11603b_a01	TrAa01ClFe	$FeCl^+ + O_2^- \rightarrow Fe^{3+} + Cl^- + HO2^- + OH^-$	1E7		de Laat and Le (2006)
A11604a_a01	TrAa01ClFe	$FeCl^{2+} + HO_2 \rightarrow Fe^{2+} + Cl^- + O_2 + H^+$	2E4		de Laat and Le (2006)
A11604b_a01	TrAa01ClFe	$FeCl_2^+ + HO_2 \rightarrow Fe^{2+} + 2\ Cl^- + O_2 + H^+$	2E4		de Laat and Le (2006)
A11604c_a01	TrAa01ClFe	$FeCl^{2+} + O_2^- \rightarrow Fe^{2+} + Cl^- + O_2$	5E7		de Laat and Le (2006)
A11604d_a01	TrAa01ClFe	$FeCl_2^+ + O_2^- \rightarrow Fe^{2+} + 2\ Cl^- + O_2$	5E7		de Laat and Le (2006)
A11605_a01	TrAa01ClFe	$FeO^{2+} + Cl^- \rightarrow Fe^{3+} + Cl + 2\ OH^-$	1E2		Jacobsen et al. (1998)*
A11901_a01	TrAa01FeS	$FeO^{2+} + SO_2 \rightarrow Fe^{3+} + SO_3^-$	4.5E5		Jacobsen et al. (1998)*
A11902_a01	TrAa01FeS	$FeO^{2+} + HSO_3^- \rightarrow Fe^{3+} + SO_3^- + OH^-$	2.5E5		Jacobsen et al. (1998)*
A11903_a01	TrAa01FeS	$FeOH^{2+} + HSO_3^- \rightarrow Fe^{2+} + SO_3^- + H_2O$	30		Ziajka et al. (1994)

Table 6: Aqueous phase reactions (...continued)

#	labels	reaction	k_0 $[M^{1-n}s^{-1}]$	$-E_a/R[K]$	reference
A11904_a01	TrAa01FeS	Fe^{2+} + SO_5^- → $FeOH^{2+}$ + HSO_5^-	8E5		Ziajka et al. (1994)*
A11905_a01	TrAa01FeS	Fe^{2+} + HSO_5^- → $FeOH^{2+}$ + SO_4^-	3.0E4		Gilbert and Stell (1990)
A11906_a01	TrAa01FeS	Fe^{2+} + SO_4^- → $FeSO_4^+$	3.6E7		McElroy and Waygood (1990)*
A11907_a01	TrAa01FeS	$FeOH^{2+}$ + SO_3^- → Fe^{2+} + HSO_4^-	3E7		Warneck (2018)
A11908_a01	TrAa01FeS	$FeSO_3^+$ + SO_3^- → Fe^{2+} + SO_4^{2-} + SO_2	2.16E6		Warneck (2018)*
A3103_a01MK	TrAa01ScN	NO + NO + O_2 → NO_2 + NO_2	6.4E6		Pires et al. (1994)
A3104_a01MK	TrAa01ScN	NO_2 + NO_2 → N_2O_4	4.5E8		Mack and Bolton (1999)
A3105_a01MK	TrAa01ScN	NO + NO_2 → N_2O_3	1.1E9		Mack and Bolton (1999)
A3106_a01MK	TrAa01ScN	NO_2 + O_2^- → O_2 + NO_2^-	1.0E8		Ervens et al. (2003)
A3107_a01MK	TrAa01ScN	NO_2 + O_2^- → NO_4^-	4.5E9		Ervens et al. (2003)
A3109_a01MK	TrAa01ScN	N_2O_3 → 2 NO_2^- + 2 H^+	1.6E3		Licht et al. (1988)
A3205_a01MK	TrAa01ScN	NO + OH → HONO	1.0E10		Mack and Bolton (1999)
A3206_a01MK	TrAa01ScN	HONO + HONO → N_2O_3 + H_2O	5.6		Schwartz (1983)
A4111_a01MK	TrAa01Sc	CH_3OO + CH_3OO → HCHO + CH_3OH + HO_2	1.7E8	-2200	Ervens et al. (2004)
A4112_a01MK	TrAa01Sc	CH_3OH + OH → $CH_2(OH)_2$ + HO_2	1.0E9		Ervens et al. (2004)
A4113_a01MK	TrAa01Sc	CH_3OOH + OH → HCOOH + HO_2	6.0E6	-1715	Ervens et al. (2004)
A4114_a01MK	TrAa01Sc	$CH_2(OH)_2$ + OH → HCOOH + HO_2	1.0E9		Ervens et al. (2004)
A4120_a01MK	TrAa01AmiScN	MMA + N_2O_4 → 0.3 MMNNO2 + 0.7 CH_2NH_2 + 1.7 HONO	4.0E7		Challis and Kyrtopoulos (1979)
A4121_a01MK	TrAa01AmiScN	MMA + OH → 0.5 CH_2NH_2 + 0.5 $CH_3NH_2^+$ + 0.5 OH^- + H_2O	6.9E9		Simić et al. (1971)
A4122_a01MK	TrAa01AmiScN	CH_2NH_2 + O_2 → $CH_2NH_2^+$ + O_2^-	3.5E9		Duff et al. (2003)
A4123_a01MK	TrAa01AmiScN	$CH_2NH_2^+$ → NH_3 + HCHO + H^+	4.0		Duff et al. (2003)
A4124_a01MK	TrAa01AmiScN	CH_2NH_2 + CH_2NH_2 → $NH_2CH_2CH_2NH_2$	2.0E9		Simić et al. (1971)
A4125_a01MK	TrAa01AmiScN	$CH_3NH_2^+$ + $CH_3NH_2^+$ → $CH_3NHNHCH_3$ + 2 H^+	2.0E9		Simić et al. (1971)
A4126_a01MK	TrAa01AmiScN	MMA^+ + OH → 0.275 $CH_2NH_3^+$ + 0.725 $CH_3NH_2^+$ + H_2O	1.1E8		Simić et al. (1971)
A4127_a01MK	TrAa01AmiScN	HNCO → NH_3 + CO_2	7.83E-4		Roberts et al. (2011)
A4128_a01MK	TrAa01AmiScN	H2NCHO + OH → HNCO + H_2O	3.7E8		Munoz et al. (2000)
A4129_a01MK	TrAa01AScN	MMNNO2 + OH → NH_4^+ + HCHO + OH^- + NO_2	5.44E8		Fuller-Rowell (1993)
A4200_a01MK	TrAa01ScC	GLYOX + OH → $HCOCO_2H$ + HO_2	1.1E9	-1516	Ervens et al. (2004)
A4201_a01MK	TrAa01ScC	$HCOCO_2H$ + OH → OXALAC + HO_2 + 2 CO_2	3.6E8	-1000	Ervens et al. (2004)
A4202_a01MK	TrAa01ScC	$HCOCOO^-$ + OH → $HC_2O_4^-$ + HO_2 + 2 CO_2	2.9E9	-4300	Ervens et al. (2004)
A4203_a01MK	TrAa01ScC	$HOCH_2CHO$ + OH → GLYOX + HO_2	1.2E9		Ervens et al. (2004)

Table 6: Aqueous phase reactions (...continued)

#	labels	reaction	k_0 $[M^{1-n}s^{-1}]$	$-E_a/R[K]$	reference
A4204_a01MK	TrAa01ScC	$HOCH_2CO_2H + OH \rightarrow HCOCO_2H + HO_2$	1.2E9		Ervens et al. (2004)
A4205_a01MK	TrAa01Sc	$C_2O_4^{2-} + OH \rightarrow OH^- + O_2^-$	1.6E8	-4300	Ervens et al. (2004)
A4206_a01MK	TrAa01Sc	$HC_2O_4^- + OH \rightarrow OH^- + HO_2$	1.9E8	-2800	Ervens et al. (2004)
A4207_a01MK	TrAa01Sc	$OXALAC + OH \rightarrow H_2O + HO_2$	1.4E6		Ervens et al. (2004)
A4208_a01MK	TrAa01ScC	$CH_3COO^- + OH \rightarrow HO_2 + OH^- + CH_3COO_2$	1.0E8	-1800	Ervens et al. (2004)
A4209_a01MK	TrAa01ScC	$CH_3COOH + OH \rightarrow HO_2 + H_2O + CH_3COO_2$	1.5E7	-1330	Ervens et al. (2004)
A4210_a01MK	TrAa01AmiScCN	$DMA + N_2O_3 \rightarrow NDMA + HONO$	4.29E7		Mirvish (1975)
A4211_a01MK	TrAa01AmiScCN	$DMA + N_2O_4 \rightarrow 0.4\ DMNNO2 + 0.4\ NDMA + 0.2\ CH_3NHCH_2 + 0.6\ HONO + 0.6\ HNO_3$	4.0E7		Challis and Kyrtopoulos (1979)
A4212_a01MK	TrAa01AmiScCN	$DMA + HONO \rightarrow NDMA + H_2O$	0.1		Anastasio and Chu (2009)
A4213_a01MK	TrAa01AmiScCN	$DMA + NO_2^- \rightarrow NDMA + OH^-$	0.1		Anastasio and Chu (2009)
A4214_a01MK	TrAa01AmiScCN	$(CH_3)_2NH^+ + NO_2 \rightarrow DMNNO2 + H^+$	1.0E4		Fell et al. (1990)
A4215_a01MK	TrAa01AmiScCN	$DMA + OH \rightarrow 0.5\ CH_3NHCH_2 + 0.5\ (CH_3)_2NH^+ + 0.5\ OH^- + H_2O$	8.9E9		Lee et al. (2007a)
A4216_a01MK	TrAa01AmiScCN	$CH_3NHCH_2 + O_2 \rightarrow CH_3NH^+CH_2 + O_2^-$	3.5E9		Duff et al. (2003)
A4217_a01MK	TrAa01AmiScCN	$CH_3NH^+CH_2 \rightarrow MMA + HCHO + H^+$	4.0		Duff et al. (2003)
A4218_a01MK	TrAa01AmiScCN	$CH_3NHCH_2 + CH_3NHCH_2 \rightarrow DMA + CH_3NH^+CH_2 + OH^-$	2.0E9		Das and von Sonntag (1986)
A4219_a01MK	TrAa01AmiScCN	$CH_3NHCH_2 + (CH_3)_2NH^+ \rightarrow DMA + CH_3NH^+CH_2$	2.0E9		Das and von Sonntag (1986)
A4220_a01MK	TrAa01AmiScCN	$DMA^+ + OH \rightarrow 0.275\ CH_3NH_2^+CH_2 + 0.725\ (CH_3)_2NH^+ + H_2O$	6.0E7		Lee et al. (2007a)
A4221_a01MK	TrAa01AmiScCN	$MEA + N_2O_4 \rightarrow 0.3\ MEANNO2 + 0.7\ NH_2CH_2CHOH + 1.7\ HONO$	4.0E7		Challis and Kyrtopoulos (1979)
A4222_a01MK	TrAa01AmiScCN	$MEA + OH \rightarrow 0.5\ NH_2CH_2CHOH + 0.5\ HOCH_2CH_2NH_2^+ + 0.5\ OH^- + H_2O$	2.9E9		Kishore et al. (2004)
A4223_a01MK	TrAa01AmiScCN	$NH_2CH_2CHOH + OH \rightarrow CH_2NH_2 + HCHO$	1.0E9		Kishore et al. (2004)
A4224_a01MK	TrAa01AmiScCN	$MEA^+ + OH \rightarrow 0.275\ HOCHCH_2NH_3^+ + 0.725\ HOCH_2CH_2NH_2^+ + H_2O$	3.0E8		Kishore et al. (2004)
A4225_a01MK	TrAa01AmiScCN	$H2NCOCH2OH + OH \rightarrow HNCO + HCHO + HO_2$	1.9E8		Roble (1995)
A4226_a01MK	TrAa01AmiScCN	$CH_3NHCHO + OH \rightarrow HNCO + HCHO + HO_2$	1.2E9		Roble (1995)
A4227_a01MK	TrAa01AmiScCN	$NDMA + OH \rightarrow MMA^+ + HCHO + OH^- + NO$	4.5E8		Lee et al. (2007b)
A4228_a01MK	TrAa01AmiScCN	$DMNNO2 + OH \rightarrow MMA^+ + HCHO + OH^- + NO_2$	5.44E8		Fuller-Rowell (1993)

Table 6: Aqueous phase reactions (...continued)

#	labels	reaction	k_0 $[M^{1-n}s^{-1}]$	$-E_a/R[K]$	reference
A4229_a01MK	TrAa01AmiScCN	MEANNO + OH → $CH_2NH_3^+$ + HCHO + OH^- + NO	4.5E8		Lee et al. (2007b)
A4230_a01MK	TrAa01AmiScCN	MEANNO2 + OH → $CH_2NH_3^+$ + HCHO + OH^- + NO_2	5.44E8		Fuller-Rowell (1993)
A4231_a01MK	TrAa01AmiScCN	CH_3NCO → MMA + CO_2	7.83E-4		Roberts et al. (2011)
A4300_a01MK	TrAa01ScC	MGLYOX + OH → MGLYOAC + HO_2	1.1E9	-1600	Ervens et al. (2004)
A4301_a01MK	TrAa01ScC	CH_3COCOO^- + OH → CH_3COO_2 + HO_2 + CO_2 + O_2^-	7.0E8		Ervens et al. (2004)
A4302_a01MK	TrAa01ScC	MGLYOAC + OH → CH_3CHO + HO_2 + H_2O + CO_2	1.2E8		Ervens et al. (2004)
A4303_a01MK	TrAa01Sc	DOC + OH → DOCO + HO_2	4.1E8		Barth (1992)
A4304_a01MK	TrAa01Sc	MALONAC + OH → OXALAC + HO_2	5.0E7		Ervens et al. (2004)
A4310_a01MK	TrAa01AmiScCN	TMA + N_2O_3 → 0.075 NDMA + 1.9 HONO + 0.95 $(CH_3)_2NCH_2$	4.0E7		Challis and Kyrtopoulos (1979)
A4311_a01MK	TrAa01AmiScCN	TMA + N_2O_4 → 0.05 DMNNO2 + 0.1 NDMA + 0.90 $(CH_3)_2NCH_2$ + 0.9 HONO + 0.9 HNO_3	4.0E7		Challis and Kyrtopoulos (1979)
A4312_a01MK	TrAa01AmiScCN	TMA + OH → 0.5 $(CH_3)_2NCH_2$ + 0.5 $(CH_3)_3N^+$ + 0.5 OH^- + H_2O	1.2E10		Das and von Sonntag (1986)
A4313_a01MK	TrAa01AmiScCN	$(CH_3)_2NCH_2$ + O_2 → $(CH_3)_2N^+CH_2$ + O_2^-	3.5E9		Duff et al. (2003)
A4314_a01MK	TrAa01AmiScCN	$(CH_3)_2N^+CH_2$ → DMA + HCHO + H^+	4.0		Duff et al. (2003)
A4315_a01MK	TrAa01AmiScCN	$(CH_3)_2NCH_2$ + $(CH_3)_2NCH_2$ → TMA + $(CH_3)_2N^+CH_2$ + OH^-	2.0E9		Das and von Sonntag (1986)
A4316_a01MK	TrAa01AmiScCN	$(CH_3)_2NCH_2$ + $(CH_3)_3N^+$ → TMA + $(CH_3)_2N^+CH_2$	2.0E9		Das and von Sonntag (1986)
A4317_a01MK	TrAa01AmiScCN	TMA^+ + OH → 0.275 $(CH_3)_2NH^+CH_2$ + 0.725 $(CH_3)_3N^+$ + H_2O	4.0E8		Das and von Sonntag (1986)
A4318_a01MK	TrAa01AmiScCN	DMNCHO + OH → CH_3NCO + HCHO + HO_2	1.7E9		Roble (1995)
A4400_a01MK	TrAa01Sc	SUCCAC + OH → MALONAC + HO_2	5.0E7		Ervens et al. (2004)
A4410_a01MK	TrAa01AmiScCN	DEA + N_2O_3 → NDELA + HONO	1.0E7		Mitch and Herckes (2012)
A4411_a01MK	TrAa01AmiScCN	DEA + N_2O_4 → 0.4 DEANNO2 + 0.4 NDELA + 0.2 $HOETNHCH_2CHOH$ + 0.6 HONO + 0.6 HNO_3	4.0E7		Challis and Kyrtopoulos (1979)
A4412_a01MK	TrAa01AmiScCN	DEA + HONO → NDELA + H_2O	0.1		Anastasio and Chu (2009)
A4413_a01MK	TrAa01AmiScCN	DEA + NO_2^- → NDELA + OH^-	0.1		Anastasio and Chu (2009)
A4414_a01MK	TrAa01AmiScCN	$(HOET)_2NH^+$ + NO_2 → DEANNO2 + H^+	0.0		Fell et al. (1990)

Table 6: Aqueous phase reactions (...continued)

#	labels	reaction	k_0 $[M^{1-n}s^{-1}]$	$-E_a/R[K]$	reference
A4415_a01MK	TrAa01AmiScCN	DEA + OH → 0.5 $HOETNHCH_2CHOH$ + 0.5 $(HOET)_2NH^+$ + 0.5 OH^- + H_2O	8.0E9		Kishore et al. (2004)
A4416_a01MK	TrAa01AmiScCN	$HOETNHCH_2CHOH$ + OH → MEA + 2 HCHO	1.0E9		Kishore et al. (2004)
A4417_a01MK	TrAa01AmiScCN	DEA^+ + OH → 0.275 $HOETNH_2CH_2CHOH^+$ + 0.725 $(HOET)_2NH^+$ + H_2O	4.5E8		Kishore et al. (2004)
A4418_a01MK	TrAa01AmiScCN	NDELA + OH → MEA^+ + 2 HCHO + OH^- + NO	6.99E8		Fuller-Rowell (1993)
A4419_a01MK	TrAa01AmiScCN	DEANNO2 + OH → MEA^+ + 2 HCHO + OH^-+ NO_2	8.67E8		Fuller-Rowell (1993)
A4500_a01MK	TrAa01Sc	GLUTARAC + OH → SUCCAC + HO_2	1.0E7		Ervens et al. (2004)
A4600_a01MK	TrAa01Sc	ADIPAC + OH → GLUTARAC + HO_2	1.0E7		Ervens et al. (2004)
A4610_a01MK	TrAa01AmiScCN	TEA + OH → 0.5 $DENCH_2CHOH$ + 0.5 $(HOET)_3N^+$ + 0.5 OH^- + H_2O	2.0E10		Kishore et al. (2004)
A4611_a01MK	TrAa01AmiScCN	$DENCH_2CHOH$ + OH → $(HOET)_2N^+CH_2CH_2OH$ + OH^-	3.18E9		Das and von Sonntag (1986)
A4612_a01MK	TrAa01AmiScCN	$(HOET)_2N^+CH_2CH_2OH$ → DEA + $HOCH_2CHO$ + H^+	3.0E6		Das and von Sonntag (1986)
A4613_a01MK	TrAa01AmiScCN	$(HOET)_3N^+$ + $DENCH_2CHOH$ → TEA + $(HOET)_2N^+CH_2CH_2OH$	2.0E9		Das and von Sonntag (1986)
A4614_a01MK	TrAa01AmiScCN	TEA^+ + OH → 0.275 $(HOET)_2NH^+CH_2CHOH$ + 0.725 $(HOET)_3N^+$ + H_2O	2.0E9		Kishore et al. (2004)
A4615_a01MK	TrAa01AmiScCN	TEA + N_2O_3 → 0.075 NDELA + 0.95 $DENCH_2CHOH$ + 1.9 HONO	2.9E5		Mitch and Herckes (2012)
A4616_a01MK	TrAa01AmiScCN	TEA + N_2O_4 → 0.05 DEANNO2 + 0.1 NDELA + 0.9 $DENCH_2CHOH$ + 0.9 HONO + 0.9 HNO_3	4.0E7		Challis and Kyrtopoulos (1979)
A9405_a01MK	TrAa01MblScS	HSO_3^- + $CH_2(OH)_2$ → $CH_2OHSO_3^-$	0.436	-2990	Ervens et al. (2004)
A9406a_a01MK	TrAa01MblScS	$CH_2OHSO_3^-$ → $CH_2(OH)_2$ + HSO_3^-	1.22E-7		Ervens et al. (2004)
A9406b_a01MK	TrAa01MblScS	$CH_2OHSO_3^-$ → $CH_2(OH)_2$ + SO_3^{2-} + H^+	3.8E-6	-5530	Ervens et al. (2004)
A9407_a01MK	TrAa01MblScS	SO_3^{2-} + $CH_2(OH)_2$ → $CH_2OHSO_3^-$ + O_2^-	1.36E5	-2450	Ervens et al. (2004)
A9408_a01MK	TrAa01MblScS	$CH_2OHSO_3^-$ + OH → HCOOH + HO_2 + HSO_3^-	3.0E8		Ervens et al. (2004)
A9409_a01MK	TrAa01MblScCS	DMS + O_3 → DMSO	8.61E8	-2600	Hoffmann et al. (2016)
A9410_a01MK	TrAa01MblScCS	DMS + OH → DMSO + HO_2	1.9E10		Hoffmann et al. (2016)
A9411_a01MK	TrAa01MblScCS	DMSO + OH → MSIA + HO_2 + CO_2	6.65E9	-1270	Hoffmann et al. (2016)
A9412_a01MK	TrAa01MblScS	MSIA + O_3 → $CH_3SO_3^-$ + H^+	3.5E7		Hoffmann et al. (2016)
A9413_a01MK	TrAa01MblScS	MSIA + OH → $CH_3SO_3^-$ + H^+ + HO_2	6.0E9		Hoffmann et al. (2016)

Specific notes

A6102_a01: Jacobi (1996) found an upper limit of 6E9 and cite an upper limit from another study of 2E9. Here, we set the rate coefficient to 1E9.

A6301_a01: There is also an earlier study by Exner et al. (1992) which found a smaller rate coefficient but did not consider the back reaction.

A9105_a01: The rate coefficient for the sum of the paths (leading to either HSO_5^- or SO_4^{2-}) is from Huie and Neta (1987), the ratio 0.28/0.72 is from Deister and Warneck (1990).

A9106_a01: See also: (Huie and Neta, 1987; Warneck, 1991). If this reaction produces a lot of SO_4^-, it will have an effect. However, we currently assume only the stable $S_2O_8^{2-}$ as product. Since $S_2O_8^{2-}$ is not treated explicitly in the mechanism, SO_4^{2-} is used as a proxy and the second sulfur atom is put into the lumped LSULFUR.

A9205_a01: D. Sedlak, pers. comm. (1993).

A9208_a01: D. Sedlak, pers. comm. (1993).

A9400_a01: Product $2.48 \times 10^7 \times 5.5 \times 10^{-4}$ considering the hydrated form of HCHO.

A9402_a01: Product $790 \times 5.5 \times 10^{-4}$ considering the hydrated form of HCHO.

A9605_a01: Assumed to be the same as for SO_3^{2-} + HOCl.

A11302_a01: value from Pikaev et al. (1974)

A11605_a01: products assumed

A11901_a01: products assumed

A11902_a01: products assumed

A11904_a01: Assumed. Note that CAPRAM 2.4 lists k=4.3E7 from Herrmann Air Pollution Research Report 57 and it also lists k= 2.65E7 from Williams PhD 1996 http://lib.leeds.ac.uk/record=b1835184~S5. Brand and van Eldik (1995) also list k=3.56E4 from Waygood EUROTRAC 1992 report.

A11906_a01: 3E8*6500/(48000+6500)

A11908_a01: Assuming that the intermediate $S_2O_6^{2-}$ dissociates quickly.

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