Jozef Peeters

List of Publications by Year in Descending Order

134 papers	5,854 citations	47 h-index	69 g-index
138 ext. papers	6,233 ext. citations	3.7 avg, IF	5.72 L-index

#	Paper	IF	Citations
134	The CH(X) + HO reaction: two transition state kinetics. *Physical Chemistry Chemical Physics*, **2021**, 23, 16142-16149	3.6	0
133	Chemistry and deposition in the Model of Atmospheric composition at Global and Regional scales using Inversion Techniques for Trace gas Emissions (MAGRITTE v1.1) Part 1: Chemical mechanism. *Geoscientific Model Development*, **2019**, 12, 2307-2356	6.3	19
132	The photolysis of hydroperoxycarbonyls. *Physical Chemistry Chemical Physics*, **2018**, 20, 6970-6979	3.6	10
131	Theoretically derived mechanisms of HPALD photolysis in isoprene oxidation. *Physical Chemistry Chemical Physics*, **2017**, 19, 9096-9106	3.6	11
130	The reaction of methyl peroxy and hydroxyl radicals as a major source of atmospheric methanol. *Nature Communications*, **2016**, 7, 13213	17.4	54
129	Atmospheric Vinyl Alcohol to Acetaldehyde Tautomerization Revisited. *Journal of Physical Chemistry Letters*, **2015**, 6, 4005-11	6.4	16
128	Fast (E)-(Z) Isomerization Mechanisms of Substituted Allyloxy Radicals in Isoprene Oxidation. *Journal of Physical Chemistry A*, **2015**, 119, 7270-6	2.8	7
127	Hydroxyl radical recycling in isoprene oxidation driven by hydrogen bonding and hydrogen tunneling: the upgraded LIM1 mechanism. *Journal of Physical Chemistry A*, **2014**, 118, 8625-43	2.8	160
126	Absolute rate coefficient of the gas-phase reaction between hydroxyl radical (OH) and hydroxyacetone: investigating the effects of temperature and pressure. *Journal of Physical Chemistry A*, **2013**, 117, 12208-15	2.8	9
125	Experimental and theoretical study of the reaction of the ethynyl radical with nitrous oxide, C2H + N2O. *Physical Chemistry Chemical Physics*, **2012**, 14, 7456-70	3.6	6
124	Unusually fast 1,6-h shifts of enolic hydrogens in peroxy radicals: formation of the first-generation C2 and C3 carbonyls in the oxidation of isoprene. *Journal of Physical Chemistry A*, **2012**, 116, 6134-41	2.8	29
123	A theoretical study of the OH-initiated gas-phase oxidation mechanism of pinene (C10H16): first generation products. *Physical Chemistry Chemical Physics*, **2012**, 14, 3802-15	3.6	44
122	Theoretical and experimental investigation of the C2H + SO2 reaction over the range T = 295800 K. *Chemical Physics Letters*, **2011**, 513, 201-207	2.5	5
121	Ozone Based Atomic Layer Deposition of Hafnium Oxide and Impact of Nitrogen Oxide Species. *Journal of the Electrochemical Society*, **2011**, 158, D259	3.9	16
120	Ozone-Based Metal Oxide Atomic Layer Deposition: Impact of N[sub 2]/O[sub 2] Supply Ratio in Ozone Generation. *Electrochemical and Solid-State Letters*, **2010**, 13, H176		14
119	Theoretical Study of the HOCH2OO + HO2 Reaction: Detailed Molecular Mechanisms of the Three Reaction Channels. *Zeitschrift Fur Physikalische Chemie*, **2010**, 224, 1081-1093	3.1	7
118	HO(x) radical regeneration in isoprene oxidation via peroxy radical isomerisations. II: experimental evidence and global impact. *Physical Chemistry Chemical Physics*, **2010**, 12, 14227-35	3.6	159

117	Kinetics of O(1D) + H2O and O(1D) + H2: absolute rate coefficients and O(3P) yields between 227 and 453 K. *Physical Chemistry Chemical Physics*, **2010**, 12, 9213-21	3.6	25
116	A structure-activity relationship for the rate coefficient of H-migration in substituted alkoxy radicals. *Physical Chemistry Chemical Physics*, **2010**, 12, 12608-20	3.6	74
115	Evaluation of a detailed model of secondary organic aerosol formation from □pinene against dark ozonolysis experiments. *Atmospheric Environment*, **2010**, 44, 5434-5442	5.3	12
114	HOx regeneration in the oxidation of isoprene III: theoretical study of the key isomerisation of the Z-□hydroxy-peroxy isoprene radicals. *ChemPhysChem*, **2010**, 11, 3996-4001	3.2	34
113	Mechanism of the catalytic deperoxidation of tert-butylhydroperoxide with cobalt(II) acetylacetonate. *Chemistry - A European Journal*, **2010**, 16, 13226-35	4.8	60
112	Decomposition of substituted alkoxy radicals--part I: a generalized structure-activity relationship for reaction barrier heights. *Physical Chemistry Chemical Physics*, **2009**, 11, 9062-74	3.6	113
111	Theoretical study of the gas-phase ozonolysis of beta-pinene (C10H16). *Physical Chemistry Chemical Physics*, **2009**, 11, 5643-56	3.6	76
110	Absolute rate coefficients over extended temperature ranges and mechanisms of the CF(X(2)Pi) reactions with F(2), Cl(2) and O(2). *Physical Chemistry Chemical Physics*, **2009**, 11, 4319-25	3.6	4
109	HOx radical regeneration in the oxidation of isoprene. *Physical Chemistry Chemical Physics*, **2009**, 11, 5935-9	3.6	384
108	The gas-phase ozonolysis of beta-caryophyllene (C(15)H(24)). Part I: an experimental study. *Physical Chemistry Chemical Physics*, **2009**, 11, 4152-72	3.6	111
107	The gas-phase ozonolysis of beta-caryophyllene (C(15)H(24)). Part II: A theoretical study. *Physical Chemistry Chemical Physics*, **2009**, 11, 4173-83	3.6	58
106	A temperature dependence kinetic study of O(1D) + CH4: overall rate coefficient and product yields. *Physical Chemistry Chemical Physics*, **2008**, 10, 5714-22	3.6	21
105	Pronounced non-Arrhenius behaviour of hydrogen-abstractions from toluene and derivatives by phthalimide-N-oxyl radicals: a theoretical study. *Physical Chemistry Chemical Physics*, **2008**, 10, 1125-32	3.6	29
104	Autoxidation Chemistry: Bridging the Gap Between Homogeneous Radical Chemistry and (Heterogeneous) Catalysis. *Topics in Catalysis*, **2008**, 48, 41-48	2.3	34
103	Autoxidation of Hydrocarbons: From Chemistry to Catalysis. *Topics in Catalysis*, **2008**, 50, 124-132	2.3	86
102	Origin of byproducts during the catalytic autoxidation of cyclohexane. *Journal of Physical Chemistry A*, **2008**, 112, 1747-53	2.8	26
101	Silica-supported chromium oxide: colloids as building blocks. *Physical Chemistry Chemical Physics*, **2007**, 9, 5382-6	3.6	7
100	Diazo chemistry controlling the selectivity of olefin ketonisation by nitrous oxide. *Physical Chemistry Chemical Physics*, **2007**, 9, 4269-74	3.6	26

99	Low-volatility poly-oxygenates in the OH-initiated atmospheric oxidation of alpha-pinene: impact of non-traditional peroxyl radical chemistry. *Physical Chemistry Chemical Physics*, **2007**, 9, 5241-8	3.6	91
98	Quantum chemical and statistical rate investigation of the CF2(a3B1)+NO(X2Pi) reaction: a fast chemical quenching process. *Journal of Physical Chemistry A*, **2007**, 111, 6628-36	2.8	1
97	Theoretical reinvestigation of the O(3P) + C6H6 reaction: quantum chemical and statistical rate calculations. *Journal of Physical Chemistry A*, **2007**, 111, 3836-49	2.8	30
96	Designing a Novel Heterogeneous Catalytic System for Cyclohexane Autoxidation. *Studies in Surface Science and Catalysis*, **2007**, 373-376	1.8	
95	The formation of byproducts in the autoxidation of cyclohexane. *Chemistry - A European Journal*, **2007**, 13, 754-61	4.8	68
94	Solvent- and Metal-Free Ketonization of Fatty Acid Methyl Esters and Triacylglycerols with Nitrous Oxide. *Advanced Synthesis and Catalysis*, **2007**, 349, 1604-1608	5.6	28
93	Mechanism of thermal toluene autoxidation. *ChemPhysChem*, **2007**, 8, 2678-88	3.2	51
92	Silica-immobilized N-hydroxyphthalimide: An efficient heterogeneous autoxidation catalyst. *Journal of Catalysis*, **2007**, 251, 204-212	7.3	52
91	Reaction of HO with glycolaldehyde, HOCH2CHO: rate coefficients (240-362 K) and mechanism. *Journal of Physical Chemistry A*, **2007**, 111, 897-908	2.8	28
90	Autoxidation of ethylbenzene: the mechanism elucidated. *Journal of Organic Chemistry*, **2007**, 72, 3057-64	4.2	83
89	Kinetic parameters for gas-phase reactions: experimental and theoretical challenges. *Physical Chemistry Chemical Physics*, **2007**, 9, 4071-84	3.6	15
88	Autoxidation catalysis with N-hydroxyimides: more-reactive radicals or just more radicals?. *Physical Chemistry Chemical Physics*, **2007**, 9, 686-90	3.6	52
87	Structure-activity relationship for the addition of OH to (poly)alkenes: site-specific and total rate constants. *Journal of Physical Chemistry A*, **2007**, 111, 1618-31	2.8	89
86	To the core of autocatalysis in cyclohexane autoxidation. *Chemistry - A European Journal*, **2006**, 12, 4229-48	4.8	127
85	Silica-immobilized chromium colloids for cyclohexane autoxidation. *Angewandte Chemie - International Edition*, **2006**, 45, 7584-8	16.4	26
84	Enhanced activity and selectivity in cyclohexane autoxidation by inert H-bond acceptor catalysts. *ChemPhysChem*, **2006**, 7, 1142-8	3.2	19
83	Silica-Immobilized Chromium Colloids for Cyclohexane Autoxidation. *Angewandte Chemie*, **2006**, 118, 7746-7750	3.6	4
82	Reaction of HO with hydroxyacetone (HOCH2C(O)CH3): rate coefficients (233-363 K) and mechanism. *Physical Chemistry Chemical Physics*, **2006**, 8, 236-46	3.6	42

81	Quantum chemical and statistical rate study of the reaction of O(3P) with allene: O-addition and H-abstraction channels. *Journal of Physical Chemistry A*, **2006**, 110, 12166-76	2.8	24
80	Absolute rate coefficient of the OH + CH(3)C(O)OH reaction at T = 287-802 K. The two faces of pre-reactive H-bonding. *Journal of Physical Chemistry A*, **2006**, 110, 12852-9	2.8	26
79	Quantum chemical and theoretical kinetics study of the O(3P) + C2H2 reaction: a multistate process. *Journal of Physical Chemistry A*, **2006**, 110, 6696-706	2.8	44
78	Theoretical study of the blue-shifting hydrogen bonds between CH2X2 and CHX3 (X=F, Cl, Br) and hydrogen peroxide. *Journal of Molecular Structure*, **2006**, 792-793, 16-22	3.4	20
77	Understanding the autoxidation of hydrocarbons at the molecular level and consequences for catalysis. *Journal of Molecular Catalysis A*, **2006**, 251, 221-228		80
76	Kinetics of alpha-hydroxy-alkylperoxyl radicals in oxidation processes. HO2*-initiated oxidation of ketones/aldehydes near the tropopause. *Journal of Physical Chemistry A*, **2005**, 109, 4303-11	2.8	81
75	The kinetics of the CF3 + CF3 and CF3 + F combination reactions at 290 K and at He-pressures of approximately 1-6 Torr. *Physical Chemistry Chemical Physics*, **2005**, 7, 1187-93	3.6	7
74	Experimental and theoretical studies of the C2F4 + O reaction: nonadiabatic reaction mechanism. *Journal of Physical Chemistry A*, **2005**, 109, 9786-94	2.8	7
73	Theoretical and experimental study of the product branching in the reaction of acetic acid with OH radicals. *Journal of Physical Chemistry A*, **2005**, 109, 2401-9	2.8	42
72	Potential energy surfaces, product distributions and thermal rate coefficients of the reaction of O(3P) with C2H4(X1Ag): a comprehensive theoretical study. *Journal of Physical Chemistry A*, **2005**, 109, 7489-99	2.8	87
71	Pulsed laser photolysis and quantum chemical-statistical rate study of the reaction of the ethynyl radical with water vapor. *Journal of Chemical Physics*, **2005**, 122, 114307	3.9	28
70	Energetics and chemical bonding of the 1,3,5-tridehydrobenzene triradical and its protonated form. *Chemical Physics*, **2005**, 316, 125-140	2.3	24
69	Theoretical study of the kinetics of hydrogen abstraction in reactions of simple hydrogen compounds with triplet difluorocarbene. *Chemical Physics Letters*, **2005**, 402, 460-467	2.5	5
68	Quantum chemical study of the electronic structure of the 1-methylene-3,5-didehydrobenzene triradical (C7H5). *Chemical Physics Letters*, **2005**, 404, 150-155	2.5	4
67	Autoxidation of cyclohexane: conventional views challenged by theory and experiment. *ChemPhysChem*, **2005**, 6, 637-45	3.2	106
66	A Generalized Structure-Activity Relationship for the Decomposition of (Substituted) Alkoxy Radicals. *Journal of Atmospheric Chemistry*, **2004**, 48, 59-80	3.2	62
65	Direct ab initio dynamics studies of the reactions of HNO with H and OH radicals. *Chemical Physics Letters*, **2004**, 388, 94-99	2.5	19
64	Computational study of the stability of β-hydroperoxyl- or β-alkylperoxyl substituted alkyl radicals. *Chemical Physics Letters*, **2004**, 393, 432-436	2.5	81

63	Absolute rate coefficients of the reactions of CF2(□B1) with C3H8, C3H6, iso-C4H8 and C3H4 between 295 and 550 K. *Physical Chemistry Chemical Physics*, **2004**, 6, 2211-2215	3.6	5
62	Theoretical study of the reaction of the ethynyl radical with ammonia (C2H + NH3): hydrogen abstraction versus condensation. *Physical Chemistry Chemical Physics*, **2004**, 6, 4111	3.6	13
61	Tropopause chemistry revisited: HO2*-initiated oxidation as an efficient acetone sink. *Journal of the American Chemical Society*, **2004**, 126, 9908-9	16.4	34
60	Nontraditional (Per)oxy Ring-Closure Paths in the Atmospheric Oxidation of Isoprene and Monoterpenes. *Journal of Physical Chemistry A*, **2004**, 108, 5197-5204	2.8	48
59	The 5-Dehydro-m-xylylene Triradical and Its Nitrogen and Phosphorus Derivatives:□Open-Shell Doublet versus Quartet Ground State. *Journal of Physical Chemistry A*, **2004**, 108, 8411-8418	2.8	11
58	Use of DFT-Based Reactivity Descriptors for Rationalizing Radical Reactions: A Critical Analysis. *Journal of Physical Chemistry A*, **2004**, 108, 484-489	2.8	36
57	No Barrier for the Gas-Phase C2H + NH3 Reaction. *Journal of Physical Chemistry A*, **2004**, 108, 3695-3698	2.8	20
56	Theoretical Study of the Interaction between Methyl Fluoride, Methyl Chloride, and Methyl Bromide with Hydrogen Peroxide. *Journal of Physical Chemistry A*, **2004**, 108, 11101-11108	2.8	28
55	Mechanism of the catalytic oxidation of hydrocarbons by N-hydroxyphthalimide: a theoretical study. *Chemical Communications*, **2004**, 1140-1	5.8	63
54	The reaction of acetaldehyde and propionaldehyde with hydroxyl radicals: experimental determination of the primary H2O yield at room temperature. *Journal of Photochemistry and Photobiology A: Chemistry*, **2003**, 157, 269-274	4.7	17
53	Temperature and pressure dependent product distribution of the addition of CN radicals to C2H4. *Physical Chemistry Chemical Physics*, **2003**, 5, 5070	3.6	12
52	Absolute rate coefficients of the reactions of CF2(□B1) with NO and H2 between 287 K and 600 K. *Physical Chemistry Chemical Physics*, **2003**, 5, 5405-5408	3.6	5
51	Reactions of chemically activated C9H9 species II: The reaction of phenyl radicals with allene and cyclopropene, and of benzyl radicals with acetylene. *Physical Chemistry Chemical Physics*, **2003**, 5, 2807	3.6	55
50	Absolute rate coefficient of the reaction of CF2(□B1) with O2 between 288 and 600 K. *Physical Chemistry Chemical Physics*, **2003**, 5, 2376-2380	3.6	5
49	The 1,5-H-shift in 1-butoxy: A case study in the rigorous implementation of transition state theory for a multirotamer system. *Journal of Chemical Physics*, **2003**, 119, 5159-5170	3.9	91
48	An experimental and theoretical study of the reaction of ethynyl radicals with nitrogen dioxide (HC?C+NO2). *Journal of Chemical Physics*, **2003**, 118, 10996-11008	3.9	24
47	A combined crossed-beam, ab initio, and RiceRamspergerKasselMarcus investigation of the reaction of carbon atoms C(3Pj) with benzene, C6H6(X 1A1g) and d6-benzene, C6D6(X 1A1g). *Journal of Chemical Physics*, **2002**, 116, 3248-3262	3.9	36
46	The reaction of C2H with H2: Absolute rate coefficient measurements and ab initio study. *Journal of Chemical Physics*, **2002**, 116, 3700-3709	3.9	36

45	Reaction of phenyl radicals with propyne. *Journal of the American Chemical Society*, **2002**, 124, 2781-9	16.4	61
44	The OH-initiated atmospheric oxidation of pinonaldehyde: Detailed theoretical study and mechanism construction. *Physical Chemistry Chemical Physics*, **2002**, 4, 5795-5805	3.6	43
43	Absolute Rate Coefficient of the HCCO + NO Reaction over the RangeT= 297802 K. *Journal of Physical Chemistry A*, **2002**, 106, 12242-12247	2.8	25
42	The acetic acid forming channel in the acetone + OH reaction: A combined experimental and theoretical investigation. *Physical Chemistry Chemical Physics*, **2002**, 4, 461-466	3.6	47
41	Enhanced H-atom abstraction from pinonaldehyde, pinonic acid, pinic acid, and related compounds: theoretical study of CH bond strengths. *Physical Chemistry Chemical Physics*, **2002**, 4, 467-472	3.6	43
40	Reactions of chemically activated C9H9 species.. *Physical Chemistry Chemical Physics*, **2002**, 4, 2019-2027	3.6	39
39	H-atom abstraction by OH-radicals from (biogenic) (poly)alkenes: CH bond strengths and abstraction rates. *Chemical Physics Letters*, **2001**, 333, 162-168	2.5	39
38	NOx reduction by reburning: theoretical study of the branching ratio of the HCCO+NO reaction. *Chemical Physics Letters*, **2001**, 344, 400-406	2.5	38
37	The reaction of C2H radicals with C2H6: Absolute rate coefficient measurements for T= 295800 K, and quantum chemical study of the molecular mechanism. *Physical Chemistry Chemical Physics*, **2001**, 3, 3070-3074	3.6	27
36	Laser-induced fluorescence of nascent CH from ultraviolet photodissociation of HCCO and the absolute rate coefficient of the HCCO+O2 reaction over the range T=296839 K. *Journal of Chemical Physics*, **2001**, 114, 10332-10341	3.9	24
35	The detailed mechanism of the OH-initiated atmospheric oxidation of pinene: a theoretical study. *Physical Chemistry Chemical Physics*, **2001**, 3, 5489-5504	3.6	107
34	Experimental and theoretical study of the gas phase reaction of ethynyl radical with methane (HC?C+CH4). *Chemical Physics Letters*, **2000**, 329, 412-420	2.5	27
33	Experimental and theoretical study of the reaction of the ethynyl radical with acetylene (HC?C+HC?CH). *Chemical Physics*, **2000**, 262, 243-252	2.3	40
32	Electronic Structure Calculations on the Reaction of Vinyl Radical with Nitric Oxide. *Journal of Physical Chemistry A*, **2000**, 104, 1905-1914	2.8	12
31	Theoretical Study of the Formation of Acetone in the OH-Initiated Atmospheric Oxidation of Pinene. *Journal of Physical Chemistry A*, **2000**, 104, 11140-11146	2.8	42
30	The Atmospheric Chemistry of the Acetonoxy Radical. *Journal of Physical Chemistry A*, **2000**, 104, 11578-11588	2.8	65
29	Theoretical Investigation of the Role of Intramolecular Hydrogen Bonding in Hydroxyethoxy and Hydroxyethylperoxy Radicals in the Tropospheric Oxidation of Ethene. *Journal of Physical Chemistry A*, **1999**, 103, 1768-1775	2.8	63
28	H-atom abstraction in reactions of cyclic polyalkenes with OH. *Chemosphere*, **1999**, 38, 1189-1195	8.4	34

27	Decomposition of β-Hydroxypropoxy Radicals in the OH-Initiated Oxidation of Propene. A Theoretical and Experimental Study. *Journal of Physical Chemistry A*, **1999**, 103, 4693-4702	2.8	50
26	Detailed Microvariational RRKM Master Equation Analysis of the Product Distribution of the C2H2 + CH(X2Π) Reaction over Extended Temperature and Pressure Ranges. *Journal of Physical Chemistry A*, **1999**, 103, 5523-5533	2.8	34
25	Theoretical studies on C2H+NO reactions: mechanism for HCN+CO and HCO+CN formation. *Chemical Physics Letters*, **1998**, 283, 91-96	2.5	21
24	Theoretical studies on the C2H+O2 reaction: mechanism for HCO+CO, HCCO+O and CH+CO2 formation. *Chemical Physics Letters*, **1998**, 287, 109-118	2.5	27
23	Reactions of Cl Atoms with Selected VOCs: Kinetics, Products and Mechanisms. *Journal of Atmospheric Chemistry*, **1998**, 31, 247-267	3.2	48
22	Theoretical analysis of reactions related to the HNO2 energy surface: OH + NO and H + NO2. *Chemical Physics*, **1998**, 230, 1-11	2.3	30
21	Determination of the rate constants for the gas-phase reactions of methyl butenol with OH radicals, ozone, NO3 radicals, and Cl atoms. *International Journal of Chemical Kinetics*, **1998**, 30, 589-594	1.4	46
20	Laboratory and Theoretical Study of the Oxy Radicals in the OH- and Cl-Initiated Oxidation of Ethene. *Journal of Physical Chemistry A*, **1998**, 102, 8116-8123	2.8	135
19	B3LYP-DFT characterization of the potential energy surface of the CH(X 2Π)+C2H2 reaction. *Journal of Chemical Physics*, **1998**, 108, 1068-1080	3.9	70
18	Stochastic simulation of chemically activated unimolecular reactions. *Journal of Chemical Physics*, **1997**, 106, 6564-6573	3.9	72
17	Direct Identification of the C2H(X2Σ+) + O(3P) → CH(A2Δ) + CO Reaction as the Source of the CH(A2Δ→X2Π) Chemiluminescence in C2H2/O/H Atomic Flames. *Journal of Physical Chemistry A*, **1997**, 101, 2546-2551	2.8	50
16	Cycloaddition Reactions of CF with Unsaturated Hydrocarbons. Correlation of Activation Energies and Frequency Factors with the Hydrocarbon Ionization Energies. *Journal of Physical Chemistry A*, **1997**, 101, 787-793	2.8	11
15	Identification of the Sequence CH(2Π) + C2H2 → C3H2 + H (and C3H + H2) Followed by C3H2 + O → C2H + HCO (or H + CO) as C2H Source in C2H2/O/H Atomic Flames. *The Journal of Physical Chemistry*, **1996**, 100, 998-1007		71
14	Absolute Rate Coefficients of the Reactions of C2H with NO and H2 between 295 and 440 K. *The Journal of Physical Chemistry*, **1996**, 100, 15124-15129		50
13	Reaction of Isocyanic Acid and Hydrogen Atom (H + HNCO): Theoretical Characterization. *The Journal of Physical Chemistry*, **1996**, 100, 1615-1621		37
12	Rate Coefficients of the Reactions of C2H with O2, C2H2, and H2O between 295 and 450 K. *The Journal of Physical Chemistry*, **1995**, 99, 16284-16289		65
11	CH(a4.SIGMA.- and/or X2.PI.) Formation in the Reaction Between Ketenyl Radicals and Oxygen Atoms. Determination of the CH-yield between 405 and 960 K. *The Journal of Physical Chemistry*, **1995**, 99, 3583-3591		45
10	Experimental Investigation of the Reaction between Nitric Oxide and Ketenyl Radicals (HCCO + NO): Rate Coefficient at T = 290-670 K and Product Distribution at 700 K. *The Journal of Physical Chemistry*, **1994**, 98, 8036-8043		56

9	Formation of CH(a4.SIGMA.- and/or X2.PI.) in the Reaction of Ketenyl Radicals with Oxygen Atoms. Determination of the Methylidyne Yield at 290 K and ab Initio Calculations. *The Journal of Physical Chemistry*, **1994**, 98, 11988-11996		25
8	Branching ratio of the C2H2 + O reaction at 290 K from kinetic modelling of relative methylene concentration versus time profiles in C2H2/O/H systems. *International Journal of Chemical Kinetics*, **1994**, 26, 869-886	1.4	51
7	Theoretical Characterization of the Reaction between Nitric Oxide and Ketenyl Radicals (HCCO + NO): CO versus CO2 Loss. *The Journal of Physical Chemistry*, **1994**, 98, 8030-8035		56
6	Oxidation of silicon in plasma afterglows: New model of oxide growth including recombination of diffusing O atoms. *Journal of Applied Physics*, **1993**, 73, 2477-2485	2.5	11
5	A model for the plasma anodization of silicon at constant voltage. *Journal of Applied Physics*, **1993**, 74, 639-644	2.5	2
4	Product distributions of the acetylene + atomic oxygen and HCCO (ketenyl) + atomic hydrogen reactions. Rate constant of methylene(.apprx.X3B1) + atomic hydrogen. *The Journal of Physical Chemistry*, **1992**, 96, 9810-9816		66
3	A new model for the plasma anodization of silicon at constant current. *Journal of Applied Physics*, **1992**, 72, 719-724	2.5	29
2	Mechanism of □prompt□no formation in hydrocarbon flames. *Proceedings of the Combustion Institute*, **1977**, 16, 1055-1064		51
1	Production of chemi-ions and formation of CH and CH2 radicals in methane-oxygen and ethylene-oxygen flames. *Proceedings of the Combustion Institute*, **1975**, 15, 969-977		30