

Craig A Taatjes

List of Publications by Year in descending order

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205
papers

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30070

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docs citations

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times ranked

4319
citing authors

#	ARTICLE	IF	CITATIONS
1	Direct Kinetic Measurements of Criegee Intermediate (CH_2OO) Formed by Reaction of CH_2I with O_2 . <i>Science</i> , 2012, 335, 204-207.	12.6	649
2	Kinetics of elementary reactions in low-temperature autoignition chemistry. <i>Progress in Energy and Combustion Science</i> , 2011, 37, 371-421.	31.2	586
3	Direct Measurements of Conformer-Dependent Reactivity of the Criegee Intermediate CH_3CHOO . <i>Science</i> , 2013, 340, 177-180.	12.6	379
4	Enols Are Common Intermediates in Hydrocarbon Oxidation. <i>Science</i> , 2005, 308, 1887-1889.	12.6	306
5	Photoionization cross sections for reaction intermediates in hydrocarbon combustion. <i>International Journal of Mass Spectrometry</i> , 2005, 247, 18-27.	1.5	289
6	Research frontiers in the chemistry of Criegee intermediates and tropospheric ozonolysis. <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 1704.	2.8	244
7	The physical chemistry of Criegee intermediates in the gas phase. <i>International Reviews in Physical Chemistry</i> , 2015, 34, 309-360.	2.3	221
8	Rate Coefficients of C1 and C2 Criegee Intermediate Reactions with Formic and Acetic Acid Near the Collision Limit: Direct Kinetics Measurements and Atmospheric Implications. <i>Angewandte Chemie - International Edition</i> , 2014, 53, 4547-4550.	13.8	219
9	Studies of a fuel-rich propane flame with photoionization mass spectrometry. <i>Proceedings of the Combustion Institute</i> , 2005, 30, 1681-1688.	3.9	210
10	The multiplexed chemical kinetic photoionization mass spectrometer: A new approach to isomer-resolved chemical kinetics. <i>Review of Scientific Instruments</i> , 2008, 79, 104103.	1.3	190
11	Direct Observation of the Gas-Phase Criegee Intermediate (CH_2OO). <i>Journal of the American Chemical Society</i> , 2008, 130, 11883-11885.	13.7	189
12	Imaging combustion chemistry via multiplexed synchrotron-photoionization mass spectrometry. <i>Physical Chemistry Chemical Physics</i> , 2008, 10, 20-34.	2.8	185
13	Measurements, Theory, and Modeling of OH Formation in Ethyl + O_2 and Propyl + O_2 Reactions. <i>Journal of Physical Chemistry A</i> , 2003, 107, 4415-4427.	2.5	160
14	Identification and Chemistry of C_4H_3 and C_4H_5 Isomers in Fuel-Rich Flames. <i>Journal of Physical Chemistry A</i> , 2006, 110, 3670-3678.	2.5	143
15	Direct measurement of Criegee intermediate (CH_2OO) reactions with acetone, acetaldehyde, and hexafluoroacetone. <i>Physical Chemistry Chemical Physics</i> , 2012, 14, 10391.	2.8	143
16	Detection and Identification of the Keto-Hydroperoxide ($\text{HOOCH}_2\text{OCHO}$) and Other Intermediates during Low-Temperature Oxidation of Dimethyl Ether. <i>Journal of Physical Chemistry A</i> , 2015, 119, 7361-7374.	2.5	143
17	Criegee intermediates and their impacts on the troposphere. <i>Environmental Sciences: Processes and Impacts</i> , 2018, 20, 437-453.	3.5	136
18	Direct observation and kinetics of a hydroperoxyalkyl radical (QOOH). <i>Science</i> , 2015, 347, 643-646.	12.6	130

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19	Role of peroxy chemistry in the high-pressure ignition of n-butanol – Experiments and detailed kinetic modelling. <i>Combustion and Flame</i> , 2011, 158, 1444-1455.	5.2	121
20	Synchrotron photoionization measurements of combustion intermediates: Photoionization efficiency and identification of C ₃ H ₂ isomers. <i>Physical Chemistry Chemical Physics</i> , 2005, 7, 806.	2.8	113
21	Uncovering the Fundamental Chemistry of Alkyl + O ₂ Reactions via Measurements of Product Formation. <i>Journal of Physical Chemistry A</i> , 2006, 110, 4299-4312.	2.5	106
22	Production of cold formaldehyde molecules for study and control of chemical reaction dynamics with hydroxyl radicals. <i>Physical Review A</i> , 2006, 73, .	2.5	106
23	Quantification of the Keto-Hydroperoxide (HOOCH ₂ OCHO) and Other Elusive Intermediates during Low-Temperature Oxidation of Dimethyl Ether. <i>Journal of Physical Chemistry A</i> , 2016, 120, 7890-7901.	2.5	104
24	Regional and global impacts of Criegee intermediates on atmospheric sulphuric acid concentrations and first steps of aerosol formation. <i>Faraday Discussions</i> , 2013, 165, 45.	3.2	103
25	Kinetics of Cl Atom Reactions with Methane, Ethane, and Propane from 292 to 800 K. <i>Journal of Physical Chemistry A</i> , 1997, 101, 1873-1880.	2.5	102
26	Initial Steps of Aromatic Ring Formation in a Laminar Premixed Fuel-Rich Cyclopentene Flame. <i>Journal of Physical Chemistry A</i> , 2007, 111, 4081-4092.	2.5	102
27	Criegee Intermediate Reactions with Carboxylic Acids: A Potential Source of Secondary Organic Aerosol in the Atmosphere. <i>ACS Earth and Space Chemistry</i> , 2018, 2, 833-842.	2.7	102
28	Criegee Intermediates: What Direct Production and Detection Can Teach Us About Reactions of Carbonyl Oxides. <i>Annual Review of Physical Chemistry</i> , 2017, 68, 183-207.	10.8	98
29	Combustion Chemistry of Enols: Possible Ethenol Precursors in Flames. <i>Journal of Physical Chemistry A</i> , 2006, 110, 3254-3260.	2.5	96
30	Additional chain-branching pathways in the low-temperature oxidation of branched alkanes. <i>Combustion and Flame</i> , 2016, 164, 386-396.	5.2	94
31	Theory, measurements, and modeling of OH and HO ₂ formation in the reaction of cyclohexyl radicals with O ₂ . <i>Physical Chemistry Chemical Physics</i> , 2007, 9, 4315.	2.8	92
32	Products of the Benzene + O(³ P) Reaction. <i>Journal of Physical Chemistry A</i> , 2010, 114, 3355-3370.	2.5	92
33	Combustion chemistry of the propanol isomers – investigated by electron ionization and VUV-photoionization molecular-beam mass spectrometry. <i>Combustion and Flame</i> , 2009, 156, 1181-1201.	5.2	91
34	The reaction of hydroxyethyl radicals with O ₂ : A theoretical analysis and experimental product study. <i>Proceedings of the Combustion Institute</i> , 2009, 32, 271-277.	3.9	90
35	Benzene precursors and formation routes in a stoichiometric cyclohexane flame. <i>Proceedings of the Combustion Institute</i> , 2007, 31, 565-573.	3.9	89
36	Absolute Photoionization Cross-Section of the Methyl Radical. <i>Journal of Physical Chemistry A</i> , 2008, 112, 9336-9343.	2.5	89

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37	Low-temperature combustion chemistry of biofuels: pathways in the initial low-temperature (550 Tj ETQq1 1 0.784314 rgBT/Overlo	2.8	88
38	Energy-Resolved Photoionization of Alkylperoxy Radicals and the Stability of Their Cations. Journal of the American Chemical Society, 2006, 128, 13559-13567.	13.7	87
39	Direct Measurements of Unimolecular and Bimolecular Reaction Kinetics of the Criegee Intermediate (CH ₃) ₂ COO. Journal of Physical Chemistry A, 2017, 121, 4-15.	2.5	87
40	Infrared frequency-modulation probing of product formation in alkyl + O ₂ reactions. Part IV. For Part III see ref. 12. Reactions of propyl and butyl radicals with O ₂ Electronic Supplementary Information available. See http://www.rsc.org/suppdata/fd/b1/b102237g/ . Faraday Discussions, 2001, 119, 101-120.	3.2	86
41	Absolute photoionization cross-section of the propargyl radical. Journal of Chemical Physics, 2012, 136, 134307.	3.0	86
42	The reaction of Criegee intermediate CH ₂ OO with water dimer: primary products and atmospheric impact. Physical Chemistry Chemical Physics, 2017, 19, 21970-21979.	2.8	83
43	Direct detection of polyynes formation from the reaction of ethynyl radical (C ₂ H) with propyne (CH ₃ C≡CH) and allene (CH ₂ C=CH ₂). Physical Chemistry Chemical Physics, 2007, 9, 4291-4300.	2.8	79
44	Cyclic Versus Linear Isomers Produced by Reaction of the Methylidyne Radical (CH) with Small Unsaturated Hydrocarbons. Journal of the American Chemical Society, 2009, 131, 993-1005.	13.7	77
45	Formally direct pathways and low-temperature chain branching in hydrocarbon autoignition: the cyclohexyl + O ₂ reaction at high pressure. Physical Chemistry Chemical Physics, 2009, 11, 1320.	2.8	76
46	Infrared Frequency-Modulation Probing of Product Formation in Alkyl + O ₂ Reactions: I. The Reaction of C ₂ H ₅ with O ₂ between 295 and 698 K. Journal of Physical Chemistry A, 2000, 104, 11549-11560.	2.5	73
47	VUV Photoionization Cross Sections of HO ₂ , H ₂ O ₂ , and H ₂ CO. Journal of Physical Chemistry A, 2015, 119, 1279-1291.	2.5	66
48	The influence of ethanol addition on premixed fuel-rich propene-oxygen-argon flames. Proceedings of the Combustion Institute, 2007, 31, 1119-1127.	3.9	64
49	Photoionization mass spectrometric studies and modeling of fuel-rich allene and propyne flames. Proceedings of the Combustion Institute, 2007, 31, 1157-1164.	3.9	63
50	Direct kinetic measurements and theoretical predictions of an isoprene-derived Criegee intermediate. Proceedings of the National Academy of Sciences of the United States of America, 2020, 117, 9733-9740.	7.1	63
51	Quantum Interference as the Source of Steric Asymmetry and Parity Propensity Rules in NO ⁺ Rare Gas Inelastic Scattering. Journal of the American Chemical Society, 2006, 128, 8777-8789.	13.7	61
52	Synchrotron-based double imaging photoelectron/photoion coincidence spectroscopy of radicals produced in a flow tube: OH and OD. Journal of Chemical Physics, 2015, 142, 164201.	3.0	60
53	Directly measuring reaction kinetics of ¹ QOOH - a crucial but elusive intermediate in hydrocarbon autoignition. Physical Chemistry Chemical Physics, 2013, 15, 10753.	2.8	58
54	Influence of functional groups on low-temperature combustion chemistry of biofuels. Progress in Energy and Combustion Science, 2021, 86, 100925.	31.2	58

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55	Reaction of the C ₂ H Radical with 1-Butyne (C ₄ H ₆): Low-Temperature Kinetics and Isomer-Specific Product Detection. <i>Journal of Physical Chemistry A</i> , 2010, 114, 3340-3354.	2.5	57
56	Reactions of the CN Radical with Benzene and Toluene: Product Detection and Low-Temperature Kinetics. <i>Journal of Physical Chemistry A</i> , 2010, 114, 1749-1755.	2.5	56
57	Criegee intermediates: production, detection and reactivity. <i>International Reviews in Physical Chemistry</i> , 2020, 39, 385-424.	2.3	56
58	Atmospheric transformation of enols: A potential secondary source of carboxylic acids in the urban troposphere. <i>Geophysical Research Letters</i> , 2007, 34, .	4.0	55
59	Diode laser probing of I*(2P _{1/2}) Doppler profiles: Time evolution of a fast, anisotropic velocity distribution in a thermal bath. <i>Journal of Chemical Physics</i> , 1990, 93, 6543-6553.	3.0	54
60	Absolute and Site-Specific Abstraction Rate Coefficients for Reactions of Cl with CH ₃ CH ₂ OH, CH ₃ CD ₂ OH, and CD ₃ CH ₂ OH between 295 and 600 K. <i>Journal of Physical Chemistry A</i> , 1999, 103, 9805-9814.	2.5	54
61	Spectroscopy of the Simplest Criegee Intermediate CH ₂ OO: Simulation of the First Bands in Its Electronic and Photoelectron Spectra. <i>Chemistry - A European Journal</i> , 2012, 18, 12411-12423.	3.3	54
62	A coordinated investigation of the combustion chemistry of diisopropyl ketone, a prototype for biofuels produced by endophytic fungi. <i>Combustion and Flame</i> , 2014, 161, 711-724.	5.2	54
63	Infrared Frequency-Modulation Probing of Product Formation in Alkyl + O ₂ Reactions: II. The Reaction of C ₃ H ₇ with O ₂ between 296 and 683 K. <i>Journal of Physical Chemistry A</i> , 2001, 105, 3205-3213.	2.5	53
64	Temperature controlled multiple pass absorption cell for gas phase chemical kinetics studies. <i>Review of Scientific Instruments</i> , 1997, 68, 1875-1878.	1.3	52
65	New mechanistic insights to the O(3P) + propene reaction from multiplexed photoionization mass spectrometry. <i>Physical Chemistry Chemical Physics</i> , 2012, 14, 10410.	2.8	51
66	Time-resolved infrared absorption measurements of product formation in Cl atom reactions with alkenes and alkynes. <i>International Reviews in Physical Chemistry</i> , 1999, 18, 419-458.	2.3	50
67	Direct detection of pyridine formation by the reaction of CH (CD) with pyrrole: a ring expansion reaction. <i>Physical Chemistry Chemical Physics</i> , 2010, 12, 8750.	2.8	49
68	Probing the low-temperature chain-branching mechanism of n-butane autoignition chemistry via time-resolved measurements of ketohydroperoxide formation in photolytically initiated n-C ₄ H ₁₀ oxidation. <i>Proceedings of the Combustion Institute</i> , 2015, 35, 291-298.	3.9	48
69	Infrared Laser Absorption Measurements of the Kinetics of Propargyl Radical Self-Reaction and the 193 nm Photolysis of Propyne. <i>Journal of Physical Chemistry A</i> , 2003, 107, 4843-4850.	2.5	47
70	Infrared Absorption Probing of the Cl + C ₃ H ₆ Reaction: Rate Coefficients for HCl Production between 290 and 800 K. <i>Journal of Physical Chemistry A</i> , 1997, 101, 5776-5782.	2.5	46
71	Pressure-Dependent I-Atom Yield in the Reaction of CH ₂ I with O ₂ Shows a Remarkable Apparent Third-Body Efficiency for O ₂ . <i>Journal of Physical Chemistry Letters</i> , 2012, 3, 3399-3403.	4.6	46
72	Low-temperature combustion chemistry of biofuels: Pathways in the low-temperature (550-700K) oxidation chemistry of isobutanol and tert-butanol. <i>Proceedings of the Combustion Institute</i> , 2013, 34, 493-500.	3.9	46

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73	Experimental and computational studies of Criegee intermediate reactions with NH ₃ and CH ₃ NH ₂ . Physical Chemistry Chemical Physics, 2019, 21, 14042-14052.	2.8	46
74	The vinyl radical ($\tilde{\chi}^2A''$) spectrum between 530 and 415 nm measured by cavity ring-down spectroscopy. Journal of Chemical Physics, 1999, 110, 1841-1843.	3.0	45
75	A general method for Doppler determination of cylindrically symmetric velocity distributions: An application of Fourier transform Doppler spectroscopy. Journal of Chemical Physics, 1990, 93, 6554-6559.	3.0	44
76	High-resolution diode laser absorption spectroscopy of the O-H stretch overtone band (2,0,0) \rightarrow (0,0,0) of the HO ₂ radical. Journal of Molecular Spectroscopy, 2003, 219, 163-169.	1.2	43
77	Deuterium Kinetic Isotope Effect and Temperature Dependence in the Reactions of CH ₂ with Methane and Acetylene. Journal of Physical Chemistry A, 1997, 101, 1881-1886.	2.5	42
78	Synchrotron Photoionization Mass Spectrometry Measurements of Kinetics and Product Formation in the Allyl Radical (H ₂ CCHCH ₂) Self-Reaction. Journal of Physical Chemistry A, 2008, 112, 9366-9373.	2.5	42
79	Low-Temperature Combustion Chemistry of <i>n</i> -Butanol: Principal Oxidation Pathways of Hydroxybutyl Radicals. Journal of Physical Chemistry A, 2013, 117, 11983-12001.	2.5	40
80	Pressure-Dependent Competition among Reaction Pathways from First- and Second-O ₂ Additions in the Low-Temperature Oxidation of Tetrahydrofuran. Journal of Physical Chemistry A, 2016, 120, 6582-6595.	2.5	40
81	Time-resolved wavelength modulation spectroscopy measurements of HO ₂ kinetics. Applied Optics, 1997, 36, 5817.	2.1	39
82	Measurements and Modeling of HO ₂ Formation in the Reactions of <i>n</i> -C ₃ H ₇ and <i>i</i> -C ₃ H ₇ Radicals with O ₂ . Journal of Physical Chemistry B, 2005, 109, 8374-8387.	2.6	38
83	Photoionization of 1-Alkenylperoxy and Alkylperoxy Radicals and a General Rule for the Stability of Their Cations. Journal of the American Chemical Society, 2007, 129, 14019-14025.	13.7	38
84	Unconventional Peroxy Chemistry in Alcohol Oxidation: The Water Elimination Pathway. Journal of Physical Chemistry Letters, 2013, 4, 350-354.	4.6	38
85	Time- and Isomer-Resolved Measurements of Sequential Addition of Acetylene to the Propargyl Radical. Journal of Physical Chemistry Letters, 2015, 6, 4153-4158.	4.6	38
86	Products of Criegee intermediate reactions with NO ₂ : experimental measurements and tropospheric implications. Faraday Discussions, 2017, 200, 313-330.	3.2	38
87	Infrared Absorption Probing of the Cl + C ₂ H ₄ Reaction: Direct Measurement of Arrhenius Parameters for Hydrogen Abstraction. Journal of Physical Chemistry A, 1997, 101, 4172-4177.	2.5	37
88	RECENT PROGRESS IN INFRARED ABSORPTION TECHNIQUES FOR ELEMENTARY GAS-PHASE REACTION KINETICS. Annual Review of Physical Chemistry, 2001, 52, 41-70.	10.8	37
89	Multiscale Informatics for Low-Temperature Propane Oxidation: Further Complexities in Studies of Complex Reactions. Journal of Physical Chemistry A, 2015, 119, 7095-7115.	2.5	37
90	Synchrotron photoionization measurements of combustion intermediates: the photoionization efficiency of HONO. Chemical Physics Letters, 2004, 394, 19-24.	2.6	36

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91	Research needs for future internal combustion engines. <i>Physics Today</i> , 2008, 61, 47-52.	0.3	36
92	New experiments and validated master-equation modeling for OH production in propyl+O ₂ reactions. <i>Proceedings of the Combustion Institute</i> , 2011, 33, 293-299.	3.9	35
93	Steric asymmetry and lambda-doublet propensities in state-to-state rotationally inelastic scattering of NO(2 $\hat{1}$ /2) with He. <i>Journal of Chemical Physics</i> , 2004, 121, 11691-11701.	3.0	34
94	Isomer-specific product detection of CN radical reactions with ethene and propene by tunable VUV photoionization mass spectrometry. <i>International Journal of Mass Spectrometry</i> , 2009, 280, 113-118.	1.5	34
95	Laser double-resonance measurements of rotational relaxation rates of HF($\hat{\nu}_0$ =13) with rare gases, H ₂ , and D ₂ . <i>Journal of Chemical Physics</i> , 1988, 89, 302-308.	3.0	33
96	Two-Dimensional Imaging of the Photolysis of Oriented Molecules. <i>The Journal of Physical Chemistry</i> , 1995, 99, 4360-4363.	2.9	33
97	Synchrotron Photoionization Mass Spectrometry Measurements of Product Formation in Low-Temperature <i>n</i> -Butane Oxidation: Toward a Fundamental Understanding of Autoignition Chemistry and <i>n</i> -C ₄ H ₉ + O ₂ Reactions. <i>Journal of Physical Chemistry A</i> , 2013, 117, 12216-12225.	2.5	33
98	Influence of oxygenation in cyclic hydrocarbons on chain-termination reactions from R + O ₂ : tetrahydropyran and cyclohexane. <i>Proceedings of the Combustion Institute</i> , 2017, 36, 597-606.	3.9	33
99	Rapid Discovery and Functional Characterization of Terpene Synthases from Four Endophytic Xylariaceae. <i>PLoS ONE</i> , 2016, 11, e0146983.	2.5	33
100	Infrared laser absorption measurements of HCl(ν =1) production in reactions of Cl atoms with isobutane, methanol, acetaldehyde, and toluene at 295 K. <i>Chemical Physics Letters</i> , 2002, 366, 417-425.	2.6	32
101	Time-resolved measurements of OH and HO ₂ product formation in pulsed-photolytic chlorine atom initiated oxidation of neopentane. <i>Physical Chemistry Chemical Physics</i> , 2003, 5, 1584-1592.	2.8	32
102	New Insights into Low-Temperature Oxidation of Propane from Synchrotron Photoionization Mass Spectrometry and Multiscale Informatics Modeling. <i>Journal of Physical Chemistry A</i> , 2015, 119, 7116-7129.	2.5	32
103	The reaction of hydroxyl and methylperoxy radicals is not a major source of atmospheric methanol. <i>Nature Communications</i> , 2018, 9, 4343.	12.8	32
104	Measurements and Modeling of DO ₂ Formation in the Reactions of C ₂ D ₅ and C ₃ D ₇ Radicals with O ₂ . <i>Journal of Physical Chemistry A</i> , 2007, 111, 4015-4030.	2.5	31
105	Infrared Frequency-Modulation Probing of Cl + C ₃ H ₄ (Allene, Propyne) Reactions: Kinetics of HCl Production from 292 to 850 K. <i>Journal of Physical Chemistry A</i> , 1998, 102, 4846-4856.	2.5	30
106	Experimental Evidence of Dioxole Unimolecular Decay Pathway for Isoprene-Derived Criegee Intermediates. <i>Journal of Physical Chemistry A</i> , 2020, 124, 3542-3554.	2.5	30
107	Polarization-resolved (2+1) resonance-enhanced multiphoton ionization spectroscopy of CF ₃ I (6s) Rydberg states. <i>Journal of Chemical Physics</i> , 1993, 98, 4355-4371.	3.0	29
108	Measurements and Automated Mechanism Generation Modeling of OH Production in Photolytically Initiated Oxidation of the Neopentyl Radical. <i>Journal of Physical Chemistry A</i> , 2007, 111, 3891-3900.	2.5	29

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109	Product Detection of the CH Radical Reaction with Acetaldehyde. Journal of Physical Chemistry A, 2012, 116, 6091-6106.	2.5	29
110	Temperature Dependence and Deuterium Kinetic Isotope Effects in the CH (CD) + C ₂ H ₄ (C ₂ D ₄) Reaction between 295 and 726 K. Journal of Physical Chemistry A, 2001, 105, 5393-5401.	2.5	28
111	How does the molecular velocity distribution affect kinetics measurements by time-resolved mass spectrometry?. International Journal of Chemical Kinetics, 2007, 39, 565-570.	1.6	28
112	Infrared Frequency-Modulation Probing of Product Formation in Alkyl + O ₂ Reactions: III. The Reaction of Cyclopentyl Radical (c-C ₅ H ₉) with O ₂ between 296 and 723 K. Journal of Physical Chemistry A, 2001, 105, 6646-6654.	2.5	27
113	Temperature dependence and deuterium kinetic isotope effects in the HCO(DCO)+O ₂ reaction between 296 and 673 K. Chemical Physics Letters, 2001, 347, 79-86.	2.6	27
114	Vinyl radical visible spectroscopy and excited state dynamics. Journal of Chemical Physics, 2002, 116, 8343.	3.0	27
115	Measurements and Quasi-Quantum Modeling of the Steric Asymmetry and Parity Propensities in State-to-State Rotationally Inelastic Scattering of NO (2 $\hat{1}$ 1/2) with D ₂ . Journal of Physical Chemistry A, 2007, 111, 7631-7639.	2.5	27
116	Temperature-Dependent Kinetics of the Vinyl Radical (C ₂ H ₃) Self-Reaction. Journal of Physical Chemistry A, 2009, 113, 1278-1286.	2.5	27
117	Hydroxyacetone Production From C ₃ Criegee Intermediates. Journal of Physical Chemistry A, 2017, 121, 16-23.	2.5	27
118	Rotational excitation in scattering of hyperthermal NO from Pt(111). Journal of Chemical Physics, 1995, 102, 3835-3847.	3.0	26
119	Infrared frequency-modulation measurements of absolute rate coefficients for Cl+HD $\hat{+}$ HCl(DCl)+D(H) between 295 and 700 K. Chemical Physics Letters, 1999, 306, 33-40.	2.6	26
120	Kinetics of the reaction of vinyl radicals with NO: Ab initio theory, master equation predictions, and laser absorption measurements. Physical Chemistry Chemical Physics, 2004, 6, 2216-2223.	2.8	25
121	Low-temperature combustion chemistry of novel biofuels: resonance-stabilized QOOH in the oxidation of diethyl ketone. Physical Chemistry Chemical Physics, 2014, 16, 13027-13040.	2.8	25
122	Kinetic Isotope Effects and Variable Reaction Coordinates in Barrierless Recombination Reactions. Journal of Physical Chemistry A, 2001, 105, 8567-8578.	2.5	24
123	Formation and stability of gas-phase o-benzoquinone from oxidation of ortho-hydroxyphenyl: a combined neutral and distonic radical study. Physical Chemistry Chemical Physics, 2016, 18, 4320-4332.	2.8	24
124	Influence of the Ether Functional Group on Ketohydroperoxide Formation in Cyclic Hydrocarbons: Tetrahydropyran and Cyclohexane. Journal of Physical Chemistry A, 2019, 123, 3634-3646.	2.5	23
125	Dynamical information in angular distributions of fragments from photolysis of oriented molecules. Chemical Physics Letters, 1993, 203, 363-370.	2.6	22
126	Temperature dependence of the reaction C ₂ H (C ₂ D) + O ₂ between 295 and 700 K. Chemical Physics Letters, 1997, 270, 580-586.	2.6	22

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127	Isomer-Selective Study of the OH Initiated Oxidation of Isoprene in the Presence of O ₂ and NO. I. The Minor Inner OH-Addition Channel. Journal of Physical Chemistry A, 2010, 114, 904-912.	2.5	22
128	Formation of dimethylketene and methacrolein by reaction of the CH radical with acetone. Physical Chemistry Chemical Physics, 2013, 15, 4049.	2.8	22
129	Product Branching Fractions of the CH + Propene Reaction from Synchrotron Photoionization Mass Spectrometry. Journal of Physical Chemistry A, 2013, 117, 6450-6457.	2.5	22
130	Kinetic Isotope Effect in the CH[2 ¹] + O ₂ Reaction. The Journal of Physical Chemistry, 1996, 100, 17840-17845.	2.9	21
131	HO ₂ OOH-mediated reactions in cyclohexene oxidation. Proceedings of the Combustion Institute, 2019, 37, 323-335.	3.9	21
132	Reaction of Perfluorooctanoic Acid with Criegee Intermediates and Implications for the Atmospheric Fate of Perfluorocarboxylic Acids. Environmental Science & Technology, 2019, 53, 1245-1251.	10.0	21
133	Product Formation in the Cl-Initiated Oxidation of Cyclopropane. Journal of Physical Chemistry A, 2003, 107, 1992-2002.	2.5	20
134	Pressure and Temperature Dependence of the Reaction of Vinyl Radical with Ethylene. Journal of Physical Chemistry A, 2007, 111, 6843-6851.	2.5	20
135	Absolute photoionization cross-sections of selected furanic and lactonic potential biofuels. International Journal of Mass Spectrometry, 2013, 348, 39-46.	1.5	20
136	Time-resolved measurements of product formation in the low-temperature (550-675 K) oxidation of neopentane: a probe to investigate chain-branching mechanism. Physical Chemistry Chemical Physics, 2017, 19, 13731-13745.	2.8	20
137	Direct kinetics study of CH ₂ OO + methyl vinyl ketone and CH ₂ OO + methacrolein reactions and an upper limit determination for CH ₂ OO + CO reaction. Physical Chemistry Chemical Physics, 2018, 20, 19373-19381.	2.8	20
138	Vacuum ultraviolet photoionization cross section of the hydroxyl radical. Journal of Chemical Physics, 2018, 148, 184302.	3.0	20
139	Reaction mechanisms of a cyclic ether intermediate: Ethyloxirane. International Journal of Chemical Kinetics, 2021, 53, 43-59.	1.6	20
140	Rotational rainbows in NO scattering from Pt(111). Faraday Discussions, 1993, 96, 297.	3.2	19
141	Ultraviolet photodissociation of vinyl iodide: understanding the halogen dependence of photodissociation mechanisms in vinyl halides. Physical Chemistry Chemical Physics, 2008, 10, 713-728.	2.8	19
142	Photoionization Mass Spectrometric Measurements of Initial Reaction Pathways in Low-Temperature Oxidation of 2,5-Dimethylhexane. Journal of Physical Chemistry A, 2014, 118, 10188-10200.	2.5	19
143	Investigating the Tropospheric Chemistry of Acetic Acid Using the Global 3D Chemistry Transport Model, STOCHEM-CRI. Journal of Geophysical Research D: Atmospheres, 2018, 123, 6267-6281.	3.3	19
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