

# Jim J Lin

## List of Publications by Citations

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#	Paper	IF	Citations
138	Application of time-sliced ion velocity imaging to crossed molecular beam experiments. <i>Review of Scientific Instruments</i> , <b>2003</b> , 74, 2495-2500	1.7	282
137	Atmospheric chemistry. Direct kinetic measurement of the reaction of the simplest Criegee intermediate with water vapor. <i>Science</i> , <b>2015</b> , 347, 751-4	33.3	218
136	State-specific correlation of coincident product pairs in the F + CD <sub>4</sub> reaction. <i>Science</i> , <b>2003</b> , 300, 966-9	33.3	216
135	Kinetics of a Criegee intermediate that would survive high humidity and may oxidize atmospheric SO <sub>2</sub> . <i>Proceedings of the National Academy of Sciences of the United States of America</i> , <b>2015</b> , 112, 10857-62	11.5	180
134	Photodissociation of H <sub>2</sub> O at 121.6 nm: A state-to-state dynamical picture. <i>Journal of Chemical Physics</i> , <b>2000</b> , 113, 10073-10090	3.9	163
133	Reactive resonance in a polyatomic reaction. <i>Physical Review Letters</i> , <b>2004</b> , 92, 103201	7.4	133
132	A Quantum State-Resolved Insertion Reaction: O((1)D) + H <sub>2</sub> (J = 0) → OH((2) product operator product operator, v, N) + H((2)S). <i>Science</i> , <b>2000</b> , 289, 1536-1538	33.3	123
131	Competition between H <sub>2</sub> O and (H <sub>2</sub> O) <sub>2</sub> reactions with CH <sub>2</sub> OO/CH <sub>3</sub> CHOO. <i>Physical Chemistry Chemical Physics</i> , <b>2016</b> , 18, 4557-68	3.6	110
130	The UV absorption spectrum of the simplest Criegee intermediate CH <sub>2</sub> OO. <i>Physical Chemistry Chemical Physics</i> , <b>2014</b> , 16, 10438-43	3.6	106
129	Strong Negative Temperature Dependence of the Simplest Criegee Intermediate CH <sub>2</sub> OO Reaction with Water Dimer. <i>Journal of Physical Chemistry Letters</i> , <b>2015</b> , 6, 2708-13	6.4	105
128	Crossed-beam scattering of F+CD <sub>4</sub> →DF+CD <sub>3</sub> (NK): The integral cross sections. <i>Journal of Chemical Physics</i> , <b>2003</b> , 119, 2538-2544	3.9	103
127	Structure-dependent reactivity of Criegee intermediates studied with spectroscopic methods. <i>Chemical Society Reviews</i> , <b>2017</b> , 46, 7483-7497	58.5	99
126	Probing the effect of the H <sub>2</sub> rotational state in O(1D)+H <sub>2</sub> →OH+H: Theoretical dynamics including nonadiabatic effects and a crossed molecular beam study. <i>Journal of Chemical Physics</i> , <b>2000</b> , 113, 7330-7344	3.9	92
125	Imaging the pair-correlated excitation function: The F+CH <sub>4</sub> →HF(v=0)+CH <sub>3</sub> (nu=0) reaction. <i>Journal of Chemical Physics</i> , <b>2004</b> , 120, 117-22	3.9	81
124	On the Cl*(2P <sub>1/2</sub> ) Reactivity and the Effect of Bend Excitation in the Cl + CH <sub>4</sub> /CD <sub>4</sub> Reactions. <i>Journal of Physical Chemistry A</i> , <b>2004</b> , 108, 7832-7836	2.8	77
123	Detailed mechanism of the CH <sub>2</sub> + O <sub>3</sub> reaction: yield and self-reaction of the simplest Criegee intermediate CH <sub>2</sub> OO. <i>Journal of Chemical Physics</i> , <b>2014</b> , 141, 104308	3.9	76
122	Observation of a reactive resonance in the integral cross section of a six-atom reaction: F+CHD <sub>3</sub> . <i>Journal of Chemical Physics</i> , <b>2004</b> , 121, 813-8	3.9	74

121	The O(1D)+H <sub>2</sub> reaction at 56 meV collision energy: A comparison between quantum mechanical, quasiclassical trajectory, and crossed beam results. <i>Journal of Chemical Physics</i> , <b>2002</b> , 116, 10692-10703	3.9	72
120	Insights into dynamics of the F+CD <sub>4</sub> reaction via product pair correlation. <i>Journal of Chemical Physics</i> , <b>2003</b> , 119, 4997-5000	3.9	67
119	UV absorption spectrum of the C <sub>2</sub> Criegee intermediate CH <sub>3</sub> CHOO. <i>Journal of Chemical Physics</i> , <b>2014</b> , 141, 074302	3.9	66
118	Unimolecular Decomposition Rate of the Criegee Intermediate (CH <sub>3</sub> ) <sub>2</sub> COO Measured Directly with UV Absorption Spectroscopy. <i>Journal of Physical Chemistry A</i> , <b>2016</b> , 120, 4789-98	2.8	65
117	Mode-correlated product pairs in the F+CHD <sub>3</sub> ->DF+CHD <sub>2</sub> reaction. <i>Journal of Chemical Physics</i> , <b>2003</b> , 119, 8289-8296	3.9	65
116	Multiple dynamical pathways in the O(1D)+CH <sub>4</sub> reaction: A comprehensive crossed beam study. <i>Journal of Chemical Physics</i> , <b>2000</b> , 113, 5287	3.9	64
115	New low background crossed molecular beam apparatus: Low background detection of H <sub>2</sub> . <i>Review of Scientific Instruments</i> , <b>1998</b> , 69, 1642-1646	1.7	64
114	Imaging the isotope effects in the ground state reaction of Cl CH <sub>4</sub> and CD <sub>4</sub> . <i>Molecular Physics</i> , <b>2005</b> , 103, 1757-1763	1.7	57
113	Dissociation dynamics of the water molecule on the $\sigma^*_{B1}$ electronic surface. <i>Journal of Chemical Physics</i> , <b>2000</b> , 113, 10597-10604	3.9	54
112	Photodissociation of D <sub>2</sub> O at 121.6 nm: A state-to-state dynamical picture. <i>Journal of Chemical Physics</i> , <b>2001</b> , 114, 7830-7837	3.9	53
111	Rotationally selected product pair correlation in F+CD(4)->DF(nu(0)+CD(3)(nu=0,N). <i>Journal of Chemical Physics</i> , <b>2004</b> , 120, 5863-6	3.9	51
110	Photodissociation dynamics of CH <sub>3</sub> Cl at 157.6 nm: Evidence for CH <sub>2</sub> (X 3B <sub>1</sub> / $\sigma^*_{A1}$ )+HCl product channels. <i>Chemical Physics Letters</i> , <b>2002</b> , 361, 374-382	2.5	51
109	UV absorption cross sections of ClOOCl are consistent with ozone degradation models. <i>Science</i> , <b>2009</b> , 324, 781-4	33.3	49
108	Mode correlation of product pairs in the reaction OH+CD <sub>4</sub> ->HOD+CD <sub>3</sub> . <i>Journal of Chemical Physics</i> , <b>2005</b> , 122, 131102	3.9	47
107	Dynamics of the O(1D)+CH <sub>4</sub> reaction: Atomic hydrogen channel vs molecular hydrogen channel. <i>Journal of Chemical Physics</i> , <b>1999</b> , 110, 10821-10829	3.9	47
106	The low-lying bending vibrational levels of the CCH (X 2 $\Pi$ ) radical studied by laser-induced fluorescence. <i>Journal of Chemical Physics</i> , <b>1993</b> , 98, 6690-6696	3.9	47
105	Temperature dependence of the reaction of anti-CHCHOO with water vapor. <i>Physical Chemistry Chemical Physics</i> , <b>2016</b> , 18, 28189-28197	3.6	46
104	Nonstatistical behavior of reactive scattering in the (18)O+(32)O(2) isotope exchange reaction. <i>Journal of the American Chemical Society</i> , <b>2007</b> , 129, 2866-70	16.4	44

103	State-to-state dynamics for O(1D)+D <sub>2</sub> →OD+D: evidence for a collinear abstraction mechanism. <i>Physical Review Letters</i> , <b>2001</b> , 86, 408-11	7.4	44
102	Rotationally selected product pair correlation: F+CD <sub>4</sub> → DF( $\nu_1$ )+CD <sub>3</sub> ( $\nu_2 = 0$ and 2, N). <i>Journal of Chemical Physics</i> , <b>2006</b> , 124, 104309	3.9	43
101	Unimolecular dissociation of the CH <sub>3</sub> OCO radical: an intermediate in the CH <sub>3</sub> O + CO reaction. <i>Journal of Physical Chemistry A</i> , <b>2006</b> , 110, 1625-34	2.8	42
100	Site-specific dissociation dynamics of ethylene at 157 nm: Atomic and molecular hydrogen elimination. <i>Journal of Chemical Physics</i> , <b>2000</b> , 113, 9668-9677	3.9	37
99	Two photoionization thresholds of N <sub>3</sub> produced by ClN <sub>3</sub> photodissociation at 248 nm: further evidence for cyclic N <sub>3</sub> . <i>Journal of Chemical Physics</i> , <b>2005</b> , 123, 051101	3.9	36
98	Quantum state specific dynamics for the O(1D)+HD→OD+H reaction. <i>Journal of Chemical Physics</i> , <b>2000</b> , 113, 1325-1328	3.9	36
97	State-correlation matrix of the product pair from F + CD <sub>4</sub> → DF( $\nu_1$ )+ CD <sub>3</sub> (0 $\nu_2$ 0 0). <i>Physical Chemistry Chemical Physics</i> , <b>2006</b> , 8, 3000-6	3.6	35
96	Energy Dependence of Oxygen Isotope Exchange and Quenching in the O(1D) + CO <sub>2</sub> Reaction: A Crossed Molecular Beam Study. <i>Journal of Physical Chemistry A</i> , <b>2004</b> , 108, 7995-8001	2.8	35
95	Photodissociation dynamics of cyclopropane at 157 nm. <i>Journal of Chemical Physics</i> , <b>2002</b> , 117, 153-160	3.9	34
94	Crossed molecular beam studies of the O(1D)+CH <sub>4</sub> reaction: Evidences for the CH <sub>2</sub> OH+H channel. <i>Journal of Chemical Physics</i> , <b>1998</b> , 109, 2975-2978	3.9	34
93	Direct kinetic measurements and theoretical predictions of an isoprene-derived Criegee intermediate. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , <b>2020</b> , 117, 9733-9740	11.5	33
92	Photodissociation dynamics of H <sub>2</sub> O at 121.6 nm: Effect of parent rotational excitation on reaction pathways. <i>Journal of Chemical Physics</i> , <b>1999</b> , 110, 4123-4126	3.9	32
91	Photodissociation of O <sub>2</sub> at 157 nm: Experimental observation of anisotropy mixing in the O <sub>2</sub> +h $\nu$ →O(3P)+O(3P) channel. <i>Journal of Chemical Physics</i> , <b>1998</b> , 109, 1758-1762	3.9	31
90	Products of Criegee intermediate reactions with NO: experimental measurements and tropospheric implications. <i>Faraday Discussions</i> , <b>2017</b> , 200, 313-330	3.6	27
89	Observation of photochemical C-N bond cleavage in CH <sub>3</sub> N <sub>3</sub> : a new photochemical route to cyclic N <sub>3</sub> . <i>Journal of Physical Chemistry A</i> , <b>2008</b> , 112, 1105-11	2.8	27
88	Dynamics of the O(1D)+CO <sub>2</sub> oxygen isotope exchange reaction. <i>Journal of Chemical Physics</i> , <b>2003</b> , 119, 8213-8216	3.9	27
87	Criegee Intermediate Reaction with Alcohol Is Enhanced by a Single Water Molecule. <i>Journal of Physical Chemistry Letters</i> , <b>2018</b> , 9, 7040-7044	6.4	27
86	Photodissociation cross section of ClOOCl at 330 nm. <i>Journal of Physical Chemistry A</i> , <b>2010</b> , 114, 4791-7	2.8	26

85	Photodissociation dynamics of 1,1-difluoroethylene at 157 nm excitation. <i>Journal of Chemical Physics</i> , <b>1998</b> , 109, 10838-10846	3.9	26
84	Site and isotope effects on the molecular hydrogen elimination from ethylene at 157 nm excitation. <i>Journal of Chemical Physics</i> , <b>1998</b> , 109, 2979-2982	3.9	26
83	Photodissociation Dynamics of Methanol at 157 nm. <i>Journal of Physical Chemistry A</i> , <b>1999</b> , 103, 10324-10332	2.6	26
82	Competing atomic and molecular hydrogen pathways in the photodissociation of methanol at 157 nm. <i>Journal of Chemical Physics</i> , <b>1999</b> , 111, 5-9	3.9	26
81	Temperature-Dependent Rate Coefficients for the Reaction of CHOO with Hydrogen Sulfide. <i>Journal of Physical Chemistry A</i> , <b>2017</b> , 121, 938-945	2.8	25
80	Photodissociation cross sections of CLOOCl at 248.4 and 266 nm. <i>Journal of Chemical Physics</i> , <b>2009</b> , 131, 174301	3.9	25
79	Multiple pathway dynamics of the O(1D)+C <sub>2</sub> H <sub>6</sub> reaction: A crossed beam study. <i>Journal of Chemical Physics</i> , <b>2001</b> , 115, 849-857	3.9	25
78	Absolute UV absorption cross sections of dimethyl substituted Criegee intermediate (CH <sub>3</sub> ) <sub>2</sub> COO. <i>Chemical Physics Letters</i> , <b>2016</b> , 653, 155-160	2.5	23
77	Product branching from the CH <sub>2</sub> CH <sub>2</sub> OH radical intermediate of the OH + ethene reaction. <i>Journal of Physical Chemistry A</i> , <b>2011</b> , 115, 9097-110	2.8	23
76	A crossed molecular beam study of the O(1D) + C(3)H(8) reaction: multiple reaction pathways. <i>Journal of the American Chemical Society</i> , <b>2001</b> , 123, 322-30	16.4	23
75	Photodissociation of hydrogen sulfide at 157.6 nm: Observation of SH bimodal rotational distribution. <i>Journal of Chemical Physics</i> , <b>1999</b> , 111, 3940-3945	3.9	22
74	The simplest all-nitrogen ring: photolytically filling the cyclic-N <sub>3</sub> well. <i>Journal of Chemical Physics</i> , <b>2007</b> , 126, 041101	3.9	21
73	Photodissociation dynamics of propyne at 157 nm. <i>Journal of Chemical Physics</i> , <b>2000</b> , 112, 6656-6665	3.9	21
72	Investigation of the O+allyl addition/elimination reaction pathways from the OCH(2)CHCH(2) radical intermediate. <i>Journal of Chemical Physics</i> , <b>2008</b> , 129, 084301	3.9	20
71	A complete look at a multiple pathway reaction: The reaction of O( <sup>1</sup> D) with ethane. <i>Journal of Chemical Physics</i> , <b>2001</b> , 114, 4	3.9	20
70	High resolution quantum cascade laser spectroscopy of the simplest Criegee intermediate, CHOO, between 1273 cm and 1290 cm. <i>Journal of Chemical Physics</i> , <b>2017</b> , 146, 244302	3.9	19
69	Primary photodissociation pathways of epichlorohydrin and analysis of the C-C bond fission channels from an O( <sup>3</sup> P)+allyl radical intermediate. <i>Journal of Chemical Physics</i> , <b>2010</b> , 133, 094306	3.9	19
68	Deciphering the nature of the reactive resonance in F + CHD <sub>3</sub> : correlated differential cross-sections of the two isotopic channels. <i>Molecular Physics</i> , <b>2010</b> , 108, 957-968	1.7	19

67	Determining the partial photoionization cross-sections of ethyl radicals. <i>Journal of Physical Chemistry A</i> , <b>2007</b> , 111, 12417-22	2.8	19
66	Collision-free photochemistry of methylazide: observation of unimolecular decomposition of singlet methylnitrene. <i>Journal of Chemical Physics</i> , <b>2006</b> , 125, 133302	3.9	19
65	Product angular anisotropy in CO <sub>2</sub> photodissociation at 157 nm. <i>Chemical Physics Letters</i> , <b>2003</b> , 382, 665-670	2.5	19
64	The non-statistical dynamics of the <sup>13</sup> C + <sup>18</sup> O isotope exchange reaction at two energies. <i>Journal of Chemical Physics</i> , <b>2014</b> , 141, 064311	3.9	18
63	Unimolecular dissociation of the propargyl radical intermediate of the CH+C <sub>2</sub> H <sub>2</sub> and C+C <sub>2</sub> H <sub>3</sub> reactions. <i>Journal of Chemical Physics</i> , <b>2006</b> , 125, 133306	3.9	18
62	Kinetics of the reaction of the simplest Criegee intermediate with ammonia: a combination of experiment and theory. <i>Physical Chemistry Chemical Physics</i> , <b>2018</b> , 20, 29669-29676	3.6	18
61	The bending vibrational levels of the acetylene cation: a case study of the Renner-Teller effect in a molecule with two degenerate bending vibrations. <i>Journal of Chemical Physics</i> , <b>2006</b> , 125, 133201	3.9	17
60	Kinetics of the simplest Criegee intermediate reaction with ozone studied using a mid-infrared quantum cascade laser spectrometer. <i>Physical Chemistry Chemical Physics</i> , <b>2017</b> , 20, 97-102	3.6	17
59	Dynamics of Atomic and Molecular Hydrogen Elimination from Small Alkanes Following 157-nm Excitation. <i>Journal of Physical Chemistry A</i> , <b>2000</b> , 104, 7189-7199	2.8	16
58	Reactivity of Criegee Intermediates toward Carbon Dioxide. <i>Journal of Physical Chemistry Letters</i> , <b>2018</b> , 9, 184-188	6.4	15
57	Barrierless reactions between two closed-shell molecules. I. Dynamics of F(2)+CH(3)SCH(3) reaction. <i>Journal of Chemical Physics</i> , <b>2008</b> , 128, 104317	3.9	15
56	Dissociation rates of benzene at VUV excitation. <i>Journal of Chemical Physics</i> , <b>2001</b> , 115, 9623-9626	3.9	15
55	Multiple dynamical pathways in the O+SiH <sub>4</sub> reaction studied by the crossed molecular beam method. <i>Journal of Chemical Physics</i> , <b>2000</b> , 113, 1831-1842	3.9	15
54	Temperature and isotope effects in the reaction of CHCHO with methanol. <i>Physical Chemistry Chemical Physics</i> , <b>2019</b> , 21, 13633-13640	3.6	13
53	Unimolecular decomposition rates of a methyl-substituted Criegee intermediate -CHCHO. <i>RSC Advances</i> , <b>2020</b> , 10, 8518-8524	3.7	13
52	Does ozone-water complex produce additional OH radicals in the atmosphere?. <i>Journal of Physical Chemistry A</i> , <b>2012</b> , 116, 12082-8	2.8	13
51	UV photolysis of ClOOCl and the ozone hole. <i>Chemistry - an Asian Journal</i> , <b>2011</b> , 6, 1664-78	4.5	13
50	Photodissociation dynamics of ClOOCl at 248.4 and 308.4 nm. <i>Physical Chemistry Chemical Physics</i> , <b>2011</b> , 13, 8195-203	3.6	13

49	Photodissociation dynamics of OCLO at 157 nm. <i>Journal of Chemical Physics</i> , <b>1998</b> , 108, 10061-10069	3.9	13
48	O(1D) reaction with cyclopropane: Evidence of O atom insertion into the C-C bond. <i>Journal of Chemical Physics</i> , <b>2001</b> , 115, 7-10	3.9	12
47	Water Vapor Does Not Catalyze the Reaction between Methanol and OH Radicals. <i>Angewandte Chemie - International Edition</i> , <b>2019</b> , 58, 5013-5017	16.4	12
46	Dynamics of the F2 reaction with the simplest pi-bonding molecule. <i>Journal of Chemical Physics</i> , <b>2008</b> , 128, 184302	3.9	11
45	Dynamics of the F2+CH3SCH3 reaction: a molecule-molecule reaction without entrance barrier. <i>Journal of Chemical Physics</i> , <b>2007</b> , 127, 101101	3.9	11
44	Multiple channel dynamics of the O+CH3F reaction. <i>Journal of Chemical Physics</i> , <b>2000</b> , 113, 9678-9685	3.9	11
43	Effects of water vapor on the reaction of CHOO with NH. <i>Physical Chemistry Chemical Physics</i> , <b>2019</b> , 21, 22589-22597	3.6	11
42	Surprisingly long lifetime of methacrolein oxide, an isoprene derived Criegee intermediate, under humid conditions. <i>Communications Chemistry</i> , <b>2021</b> , 4,	6.3	11
41	Temperature-Dependent Rate Coefficient for the Reaction of CHSH with the Simplest Criegee Intermediate. <i>Journal of Physical Chemistry A</i> , <b>2019</b> , 123, 4096-4103	2.8	10
40	Selective ionization of photofragments using tunable radiation from a synchrotron. <i>Journal of Electron Spectroscopy and Related Phenomena</i> , <b>2005</b> , 144-147, 135-138	1.7	10
39	A complete look at the dissociation dynamics of vinylfluoride at 157 nm. <i>Journal of Chemical Physics</i> , <b>2002</b> , 116, 6982-6989	3.9	10
38	UV Spectrum of the Simplest Deuterated Criegee Intermediate CD2OO. <i>Journal of the Chinese Chemical Society</i> , <b>2017</b> , 64, 360-368	1.5	9
37	Hydrogen-Bonding Mediated Reactions of Criegee Intermediates in the Gas Phase: Competition between Bimolecular and Termolecular Reactions and the Catalytic Role of Water. <i>Journal of Physical Chemistry A</i> , <b>2019</b> , 123, 8336-8348	2.8	9
36	Synergy of Water and Ammonia Hydrogen Bonding in a Gas-Phase Reaction. <i>Journal of Physical Chemistry A</i> , <b>2019</b> , 123, 1337-1342	2.8	9
35	Radical intermediates in the addition of OH to propene: photolytic precursors and angular momentum effects. <i>Journal of Physical Chemistry A</i> , <b>2014</b> , 118, 3211-29	2.8	9
34	Photofragmentation translational spectroscopy of methyl azide (CH3N3) photolysis at 193 nm: molecular and radical channel product branching ratio. <i>Journal of Physical Chemistry A</i> , <b>2012</b> , 116, 4695-704	2.8	9
33	Dynamics of the F2 reaction with propene: the effect of methyl substitution. <i>Journal of Physical Chemistry A</i> , <b>2009</b> , 113, 4381-6	2.8	9
32	Investigations of oxysilanes from the crossed-beam reaction of atomic oxygen with silane using tunable vacuum-ultraviolet ionization. <i>Chemical Physics Letters</i> , <b>2007</b> , 444, 237-241	2.5	9



31	Experimental and theoretical investigations of the O(1D) reaction with cyclopropane. <i>Journal of Chemical Physics</i> , <b>2002</b> , 116, 8292	3.9	9
30	Photodissociation dynamics of trifluoroethylene at 157 nm excitation. <i>Journal of Chemical Physics</i> , <b>1998</b> , 109, 10719-10726	3.9	9
29	Photoproduct channels from BrCD <sub>2</sub> CD <sub>2</sub> OH at 193 nm and the HDO + vinyl products from the CD <sub>2</sub> CD <sub>2</sub> OH radical intermediate. <i>Journal of Physical Chemistry A</i> , <b>2012</b> , 116, 6394-407	2.8	8
28	Note: a transient absorption spectrometer using an ultra bright laser-driven light source. <i>Review of Scientific Instruments</i> , <b>2013</b> , 84, 086106	1.7	8
27	Site specificity in molecular hydrogen elimination from photodissociation of propane at 157 nm. <i>Journal of Chemical Physics</i> , <b>1999</b> , 111, 1793-1796	3.9	8
26	Dynamics of reactions O((1)D)+C(6)H(6) and C(6)D(6). <i>Journal of Chemical Physics</i> , <b>2008</b> , 129, 174303	3.9	7
25	Crossed molecular beam studies of the O(1D)+NH <sub>3</sub> reaction. <i>Journal of Chemical Physics</i> , <b>2001</b> , 115, 842-848	3.9	7
24	Site specific dissociation dynamics of propane at 157 nm excitation. <i>Journal of Chemical Physics</i> , <b>2000</b> , 112, 8027-8037	3.9	7
23	Kinetics of Unimolecular Decay of Methyl Vinyl Ketone Oxide, an Isoprene-Derived Criegee Intermediate, under Atmospherically Relevant Conditions. <i>Journal of Physical Chemistry A</i> , <b>2020</b> , 124, 9375-9381	2.8	6
22	The role of the iodine-atom adduct in the synthesis and kinetics of methyl vinyl ketone oxide-a resonance-stabilized Criegee intermediate. <i>Physical Chemistry Chemical Physics</i> , <b>2020</b> , 22, 13603-13612	3.6	6
21	Dynamics of reactions between two closed-shell molecules. <i>Physical Chemistry Chemical Physics</i> , <b>2011</b> , 13, 19206-13	3.6	6
20	Barrierless reactions between two closed-shell molecules. II. Dynamics of F <sub>2</sub> + CH <sub>3</sub> SSCH <sub>3</sub> reaction. <i>Journal of Chemical Physics</i> , <b>2009</b> , 130, 014301	3.9	6
19	Exploring the dynamics of reaction N+SiH(4) with crossed molecular-beam experiments and quantum-chemical calculations. <i>Journal of Chemical Physics</i> , <b>2008</b> , 129, 174304	3.9	5
18	Absolute Infrared Absorption Cross Section of the Simplest Criegee Intermediate Near 1285.7 cm. <i>Journal of Physical Chemistry A</i> , <b>2018</b> , 122, 8874-8881	2.8	5
17	A crossed beam study of 18O(3P)+NO <sub>2</sub> and 18O(1D)+NO <sub>2</sub> : isotope exchange and O <sub>2</sub> +NO formation channels. <i>Journal of Chemical Physics</i> , <b>2012</b> , 137, 044302	3.9	4
16	Dissociative photoionization of ClN <sub>3</sub> using high-resolution synchrotron radiation: The N-Cl bond energy in ClN <sub>3</sub> . <i>International Journal of Mass Spectrometry</i> , <b>2007</b> , 265, 261-266	1.9	4
15	Crossed molecular beam studies on the reaction dynamics of O(1D)+N <sub>2</sub> O. <i>Journal of Chemical Physics</i> , <b>2006</b> , 125, 133121	3.9	4
14	Dynamics of Atomic and Molecular Hydrogen Elimination from Hydrocarbons at VUV Excitation. <i>Journal of the Chinese Chemical Society</i> , <b>1999</b> , 46, 435-444	1.5	4



13	Water Vapor Does Not Catalyze the Reaction between Methanol and OH Radicals. <i>Angewandte Chemie</i> , <b>2019</b> , 131, 5067-5071	3.6	3
12	Electronic quenching of O( <sup>1</sup> D) by Xe: Oscillations in the product angular distribution and their dependence on collision energy. <i>Journal of Chemical Physics</i> , <b>2015</b> , 143, 054307	3.9	3
11	Ion dissociation of hydrazoic acid investigated by synchrotron-radiation-based photoionization mass spectrometry. <i>Journal of Physical Chemistry A</i> , <b>2009</b> , 113, 3822-9	2.8	3
10	Double methyl substitution in simple alkenes leads to low-barrier reactions towards molecular fluorine. <i>Chemical Physics Letters</i> , <b>2011</b> , 510, 42-47	2.5	3
9	Photodissociation Dynamics of Benzene at 193 nm. <i>Journal of the Chinese Chemical Society</i> , <b>2002</b> , 49, 1-6	1.5	3
8	Reactions of Criegee Intermediates are Enhanced by Hydrogen-Atom Relay Through Molecular Design. <i>ChemPhysChem</i> , <b>2020</b> , 21, 2056-2059	3.2	3
7	Temperature-dependent kinetics of the simplest Criegee intermediate reaction with dimethyl sulfoxide. <i>Journal of the Chinese Chemical Society</i> , <b>2020</b> , 67, 1563-1570	1.5	3
6	Further studies into the photodissociation pathways of 2-bromo-2-nitropropane and the dissociation channels of the 2-nitro-2-propyl radical intermediate. <i>Journal of Physical Chemistry A</i> , <b>2014</b> , 118, 4707-22	2.8	2
5	Formation of NO <sub>3</sub> in the Photolysis of N <sub>2</sub> O <sub>5</sub> <b>2013</b> , 1,		2
4	Kinetics of dimethyl sulfide (DMS) reactions with isoprene-derived Criegee intermediates studied with direct UV absorption. <i>Atmospheric Chemistry and Physics</i> , <b>2020</b> , 20, 12983-12993	6.8	2
3	Substituent Effect in the Reactions between Criegee Intermediates and 3-Aminopropanol. <i>Journal of Physical Chemistry A</i> , <b>2021</b> , 125, 6580-6590	2.8	2
2	Photolysis cross-section of ozone dimer. <i>Chemistry - an Asian Journal</i> , <b>2011</b> , 6, 2925-30	4.5	1
1	The heat of formation of chlorine-isocyanate and the relative stability of isoelectronic molecules: an experimental and theoretical study. <i>Journal of Chemical Physics</i> , <b>2006</b> , 124, 241106	3.9	1