Scott L Anderson

List of Publications by Year
in descending order

196 papers	8,184 citations	34105 52 h-index	62596 80 g-index
199 all docs	199 docs citations	199 times ranked	5359 citing authors

#	Article	IF	Citations
1	Electronic Structure Controls Reactivity of Size-Selected Pd Clusters Adsorbed on TiO ₂ Surfaces. Science, 2009, 326, 826-829.	12.6	552
2	CO Oxidation on Aun/TiO2Catalysts Produced by Size-Selected Cluster Deposition. Journal of the American Chemical Society, 2004, 126, 5682-5683.	13.7	338
3	Scalable and safe synthetic organic electroreduction inspired by Li-ion battery chemistry. Science, 2019, 363, 838-845.	12.6	305
4	Collision-induced dissociation and ab initio studies of boron cluster ions: determination of structures and stabilities. The Journal of Physical Chemistry, 1988, 92, 5803-5812.	2.9	244
5	Collisionâ€induced dissociation of aluminum cluster ions: Fragmentation patterns, bond energies, and structures for Al+2â€“Al+7. Journal of Chemical Physics, 1987, 87, 260-268.	3.0	160
6	Oxide-Free, Catalyst-Coated, Fuel-Soluble, Air-Stable Boron Nanopowder as Combined Combustion Catalyst and High Energy Density Fuel. Energy & Fuels, 2009, 23, 6111-6120.	5.1	132
7	Chemistry of metal and semimetal cluster ions. Chemical Reviews, 1992, 92, 1541-1565.	47.7	127
8	Ethylene Dehydrogenation on Pt_{4,7,8} Clusters on Al₂O₃: Strong Cluster Size Dependence Linked to Preferred Catalyst Morphologies. ACS Catalysis, 2017, 7, 3322-3335.	11.2	124
9	Breakdown and Combustion of JP-10 Fuel Catalyzed by Nanoparticulate CeO2and Fe2O3. Energy & Fuels, 2006, 20, 1886-1894.	5.1	122
10	Collision ofLi+andNa+withC60: Insertion, fragmentation, and thermionic emission. Physical Review Letters, 1992, 69, 1352-1355.	7.8	110
11	Strong Effects of Cluster Size and Air Exposure on Oxygen Reduction and Carbon Oxidation Electrocatalysis by Size-Selected Pt_n(nâ‰ˆ 11) on Glassy Carbon Electrodes. Journal of the American Chemical Society, 2013, 135, 3073-3086.	13.7	109
12	Boron Switch for Selectivity of Catalytic Dehydrogenation on Size-Selected Pt Clusters on Al₂O₃. Journal of the American Chemical Society, 2017, 139, 11568-11575.	13.7	103
13	Ne++C60: Collision energy and impact parameter dependence for endohedral complex formation, fragmentation, and charge transfer. Journal of Chemical Physics, 1992, 96, 3344-3347.	3.0	98
14	Interaction of boron cluster ions with water: Single collision dynamics and sequential etching. Journal of Chemical Physics, 1990, 92, 292-303.	3.0	97
15	Resonance-enhanced multiphoton ionization of molecular hydrogen via the E,F1Î£g+ state: Photoelectron energy and angular distributions. Chemical Physics Letters, 1984, 105, 22-27.	2.6	96
16	Agglomeration, support effects, and CO adsorption on Au/TiO2(110) prepared by ion beam deposition. Surface Science, 2005, 578, 5-19.	1.9	92
17	Production and collision-induced dissociation of small boron cluster ions. The Journal of Physical Chemistry, 1987, 91, 5161-5163.	2.9	91
18	Oxidation of small boron cluster ions (B+1â€“13) by oxygen. Journal of Chemical Physics, 1988, 89, 2848-2860.	3.0	91

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19	Collision of alkali ions with C60/C70: Insertion, thermionic emission, and fragmentation. Journal of Chemical Physics, 1993, 99, 5858-5870.	3.0	90
20	Fragmentation of nitrous oxide by monochromatic soft x rays. Journal of Chemical Physics, 1986, 85, 5755-5762.	3.0	89
21	Cluster size effects on CO oxidation activity, adsorbate affinity, and temporal behavior of model Aunâ^•TiO2 catalysts. Journal of Chemical Physics, 2005, 123, 124710.	3.0	87
22	Reactions of boron cluster ions (B+n,n=2â€“24) with N2O: NO versus NN bond activation as a function of size. Journal of Chemical Physics, 1991, 94, 6446-6458.	3.0	81
23	Size-dependent oxidation of Pdn (nâ©½13) on alumina/NiAl(110): Correlation with Pd core level binding energies. Surface Science, 2009, 603, 2764-2770.	1.9	81
24	Size-Dependent Oxygen Activation Efficiency over Pd_n/TiO₂(110) for the CO Oxidation Reaction. Journal of the American Chemical Society, 2010, 132, 13097-13099.	13.7	79
25	Electrocatalysis by Mass-Selected Pt_n Clusters. Accounts of Chemical Research, 2016, 49, 2632-2639.	15.6	79
26	The effect of vibration and translational energy on the reaction dynamics of the H+2 +H2 system. Journal of Chemical Physics, 1981, 75, 2153-2162.	3.0	76
27	Multiphoton ionization photoelectron spectroscopy: a new method for determining vibrational structure of molecular ions. Chemical Physics Letters, 1982, 93, 11-15.	2.6	73
28	Hypergolic ionic liquids to mill, suspend, and ignite boron nanoparticles. Chemical Communications, 2012, 48, 4311.	4.1	72
29	Ne++C60 collisions: The dynamics of charge and energy transfer, fragmentation, and endohedral complex formation. Journal of Chemical Physics, 1993, 99, 3468-3479.	3.0	71
30	Dissociation Energies for Small Carbon Cluster Ions (C2-19+) Measured by Collision-Induced Dissociation. The Journal of Physical Chemistry, 1995, 99, 10736-10741.	2.9	69
31	Boron cluster ion oxidation: Reactions with CO2, dissociation of boron cluster oxide (BnO+) ions, and sequential oxidation. Journal of Chemical Physics, 1991, 94, 2833-2847.	3.0	68
32	Cluster Size Controls Branching between Water and Hydrogen Peroxide Production in Electrochemical Oxygen Reduction at Pt_n/ITO. Journal of Physical Chemistry C, 2015, 119, 11160-11170.	3.1	68
33	Multiphoton ionization photoelectron spectroscopy of phenol: Vibrational frequencies and harmonic force field for the 2B1 cation. Journal of Chemical Physics, 1985, 82, 5329-5339.	3.0	66
34	GeradeRydberg states of acetylene studied by multiphoton ionization and photoelectron spectroscopy. Journal of Chemical Physics, 1987, 87, 5105-5115.	3.0	65
35	Fullerene (C61+) production and decomposition in carbon-13 (1+) + fullerene (C60) collisions: carbon-atom exchange and the fragmentation pattern as a function of energy. The Journal of Physical Chemistry, 1992, 96, 3574-3576.	2.9	65
36	O++C60Â· C60O+ production and decomposition, charge transfer, and formation of C59O+. Dopeyball or [CO@C58+. Chemical Physics Letters, 1992, 199, 373-378.	2.6	65

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37	Vibrational mode effects, scattering dynamics, and energy disposal in reaction of C2H+2with methane. Journal of Chemical Physics, 1995, 102, 1199-1216.	3.0	63
38	MPI photoelectron spectroscopy of ungerade excited states of acetylene: Intermediate state mixing and ion state selection. Journal of Chemical Physics, 1987, 87, 852-860.	3.0	62
39	Deposition dynamics and chemical properties of size-selected Ir clusters on TiO2. Surface Science, 2003, 542, 253-275.	1.9	62
40	Modular terpene synthesis enabled by mild electrochemical couplings. Science, 2022, 375, 745-752.	12.6	62
41	Thermal decomposition of JP-10 studied by micro-flowtube pyrolysis-mass spectrometry. Combustion and Flame, 2006, 144, 662-674.	5.2	61
42	The effects of collision energy and ion vibrational excitation on proton and charge transfer in H2++N2, CO, O2. Journal of Chemical Physics, 1982, 77, 1842-1854.	3.0	58
43	CO adsorption and desorption on size-selected Pdn/TiO2(110) model catalysts: Size dependence of binding sites and energies, and support-mediated adsorption. Journal of Chemical Physics, 2012, 136, 204705.	3.0	58
44	Nitrogen ion (N+) + C60 fullerene reactive scattering: substitution, charge transfer, and fragmentation. The Journal of Physical Chemistry, 1992, 96, 10597-10600.	2.9	57
45	Interaction of small boron cluster ions with HF. Journal of Chemical Physics, 1997, 106, 9511-9522.	3.0	56
46	Initial and Final State Effects in the Ultraviolet and X-ray Photoelectron Spectroscopy (UPS and XPS) of Size-Selected Pd_n Clusters Supported on TiO₂(110). Journal of Physical Chemistry C, 2015, 119, 6033-6046.	3.1	56
47	Proton affinities of hydrogen halides determined by the molecular beam photoionization method. Journal of Chemical Physics, 1979, 71, 605-609.	3.0	55
48	Direct dynamics study of energy transfer and collision-induced dissociation: Effects of impact energy, geometry, and reactant vibrational mode in H2CO+â€“Ne collisions. Journal of Chemical Physics, 2003, 119, 3040-3050.	3.0	55
49	Dissociation energies for carbon cluster ions (C+2â€“15): A system where photodissociation is misleading. Journal of Chemical Physics, 1991, 95, 4719-4720.	3.0	54
50	Air-stable, unoxidized, hydrocarbon-dispersible boron nanoparticles. Journal of Materials Research, 2009, 24, 3462-3464.	2.6	54
51	Size-dependent electronic structure controls activity for ethanol electro-oxidation at Pt_n/indium tin oxide (n = 1 to 14). Physical Chemistry Chemical Physics, 2015, 17, 17601-17610.	2.8	54
52	Dynamics of boron cluster ion reactions with deuterium: Adduct formation and decay. Journal of Chemical Physics, 1989, 91, 226-239.	3.0	53
53	Inherent Size Effects on XANES of Nanometer Metal Clusters: Size-Selected Platinum Clusters on Silica. Journal of Physical Chemistry C, 2017, 121, 361-374.	3.1	52
54	Large, modeâ€selective vibrational effect on the reaction of C2H+2 with methane. Journal of Chemical Physics, 1994, 101, 5410-5412.	3.0	49

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55	Proton transfer in the [phenol-NH3]+ system: An experimental and ab initio study. Journal of Chemical Physics, 2000, 112, 5717-5721.	3.0	49
56	Photoionization of (H2)2 and the clusters of O2 molecules. Journal of Chemical Physics, 1980, 73, 4779-4783.	3.0	48
57	Simple radio-frequency power source for ion guides and ion traps. Review of Scientific Instruments, 1997, 68, 3357-3362.	1.3	48
58	Sintering, oxidation, and chemical properties of size-selected nickel clusters on TiO2(110). Journal of Chemical Physics, 2002, 117, 5001-5011.	3.0	48
59	Multiphoton Ionization State Selection: Vibrational-Mode and Rotational-State Control. Advances in Chemical Physics, 2007, , 177-212.	0.3	47
60	The effects of bending and stretching vibration on the reaction of acetylene cations with methane. Journal of Chemical Physics, 1989, 90, 1577-1587.	3.0	46
61	Exploring the Structure of Nitrogen-Rich Ionic Liquids and Their Binding to the Surface of Oxide-Free Boron Nanoparticles. Journal of Physical Chemistry C, 2013, 117, 5693-5707.	3.1	45
62	Nonadiabaticity in ionâ€“molecule reactions: Coupling of proton and charge transfer in the H+2 and D+2+Ar system. Journal of Chemical Physics, 1982, 77, 748-755.	3.0	44
63	Interaction of Mn+ and Mn2+ with C60. Exohedral and endohedral metal-fullerene bonding. Chemical Physics Letters, 1995, 243, 45-48.	2.6	42
64	Chemistry and cooling of transition metal cluster ions. Chemical Physics Letters, 1985, 122, 410-414.	2.6	40
65	Methane ignition catalyzed by in situ generated palladium nanoparticles. Combustion and Flame, 2010, 157, 421-435.	5.2	40
66	Oxygen activation and CO oxidation over size-selected Pt_n/alumina/Re(0001) model catalysts: correlations with valence electronic structure, physical structure, and binding sites. Physical Chemistry Chemical Physics, 2014, 16, 26443-26457.	2.8	40
67	Coking-Resistant Sub-Nano Dehydrogenation Catalysts: Pt_nSn_x/SiO₂ (n = 4, 7). ACS Catalysis, 2020, 10, 4543-4558.	11.2	40
68	Multiphoton ionization photoelectron spectroscopy study of OCS: Rydberg vibronic structure and ion state selection. Journal of Chemical Physics, 1988, 89, 5527-5534.	3.0	39
69	Multiphoton ionization and photoelectron spectroscopy of formaldehyde via its 3p Rydberg states. Journal of Chemical Physics, 2001, 114, 9797-9806.	3.0	39
70	Pyrolysis Chemistry of Cubane and Methylcubane:â€‰ The Effect of Methyl Substitution on Stability and Product Branching. Journal of Physical Chemistry A, 2003, 107, 1162-1174.	2.5	39
71	Single CdSe/ZnS Nanocrystals in an Ion Trap: Charge and Mass Determination and Photophysics Evolution with Changing Mass, Charge, and Temperature. ACS Nano, 2014, 8, 2387-2398.	14.6	39
72	Observation of circular dichroism in photoelectron angular distributions. Journal of Chemical Physics, 1986, 85, 6803-6804.	3.0	38

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73	Collision-induced dissociation of formaldehyde cations: The effects of vibrational mode, collision energy, and impact parameter. Journal of Chemical Physics, 2002, 116, 5530-5543.	3.0	38
74	Direct Dynamics Trajectory Study of Vibrational Effects:Â Can Polanyi Rules Be Generalized to a Polyatomic System?. Journal of the American Chemical Society, 2004, 126, 8602-8603.	13.7	38
75	Agglomeration, Sputtering, and Carbon Monoxide Adsorption Behavior for Au/Al2O3Prepared by Aun+Deposition on Al2O3/NiAl(110). Journal of Physical Chemistry B, 2005, 109, 11340-11347.	2.6	38
76	Fragmentation of acetone following excitation in the region of the oxygen K edge. Journal of Chemical Physics, 1987, 86, 4442-4445.	3.0	37
77	Reaction of aluminum cluster ions with oxygen and nitrous oxide: Energetics and dynamics of cluster oxidation. Journal of Chemical Physics, 1988, 89, 273-286.	3.0	37
78	Reaction mechanisms and energy disposal in the [C2H2:OCS]+ system: A modeâ€selective differential cross section study. Journal of Chemical Physics, 1996, 105, 3089-3107.	3.0	37
79	Simplified radio-frequency generator for driving ion guides, traps, and other capacitive loads. Review of Scientific Instruments, 2000, 71, 4335.	1.3	36
80	Hydrazine Decomposition over Irn/Al2O3Model Catalysts Prepared by Size-Selected Cluster Deposition. Journal of Physical Chemistry B, 2005, 109, 381-388.	2.6	36
81	In Situ Generation of Pd/PdO Nanoparticle Methane Combustion Catalyst: Correlation of Particle Surface Chemistry with Ignition. Journal of Physical Chemistry C, 2009, 113, 20632-20639.	3.1	36
82	Water on Rutile TiO2(110) and Au/TiO2(110): Effects on Au Mobility and the Isotope Exchange Reaction. Journal of Physical Chemistry C, 2008, 112, 9006-9015.	3.1	35
83	Alumina support and Pdn cluster size effects on activity of Pdn for catalytic oxidation of CO. Faraday Discussions, 2013, 162, 323.	3.2	35
84	A phase-space-compressing, mass-selecting beamline for hyperthermal, focused ion beam deposition. Review of Scientific Instruments, 1998, 69, 4106-4115.	1.3	34
85	Mode-Selective Differential Scattering as a Probe of Polyatomic Ion Reaction Mechanisms. Accounts of Chemical Research, 1997, 30, 28-36.	15.6	33
86	The effects of different vibrational modes and collision energy on the reaction of acetylene cations with carbonyl sulfide. Journal of Chemical Physics, 1990, 92, 7356-7364.	3.0	32
87	Complex formation and decay in ion-molecule reactions: Mode-selective scattering as a dynamical probe. International Reviews in Physical Chemistry, 2001, 20, 165-188.	2.3	30
88	Synthesis of Nanoparticles from Malleable and Ductile Metals Using Powder-Free, Reactant-Assisted Mechanical Attrition. ACS Applied Materials & Interfaces, 2014, 6, 19579-19591.	8.0	30
89	Boron Nanoparticles with High Hydrogen Loading: Mechanism for Bâ€“H Binding and Potential for Improved Combustibility and Specific Impulse. ACS Applied Materials & Interfaces, 2014, 6, 8513-8525.	8.0	30
90	Binding of Alkenes and Ionic Liquids to Bâ€“H-Functionalized Boron Nanoparticles: Creation of Particles with Controlled Dispersibility and Minimal Surface Oxidation. ACS Applied Materials & Interfaces, 2015, 7, 9991-10003.	8.0	29

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91	Vibrational mode-selected differential scattering of NH3+ methanol (d1, d3, d4): Control of product branching by hydrogen-bonded complex formation. Journal of Chemical Physics, 1998, 108, 2395-2407.	3.0	27
92	Cluster size effects on hydrazine decomposition on Irn/Al2O3/NiAl(110). Surface Science, 2006, 600, 461-467.	1.9	27
93	Functionalization and Passivation of Boron Nanoparticles with a Hypergolic Ionic Liquid. Journal of Propulsion and Power, 2013, 29, 489-495.	2.2	27
94	Dynamics of the C2H2++ ND3Reaction:Â A Vibrational-Mode-Selective Scattering Study. Journal of Physical Chemistry A, 1997, 101, 6504-6512.	2.5	26
95	Boron Oxide Oligomer Collision-Induced Dissociation:Â Thermochemistry, Structure, and Implications for Boron Combustion. Journal of Physical Chemistry A, 1997, 101, 9935-9941.	2.5	26
96	The effects of collision energy, vibrational mode, and vibrational angular momentum on energy transfer and dissociation in NO2+â€“rare gas collisions: An experimental and trajectory study. Journal of Chemical Physics, 2006, 125, 133115.	3.0	26
97	Cluster size effects on sintering, CO adsorption, and implantation in Ir/SiO2. Journal of Chemical Physics, 2009, 131, 114701.	3.0	26
98	Reaction of magnetically state selected NO with O3: Effect of fs states and rotational states on reactivity. Journal of Chemical Physics, 1980, 72, 6521-6528.	3.0	25
99	The effects of reactant vibrational, fine structure, and collision energy on the reactions of OCS+with C2H2: Complementary studies of reactions in the [C2H2+OCS]+system. Journal of Chemical Physics, 1991, 94, 6459-6468.	3.0	25
100	Comparison of bending, Câ€“C stretching, and collision energy effects on the reaction of C2H+2with D2. Journal of Chemical Physics, 1992, 96, 5781-5788.	3.0	25
101	Optically detected, single nanoparticle mass spectrometer with pre-filtered electrospray nanoparticle source. Review of Scientific Instruments, 2014, 85, 014104.	1.3	25
102	Preparation of Size- and Composition-Controlled Pt_nSn_x/SiO₂ (n = 4, 7, 24) Bimetallic Model Catalysts with Atomic Layer Deposition. Journal of Physical Chemistry C, 2019, 123, 16194-16209.	3.1	25
103	Sn-modification of Pt7/alumina model catalysts: Suppression of carbon deposition and enhanced thermal stability. Journal of Chemical Physics, 2020, 152, 024702.	3.0	25
104	The effects of vibrational mode, spinâ€“orbit state, and collision energy on collisionâ€induced dissociation and predissociation of OCS+. Journal of Chemical Physics, 1991, 95, 3275-3282.	3.0	24
105	Complex formation, rearrangement, and reaction in PhOH++ND3: Vibrational mode effects, recoil velocities, andab initiostudies. Journal of Chemical Physics, 2000, 113, 4158-4170.	3.0	24
106	Spectroscopic Study on the Intermediates and Reaction Rates in the Oxidation of Levitated Droplets of Energetic Ionic Liquids by Nitrogen Dioxide. Journal of Physical Chemistry A, 2018, 122, 7351-7377.	2.5	24
107	Third harmonic interference effects in the mpi spectrum of acetylene. Chemical Physics Letters, 1986, 129, 31-35.	2.6	23
108	A mode-selective differential scattering study of the C2H2++methanol reaction: Influence of collision intermediates, collision times, and transition states. Journal of Chemical Physics, 1998, 108, 7173-7184.	3.0	23

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109	Reaction of formaldehyde cation with methane: Effects of collision energy and H2CO+ and methane vibrations. Journal of Chemical Physics, 2003, 119, 200-214.	3.0	23
110	Effects of Alumina Thickness on CO Oxidation Activity over Pd₂₀/Alumina/Re(0001): Correlated Effects of Alumina Electronic Properties and Pd₂₀ Geometry on Activity. Journal of Physical Chemistry C, 2015, 119, 1359-1375.	3.1	23
111	Vibrational effects in proton and charge transfer in the H+2 + Ar system. Chemical Physics Letters, 1981, 82, 392-395.	2.6	22
112	Hydride abstraction by NO+ from ethanol: Effects of collision energy and ion rotational state. Journal of Chemical Physics, 2000, 113, 3002-3010.	3.0	22
113	Cluster ion beam study of a system with structural isomers: C+n+D2 (n=2â€“12). Chemical Physics Letters, 1991, 177, 146-152.	2.6	21
114	Pyrolysis and Isomerization of Quadricyclane, Norbornadiene, and Toluene. Journal of Physical Chemistry A, 1998, 102, 9202-9212.	2.5	21
115	Rapid Aluminum Nanoparticle Production by Milling in NH₃ and CH₃NH₂ Atmospheres: An Experimental and Theoretical Study. ACS Applied Materials & Interfaces, 2015, 7, 16101-16116.	8.0	21
116	Reactions of bare aluminum cluster ions. Chemical Physics Letters, 1986, 129, 429-432.	2.6	20
117	Nâ€“O versus Nâ€“N bond activation in reaction of N2O with carbon cluster ions: Experimental andabinitiostudies of the effects of geometric and electronic structure. Journal of Chemical Physics, 1994, 100, 8784-8794.	3.0	20
118	Collisions of rare gas ions with C60: Endohedral formation, energy transfer, and scattering dynamics. Journal of Chemical Physics, 1997, 107, 8370-8379.	3.0	20
119	The effects of vibrational mode and collision energy on the reaction of formaldehyde cation with carbonyl sulfide. Journal of Chemical Physics, 2002, 117, 8292-8307.	3.0	20
120	Dynamical control of â€˜statisticalâ€™ ionâ€“molecule reactions. International Journal of Mass Spectrometry, 2005, 241, 173-184.	1.5	20
121	Thermal and adsorbate effects on the activity and morphology of size-selected Pdn/TiO2 model catalysts. Surface Science, 2014, 621, 40-50.	1.9	20
122	Aluminum Nanoparticle Production by Acetonitrile-Assisted Milling: Effects of Liquid- vs Vapor-Phase Milling and of Milling Method on Particle Size and Surface Chemistry. Journal of Physical Chemistry C, 2016, 120, 19613-19629.	3.1	20
123	Ion beam studies of atomic ion collisions with C60: chemistry at surface, substitutional, and endohedral sites. International Journal of Mass Spectrometry and Ion Processes, 1994, 138, 173-185.	1.8	19
124	Effects of Composition, Structure, and H Atom Addition on the Chemistry of Boron Oxide Cluster Ions with HF. The Journal of Physical Chemistry, 1995, 99, 16276-16283.	2.9	19
125	Transition-metal C60 bonding by guided ion beam scattering. International Journal of Mass Spectrometry, 1999, 185-187, 603-615.	1.5	19
126	Alloying with Sn Suppresses Sintering of Size-Selected Subnano Pt Clusters on SiO₂ with and without Adsorbates. Chemistry of Materials, 2020, 32, 8595-8605.	6.7	19

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127	Direct Dynamics Trajectory Study of the Reaction of Formaldehyde Cation with D2:Â Vibrational and Zero-Point Energy Effects on Quasiclassical Trajectoriesâ€. Journal of Physical Chemistry A, 2005, 109, 11376-11384.	2.5	18
128	Multiphoton ionization photoelectron spectroscopy of acetaldehyde via the Ãƒâ€‰1Aâ€³, BÌƒ, CÌƒ, and DÌƒ states. Journal of Chemical Physics, 2001, 114, 3018-3028.	3.0	17
129	Spectroscopic Investigation of the Primary Reaction Intermediates in the Oxidation of Levitated Droplets of Energetic Ionic Liquids. Journal of Physical Chemistry Letters, 2017, 8, 6053-6059.	4.6	17
130	Diborane Interactions with Pt₇/Alumina: Preparation of Size-Controlled Borated Pt Model Catalysts. Journal of Physical Chemistry C, 2018, 122, 1631-1644.	3.1	17
131	Use of a quadrupole mass filter for high energy resolution ion beam production. Review of Scientific Instruments, 1995, 66, 3706-3708.	1.3	16
132	Vibrational mode and collision energy effects on a highly constrained reaction: OCS+(Î½)+OCSâ†'CS+2+CO2 and S+2+2 CO. Journal of Chemical Physics, 1995, 102, 1188-1191.	3.0	16
133	Kinetic parameters for heterogenous boron combustion reactions via the Cluster Beam approach. Combustion and Flame, 1996, 105, 68-79.	5.2	16
134	Multiphoton ionization vibrational state selection of H2O+, D2O+ and HDO+. Chemical Physics Letters, 2007, 440, 171-175.	2.6	16
135	In Situ Small-Angle X-ray Scattering from Pd Nanoparticles Formed by Thermal Decomposition of Organo-Pd Catalyst Precursors Dissolved in Hydrocarbons. Journal of Physical Chemistry C, 2013, 117, 22627-22635.	3.1	16
136	Mass-selected supported cluster catalysts: Size effects on CO oxidation activity, electronic structure, and thermal stability of Pdn/alumina (nâ‰30) model catalysts. International Journal of Mass Spectrometry, 2014, 370, 1-15.	1.5	16
137	Effects of acetonitrile-assisted ball-milled aluminum nanoparticles on the ignition of acoustically levitated exo-tetrahydrodicyclopentadiene (JP-10) droplets. Chemical Physics Letters, 2020, 754, 137679.	2.6	16
138	Oxidation of small carbon cluster ions by O2: Effects of structure on the reaction mechanism. Journal of Chemical Physics, 1992, 97, 8164-8172.	3.0	15
139	Photoluminescence of Charged CdSe/ZnS Quantum Dots in the Gas Phase: Effects of Charge and Heating on Absorption and Emission Probabilities. ACS Nano, 2014, 8, 12534-12548.	14.6	15
140	Effect of O₂ and CO Exposure on the Photoelectron Spectroscopy of Size-Selected Pd_n Clusters Supported on TiO₂(110). Journal of Physical Chemistry C, 2016, 120, 2126-2138.	3.1	15
141	Reaction of formaldehyde cation with molecular hydrogen: Effects of collision energy and H2CO+ vibrations. Journal of Chemical Physics, 2004, 120, 8528-8536.	3.0	14
142	Vibrational Mode Effects as a Probe of Inter-channel Coupling in the Reactions of Formaldehyde Cation with Ammonia and Waterâ€. Journal of Physical Chemistry A, 2004, 108, 9945-9956.	2.5	14
143	State-Selected C2H2+ Reactions with Methane at High Internal Energies. H+ and H- Transfer Reactions, Two New Channels in the C2H2+ A State Region. The Journal of Physical Chemistry, 1995, 99, 15523-15531.	2.9	13
144	Single Nanoparticle Mass Spectrometry as a High Temperature Kinetics Tool: Sublimation, Oxidation, and Emission Spectra of Hot Carbon Nanoparticles. Journal of Physical Chemistry A, 2015, 119, 12538-12550.	2.5	13

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145	Thermally Brightened CdSe/ZnS Quantum Dots as Noncontact Probes for Surface Chemistry Studies of Dark Nanoparticles Trapped in the Gas Phase. Journal of Physical Chemistry C, 2015, 119, 14561-14570.	3.1	13
146	Thermal emission spectroscopy for single nanoparticle temperature measurement: optical system design and calibration. Applied Optics, 2019, 58, 642.	1.8	13
147	Vibrational mode and collision energy effects on proton transfer in phenol cationâ€“methylamine collisions. Journal of Chemical Physics, 2000, 112, 10831-10837.	3.0	12
148	Thermal Emission Spectroscopy of Single, Isolated Carbon Nanoparticles: Effects of Particle Size, Material, Charge, Excitation Wavelength, and Thermal History. Journal of Physical Chemistry C, 2020, 124, 1704-1716.	3.1	12
149	Reactions of Boron Oxide and BnOmH+ Cluster Ions with Water. Journal of Physical Chemistry A, 1999, 103, 226-234.	2.5	11
150	Vibrational and collision energy effects on the reaction of CH3CHO+ with methanol. Journal of Chemical Physics, 2001, 115, 5843-5858.	3.0	11
151	Reaction of acetaldehyde cations with water: The effects of CH3CHO+ vibrational mode and impact parameter on reactivity and product branching. Journal of Chemical Physics, 2001, 115, 1274-1286.	3.0	11
152	Vibrational mode and collision energy effects on reaction of H2CO+ with C2D4. Journal of Chemical Physics, 2004, 121, 11746-11759.	3.0	10
153	In situ X-ray Scattering and Dynamical Modeling of Pd Catalyst Nanoparticles Formed in Flames. Journal of Physical Chemistry C, 2015, 119, 19073-19082.	3.1	10
154	Combustion Behavior of High Energy Density Boraneâ€“Aluminum Nanoparticles in Hypergolic Ionic Liquids. Energy & Fuels, 2018, 32, 7898-7908.	5.1	10
155	Selective growth of Al2O3 on size-selected platinum clusters by atomic layer deposition. Surface Science, 2020, 691, 121485.	1.9	10
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