

# Cheuk-Yiu Ng

## List of Publications by Year in Descending Order

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

262  
papers

7,576  
citations

42  
h-index

65  
g-index

273  
ext. papers

7,840  
ext. citations

4  
avg, IF

5.68  
L-index

#	Paper	IF	Citations
262	Chemical activation of oxygen molecule by quantum electronic state selected vanadium cation: observation of spin-orbit state effects. <i>Molecular Physics</i> , <b>2021</b> , 119, e1767309	1.7	3
261	Quantum electronic control on chemical activation of methane by collision with spin-orbit state selected vanadium cation. <i>Physical Chemistry Chemical Physics</i> , <b>2021</b> , 23, 273-286	3.6	6
260	What is the Bond Dissociation Energy of the Vanadium Hydride Cation?. <i>Journal of Physical Chemistry A</i> , <b>2020</b> , 124, 5306-5313	2.8	4
259	Chemical Activation of Water Molecule by Collision with Spin-Orbit-State-Selected Vanadium Cation: Quantum-Electronic-State Control of Chemical Reactivity. <i>Journal of Physical Chemistry A</i> , <b>2020</b> , 124, 8884-8896	2.8	5
258	Photodissociation branching ratios of $^{12}\text{C}^{16}\text{O}$ from 108000 $\text{cm}^{-1}$ to 113200 $\text{cm}^{-1}$ measured by two-color VUV-VUV laser pump-probe time-slice velocity-map ion imaging method: Observation of channels for producing $\text{O}(^1\text{D})$ . <i>Chinese Journal of Chemical Physics</i> , <b>2020</b> , 33, 91-100	0.9	9
257	Branching Ratios of the $\text{N}(^1\text{D})$ and $\text{N}(^3\text{D})$ Spin-Orbit States Produced in the State-Selected Photodissociation of $\text{N}_2$ Determined Using Time-Sliced Velocity-Mapped-Imaging Photoionization Mass Spectrometry (TS-VMI-PI-MS). <i>Journal of Physical Chemistry A</i> , <b>2019</b> , 123, 2289-2300	2.8	10
256	Quantum state-to-state vacuum ultraviolet photodissociation dynamics of small molecules. <i>Chinese Journal of Chemical Physics</i> , <b>2019</b> , 32, 23-34	0.9	16
255	Quantum Spin-Orbit Electronic State Selection of Atomic Transition Metal Vanadium Cation for Chemical Reactivity Studies. <i>Journal of Physical Chemistry A</i> , <b>2019</b> , 123, 2310-2319	2.8	8
254	Quantum state control on the chemical reactivity of a transition metal vanadium cation in carbon dioxide activation. <i>Physical Chemistry Chemical Physics</i> , <b>2019</b> , 21, 6868-6877	3.6	10
253	High-Level Ab Initio Predictions for the Ionization Energy, Bond Dissociation Energies, and Heats of Formation of Vanadium Methylidyne Radical and Its Cation ( $\text{VCH}/\text{VCH}^+$ ). <i>Journal of Physical Chemistry A</i> , <b>2019</b> , 123, 7454-7462	2.8	2
252	Chemical Activation of a Deuterium Molecule by Collision with a Quantum Electronic State-Selected Vanadium Cation. <i>Journal of Physical Chemistry A</i> , <b>2019</b> , 123, 5937-5944	2.8	5
251	Ab Initio Study of Ground-state $\text{CS}$ Photodissociation via Highly Excited Electronic States. <i>Astrophysical Journal</i> , <b>2019</b> , 882, 86	4.7	3
250	High-level ab initio predictions for the ionisation energy, bond dissociation energies and heats of formation of zirconium oxide and its cation ( $\text{ZrO}/\text{ZrO}^+$ ). <i>Molecular Physics</i> , <b>2018</b> , 116, 2709-2718	1.7	4
249	Quantum-vibrational-state-selected Integral Cross Sections and Product Branching Ratios for the Ion-molecule Reactions of $\text{N}_2^+ + (\text{X } ^2\Sigma_g^+; v = 0, 1) + \text{H}_2\text{O}$ and $\text{H}_2\text{O} + (\text{X } ^2\Sigma_g^+; v_1 + v_2 + v_3 = 000 \text{ and } 100) + \text{N}_2$ in the Collision Energy Range of 0.04-10.00 eV. <i>Astrophysical Journal</i> , <b>2018</b> , 861, 17	4.7	7
248	Quantum-State-Selected Integral Cross Sections and Branching Ratios for the Ion-Molecule Reaction of $\text{N}(^4\Sigma_g^-; 0-2) + \text{CH}$ in the Collision Energy Range of 0.05-10.00 eV. <i>Journal of Physical Chemistry A</i> , <b>2018</b> , 122, 6491-6499	2.8	7
247	Branching Ratio Measurements of the Predissociation of $\text{CO}$ by Time-Slice Velocity-Map Ion Imaging in the Energy Region from 106 250 to 107 800 $\text{cm}^{-1}$ . <i>Journal of Physical Chemistry A</i> , <b>2018</b> , 122, 8136-8142	2.8	14
246	Two-Color Pulsed-Field Ionization-Photoelectron Spectroscopy: A Quest to Benchmark State-of-the-Art ab initio Quantum Electronic Structure Calculations of Spectroscopic and Energetic Properties for Transition Metal-Containing Species <b>2018</b> , 195-249		5

245	Isotopic and quantum-rovibrational-state effects for the ion-molecule reaction in the collision energy range of 0.03-10.00 eV. <i>Physical Chemistry Chemical Physics</i> , <b>2017</b> , 19, 8694-8705	3.6	10
244	High-Level ab Initio Predictions for the Ionization Energies, Bond Dissociation Energies, and Heats of Formation of Titanium Oxides and Their Cations (TiO/TiO, n = 1 and 2). <i>Journal of Physical Chemistry A</i> , <b>2017</b> , 121, 669-679	2.8	15
243	A quantum-rovibrational-state-selected study of the reaction in the collision energy range of 0.05-10.00 eV: translational, rotational, and vibrational energy effects. <i>Physical Chemistry Chemical Physics</i> , <b>2017</b> , 19, 9778-9789	3.6	11
242	A vacuum ultraviolet laser pulsed field ionization-photoion study of methane (CH): determination of the appearance energy of methylum from methane with unprecedented precision and the resulting impact on the bond dissociation energies of CH and CH. <i>Physical Chemistry Chemical Physics</i> , <b>2017</b> , 19, 8500-8505	3.6	14
241	A quantum-rovibrational-state-selected study of the proton-transfer reaction $H(X\bar{v} = 1-3; N = 0-3) + Ne \rightarrow NeH + H$ using the pulsed field ionization-photoion method: observation of the rotational effect near the reaction threshold. <i>Physical Chemistry Chemical Physics</i> , <b>2017</b> , 19, 18619-18627	3.6	8
240	Quantum-state-selected integral cross sections for the charge transfer collision of $O(a\bar{v} = 1-2; J) [O(X\bar{v} = 22-23; J)] + Ar$ at center-of-mass collision energies of 0.05-10.00 eV. <i>Physical Chemistry Chemical Physics</i> , <b>2017</b> , 19, 29057-29067	3.6	9
239	Branching Ratios in Vacuum Ultraviolet Photodissociation of CO and N <sub>2</sub> : Implications for Oxygen and Nitrogen Isotopic Compositions of the Solar Nebula. <i>Astrophysical Journal</i> , <b>2017</b> , 850, 48	4.7	16
238	Rotationally Resolved State-to-State Photoelectron Study of Molybdenum Monoxide Cation (MoO <sup>+</sup> ). <i>Journal of Physical Chemistry A</i> , <b>2016</b> , 120, 4643-54	2.8	12
237	QUANTUM-STATE DEPENDENCE OF PRODUCT BRANCHING RATIOS IN VACUUM ULTRAVIOLET PHOTODISSOCIATION OF N <sub>2</sub> . <i>Astrophysical Journal</i> , <b>2016</b> , 819, 23	4.7	18
236	High-Resolution Threshold Photoelectron Spectroscopy by Vacuum Ultraviolet Laser Velocity-Map-Imaging Method. <i>Chinese Journal of Chemical Physics</i> , <b>2016</b> , 29, 59-69	0.9	
235	ABSOLUTE INTEGRAL CROSS SECTIONS FOR THE STATE-SELECTED ION-MOLECULE REACTION $\{N\}_2^+ + X^2\{\Sigma\}_g^+; v=0$ + C <sub>2</sub> H <sub>2</sub> IN THE COLLISION ENERGY RANGE OF 0.03-10.00 eV. <i>Astrophysical Journal</i> , <b>2016</b> , 827, 17	4.7	10
234	A HIGH-RESOLUTION VACUUM ULTRAVIOLET LASER PHOTOIONIZATION AND PHOTOELECTRON STUDY OF THE CO ATOM. <i>Astrophysical Journal</i> , <b>2016</b> , 833, 205	4.7	2
233	Comparison of experimental and theoretical quantum-state-selected integral cross-sections for the H <sub>2</sub> O <sup>+</sup> + H <sub>2</sub> (D <sub>2</sub> ) reactions in the collision energy range of 0.04-10.00 eV. <i>Physical Chemistry Chemical Physics</i> , <b>2016</b> , 18, 22509-15	3.6	22
232	State-to-state vacuum ultraviolet photodissociation study of CO <sub>2</sub> on the formation of state-correlated CO(X(1) $\bar{v}$ ; v) with O((1)D) and O((1)S) photoproducts at 11.95-12.22 eV. <i>Physical Chemistry Chemical Physics</i> , <b>2015</b> , 17, 11752-62	3.6	14
231	Rotationally resolved state-to-state photoelectron study of zirconium monoxide cation (ZrO <sup>+</sup> ). <i>Molecular Physics</i> , <b>2015</b> , 113, 2228-2242	1.7	10
230	VACUUM ULTRAVIOLET LASER PHOTOION AND PULSED FIELD IONIZATION PHOTOION STUDY OF RYDBERG SERIES OF CHLORINE ATOMS PREPARED IN THE $\{P\}_J$ (J= 3/2 and 1/2) FINE-STRUCTURE STATES. <i>Astrophysical Journal</i> , <b>2015</b> , 810, 132	4.7	3
229	Rotationally Selected and Resolved State-to-State Photoelectron Study of Vanadium Monoxide Cation VO <sup>+</sup> (X(3) $\bar{v}$ ; v <sup>+</sup> = 0-3). <i>Journal of Physical Chemistry A</i> , <b>2015</b> , 119, 11162-9	2.8	11
228	Rotationally resolved state-to-state photoionization and the photoelectron study of vanadium monocarbide and its cations (VC/VC <sup>+</sup> ). <i>Physical Chemistry Chemical Physics</i> , <b>2015</b> , 17, 9780-93	3.6	14

227	State-to-state spectroscopy and dynamics of ions and neutrals by photoionization and photoelectron methods. <i>Annual Review of Physical Chemistry</i> , <b>2014</b> , 65, 197-224	15.7	30
226	Photodissociation of CO <sub>2</sub> between 13.540 eV and 13.678 eV. <i>Physical Chemistry Chemical Physics</i> , <b>2014</b> , 16, 563-9	3.6	18
225	Photochemistry. Evidence for direct molecular oxygen production in CO <sub>2</sub> photodissociation. <i>Science</i> , <b>2014</b> , 346, 61-4	33.3	78
224	Communication: direct measurements of nascent O( <sup>3</sup> P <sub>0,1,2</sub> ) fine-structure distributions and branching ratios of correlated spin-orbit resolved product channels CO( <sup>1</sup> Σ <sup>+</sup> ) + O( <sup>3</sup> P <sub>0,1,2</sub> ) and CO(X <sup>1</sup> Σ <sup>+</sup> ; v) + O( <sup>3</sup> P <sub>0,1,2</sub> ) in VUV photodissociation of CO <sub>2</sub> . <i>Journal of Chemical Physics</i> , <b>2014</b> , 140, 231101	3.9	13
223	Rotationally resolved state-to-state photoionization and photoelectron study of titanium carbide and its cation (TiC/TiC <sup>+</sup> ). <i>Journal of Chemical Physics</i> , <b>2014</b> , 141, 144307	3.9	15
222	Rotationally resolved state-to-state photoelectron study of niobium carbide radical. <i>Journal of Chemical Physics</i> , <b>2014</b> , 141, 024304	3.9	12
221	Communication: the origin of rotational enhancement effect for the reaction of H <sub>2</sub> O <sup>+</sup> + H <sub>2</sub> (D <sub>2</sub> ). <i>Journal of Chemical Physics</i> , <b>2014</b> , 140, 011102	3.9	38
220	Communication: State-to-state photoionization and photoelectron study of vanadium methylidyne radical (VCH). <i>Journal of Chemical Physics</i> , <b>2014</b> , 140, 181101	3.9	8
219	High-level ab initio predictions for the ionization energy, bond dissociation energies, and heats of formation of cobalt carbide (CoC) and its cation (CoC <sup>+</sup> ). <i>Journal of Chemical Physics</i> , <b>2013</b> , 138, 094302	3.9	22
218	State-to-state Photoionization Dynamics of Vanadium Nitride by Two-color Laser Photoionization and Photoelectron Methods. <i>Chinese Journal of Chemical Physics</i> , <b>2013</b> , 26, 669-678	0.9	16
217	Branching ratio measurements for vacuum ultraviolet photodissociation of 12C16O. <i>Journal of Physical Chemistry A</i> , <b>2013</b> , 117, 6185-95	2.8	25
216	Communication: A vibrational study of titanium dioxide cation using the vacuum ultraviolet laser pulsed field ionization-photoelectron method. <i>Journal of Chemical Physics</i> , <b>2013</b> , 138, 041101	3.9	11
215	Rovibronically selected and resolved two-color laser photoionization and photoelectron study of titanium monoxide cation. <i>Journal of Chemical Physics</i> , <b>2013</b> , 138, 174309	3.9	16
214	Communication: State-to-state photodissociation study by the two-color VUV-VUV laser pump-probe time-slice velocity-map-imaging-photoion method. <i>Journal of Chemical Physics</i> , <b>2013</b> , 138, 191102	3.9	25
213	Rovibronically selected and resolved two-color laser photoionization and photoelectron study of cobalt carbide cation. <i>Journal of Chemical Physics</i> , <b>2013</b> , 138, 094301	3.9	18
212	ABSOLUTE INTEGRAL CROSS SECTIONS AND PRODUCT BRANCHING RATIOS FOR THE VIBRATIONALLY SELECTED ION-MOLECULE REACTIONS: N <sub>2</sub> <sup>+</sup> (X <sup>2</sup> Σ <sup>+</sup> ; v=0-2) + CH <sub>4</sub> . <i>Astrophysical Journal</i> , <b>2013</b> , 769, 72	4.7	17
211	The translational, rotational, and vibrational energy effects on the chemical reactivity of water cation H <sub>2</sub> O <sup>+</sup> (X <sup>2</sup> B <sub>1</sub> ) in the collision with deuterium molecule D <sub>2</sub> . <i>Journal of Chemical Physics</i> , <b>2013</b> , 139, 024203	3.9	32
210	Time-slice velocity-map ion imaging studies of the photodissociation of NO in the vacuum ultraviolet region. <i>Journal of Chemical Physics</i> , <b>2012</b> , 136, 134302	3.9	10

209	A HIGH-RESOLUTION PHOTOIONIZATION AND PHOTOELECTRON STUDY OF $^{58}\text{Ni}$ USING A VACUUM ULTRAVIOLET LASER. <i>Astrophysical Journal</i> , <b>2012</b> , 747, 20	4.7	8
208	Communication: Rovibrationally selected absolute total cross sections for the reaction $\text{H}_2\text{O}^+(\text{X}^2\text{B}_1; v_1(+), v_2(+), v_3(+)) = 000; \text{N}^+(\text{K}^+\text{K}^+) + \text{D}_2$ : observation of the rotational enhancement effect. <i>Journal of Chemical Physics</i> , <b>2012</b> , 137, 241101	3.9	42
207	Branching ratio measurements of the predissociation of $^{12}\text{C}^{16}\text{O}$ by time-slice velocity-map ion imaging in the energy region from 108,000 to 110,500 $\text{cm}^{-1}$ . <i>Journal of Chemical Physics</i> , <b>2012</b> , 137, 034305	3.9	22
206	Rovibrationally selected ion-molecule collision study using the molecular beam vacuum ultraviolet laser pulsed field ionization-photoion method: charge transfer reaction of $\text{N}_2^+(\text{X}^2\Sigma^+; v^+ = 0-2; \text{N}^+ = 0-9) + \text{Ar}$ . <i>Journal of Chemical Physics</i> , <b>2012</b> , 137, 104202	3.9	29
205	Communication: A vibrational study of propargyl cation using the vacuum ultraviolet laser velocity-map imaging photoelectron method. <i>Journal of Chemical Physics</i> , <b>2012</b> , 137, 161101	3.9	8
204	High-resolution threshold photoelectron study of the propargyl radical by the vacuum ultraviolet laser velocity-map imaging method. <i>Journal of Chemical Physics</i> , <b>2011</b> , 135, 224304	3.9	21
203	Communication: branching ratio measurements in the predissociation of $^{12}\text{C}^{16}\text{O}$ by time-slice velocity-map ion imaging in the vacuum ultraviolet region. <i>Journal of Chemical Physics</i> , <b>2011</b> , 135, 221101	3.9	24
202	Note: Accurate ab initio predictions of ionization energies of propargyl and allyl radicals: revisited. <i>Journal of Chemical Physics</i> , <b>2011</b> , 135, 246101	3.9	7
201	A vacuum-ultraviolet laser pulsed field ionization-photoelectron study of sulfur monoxide (SO) and its cation (SO <sup>+</sup> ). <i>Journal of Chemical Physics</i> , <b>2011</b> , 134, 144304	3.9	12
200	Communication: rovibrationally selected study of the $\text{N}_2^+(\text{X}; v^+=1, \text{N}^+= 0-8) + \text{Ar}$ charge transfer reaction using the vacuum ultraviolet laser pulsed field ionization-photoion method. <i>Journal of Chemical Physics</i> , <b>2011</b> , 134, 201105	3.9	24
199	Communication: vacuum ultraviolet laser photodissociation studies of small molecules by the vacuum ultraviolet laser photoionization time-sliced velocity-mapped ion imaging method. <i>Journal of Chemical Physics</i> , <b>2011</b> , 135, 071101	3.9	28
198	Time-sliced velocity-mapped imaging studies of the predissociation of single ro-vibronic energy levels of $\text{N}_2$ in the extreme ultraviolet region using vacuum ultraviolet photoionization. <i>Journal of Chemical Physics</i> , <b>2011</b> , 135, 134319	3.9	16
197	High-level ab initio predictions for the ionization energy, bond dissociation energies, and heats of formation of nickel carbide (NiC) and its cation (NiC <sup>+</sup> ). <i>Journal of Chemical Physics</i> , <b>2010</b> , 133, 114304	3.9	22
196	Rovibronically selected and resolved two-color laser photoionization and photoelectron study of nickel carbide cation. <i>Journal of Chemical Physics</i> , <b>2010</b> , 133, 054310	3.9	20
195	High-level ab initio predictions for the ionization energy, bond dissociation energies, and heats of formations of iron carbide (FeC) and its cation (FeC <sup>+</sup> ). <i>Journal of Physical Chemistry A</i> , <b>2009</b> , 113, 14321-8	2.8	30
194	Comment on "Experimental test of self-shielding in vacuum ultraviolet photodissociation of CO". <i>Science</i> , <b>2009</b> , 324, 1516; author reply 1516	3.3	12
193	A HIGH-RESOLUTION PHOTOIONIZATION STUDY OF $^{56}\text{Fe}$ USING A VACUUM ULTRAVIOLET LASER. <i>Astrophysical Journal</i> , <b>2009</b> , 693, 940-945	4.7	12
192	Rovibronically selected and resolved two-color laser photoionization and photoelectron study of the iron carbide cation. <i>Journal of Physical Chemistry A</i> , <b>2009</b> , 113, 4242-8	2.8	34

191	Spectroscopy and Dynamics of Neutrals and Ions by High-Resolution Infrared/Vacuum Ultraviolet Photoionization and Photoelectron Methods <b>2009</b> , 659-691		8
190	Infrared vacuum-ultraviolet laser pulsed field ionization-photoelectron study of CH <sub>3</sub> Br <sup>+</sup> (X(2)E(3/2)). <i>Journal of Physical Chemistry A</i> , <b>2008</b> , 112, 9277-82	2.8	17
189	3s Rydberg and cationic States of pyrazine studied by photoelectron spectroscopy. <i>Journal of Physical Chemistry A</i> , <b>2008</b> , 112, 2293-310	2.8	35
188	Infrared-vacuum ultraviolet pulsed field ionization-photoelectron study of C <sub>2</sub> H <sub>4</sub> ( <sup>+</sup> ) using a high-resolution infrared laser. <i>Journal of Physical Chemistry A</i> , <b>2008</b> , 112, 2572-8	2.8	8
187	A vacuum ultraviolet laser photoionization and pulsed field ionization study of nascent S(3P <sub>2</sub> ,1,0) and S(1D <sub>2</sub> ) formed in the 193.3 nm photodissociation of CS <sub>2</sub> . <i>Journal of Chemical Physics</i> , <b>2008</b> , 128, 014305	3.9	18
186	Rovibrationally selected and resolved pulsed field ionization-photoelectron study of propyne: ionization energy and spin-orbit interaction in propyne cation. <i>Journal of Chemical Physics</i> , <b>2008</b> , 128, 094311	3.9	21
185	Infrared-vacuum ultraviolet-pulsed field ionization-photoelectron study of CH(3)I( <sup>+</sup> ) using a high-resolution infrared laser. <i>Journal of Chemical Physics</i> , <b>2008</b> , 128, 104306	3.9	17
184	Vacuum ultraviolet pulsed field ionization-photoelectron and infrared-photoinduced Rydberg ionization study of trans-1,3-butadiene. <i>Journal of Chemical Physics</i> , <b>2008</b> , 129, 114305	3.9	4
183	High-resolution Rydberg tagging time-of-flight measurements of atomic photofragments by single-photon vacuum ultraviolet laser excitation. <i>Review of Scientific Instruments</i> , <b>2008</b> , 79, 123106	1.7	18
182	Vacuum ultraviolet excitation spectroscopy of the autoionizing Rydberg states of atomic sulfur in the 73 350-84 950 cm <sup>-1</sup> frequency range. <i>Journal of Chemical Physics</i> , <b>2008</b> , 128, 084303	3.9	8
181	Proton Energy Loss Spectroscopy as a State-to-State Probe of Molecular Dynamics. <i>Advances in Chemical Physics</i> , <b>2007</b> , 553-647		27
180	Diabatic Potential Energy Surfaces for Charge-Transfer Processes. <i>Advances in Chemical Physics</i> , <b>2007</b> , 73-134		58
179	Ultraviolet Photodissociation Studies of Organosulfur Molecules and Radicals: Energetics, Structure Identification, and Internal State Distribution. <i>Advances in Photochemistry</i> , <b>2007</b> , 1-116		3
178	State-Selected and State-to-State Ion-Molecular Reaction Dynamics by Photoionization and Differential Reactivity Methods. <i>Advances in Chemical Physics</i> , <b>2007</b> , 401-500		21
177	Rotationally resolved vacuum ultraviolet pulsed field ionization-photoelectron vibrational bands for HD <sup>+</sup> (X 2Sigma <sup>g+</sup> , v <sup>+</sup> =0-20). <i>Journal of Chemical Physics</i> , <b>2007</b> , 126, 164303	3.9	5
176	Combined vacuum ultraviolet laser and synchrotron pulsed field ionization study of CH <sub>2</sub> BrCl. <i>Journal of Chemical Physics</i> , <b>2007</b> , 126, 184304	3.9	9
175	Vacuum ultraviolet laser pulsed field ionization-photoelectron study of allyl radical CH <sub>2</sub> CHCH <sub>2</sub> . <i>Journal of Chemical Physics</i> , <b>2007</b> , 126, 171101	3.9	23
174	An experimental and quasiclassical trajectory study of the rovibrationally state-selected reactions: HD <sup>+</sup> (v=0-15, j=1)+He <sup>-&gt;</sup> HeH <sup>+</sup> (HeD <sup>+</sup> )+D(H). <i>Journal of Chemical Physics</i> , <b>2007</b> , 126, 234305	3.9	10

173	Assignment of rovibrational transitions of propyne in the region of 2934-2952 cm <sup>-1</sup> measured by two-color IR-vacuum ultraviolet laser photoion-photoelectron methods. <i>Journal of Chemical Physics</i> , <b>2007</b> , 127, 044313	3.9	10
172	A time-dependent wave packet quantum scattering study of the reaction HD <sup>+</sup> (v = 0 - 3; j <sub>0</sub> = 1) + He → HeH <sup>+</sup> (HeD <sup>+</sup> ) + D(H). <i>Journal of Chemical Physics</i> , <b>2007</b> , 127, 164318	3.9	17
171	Rotationally Resolved Vacuum Ultraviolet Pulsed Field Ionization-Photoelectron Vibrational Bands for H <sub>2</sub> <sup>+</sup> (X <sub>2</sub> g <sup>+</sup> , v <sub>+</sub> =0-18). <i>Chinese Journal of Chemical Physics</i> , <b>2007</b> , 20, 352-364	0.9	8
170	State Selected Charge Transfer and Chemical Reactions by the Tesico Technique. <i>Advances in Chemical Physics</i> , <b>2007</b> , 263-307		16
169	Multicoincidence Detection in Beam Studies of Ion-Molecule Reactions: Technique and Application to X <sup>-</sup> + h <sub>2</sub> Reactions. <i>Advances in Chemical Physics</i> , <b>2007</b> , 309-399		6
168	Crossed-Molecular Beam Studies of State-to-State Reaction Dynamics. <i>Advances in Chemical Physics</i> , <b>2007</b> , 501-552		3
167	Control of Transition-Metal Cation Reactivity by Electronic State Selection. <i>Advances in Chemical Physics</i> , <b>2007</b> , 213-262		18
166	Multiphoton Ionization State Selection: Vibrational-Mode and Rotational-State Control. <i>Advances in Chemical Physics</i> , <b>2007</b> , 177-212		43
165	Molecular Beam Photoionization Studies of Molecules and Clusters. <i>Advances in Chemical Physics</i> , <b>2007</b> , 263-362		76
164	Inhomogeneous RF Fields: A Versatile Tool for the Study of Processes with Slow Ions. <i>Advances in Chemical Physics</i> , <b>2007</b> , 1-176		314
163	Accurate ab initio predictions of ionization energies and heats of formation for the 2-propyl, phenyl, and benzyl radicals. <i>Journal of Chemical Physics</i> , <b>2006</b> , 124, 044323	3.9	25
162	A photodissociation study of CH <sub>2</sub> BrCl in the A-band using the time-sliced ion velocity imaging method. <i>Journal of Chemical Physics</i> , <b>2006</b> , 124, 034309	3.9	45
161	Infrared vibrational spectroscopy of cis-dichloroethene in Rydberg states. <i>Journal of Chemical Physics</i> , <b>2006</b> , 124, 64310	3.9	6
160	Direct identification of propargyl radical in combustion flames by vacuum ultraviolet photoionization mass spectrometry. <i>Journal of Chemical Physics</i> , <b>2006</b> , 124, 74302	3.9	29
159	The study of state-selected ion-molecule reactions using the vacuum ultraviolet pulsed field ionization-photoion technique. <i>Journal of Chemical Physics</i> , <b>2006</b> , 125, 132306	3.9	37
158	Rovibrationally selected and resolved state-to-state photoionization of ethylene using the infrared-vacuum ultraviolet pulsed field ionization-photoelectron method. <i>Journal of Chemical Physics</i> , <b>2006</b> , 125, 133304	3.9	19
157	Vacuum ultraviolet laser pulsed field ionization-photoelectron study of cis-dichloroethene. <i>Journal of Chemical Physics</i> , <b>2006</b> , 124, 224311	3.9	11
156	Accurate ab initio Predictions of Ionization Energies and Heats of Formation for Cyclopropenylidene, Propargylene and Propadienylidene. <i>Chinese Journal of Chemical Physics</i> , <b>2006</b> , 19, 29-38	0.9	28

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49	Experimental and theoretical total state-selected and state-to-state absolute cross sections. I. The H <sub>2</sub> (X,v)+Ar reaction. <i>Journal of Chemical Physics</i> , <b>1990</b> , 93, 4818-4831	3.9	37
48	A photoion-photoelectron coincidence study of (CO) <sub>2</sub> and (CO) <sub>3</sub> . <i>Journal of Chemical Physics</i> , <b>1989</b> , 90, 6026-6033	3.9	15

47	A photoion-photoelectron coincidence study of Kr and Xe dimers. <i>Journal of Chemical Physics</i> , <b>1989</b> , 90, 4689-4696	3.9	21
46	A state-selected study of the unimolecular decomposition of C <sub>2</sub> H <sub>2</sub> (A, B) using the photoion-photoelectron coincidence method. <i>Journal of Chemical Physics</i> , <b>1989</b> , 91, 2898-2900	3.9	34
45	A photoion-photoelectron coincidence study of (N <sub>2</sub> ) <sub>2</sub> and (N <sub>2</sub> ) <sub>3</sub> . <i>Journal of Chemical Physics</i> , <b>1989</b> , 91, 849-856	3.9	30
44	A photoion-photoelectron coincidence study of Ar <sub>n</sub> (n=2-4). <i>Journal of Chemical Physics</i> , <b>1989</b> , 90, 2995-3003	3.9	73
43	A 193 nm laser photofragmentation time-of-flight mass spectrometric study of CS <sub>2</sub> and CS <sub>2</sub> clusters. <i>Journal of Chemical Physics</i> , <b>1988</b> , 88, 1658-1669	3.9	71
42	A photoion-photoelectron coincidence study of (CO) <sub>2</sub> . <i>Journal of Chemical Physics</i> , <b>1988</b> , 88, 4098-4099	3.9	13
41	Observation of the formation of C <sup>+</sup> , O <sup>+</sup> , and ArC <sup>+</sup> in the collisions of Ar <sup>+</sup> (2P <sub>3/2</sub> , 1/2) with CO. <i>Journal of Chemical Physics</i> , <b>1988</b> , 89, 3381-3382	3.9	19
40	Absolute state-to-state total cross sections for the reactions N <sub>2</sub> (X, v=0, 1) + Ar(1S <sub>0</sub> ) → N <sub>2</sub> (X, v) + Ar(2P <sub>3/2</sub> , 1/2). <i>Journal of Chemical Physics</i> , <b>1987</b> , 86, 170-175	3.9	42
39	A state-to-state study of the electron transfer reactions Ar <sup>+</sup> (2P <sub>3/2</sub> , 1/2) + N <sub>2</sub> (X, v=0) → Ar(1S <sub>0</sub> ) + N <sub>2</sub> (X, v). <i>Journal of Chemical Physics</i> , <b>1986</b> , 85, 3874-3890	3.9	53
38	A state-to-state study of the electron transfer reactions N <sub>2</sub> (X, v=0, 1) + Ar(1S <sub>0</sub> ) → N <sub>2</sub> (X, v) + Ar(2P <sub>3/2</sub> , 1/2). <i>Journal of Chemical Physics</i> , <b>1986</b> , 85, 7136-7145	3.9	18
37	A vibrational state-selected study of the reaction H <sub>2</sub> (v=1) + H <sub>2</sub> (v=0) → H <sub>3</sub> + H using the tandem photoionization mass spectrometry and radio frequency ion guide methods. <i>Journal of Chemical Physics</i> , <b>1986</b> , 84, 4317-4326	3.9	45
36	Photoionization studies of HgKr and HgXe. <i>Journal of Chemical Physics</i> , <b>1986</b> , 84, 1142-1145	3.9	12
35	Vibrational state distributions of H <sub>2</sub> (v) resulting from the electron transfer reactions H <sub>2</sub> (v=0, 1) + H <sub>2</sub> (v=0) → H <sub>2</sub> (v) + H <sub>2</sub> (v) in the collisional energy range of 2-6 eV. <i>Journal of Chemical Physics</i> , <b>1986</b> , 84, 197-200	3.9	15
34	Molecular beam photoionization study of S <sub>2</sub> . <i>Journal of Chemical Physics</i> , <b>1986</b> , 84, 778-782	3.9	16
33	Photoionization study of HgAr. <i>Journal of Chemical Physics</i> , <b>1985</b> , 82, 648-652	3.9	20
32	A state-to-state study of the symmetric charge transfer reaction Ar <sup>+</sup> (2P <sub>3/2</sub> , 1/2) + Ar(1S <sub>0</sub> ). <i>Journal of Chemical Physics</i> , <b>1985</b> , 82, 5489-5498	3.9	32
31	A study of the unimolecular decomposition of the (C <sub>2</sub> H <sub>4</sub> ) <sub>3</sub> complex. <i>Journal of Chemical Physics</i> , <b>1985</b> , 83, 2813-2817	3.9	13
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28	A study of the unimolecular decomposition of the $(C_2H_4)_2^+$ complex. <i>Journal of Chemical Physics</i> , <b>1984</b> , 80, 1482-1489	3.9	25
27	Molecular beam photoionization study of $HgCl_2$ . <i>Journal of Chemical Physics</i> , <b>1983</b> , 78, 37-45	3.9	24
26	Higher resolution photoionization study of acetylene near the threshold. <i>Journal of Chemical Physics</i> , <b>1982</b> , 76, 3905-3907	3.9	32
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22	A study of the ion-molecule half reactions $O_2(a^4\Sigma, v) + (O_2)_m \rightarrow O_2 + 2m + 1 + O$ , $m = 1, 2$ , or $3$ , using the molecular beam photoionization method. <i>Journal of Chemical Physics</i> , <b>1981</b> , 74, 3348-3352	3.9	32
21	Photoionization study of $CO_2$ , $N_2O$ dimers and clusters. <i>Journal of Chemical Physics</i> , <b>1981</b> , 75, 4921-4926	3.9	55
20	Molecular beam photoionization study of $SO_2$ and $(SO_2)_2$ . <i>Journal of Chemical Physics</i> , <b>1981</b> , 75, 1650-1657	3.9	56
19	Molecular beam photoionization study of $OCS$ , $(OCS)_2$ , $(OCS)_3$ , and $OCS \cdot CS_2$ . <i>Journal of Chemical Physics</i> , <b>1981</b> , 74, 1645-1651	3.9	62
18	Molecular beam photoionization study of $CO$ , $N_2$ , and $NO$ dimers and clusters. <i>Journal of Chemical Physics</i> , <b>1981</b> , 74, 3342-3347	3.9	107
17	A study of the high Rydberg state and ion-molecule reactions of carbon disulfide using the molecular beam photoionization method. <i>Journal of Chemical Physics</i> , <b>1981</b> , 74, 1125-1132	3.9	25
16	Molecular beam photoionization study of carbon disulfide, carbon disulfide dimer and clusters. <i>Journal of Chemical Physics</i> , <b>1980</b> , 73, 2523-2533	3.9	84
15	A study of the chemiionization process $CS_2^*(n) + CS_2 \rightarrow CS_3^+ + CS + e^-$ using the molecular beam photoionization method. <i>Journal of Chemical Physics</i> , <b>1980</b> , 72, 4242-4244	3.9	17
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