

Mark A Johnson

List of Publications by Year in descending order

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179
papers

10,423
citations

28242

55
h-index

38368

95
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180
all docs

180
docs citations

180
times ranked

4792
citing authors

#	ARTICLE	IF	CITATIONS
1	Spectral Signatures of Hydrated Proton Vibrations in Water Clusters. <i>Science</i> , 2005, 308, 1765-1769.	6.0	712
2	MOLECULAR ASPECTS OF HALIDE ION HYDRATION: The Cluster Approach. <i>Annual Review of Physical Chemistry</i> , 2003, 54, 173-213.	4.8	449
3	Spectroscopic Determination of the OH- Solvation Shell in the OH·(H ₂ O) _n Clusters. <i>Science</i> , 2003, 299, 1367-1372.	6.0	355
4	How Do Small Water Clusters Bind an Excess Electron?. <i>Science</i> , 2004, 306, 675-679.	6.0	276
5	Cryogenic Ion Chemistry and Spectroscopy. <i>Accounts of Chemical Research</i> , 2014, 47, 202-210.	7.6	256
6	The vibrational predissociation spectra of the H ₅ O ₂ ⁺ ·M ⁿ (RG=Ar,Ne) clusters: Correlation of the solvent perturbations in the free OH and shared proton transitions of the Zundel ion. <i>Journal of Chemical Physics</i> , 2005, 122, 244301.	1.2	228
7	Vibrational Spectroscopy of the Ionic Hydrogen Bond: Fermi Resonances and Ion·Molecule Stretching Frequencies in the Binary X·H ₂ O (X = Cl, Br, I) Complexes via Argon Predissociation Spectroscopy. <i>Journal of the American Chemical Society</i> , 1998, 120, 12361-12362.	6.6	226
8	Spectroscopic snapshots of the proton-transfer mechanism in water. <i>Science</i> , 2016, 354, 1131-1135.	6.0	213
9	Electronic absorption spectra of size-selected hydrated electron clusters: (H ₂ O) _n ⁻ , n=6-50. <i>Journal of Chemical Physics</i> , 1997, 106, 811-814.	1.2	188
10	Fundamental Excitations of the Shared Proton in the H ₃ O ₂ ⁻ and H ₅ O ₂ ⁺ Complexes. <i>Journal of Physical Chemistry A</i> , 2005, 109, 1487-1490.	1.1	168
11	Precursor of the ν_1 charge-transfer-to-solvent (CTTS) band in $\text{H}^+\cdots(\text{H}_2\text{O})_n$ clusters. <i>Journal of Chemical Physics</i> , 1996, 105, 7231-7234.	1.2	162
12	Vibrational Spectroscopy of Small Br·(H ₂ O) _n and I·(H ₂ O) _n Clusters: Infrared Characterization of the Ionic Hydrogen Bond. <i>Journal of Physical Chemistry A</i> , 1998, 102, 3067-3071.	1.1	158
13	Isolating the Spectroscopic Signature of a Hydration Shell With the Use of Clusters: Superoxide Tetrahydrate. <i>Science</i> , 2000, 287, 2461-2463.	6.0	157
14	Prying Apart a Water Molecule with Anionic H-Bonding: A Comparative Spectroscopic Study of the X·H ₂ O (X = OH, O, F, Cl, and Br) Binary Complexes in the 600-3800 cm ⁻¹ Region. <i>Journal of Physical Chemistry A</i> , 2006, 110, 4943-4952.	1.1	157
15	Demonstration of a pulsed photoelectron spectrometer on mass-selected negative ions: O ⁻ , O ₂ ⁻ , and O ₄ ⁻ . <i>Chemical Physics Letters</i> , 1986, 131, 170-174.	1.2	152
16	Photoelectron spectroscopy of the 'missing' hydrated electron clusters (H ₂ O) _n ⁻ , n=3, 5, 8 and 9: Isomers and continuity with the dominant clusters n=6, 7 and 411. <i>Chemical Physics Letters</i> , 1998, 297, 90-96.	1.2	142
17	Vibrational Characterization of Simple Peptides Using Cryogenic Infrared Photodissociation of H ₂ -Tagged, Mass-Selected Ions. <i>Journal of the American Chemical Society</i> , 2011, 133, 6440-6448.	6.6	139
18	Vibrational predissociation spectroscopy of the H ₂ -tagged mono- and dicarboxylate anions of dodecanedioic acid. <i>International Journal of Mass Spectrometry</i> , 2011, 300, 91-98.	0.7	136

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19	Mass-selected $\hat{\text{A}}$ -matrix isolation infrared spectroscopy of the $\hat{\text{A}}\cdot(\text{H}_2\text{O})_2$ complex: making and breaking the inter-water hydrogen-bond. <i>Chemical Physics</i> , 1998, 239, 485-491.	0.9	131
20	Spectroscopic Observation of Ion-Induced Water Dimer Dissociation in the $\text{X}\cdot\hat{\text{A}}\cdot(\text{H}_2\text{O})_2$ ($\text{X} = \text{F}, \text{Cl}, \text{Br}, \text{I}$) Clusters. <i>Journal of Physical Chemistry A</i> , 1999, 103, 10665-10669.	1.1	128
21	Determination of Noncovalent Docking by Infrared Spectroscopy of Cold Gas-Phase Complexes. <i>Science</i> , 2012, 335, 694-698.	6.0	127
22	Predissociation spectroscopy of the argon-solvated H_5O_2^+ $\hat{\text{A}}$ -cation in the 1000-1900 cm^{-1} region. <i>Journal of Chemical Physics</i> , 2004, 121, 11523-11526.	1.2	126
23	Photoelectron spectroscopy of $(\text{CO}_2)_n$ clusters with $n=13$: Cluster size dependence of the core molecular ion. <i>Journal of Chemical Physics</i> , 1988, 88, 5857-5863.	1.2	124
24	An infrared study of the competition between hydrogen-bond networking and ionic solvation: Halide-dependent distortions of the water trimer in the $\text{X}\hat{\text{A}}\cdot\text{H}_2\text{O}_3$, ($\text{X}=\text{Cl}, \text{Br}, \text{I}$) systems. <i>Journal of Chemical Physics</i> , 1999, 110, 7129-7132.	1.2	116
25	Vibrational spectral signature of the proton defect in the three-dimensional H^+ H_2O network. <i>Journal of Chemical Physics</i> , 2011, 134, 074701.	6.0	111
26	Snapshots of Proton Accommodation at a Microscopic Water Surface: Understanding the Vibrational Spectral Signatures of the Charge Defect in Cryogenically Cooled H^+ H_2O Clusters. <i>Journal of Physical Chemistry A</i> , 2015, 119, 9425-9440.	1.1	111
27	How the Shape of an H-Bonded Network Controls Proton-Coupled Water Activation in HONO Formation. <i>Science</i> , 2010, 327, 308-312.	6.0	99
28	Photochemistry of hydrated electron clusters $(\text{H}_2\text{O})_n$ (15-40) at 1064 nm: Size dependent competition between photofragmentation and photodetachment. <i>Journal of Chemical Physics</i> , 1988, 89, 4807-4814.	1.2	97
29	Photoelectron spectroscopy of $(\text{CO}_2)_n$ revisited: core switching in the $2n-16$ range. <i>Chemical Physics Letters</i> , 1997, 268, 429-433.	1.2	96
30	Hydration of a structured excess charge distribution: Infrared spectroscopy of the $\text{O}_2^+ \cdot (\text{H}_2\text{O})_n$ clusters. <i>Journal of Chemical Physics</i> , 2001, 114, 2698-2706.	1.2	93
31	An H/D Isotopic Substitution Study of the H_5O_2^+ $\hat{\text{A}}$ -Ar Vibrational Predissociation Spectra: Exploring the Putative Role of Fermi Resonances in the Bridging Proton Fundamentals. <i>Journal of Physical Chemistry B</i> , 2008, 112, 321-327.	1.2	92
32	Spectroscopic Study of the Ion-Radical H-Bond in H_4O_2^+ . <i>Journal of Physical Chemistry A</i> , 2009, 113, 4772-4779.	1.1	91
33	Vibrational predissociation spectroscopy of the $(\text{H}_2\text{O})_6 \hat{\text{A}} \cdot \text{Ar}_n$, $n=3-6$, clusters. <i>Journal of Chemical Physics</i> , 1998, 108, 444-449.	1.2	88
34	Isomer-Specific IR-IR Double Resonance Spectroscopy of D_2 -Tagged Protonated Dipeptides Prepared in a Cryogenic Ion Trap. <i>Journal of Physical Chemistry Letters</i> , 2012, 3, 1099-1105.	2.1	88
35	Vibrational manifestations of strong non-Condon effects in the $\text{H}_3\text{O}^+ \cdot \text{X}_3$ ($\text{X} = \text{Ar}, \text{N}_2, \text{CH}_4, \text{H}_2\text{O}$) complexes: A possible explanation for the intensity in the association band in the vibrational spectrum of water. <i>Physical Chemistry Chemical Physics</i> , 2012, 14, 7205.	1.3	82
36	Vibrational predissociation spectra of $\hat{\text{A}}\cdot(\text{H}_2\text{O})$: isotopic labels and weakly bound complexes with Ar and N_2 . <i>Chemical Physics Letters</i> , 1997, 269, 122-127.	1.2	81

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37	Gas-Phase Infrared Spectroscopy and Multidimensional Quantum Calculations of the Protonated Ammonia Dimer $N_2H_7^+$. <i>Angewandte Chemie - International Edition</i> , 2007, 46, 8691-8694.	7.2	77
38	Unraveling Anharmonic Effects in the Vibrational Predissociation Spectra of $H_5O_2^+$ and Its Deuterated Analogues. <i>Journal of Physical Chemistry A</i> , 2011, 115, 5847-5858.	1.1	75
39	Photochemistry of Halide Ion-Molecule Clusters: Dipole-Bound Excited States and the Case for Asymmetric Solvation. <i>Accounts of Chemical Research</i> , 1998, 31, 527-534.	7.6	74
40	Infrared Signatures of a Water Molecule Attached to Triatomic Domains of Molecular Anions: Evolution of the H-bonding Configuration with Domain Length. <i>Journal of Physical Chemistry A</i> , 2003, 107, 6527-6532.	1.1	74
41	Dipole-bound excited states of the $l^{\sim}CH_3CN$ and $l^{\sim}(CH_3CN)_2$ ion-molecule complexes: Evidence for asymmetric solvation. <i>Journal of Chemical Physics</i> , 1995, 103, 2006-2015.	1.2	71
42	The Vibrational Spectrum of the Neutral $(H_2O)_6$ Precursor to the "Magic" $(H_2O)_6^-$ Cluster Anion by Argon-Mediated, Population-Modulated Electron Attachment Spectroscopy. <i>Journal of Physical Chemistry A</i> , 2004, 108, 64-68.	1.1	71
43	Capture of CO_2 by a Cationic Nickel(I) Complex in the Gas Phase and Characterization of the Bound, Activated CO_2 Molecule by Cryogenic Ion Vibrational Predissociation Spectroscopy. <i>Angewandte Chemie - International Edition</i> , 2016, 55, 1282-1285.	7.2	68
44	Thermal Energy Reactions of Size-Selected Hydrated Electron Clusters $(H_2O)_n^-$. <i>The Journal of Physical Chemistry</i> , 1996, 100, 2900-2906.	2.9	67
45	Communication: He-tagged vibrational spectra of the SarGlyH ⁺ and H ⁺ (H ₂ O) _{2,3} ions: Quantifying tag effects in cryogenic ion vibrational predissociation (CIVP) spectroscopy. <i>Journal of Chemical Physics</i> , 2014, 140, 221101.	1.2	67
46	Isolating the spectra of cluster ion isomers using Ar-tag-mediated IR-IR double resonance within the vibrational manifolds: Application to $NO_2^{\sim}H_2O$. <i>Journal of Chemical Physics</i> , 2008, 129, 094303.	1.2	65
47	Pulsed photoelectron spectroscopy of negative cluster ions: Isolation of three distinguishable forms of $N_2O^{\sim}2$. <i>Journal of Chemical Physics</i> , 1988, 88, 5383-5395.	1.2	64
48	Reactions of hydrated electron clusters $(H_2O)_n^-$: scavenging the excess electron. <i>The Journal of Physical Chemistry</i> , 1989, 93, 1178-1181.	2.9	63
49	Vibrational Predissociation Spectrum of the Carbamate Radical Anion, $C_5H_5N-CO_2^{\sim}$, Generated by Reaction of Pyridine with $(CO_2)^{\sim}$. <i>Journal of the American Chemical Society</i> , 2010, 132, 15508-15511.	6.6	62
50	A Cluster Study of Anionic Hydration: Spectroscopic Characterization of the $l^{\sim}W_n$, $1 \leq n \leq 3$, Supramolecular Complexes at the Primary Steps of Solvation. <i>Journal of Physical Chemistry A</i> , 1999, 103, 443-447.	1.1	61
51	Anharmonicities and Isotopic Effects in the Vibrational Spectra of $X^{\sim}H_2O$, $^{\sim}HDO$, and $^{\sim}D_2O$ [X = Cl, Br, and I] Binary Complexes. <i>Journal of Physical Chemistry A</i> , 2010, 114, 1556-1568.	1.1	61
52	Site-specific vibrational spectral signatures of water molecules in the magic H_3O^+ (H_2O) ₂₀ and Cs ⁺ (H_2O) ₂₀ clusters. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2014, 111, 18132-18137.	3.3	59
53	Large anharmonic effects in the infrared spectra of the symmetrical $CH_3NO_2^{\sim}(H_2O)$ and $CH_3CO_2^{\sim}(H_2O)$ complexes. <i>Journal of Chemical Physics</i> , 2003, 119, 10138-10145.	1.2	57
54	Modes of Activation of Organometallic Iridium Complexes for Catalytic Water and C-H Oxidation. <i>Inorganic Chemistry</i> , 2014, 53, 423-433.	1.9	57

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55	Photoactivation of the Cl ⁻ + CH ₃ Br S _N 2 Reaction via Rotationally Resolved C- ¹³ H Stretch Excitation of the Cl- ¹³ CH ₃ Br Entrance Channel Complex. <i>Journal of the American Chemical Society</i> , 1999, 121, 6950-6951.	6.6	56
56	Argon Predissociation Spectroscopy of the OH- ¹⁸ O and Cl- ¹⁸ O Complexes in the 1000-1900 cm ⁻¹ Region: Intramolecular Bending Transitions and the Search for the Shared-Proton Fundamental in the Hydroxide Monohydrate. <i>Journal of Physical Chemistry A</i> , 2005, 109, 571-575.	1.1	56
57	Gas phase vibrational spectroscopy of the protonated water pentamer: the role of isomers and nuclear quantum effects. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 26743-26754.	1.3	53
58	Deconstructing water's diffuse OH stretching vibrational spectrum with cold clusters. <i>Science</i> , 2019, 364, 275-278.	6.0	53
59	Photoabsorption of negative cluster ions near the electron detachment threshold: A study of the (O ₂) ⁻ⁿ system. <i>Journal of Chemical Physics</i> , 1990, 93, 268-275.	1.2	52
60	Microhydration of Contact Ion Pairs in M ²⁺ OH ⁻ (H ₂ O) _n (M = Mg, Ca) Clusters: Spectral Manifestations of a Mobile Proton Defect in the First Hydration Shell. <i>Journal of Physical Chemistry A</i> , 2014, 118, 7590-7597.	1.1	52
61	Collisional activation of captured intermediates in the gas-phase S _N 2 reaction chloride + bromomethane. <i>Journal of the American Chemical Society</i> , 1991, 113, 9697-9699.	6.6	51
62	Spectroscopic Evidence for an Attractive Cation-Cation Interaction in Hydroxy-Functionalized Ionic Liquids: A Hydrogen-Bonded Chain-Like Trimer. <i>Angewandte Chemie - International Edition</i> , 2018, 57, 15364-15368.	7.2	51
63	Unraveling the Anomalous Solvatochromic Response of the Formate Ion Vibrational Spectrum: An Infrared, Ar-Tagging Study of the HCO ₂ ⁻ , DCO ₂ ⁻ , and HCO ₂ ⁻ -H ₂ O Ions. <i>Journal of Physical Chemistry Letters</i> , 2011, 2, 2437-2441.	2.1	49
64	Understanding the ionic liquid [NC ₄][NTf ₂] from individual building blocks: an IR-spectroscopic study. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 8518-8529.	1.3	48
65	Solvation of the Cl- ¹⁸ O Complex in CCl ₄ Clusters: The Effect of Solvent-Mediated Charge Redistribution on the Ionic H-Bond. <i>Journal of Physical Chemistry A</i> , 2002, 106, 1205-1209.	1.1	47
66	Hidden role of intermolecular proton transfer in the anomalously diffuse vibrational spectrum of a trapped hydronium ion. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2017, 114, E4706-E4713.	3.3	47
67	Structural Motifs in Cold Ternary Ion Complexes of Hydroxyl-Functionalized Ionic Liquids: Isolating the Role of Cation-Cation Interactions. <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 2979-2984.	2.1	47
68	Controlling the internal energy content of size-selected cluster ions: An experimental comparison of the metastable decay rate and photofragmentation methods of quantifying the internal excitation of (H ₂ O) ⁿ . <i>Journal of Chemical Physics</i> , 1991, 95, 7998-8004.	1.2	45
69	A Cluster Study of Cl ₂ -Microhydration: Size-Dependent Competition between Symmetrical H-Bonding to the Anion and the Formation of Cyclic Water Networks in the Cl ₂ -1 ⁵ (H ₂ O) Series. <i>Journal of Physical Chemistry A</i> , 2004, 108, 3910-3915.	1.1	45
70	Size-dependent collisional incorporation of D ₂ O into (H ₂ O) ⁿ around n = 15: implications on the origin of magic numbers in the hydrated electron cluster distribution. <i>Chemical Physics Letters</i> , 1991, 181, 206-212.	1.2	44
71	Disentangling the Complex Vibrational Spectrum of the Protonated Water Trimer, H ⁺ (H ₂ O) ₃ , with Two-Color IR-IR Photodissociation of the Bare Ion and Anharmonic VSCF/VCI Theory. <i>Journal of Physical Chemistry Letters</i> , 2017, 8, 3782-3789.	2.1	44
72	Spectroscopic observation of vibrational Feshbach resonances in near-threshold photoexcitation of X- ¹³ CH ₃ NO ₂ (X=I and Br). <i>Faraday Discussions</i> , 2000, 115, 395-406.	1.6	43

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73	Hiding in Plain Sight: Unmasking the Diffuse Spectral Signatures of the Protonated N-Terminus in Isolated Dipeptides Cooled in a Cryogenic Ion Trap. <i>Journal of Physical Chemistry Letters</i> , 2013, 4, 3450-3457.	2.1	43
74	Photodestruction spectra of the anionic water clusters, $(\text{H}_2\text{O})_n^-$, $n=18$ and 30: Absorption to the red of aq^- . <i>Journal of Chemical Physics</i> , 1991, 94, 5240-5242.	1.2	42
75	Observation of the dipole-bound excited state of the $\text{I}^- \cdots \text{acetone}$ ion-molecule complex. <i>Journal of Chemical Physics</i> , 1995, 102, 6335-6338.	1.2	42
76	Vibrationally Induced Proton Transfer in $\text{F}^-(\text{H}_2\text{O})$ and $\text{F}^-(\text{D}_2\text{O})$. <i>Journal of Physical Chemistry A</i> , 2008, 112, 12337-12344.	1.1	42
77	Alkali-Controlled C-H Cleavage or N-C Bond Formation by N-Derived Iron Nitrides and Imides. <i>Journal of the American Chemical Society</i> , 2016, 138, 11185-11191.	6.6	42
78	Optical-optical double resonance on cooled molecular ions: rotational assignments in the perturbed CO_2^- system. <i>Chemical Physics Letters</i> , 1982, 92, 225-231.	1.2	41
79	Infrared predissociation spectroscopy of $\text{I}^-(\text{CH}_3\text{OH})_n$, $n=1,2$: Cooperativity in asymmetric solvation. <i>Journal of Chemical Physics</i> , 2002, 116, 4853.	1.2	41
80	Dominant structural motifs of $\text{NO}^-(\text{H}_2\text{O})_n$ complexes: Infrared spectroscopic and ab initio studies. <i>Journal of Chemical Physics</i> , 2003, 118, 4945-4953.	1.2	41
81	overflow="scroll" xmlns:xocs="http://www.elsevier.com/xml/xocs/dtd" xmlns:xs="http://www.w3.org/2001/XMLSchema" xmlns:xsi="http://www.w3.org/2001/XMLSchema-instance" xmlns="http://www.elsevier.com/xml/ja/dtd" xmlns:ja="http://www.elsevier.com/xml/ja/dtd" xmlns:mml="http://www.w3.org/1998/Math/MathML" xmlns:tb="http://www.elsevier.com/xml/common/table/dtd" xmlns:sb="http://www.elsevier.com/xml/co	1.2	41
82	Caught in the Act of Dissolution. <i>Science</i> , 2002, 298, 69-69.	6.0	40
83	Generating Spectra from Ground-State Wave Functions: Unraveling Anharmonic Effects in the $\text{OH}^-(\text{H}_2\text{O})$ Vibrational Predissociation Spectrum. <i>Journal of Physical Chemistry A</i> , 2009, 113, 7346-7352.	1.1	39
84	Ionic liquids from the bottom up: Local assembly motifs in [EMIM][BF ₄] through cryogenic ion spectroscopy. <i>Journal of Chemical Physics</i> , 2013, 139, 224305.	1.2	39
85	Persistence of Dual Free Internal Rotation in $\text{NH}_4^+(\text{H}_2\text{O}) \cdots \text{He}^-(\text{H}_2\text{O})_3$ Ion-Molecule Complexes: Expanding the Case for Quantum Delocalization in He Tagging. <i>Journal of Physical Chemistry A</i> , 2015, 119, 4170-4176.	1.1	38
86	Isolation of site-specific anharmonicities of individual water molecules in the $\text{I}^-(\text{H}_2\text{O})_2$ complex using tag-free, isotopomer selective IR-IR double resonance. <i>Chemical Physics Letters</i> , 2017, 690, 159-171.	1.2	38
87	Infrared spectroscopic observation of the argon isomer distribution in evaporative ensembles of $\text{I}^-(\text{ROH})_n$ (R=methyl, ethyl, isopropyl) clusters. <i>Journal of Chemical Physics</i> , 1999, 111, 9593-9599.	1.2	37
88	Characterizing the Intramolecular H-bond and Secondary Structure in Methylated GlyGlyH ⁺ with H ₂ O Predissociation Spectroscopy. <i>Journal of the American Society for Mass Spectrometry</i> , 2011, 22, 1941-52.	1.2	37
89	Thermodynamics of Water Dimer Dissociation in the Primary Hydration Shell of the Iodide Ion with Temperature-Dependent Vibrational Predissociation Spectroscopy. <i>Journal of Physical Chemistry A</i> , 2015, 119, 1859-1866.	1.1	37
90	Molecular-level origin of the carboxylate head group response to divalent metal ion complexation at the air-water interface. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2019, 116, 14874-14880.	3.3	37

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91	Autodetachment from vibrational levels of the O_2^+ resonance across its dissociation limit by photoexcitation from O_2^+ . Journal of Chemical Physics, 1996, 105, 1807-1814.	1.2	36
92	Comparison of the local binding motifs in the imidazolium-based ionic liquids [EMIM][BF ₄] and [EMMIM][BF ₄] through cryogenic ion vibrational predissociation spectroscopy: Unraveling the roles of anharmonicity and intermolecular interactions. Journal of Chemical Physics, 2015, 142, 064306.	1.2	35
93	The angular distribution of photoelectrons ejected from the hydrated electron cluster (H ₂ O) ⁺ ₁₈ . Journal of Chemical Physics, 1990, 92, 3243-3245.	1.2	33
94	Why Does Argon Bind to Deuterium? Isotope Effects and Structures of ArH ₅ O ₂ ⁺ Complexes. Journal of Physical Chemistry A, 2008, 112, 6074-6078.	1.1	33
95	On the vibrational fine structure in the near-threshold photofragmentation spectrum of the $I^+ \cdots CH_3I$ complex: Spectroscopic observation of nonadiabatic effects in electron-molecule scattering. Journal of Chemical Physics, 1996, 105, 10416-10423.	1.2	32
96	Capturing intrinsic site-dependent spectral signatures and lifetimes of isolated OH oscillators in extended water networks. Nature Chemistry, 2020, 12, 159-164.	6.6	32
97	Isolating the Spectral Signatures of Individual Sites in Water Networks Using Vibrational Double-Resonance Spectroscopy of Cluster Isotopomers. Journal of Physical Chemistry Letters, 2010, 1, 2396-2401.	2.1	31
98	NH ⁺ ⋯F Hydrogen Bonding in a Fluorinated α -Proton Sponge Derivative: Integration of Solution, Solid-State, Gas-Phase, and Computational Studies. Journal of Organic Chemistry, 2011, 76, 7975-7984.	1.7	31
99	Photoelectron spectroscopy of the gas-phase S _N 2 reaction intermediates $I^+ \cdots CH_3I$ and $I^+ \cdots CD_3I$: Distortion of the CH ₃ I at the σ -ion dipole complex. Journal of Chemical Physics, 1993, 99, 4869-4872.	1.2	30
100	Vibrational Spectra and Fragmentation Pathways of Size-Selected, D ₂ -Tagged Ammonium/Methylammonium Bisulfate Clusters. Journal of Physical Chemistry A, 2013, 117, 13265-13274.	1.1	30
101	The charge transfer excited state of the $I^+ \cdots CH_3I$ S _N 2 reaction intermediate: Photoinduced intracluster dissociative attachment. Journal of Chemical Physics, 1994, 101, 10507-10520.	1.2	29
102	Diffuse Vibrational Signature of a Single Proton Embedded in the Oxalate Scaffold, HO ₂ CCO ₂ ⁺ . Journal of Physical Chemistry A, 2015, 119, 13018-13024.	1.1	29
103	Cooperatively enhanced hydrogen bonds in ionic liquids: closing the loop with molecular mimics of hydroxy-functionalized cations. Physical Chemistry Chemical Physics, 2019, 21, 18092-18098.	1.3	29
104	Integration of High-Resolution Mass Spectrometry with Cryogenic Ion Vibrational Spectroscopy. Journal of the American Society for Mass Spectrometry, 2019, 30, 1551-1557.	1.2	28
105	Vibronic structure of the CO ₂ ⁺ ion: reinvestigation of the antisymmetric stretch vibration in the [X̃], \tilde{A} , and [B̃] states. Molecular Physics, 1995, 85, 839-868.	0.8	27
106	Origin of the diffuse vibrational signature of a cyclic intramolecular proton bond: Anharmonic analysis of protonated 1,8-disubstituted naphthalene ions. Journal of Chemical Physics, 2013, 139, 024301.	1.2	27
107	Tag-Free and Isotopomer-Selective Vibrational Spectroscopy of the Cryogenically Cooled H ₉ O ₄ ⁺ Cation with Two-Color, IR-IR Double-Resonance Photoexcitation: Isolating the Spectral Signature of a Single OH Group in the Hydronium Ion Core. Journal of Physical Chemistry A, 2018, 122, 9275-9284.	1.1	27
108	Linking the photoelectron and infrared spectroscopies of the (H ₂ O) ₆ ⁺ isomers. Journal of Chemical Physics, 2002, 116, 1201-1203.	1.2	26

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109	Characterization of Highly Unusual NH ⁺ â€”O Hydrogen Bonding to Ester Ether Oxygen Atoms through Spectroscopic and Computational Studies. <i>Journal of Physical Chemistry A</i> , 2012, 116, 3556-3560.	1.1	26
110	Characterization of the primary hydration shell of the hydroxide ion with H ₂ tagging vibrational spectroscopy of the OH ⁻ â€”(H ₂ O) _n =2,3 and OD ⁻ â€”(D ₂ O) _n =2,3 clusters. <i>Journal of Chemical Physics</i> , 2016, 145, 134304.	1.2	26
111	Ground-State Structure of the Proton-Bound Formate Dimer by Cold-Ion Infrared Action Spectroscopy. <i>Angewandte Chemie - International Edition</i> , 2018, 57, 10615-10619.	7.2	26
112	Double-contact ion-molecule binding: Infrared characterization of the ionic H bonds to formic acid in the H ₂ Oâ€”HCOOH complex. <i>Journal of Chemical Physics</i> , 2000, 113, 7879-7884.	1.2	25
113	Synthesis, Characterization, and Nitrogenase-Relevant Reactions of an Iron Sulfide Complex with a Bridging Hydride. <i>Journal of the American Chemical Society</i> , 2015, 137, 13220-13223.	6.6	25
114	Coordination-Dependent Spectroscopic Signatures of Divalent Metal Ion Binding to Carboxylate Head Groups: H ₂ - and He-Tagged Vibrational Spectra of M ²⁺ â€”RCO ₂ ⁻ (M =) Tj ETQq0 0 0 rgBT /Overl Chemistry Letters, 2017, 8, 484-488.	2.1	25
115	Isolating the Charge-Transfer Component of the Anionic H Bond Via Spin Suppression of the Intracuster Proton Transfer Reaction in the NOâ€”H ₂ O Entrance Channel Complex. <i>Journal of Physical Chemistry A</i> , 2002, 106, 10010-10014.	1.1	24
116	Argon cluster-mediated isolation and vibrational spectra of peroxy and nominally D _{3h} isomers of CO ₃ ⁻ and NO ₃ ⁻ . <i>Journal of Chemical Physics</i> , 2008, 129, 064305.	1.2	23
117	Isotopomer-selective spectra of a single intact H ₂ O molecule in the Cs+(D ₂ O) ₅ H ₂ O isotopologue: Going beyond pattern recognition to harvest the structural information encoded in vibrational spectra. <i>Journal of Chemical Physics</i> , 2016, 144, 074305.	1.2	23
118	Survey of Ar-Tagged Predissociation and Vibrationally Mediated Photodetachment Spectroscopies of the Vinylidene Anion, C ₂ H ₂ ⁺ . <i>Journal of Physical Chemistry A</i> , 2010, 114, 1592-1601.	1.1	22
119	Vibrational spectroscopy of the cryogenically cooled O- and N-protomers of 4-aminobenzoic acid: Tag effects, isotopic labels, and identification of the E,Z isomer of the O-protomer. <i>International Journal of Mass Spectrometry</i> , 2020, 457, 116427.	0.7	22
120	Photoelectron imaging study of vibrationally mediated electron autodetachment in the type I isomer of the water hexamer anion. <i>Chemical Physics Letters</i> , 2008, 467, 32-36.	1.2	21
121	Characterization of an activated iridium water splitting catalyst using infrared photodissociation of H ₂ tagged ions. <i>Physical Chemistry Chemical Physics</i> , 2012, 14, 10109.	1.3	21
122	Identification and Partial Structural Characterization of Mass Isolated Valsartan and Its Metabolite with Messenger Tagging Vibrational Spectroscopy. <i>Journal of the American Society for Mass Spectrometry</i> , 2017, 28, 2414-2422.	1.2	21
123	Vibrational Signatures of Solvent-Mediated Deformation of the Ternary Core Ion in Size-Selected [MgSO ₄ Mg(H ₂ O) _n =4â€”11] ²⁺ Clusters. <i>Journal of Physical Chemistry A</i> , 2015, 119, 8294-8302.	1.1	20
124	Structural Evolution of the [(CO ₂) _n (H ₂ O)] ⁺ Cluster Anions: Quantifying the Effect of Hydration on the Excess Charge Accommodation Motif. <i>Journal of Physical Chemistry A</i> , 2009, 113, 8942-8948.	1.1	19
125	Bottom-Up View of Water Network-Mediated CO ₂ Reduction Using Cryogenic Cluster Ion Spectroscopy and Direct Dynamics Simulations. <i>Journal of Physical Chemistry A</i> , 2012, 116, 903-912.	1.1	19
126	Preparation and photoelectron spectrum of the CH ₃ I ⁻ anion: Rare gas cluster mediated synthesis of an ion-radical complex. <i>Journal of the American Society for Mass Spectrometry</i> , 1999, 10, 810-814.	1.2	18

#	ARTICLE	IF	CITATIONS
127	Vibrational predissociation spectroscopy and theory of Ar-tagged, protonated Imidazole (Im) $\text{Im}^+\text{H}^+\text{Ar}$ clusters. <i>Chemical Physics Letters</i> , 2011, 501, 172-178.	1.2	18
128	N_2O at water surfaces: binding forces, charge separation, energy accommodation and atmospheric implications. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 17961-17976.	1.3	18
129	Capture of CO_2 by a Cationic Nickel(I) Complex in the Gas Phase and Characterization of the Bound, Activated CO_2 Molecule by Cryogenic Ion Vibrational Predissociation Spectroscopy. <i>Angewandte Chemie</i> , 2016, 128, 1304-1307.	1.6	17
130	Preparation of Labile Ni^+ (cyclam) Cations in the Gas Phase Using Electron-Transfer Reduction through Ion-Ion Recombination in an Ion Trap and Structural Characterization with Vibrational Spectroscopy. <i>Journal of Physical Chemistry Letters</i> , 2017, 8, 5047-5052.	2.1	17
131	Vibrational Predissociation Spectroscopy of Cold Protonated Tryptophan with Different Messenger Tags. <i>Journal of Physical Chemistry A</i> , 2018, 122, 8037-8046.	1.1	17
132	Application of gas phase cryogenic vibrational spectroscopy to characterize the CO_2 , CO , N_2 and N_2O interactions with the open coordination site on a Ni(I) macrocycle using dual cryogenic ion traps. <i>Journal of Molecular Spectroscopy</i> , 2017, 332, 117-123.	0.4	16
133	Disentangling the Complex Vibrational Mechanics of the Protonated Water Trimer by Rational Control of Its Hydrogen Bonds. <i>Journal of Physical Chemistry A</i> , 2019, 123, 7965-7972.	1.1	16
134	Mechanisms and competition of halide substitution and hydrolysis in reactions of N_2O with seawater. <i>Science Advances</i> , 2019, 5, eaav6503.	4.7	16
135	Isolating the Contributions of Specific Network Sites to the Diffuse Vibrational Spectrum of Interfacial Water with Isotopomer-Selective Spectroscopy of Cold Clusters. <i>Journal of Physical Chemistry A</i> , 2020, 124, 10393-10406.	1.1	16
136	Demystifying the Diffuse Vibrational Spectrum of Aqueous Protons Through Cold Cluster Spectroscopy. <i>Annual Review of Physical Chemistry</i> , 2021, 72, 667-691.	4.8	16
137	Infrared spectra of hydrogen-bonded ion-radical complexes: H^+HCH_2 and Br^+HCHBr . <i>Journal of Chemical Physics</i> , 1999, 111, 10464-10468.	1.2	15
138	Vibrationally Induced Interconversion of H-Bonded $\text{NO}_2^+\text{H}_2\text{O}$ Isomers within $\text{NO}_2^+\text{H}_2\text{O}^+\text{Ar}$ Clusters Using IR Pump-Probe through the OH and NO Stretching Vibrations. <i>Journal of Physical Chemistry A</i> , 2009, 113, 975-981.	1.1	15
139	Vibrational predissociation spectra of the Ar-tagged $[\text{CH}_4\text{H}_3\text{O}^+]$ binary complex: spectroscopic signature of hydrogen bonding to an alkane. <i>Molecular Physics</i> , 2010, 108, 1191-1197.	0.8	15
140	Isolation and characterization of a peroxo manganese (III) dioxygen reaction intermediate using cryogenic ion vibrational predissociation spectroscopy. <i>International Journal of Mass Spectrometry</i> , 2013, 354-355, 33-38.	0.7	15
141	Autodetachment from Vibrationally Excited Vinylidene Anions. <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 1058-1063.	2.1	15
142	Spectroscopic Assessment of Intra- and Intermolecular Hydrogen Bonding in Ether-Functionalized Imidazolium Ionic Liquids. <i>Journal of Physical Chemistry A</i> , 2019, 123, 8370-8376.	1.1	15
143	Introductory lecture: advances in ion spectroscopy: from astrophysics to biology. <i>Faraday Discussions</i> , 2019, 217, 8-33.	1.6	15
144	Mapping the temperature-dependent and network site-specific onset of spectral diffusion at the surface of a water cluster cage. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2020, 117, 26047-26052.	3.3	15

#	ARTICLE	IF	CITATIONS
145	Determination of the SmO ⁺ bond energy by threshold photodissociation of the cryogenically cooled ion. <i>Journal of Chemical Physics</i> , 2021, 155, 174303.	1.2	15
146	Chain Length Dependence of Hydrogen Bond Linkages between Cationic Constituents in Hydroxy-Functionalized Ionic Liquids: Tracking Bulk Behavior to the Molecular Level with Cold Cluster Ion Spectroscopy. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 683-688.	2.1	13
147	Isotopic fractionation in low temperature ion-molecule exchange reactions: Enrichment of ²² Ne in Ne _n clusters formed by association in an ionized free jet. <i>Journal of Chemical Physics</i> , 1990, 92, 7349-7355.	1.2	11
148	Site-specific addition of D ₂ O to the (H ₂ O) ₆ ⁺ hydrated electron cluster: isomer interconversion and substitution at the double H-bond acceptor (AA) electron-binding site. <i>Physical Chemistry Chemical Physics</i> , 2008, 10, 3118.	1.3	11
149	Trapping and Structural Characterization of the XNO ₂ ⁺ NO ₃ ⁺ (X =) Tj ETQq1 1 0.784314 rg BT Reactions with Cryogenic Vibrational Spectroscopy. <i>Journal of Physical Chemistry Letters</i> , 2017, 8, 4710-4715.	2.1	11
150	Spectroscopic Signatures of Mode-Dependent Tunnel Splitting in the Iodide-Water Binary Complex. <i>Journal of Physical Chemistry A</i> , 2020, 124, 2991-3001.	1.1	11
151	Characterization of the alkali metal oxalates (MCO ₂ O ₄) ⁺ and their formation by CO ₂ reduction via the alkali metal carbonites (MCO ₂) ⁺ . <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 7460-7473.	1.3	11
152	Electronic and mechanical anharmonicities in the vibrational spectra of the H-bonded, cryogenically cooled X ⁺ HOCl (X=Cl, Br, I) complexes: Characterization of the strong anionic H-bond to an acidic OH ₂ group. <i>Journal of Chemical Physics</i> , 2022, 156, 174303.		11
153	Electron Transfer and Charge Separation in Clusters. <i>Advances in Chemical Physics</i> , 2007, , 265-302.	0.3	9
154	Spektroskopischer Nachweis einer attraktiven Kation-Kation-Wechselwirkung in OH-funktionalisierten ionischen Flüssigkeiten: ein H ₂ Br ₂ -gebundenes kettenförmiges Trimer. <i>Angewandte Chemie</i> , 2018, 130, 15590-15594.	1.6	9
155	Unmasking Rare, Large-Amplitude Motions in D ₂ -Tagged I ⁺ (H ₂ O) ₂ Isotopomers with Two-Color, Infrared-Infrared Vibrational Predissociation Spectroscopy. <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 3744-3750.	2.1	9
156	Ionic Liquid Clusters Generated from Electrospray Thrusters: Cold Ion Spectroscopic Signatures of Size-Dependent Acid-Base Interactions. <i>Journal of Physical Chemistry A</i> , 2020, 124, 10507-10516.	1.1	9
157	Isomer-specific cryogenic ion vibrational spectroscopy of the D ₂ tagged Cs ⁺ (HNO ₃)(H ₂ O) _{n=0-2} complexes: ion-driven enhancement of the acidic H-bond to water. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 4501-4507.	1.3	9
158	Preparation and Characterization of the Halogen-Bonding Motif in the Isolated Cl ⁺ IOH Complex with Cryogenic Ion Vibrational Spectroscopy. <i>Journal of Physical Chemistry Letters</i> , 2022, 13, 2750-2756.	2.1	9
159	Strong Similarities in the Local Hydration Environments of the Bromide Ion and the Cl-CCl ₃ Radical Complex: Factors Contributing to Intramolecular Distortions in the Primary Hydration Shell. <i>Journal of Physical Chemistry A</i> , 2003, 107, 9312-9318.	1.1	8
160	Experiment and theory in harmony. <i>Nature Chemistry</i> , 2009, 1, 8-9.	6.6	8
161	Water network-mediated, electron-induced proton transfer in [C ₅ H ₅ N... (H ₂ O) _n] ⁺ clusters. <i>Journal of Chemical Physics</i> , 2015, 143, 144305.	1.2	8
162	Anharmonic Densities of States for Vibrationally Excited I ⁺ (H ₂ O) ₂ , (H ₂ O) ₂ , and I ⁺ (H ₂ O) ₂ . <i>Journal of Chemical Theory and Computation</i> , 2018, 14, 3986-3997.	2.3	8

#	ARTICLE	IF	CITATIONS
163	Argon Cluster-Mediated Trapping and Vibrational Spectroscopic Characterization of an OH- \dot{A} -HCH ₂ -Intermediate in the $O^{\ominus}C^{\ominus}+CH_4$ Reaction. <i>Journal of Physical Chemistry A</i> , 2004, 108, 10116-10121.	1.1	7
164	Communication: Spectroscopic characterization of a strongly interacting C(2)H group on the EMIM ⁺ cation in the (EMIM ⁺) ₂ X ⁻ (X = BF ₄ , Cl, Br, and I) ternary building blocks of ionic liquids. <i>Journal of Chemical Physics</i> , 2017, 147, 231101.	1.2	7
165	Size-Dependent Onset of Nitric Acid Dissociation in Cs ⁺ \dot{A} ·(HNO ₃)(H ₂ O) _n =0-11 Clusters at 20 K. <i>Journal of Physical Chemistry Letters</i> , 2021, 12, 3335-3342.	2.1	7
166	Vibrational predissociation spectroscopy of Ar-tagged, trisubstituted silyl cations. <i>Chemical Physics Letters</i> , 2013, 568-569, 9-13.	1.2	5
167	Ground-State Structure of the Proton-Bound Formate Dimer by Cold Ion Infrared Action Spectroscopy. <i>Angewandte Chemie</i> , 2018, 130, 10775-10779.	1.6	5
168	Characterization of the non-covalent docking motif in the isolated reactant complex of a double proton-coupled electron transfer reaction with cryogenic ion spectroscopy. <i>Journal of Chemical Physics</i> , 2020, 152, 234309.	1.2	5
169	Vibrational Characterization of Radical Ion Adducts between Imidazole and CO ₂ . <i>Journal of Physical Chemistry A</i> , 2018, 122, 3805-3810.	1.1	4
170	Vibrational Signatures of HNO ₃ Acidity When Complexed with Microhydrated Alkali Metal Ions, M ⁺ \dot{A} ·(HNO ₃)(H ₂ O) _n =5 (M = Li, K, Na, Rb, Cs), at 20 K. <i>Journal of Physical Chemistry A</i> , 2022, 126, 1640-1647.	1.1	4
171	On the Hydrogen Oxalate Binding Motifs onto Dinuclear Cu and Ag Metal Phosphine Complexes. <i>Chemistry - A European Journal</i> , 2021, 27, 15136-15146.	1.7	3
172	Gas-Phase Reactivity of Ozone with Lanthanide Ions (Sm ⁺ , Nd ⁺) and Their Higher Oxides. <i>Journal of the American Society for Mass Spectrometry</i> , 2022, , .	1.2	3
173	Applications of Lasers and Mass Spectrometry in Molecular Spectroscopy and Molecular Structure Determination. <i>ACS Symposium Series</i> , 0, , 277-306.	0.5	3
174	Comment on ^{13}C Vibration at C2 Position of Imidazolium Cation as a Probe of the Ionic Liquid Microenvironment. <i>Journal of Physical Chemistry A</i> , 2020, 124, 755-756.	1.1	2
175	Water Network Shape-Dependence of Local Interactions with the Microhydrated \dot{A} ⁻ NO ₂ ⁻ and \dot{A} ⁻ CO ₂ ⁻ Anionic Head Groups by Cold Ion Vibrational Spectroscopy. <i>Journal of Physical Chemistry A</i> , 2022, 126, 2471-2479.	1.1	2
176	Frontispiz: Capture of CO ₂ by a Cationic Nickel(I) Complex in the Gas Phase and Characterization of the Bound, Activated CO ₂ Molecule by Cryogenic Ion Vibrational Predissociation Spectroscopy. <i>Angewandte Chemie</i> , 2016, 128, .	1.6	0
177	Frontispiece: Capture of CO ₂ by a Cationic Nickel(I) Complex in the Gas Phase and Characterization of the Bound, Activated CO ₂ Molecule by Cryogenic Ion Vibrational Predissociation Spectroscopy. <i>Angewandte Chemie - International Edition</i> , 2016, 55, .	7.2	0
178	Chemical Reduction of Nill Cyclam and Characterization of Isolated Nil Cyclam with Cryogenic Vibrational Spectroscopy and Inert-Gas-Mediated High-Resolution Mass Spectrometry. <i>Journal of Physical Chemistry A</i> , 2021, 125, 6715-6721.	1.1	0
179	Tribute to Charles A. Schmuttenmaer. <i>Journal of Physical Chemistry C</i> , 2020, 124, 22333-22334.	1.5	0