

Philippe Maitre

List of Publications by Year in descending order

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145
papers

5,837
citations

57758

44
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88630

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all docs

151
docs citations

151
times ranked

3011
citing authors

#	ARTICLE	IF	CITATIONS
1	Ligation Motifs in Zinc-Bound Sulfonamide Drugs Assayed by IR Ion Spectroscopy. <i>Molecules</i> , 2022, 27, 3144.	3.8	0
2	Gas-Phase Dissociation Chemistry of Deprotonated RGD. <i>Journal of the American Society for Mass Spectrometry</i> , 2021, 32, 55-63.	2.8	4
3	Binding Motifs in the Naked Complexes of Target Amino Acids with an Excerpt of Antitumor Active Biomolecule: An Ion Vibrational Spectroscopy Assay. <i>Chemistry - A European Journal</i> , 2021, 27, 2348-2360.	3.3	3
4	Infrared Multiple Photon Dissociation Spectroscopy of Protonated Cyameluric Acid. <i>Journal of Physical Chemistry A</i> , 2021, 125, 607-614.	2.5	0
5	Structural Insights from Tandem Mass Spectrometry, Ion Mobility-Mass Spectrometry, and Infrared/Ultraviolet Spectroscopy on Sphingonodin I: Lasso vs Branched-Cyclic Topoisomers. <i>Journal of the American Society for Mass Spectrometry</i> , 2021, 32, 1096-1104.	2.8	4
6	Guanine Tautomerism in Ionic Complexes with Ag ⁺ Investigated by IRMPD Spectroscopy and Mass Spectrometry. <i>Journal of Physical Chemistry B</i> , 2021, 125, 7137-7146.	2.6	2
7	From Preassociation to Chelation: A Survey of Cisplatin Interaction with Methionine at Molecular Level by IR Ion Spectroscopy and Computations. <i>Journal of the American Society for Mass Spectrometry</i> , 2021, 32, 2206-2217.	2.8	7
8	Applications of Infrared Multiple Photon Dissociation (IRMPD) to the Detection of Posttranslational Modifications. <i>Chemical Reviews</i> , 2020, 120, 3261-3295.	47.7	51
9	On the Interaction between Deprotonated Cytosine [C(=NH)] ⁻ and Ba ²⁺ : Infrared Multiphoton Spectroscopy and Dynamics. <i>ChemPhysChem</i> , 2020, 21, 2571-2582.	2.1	5
10	Identification and quantification of amino acids and related compounds based on Differential Mobility Spectrometry. <i>Analyst</i> , 2020, 145, 4889-4900.	3.5	5
11	Gas phase dynamics, conformational transitions and spectroscopy of charged saccharides: the oxocarbenium ion, protonated anhydrogalactose and protonated methyl galactopyranoside. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 4144-4157.	2.8	8
12	Structure of Proton-Bound Methionine and Tryptophan Dimers in the Gas Phase Investigated with IRMPD Spectroscopy and Quantum Chemical Calculations. <i>Journal of Physical Chemistry A</i> , 2020, 124, 2408-2415.	2.5	11
13	IRMPD Spectra of Protonated Hydroxybenzaldehydes: Evidence of Torsional Barriers in Carboxonium Ions. <i>ChemPhysChem</i> , 2020, 21, 749-761.	2.1	1
14	Structural and Energetic Effects of O ² -Ribose Methylation of Protonated Pyrimidine Nucleosides. <i>Journal of the American Society for Mass Spectrometry</i> , 2019, 30, 2318-2334.	2.8	5
15	Reactions of Thiiranium and Sulfonium Ions with Alkenes in the Gas Phase. <i>Journal of Organic Chemistry</i> , 2019, 84, 10076-10087.	3.2	9
16	Evaluation of the Katsuki-Sharpless Epoxidation Precatalysts by ESI-FTMS, CID, and IRMPD Spectroscopy. <i>Journal of Physical Chemistry A</i> , 2019, 123, 1022-1029.	2.5	11
17	Probing the gas-phase structure of charge-tagged intermediates of a proline catalyzed aldol reaction – vibrational spectroscopy distinguishes oxazolidinone from enamine species. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 2578-2586.	2.8	0
18	Infrared multiple photon dissociation action spectroscopy of protonated glycine, histidine, lysine, and arginine complexed with 18-crown-6 ether. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 12625-12639.	2.8	9

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19	Infrared isomer-specific fragmentation for the identification of aminobutyric acid isomers separated by differential mobility spectrometry. <i>International Journal of Mass Spectrometry</i> , 2019, 443, 16-21.	1.5	5
20	Vibrational signatures of curcumin's chelation in copper(II) complexes: An appraisal by IRMPD spectroscopy. <i>Journal of Chemical Physics</i> , 2019, 150, 165101.	3.0	8
21	The Intermediates in Lewis Acid Catalysis with Lanthanide Triflates. <i>European Journal of Organic Chemistry</i> , 2019, 2019, 3560-3566.	2.4	12
22	Short-lived intermediates (encounter complexes) in cisplatin ligand exchange elucidated by infrared ion spectroscopy. <i>International Journal of Mass Spectrometry</i> , 2019, 435, 7-17.	1.5	20
23	Resolution and Assignment of Differential Ion Mobility Spectra of Sarcosine and Isomers. <i>Journal of the American Society for Mass Spectrometry</i> , 2018, 29, 752-760.	2.8	17
24	Insights from ion mobility-mass spectrometry, infrared spectroscopy, and molecular dynamics simulations on nicotinamide adenine dinucleotide structural dynamics: NAD ⁺ vs. NADH. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 7043-7052.	2.8	14
25	Ligand-induced decarbonylation in diphosphine-ligated palladium acetates [CH ₃ CO ₂ Pd(PR ₂ CH ₂) ₂] ⁺ (R) Tj ETQ1 1 0.784314 rB	2.8	19
26	Deprotonated carbohydrate anion fragmentation chemistry: structural evidence from tandem mass spectrometry, infra-red spectroscopy, and theory. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 27897-27909.	2.8	19
27	Sequence Ion Structures and Dissociation Chemistry of Deprotonated Sucrose Anions. <i>Journal of the American Society for Mass Spectrometry</i> , 2018, 29, 2380-2393.	2.8	15
28	Evaluation of Ca ²⁺ Binding Sites in Tacrolimus by Infrared Multiple Photon Dissociation Spectroscopy. <i>Journal of Physical Chemistry B</i> , 2018, 122, 9860-9868.	2.6	6
29	Structural and Energetic Effects of O ²⁻ -Ribose Methylation of Protonated Purine Nucleosides. <i>Journal of Physical Chemistry B</i> , 2018, 122, 9147-9160.	2.6	16
30	Mimicking the Regulation Step of Fe Monooxygenases: Allosteric Modulation of Fe ^{IV} =Oxo Formation by Guest Binding in a Dinuclear Zn ^{II} =Fe ^{II} Calix[6]arene-Based Funnel Complex. <i>Chemistry - A European Journal</i> , 2017, 23, 2894-2906.	3.3	4
31	Seleniranium Ions Undergo π -Ligand Exchange via an Associative Mechanism in the Gas Phase. <i>Journal of Organic Chemistry</i> , 2017, 82, 6289-6297.	3.2	10
32	Cytosine Radical Cations: A Gas-Phase Study Combining IRMPD Spectroscopy, UVPD Spectroscopy, Ion-Molecule Reactions, and Theoretical Calculations. <i>ChemPhysChem</i> , 2017, 18, 1293-1301.	2.1	29
33	Fingerprints of Both Watson-Crick and Hoogsteen Isomers of the Isolated (Cytosine-Guanine)H ⁺ Pair. <i>Journal of Physical Chemistry Letters</i> , 2017, 8, 5501-5506.	4.6	22
34	Pterin determination in cerebrospinal fluid: state of the art. <i>Pteridines</i> , 2017, 28, 83-89.	0.5	11
35	Cisplatin and transplatin interaction with methionine: bonding motifs assayed by vibrational spectroscopy in the isolated ionic complexes. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 26697-26707.	2.8	26
36	Watson-Crick Base Pair Radical Cation as a Model for Oxidative Damage in DNA. <i>Journal of Physical Chemistry Letters</i> , 2017, 8, 3159-3165.	4.6	22

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55	Infrared spectroscopy of nucleotides in the gas phase 2. The protonated cyclic 3',5'-adenosine monophosphate. <i>RSC Advances</i> , 2013, 3, 12711.	3.6	25
56	Gas-phase infrared spectrum and acidity of the radical cation of 9-methylguanine. <i>Chemical Communications</i> , 2013, 49, 7343.	4.1	27
57	IR Spectroscopy of b_4 Fragment Ions of Protonated Pentapeptides in the X^{H} (X = C, N, O) Region. <i>Journal of Physical Chemistry A</i> , 2013, 117, 2508-2516.	2.5	16
58	IR Signature of NO Binding to a Ferrous Heme Center. <i>Journal of Physical Chemistry Letters</i> , 2013, 4, 2414-2417.	4.6	24
59	Cyanide-Arene Meisenheimer Complex Generated in Electrospray Ionization Mass Spectrometry Using Acetonitrile as a Solvent. <i>Journal of the American Society for Mass Spectrometry</i> , 2013, 24, 1603-1607.	2.8	13
60	Binding motifs of silver in prion octarepeat model peptides: a joint ion mobility, IR and UV spectroscopies, and theoretical approach. <i>Physical Chemistry Chemical Physics</i> , 2012, 14, 11433.	2.8	28
61	S-nitrosation of cysteine as evidenced by IRMPD spectroscopy. <i>International Journal of Mass Spectrometry</i> , 2012, 330-332, 160-167.	1.5	31
62	Mechanistic Investigation of the Generation of a Palladium(0) Catalyst from a Palladium(II) Allyl Complex: A Combined Experimental and DFT Study. <i>Organometallics</i> , 2012, 31, 5975-5978.	2.3	30
63	H-bonded network rearrangements in the S0, S1 and D0 states of neutral and cationic p-cresol(H2O)(NH3) complexes. <i>Physical Chemistry Chemical Physics</i> , 2012, 14, 8945.	2.8	8
64	Benzylum versus Tropylium Ion Dichotomy: Vibrational Spectroscopy of Gaseous $C_8H_9^+$ Ions. <i>Angewandte Chemie - International Edition</i> , 2012, 51, 4947-4949.	13.8	38
65	Rearrangement Pathways of the a_4 Ion of Protonated YGGFL Characterized by IR Spectroscopy and Modeling. <i>Journal of the American Society for Mass Spectrometry</i> , 2012, 23, 664-675.	2.8	29
66	IR spectroscopy of gaseous fluorocarbon ions: The perfluoroethyl anion. <i>Chemical Physics</i> , 2012, 398, 118-123.	1.9	9
67	Tyrosine side-chain catalyzed proton transfer in the YG a_2 ion revealed by theory and IR spectroscopy in the $\tilde{\nu}$ -fingerprint TM and XH (X=C, N, O) stretching regions. <i>International Journal of Mass Spectrometry</i> , 2012, 316-318, 227-234.	1.5	12
68	Direct Evidence for Tautomerization of the Uracil Moiety within the Pb^{2+} /Uridine-5'-monophosphate Complex: A Combined Tandem Mass Spectrometry and IRMPD study. <i>Inorganic Chemistry</i> , 2011, 50, 7769-7778.	4.0	35
69	Naked Five-Coordinate Fe(III)(NO) Porphyrin Complexes: Vibrational and Reactivity Features. <i>Inorganic Chemistry</i> , 2011, 50, 4445-4452.	4.0	47
70	Structure of singly hydrated, protonated phospho-tyrosine. <i>International Journal of Mass Spectrometry</i> , 2011, 308, 338-347.	1.5	22
71	Diagnosing the Protonation Site of b_1 Peptide Fragment Ions using IRMPD in the X^{H} (X = O, N, and C) Stretching Region. <i>Journal of the American Society for Mass Spectrometry</i> , 2011, 22, 1645-50.	2.8	40
72	Spectroscopic Signatures of Peptides Containing Tryptophan Radical Cations. <i>Angewandte Chemie - International Edition</i> , 2011, 50, 11430-11432.	13.8	22

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73	Structure of Pb ²⁺ /dCMP and Pb ²⁺ /CMP complexes as characterized by tandem mass spectrometry and IRMPD spectroscopy. <i>International Journal of Mass Spectrometry</i> , 2011, 304, 154-164.	1.5	23
74	Gas phase structure of micro-hydrated [Mn(CIO ₄) ⁺ and [Mn ₂ (CIO ₄) ₃] ⁺ ions probed by infrared spectroscopy. <i>Journal of the American Society for Mass Spectrometry</i> , 2010, 21, 758-772.	2.8	37
75	Protonated Sulfuric Acid: Vibrational Signatures of the Naked Ion in the Near- and Mid-IR. <i>Journal of Physical Chemistry Letters</i> , 2010, 1, 1721-1724.	4.6	12
76	Cyclization and Rearrangement Reactions of an Fragment Ions of Protonated Peptides. <i>Journal of the American Chemical Society</i> , 2010, 132, 14766-14779.	13.7	84
77	Structure of Zirconocene Complexes Relevant for Olefin Catalysis: Infrared Fingerprint of the Zr(C ₅ H ₅) ₂ (OH)(CH ₃ CN) ⁺ Cation in the Gas Phase. <i>Journal of Physical Chemistry A</i> , 2010, 114, 2073-2079.	2.5	17
78	Chiral Recognition in Cinchona Alkaloid Protonated Dimers: Mass Spectrometry and UV Photodissociation Studies. <i>Journal of Physical Chemistry A</i> , 2010, 114, 3306-3312.	2.5	39
79	Structure and Infrared Spectrum of the Ag ⁺ Phenol Ionic Complex. <i>Journal of Physical Chemistry A</i> , 2010, 114, 11053-11059.	2.5	74
80	Cysteine radical cation: A distonic structure probed by gas phase IR spectroscopy. <i>Physical Chemistry Chemical Physics</i> , 2010, 12, 9794.	2.8	55
81	IRMPD spectroscopy of protonated S-nitrosocaptopril, a biologically active, synthetic amino acid. <i>Physical Chemistry Chemical Physics</i> , 2010, 12, 13455.	2.8	20
82	Molecular Complexes of Simple Anions with Electron-Deficient Arenes: Spectroscopic Evidence for Two Types of Structural Motifs for Anion-Arene Interactions. <i>Chemistry - A European Journal</i> , 2009, 15, 8185-8195.	3.3	44
83	Mid-IR Spectroscopy and Structural Features of Protonated Carbonic Acid in the Gas Phase. <i>ChemPhysChem</i> , 2009, 10, 520-522.	2.1	10
84	Infrared Spectroscopy of Fragments from Doubly Protonated Tryptic Peptides. <i>ChemPhysChem</i> , 2009, 10, 883-885.	2.1	74
85	Structural Characterization by IRMPD Spectroscopy and DFT Calculations of Deprotonated Phosphorylated Amino Acids in the Gas Phase. <i>ChemPhysChem</i> , 2009, 10, 1630-1641.	2.1	41
86	Room-Temperature Infrared Spectroscopy Combined with Mass Spectrometry Distinguishes Gas-Phase Protein Isomers. <i>Angewandte Chemie - International Edition</i> , 2009, 48, 8340-8342.	13.8	31
87	Tautomerism of cytosine probed by gas phase IR spectroscopy. <i>International Journal of Mass Spectrometry</i> , 2009, 283, 214-221.	1.5	47
88	Infrared Spectroscopy of Fragments of Protonated Peptides: Direct Evidence for Macrocyclic Structures of <i>b₅</i> Ions. <i>Journal of the American Chemical Society</i> , 2009, 131, 11503-11508.	13.7	92
89	The Effect of a Fourth Binding Site on the Stabilization of Cationic SPS Pincer Palladium Complexes: Experimental, DFT, and Mass Spectrometric Studies. <i>Organometallics</i> , 2009, 28, 2020-2027.	2.3	11
90	Molecular Dynamics and Room Temperature Vibrational Properties of Deprotonated Phosphorylated Serine. <i>Journal of Chemical Theory and Computation</i> , 2009, 5, 2388-2400.	5.3	35

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91	Direct Probe of NO Vibration in the Naked Ferric Heme Nitrosyl Complex. <i>ChemPhysChem</i> , 2008, 9, 826-828.	2.1	33
92	Infrared spectroscopy of isolated nucleotides. 1. The cyclic 3',5'-adenosine monophosphate anion. <i>International Journal of Mass Spectrometry</i> , 2008, 270, 111-117.	1.5	43
93	Structure of Electron-Capture Dissociation Fragments from Charge-Tagged Peptides Probed by Tunable Infrared Multiple Photon Dissociation. <i>Journal of the American Chemical Society</i> , 2008, 130, 14916-14917.	13.7	59
94	Tautomerism of Uracil Probed via Infrared Spectroscopy of Singly Hydrated Protonated Uracil. <i>Journal of Physical Chemistry A</i> , 2008, 112, 12393-12400.	2.5	96
95	Vibrational Signatures of Protonated, Phosphorylated Amino Acids in the Gas Phase. <i>Journal of the American Chemical Society</i> , 2008, 130, 3359-3370.	13.7	104
96	Structural characterization under tandem mass spectrometry conditions: infrared spectroscopy of gas phase ions. <i>Physica Scripta</i> , 2008, 78, 058111.	2.5	15
97	Gas-Phase Structure of a π -Allyl Palladium Complex: Efficient Infrared Spectroscopy in a 7 T Fourier Transform Mass Spectrometer. <i>Journal of Physical Chemistry A</i> , 2007, 111, 13415-13424.	2.5	152
98	Fingerprint Vibrational Spectra of Protonated Methyl Esters of Amino Acids in the Gas Phase. <i>Journal of the American Chemical Society</i> , 2007, 129, 2829-2840.	13.7	81
99	Meisenheimer Complexes Positively Characterized as Stable Intermediates in the Gas Phase. <i>Angewandte Chemie - International Edition</i> , 2007, 46, 1995-1998.	13.8	68
100	Infrared Spectra of Isolated Protonated Polycyclic Aromatic Hydrocarbons: Protonated Naphthalene. <i>Angewandte Chemie - International Edition</i> , 2007, 46, 6714-6716.	13.8	86
101	Micro-Hydration of the $MgNO_3^+$ Cation in the Gas Phase. <i>ChemPhysChem</i> , 2007, 8, 1629-1639.	2.1	70
102	Infrared Spectra of Protonated Uracil, Thymine and Cytosine. <i>ChemPhysChem</i> , 2007, 8, 2235-2244.	2.1	128
103	Infrared spectroscopy of organometallic ions in the gas phase: From model to real world complexes. <i>Mass Spectrometry Reviews</i> , 2007, 26, 583-605.	5.4	278
104	Protonation of heterocyclic aromatic molecules: IR signature of the protonation site of furan and pyrrole. <i>International Journal of Mass Spectrometry</i> , 2007, 267, 43-53.	1.5	43
105	Gas phase infrared multiple-photon dissociation spectra of methanol, ethanol and propanol proton-bound dimers, protonated propanol and the propanol/water proton-bound dimer. <i>Physical Chemistry Chemical Physics</i> , 2006, 8, 955.	2.8	80
106	Experimental infrared spectra of $Cl^+(ROH)$ (R = H, CH ₃ , CH ₃ CH ₂) complexes in the gas-phase. <i>Physical Chemistry Chemical Physics</i> , 2006, 8, 2483-2490.	2.8	15
107	IR Spectroscopic Features of Gaseous $C_7H_7O^+$ Ions: Benzylum versus Tropylium Ion Structures. <i>Journal of Physical Chemistry A</i> , 2006, 110, 9352-9360.	2.5	50
108	IR spectroscopy of protonated toluene: Probing ring hydrogen shifts in gaseous arenium ions. <i>International Journal of Mass Spectrometry</i> , 2006, 249-250, 149-154.	1.5	49

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109	Mid-IR spectroscopy of protonated leucine methyl ester performed with an FTICR or a Paul type ion-trap. <i>International Journal of Mass Spectrometry</i> , 2006, 249-250, 14-20.	1.5	123
110	π-Complex Structure of Gaseous Benzene ⁺ NO Cations Assayed by IR Multiple Photon Dissociation Spectroscopy. <i>Journal of the American Chemical Society</i> , 2006, 128, 12553-12561.	13.7	55
111	Infrared multiphoton dissociation spectroscopy of protonated N-acetyl-alanine and alanyl-histidine. <i>International Journal of Mass Spectrometry</i> , 2005, 243, 105-113.	1.5	69
112	Infrared Spectroscopy of Protonated Phenylsilane in the Gas Phase. <i>ChemPhysChem</i> , 2005, 6, 437-440.	2.1	32
113	Protonation Sites of Isolated Fluorobenzene Revealed by IR Spectroscopy in the Fingerprint Range. <i>Journal of Physical Chemistry A</i> , 2005, 109, 7881-7887.	2.5	57
114	Infrared spectra of homogeneous and heterogeneous proton-bound dimers in the gas phase. <i>Physical Chemistry Chemical Physics</i> , 2005, 7, 2747.	2.8	77
115	Infrared Absorption Features of Gaseous Isopropyl Carbocations. <i>ChemPhysChem</i> , 2004, 5, 1679-1685.	2.1	21
116	Investigation of the protonation site in the dialanine peptide by infrared multiphoton dissociation spectroscopy. <i>Physical Chemistry Chemical Physics</i> , 2004, 6, 2659-2663.	2.8	85
117	Infrared Multiphoton Dissociation Spectroscopy of Gas-Phase Mass-Selected Hydrocarbon ⁺ Fe ⁺ Complexes. <i>Journal of the American Chemical Society</i> , 2004, 126, 11666-11674.	13.7	47
118	Reductive Nitrile Coupling in Niobium ⁺ Acetonitrile Complexes Probed by Free Electron Laser IR Multiphoton Dissociation Spectroscopy. <i>Journal of Physical Chemistry A</i> , 2004, 108, 3350-3355.	2.5	47
119	Infrared Spectrum of the Protonated Water Dimer in the Gas Phase. <i>Journal of Physical Chemistry A</i> , 2004, 108, 9008-9010.	2.5	169
120	Vibrational Signature of Charge Solvation vs Salt Bridge Isomers of Sodiated Amino Acids in the Gas Phase. <i>Journal of the American Chemical Society</i> , 2004, 126, 1836-1842.	13.7	260
121	Infrared Fingerprint of Protonated Benzene in the Gas Phase. <i>Angewandte Chemie</i> , 2003, 115, 2103-2105.	2.0	15
122	Infrared Fingerprint of Protonated Benzene in the Gas Phase. <i>Angewandte Chemie - International Edition</i> , 2003, 42, 2057-2059.	13.8	87
123	Accurate measurement of the relative bond energies of CO and H ₂ O ligands in Fe ⁺ mono- and bis-ligated complexes. <i>Rapid Communications in Mass Spectrometry</i> , 2003, 17, 351-357.	1.5	13
124	Ultrasensitive spectroscopy of ionic reactive intermediates in the gas phase performed with the first coupling of an IR FEL with an FTICR-MS. , 2003, , 541-546.		2
125	Gas Phase Infrared Spectroscopy of Selectively Prepared Ions. <i>Physical Review Letters</i> , 2002, 89, 273002.	7.8	285
126	Infrared Spectra of Gas-Phase V ⁺ (Benzene) and V ⁺ (Benzene) ₂ Complexes. <i>Journal of the American Chemical Society</i> , 2002, 124, 1562-1563.	13.7	104

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127	<i>o</i> -, <i>m</i> -, and <i>p</i> -Diphospha-benzenes and Their P ₂ (C [∞] H) ₄ Valence Isomers. An Ab Initio Theoretical Study. <i>Journal of the American Chemical Society</i> , 1999, 121, 4215-4221.	13.7	26
128	Origin of Bonding Interactions in Cu+(H ₂) _n Clusters: An Experimental and Theoretical Investigation. <i>Journal of the American Chemical Society</i> , 1998, 120, 13494-13502.	13.7	81
129	Theoretical Study of Tungsten Carbonyl Complexes (n= 1-6): Structures, Binding Energies, and Implications for Gas Phase Reactivities. <i>Journal of Physical Chemistry A</i> , 1997, 101, 3966-3976.	2.5	19
130	Binding energies of Ti+(H ₂) ₁₋₆ clusters: Theory and experiment. <i>Journal of Chemical Physics</i> , 1997, 106, 10153-10167.	3.0	48
131	Valence Bond Analysis of the Lone Pair Bond Weakening Effect for the X-H Bonds in the Series XH _n =CH ₄ , NH ₃ , OH ₂ , FH. <i>The Journal of Physical Chemistry</i> , 1996, 100, 6463-6468.	2.9	28
132	Comment on "the origin of anomalous bond dissociation energies of V+(H ₂) _n clusters" <i>Chemical Physics Letters</i> , 1995, 242, 244-248.	2.6	6
133	Structure of V(H ₂) _n + Clusters for n = 1-6. <i>The Journal of Physical Chemistry</i> , 1995, 99, 6836-6841.	2.9	34
134	Structure of Co(H ₂) _n + Clusters, for n = 1-6. <i>The Journal of Physical Chemistry</i> , 1995, 99, 3444-3447.	2.9	46
135	The structures of small iron-carbon cluster anions. Linear to planar to three-dimensional. <i>Chemical Physics Letters</i> , 1994, 227, 601-608.	2.6	59
136	Theoretical study of the low-lying states of MgN ₂ ⁺ . <i>Chemical Physics Letters</i> , 1994, 225, 467-472.	2.6	21
137	Insertion of Sc ⁺ into H ₂ : The First Example of Cluster-Mediated σ -Bond Activation by a Transition Metal Center. <i>Journal of the American Chemical Society</i> , 1994, 116, 9710-9718.	13.7	82
138	Theoretical study of the hydrogen-metal complex (H ₂ -ML ⁺) binding energies. <i>The Journal of Physical Chemistry</i> , 1993, 97, 11912-11920.	2.9	40
139	The charge-shift bonding concept. Electron-pair bonds with very large ionic-covalent resonance energies. <i>Journal of the American Chemical Society</i> , 1992, 114, 7861-7866.	13.7	155
140	Analysis of correlation consistent wavefunctions: H ₃ X ⁺ -H bond energies (X=C, Si and Ge). <i>Chemical Physics</i> , 1992, 168, 237-247.	1.9	7
141	Covalent, ionic and resonating single bonds. <i>Computational and Theoretical Chemistry</i> , 1991, 229, 163-188.	1.5	78
142	Correlation-consistent valence bond method with purely local orbitals: application to hydrogen, lithium dimer, hydrogen fluoride, fluorine and collinear hydrogen (H ₃) and lithium (Li ₃). <i>The Journal of Physical Chemistry</i> , 1990, 94, 4082-4089.	2.9	21
143	Quantitative valence bond computations of curve-crossing diagrams for model atom exchange reactions. <i>The Journal of Physical Chemistry</i> , 1990, 94, 4089-4093.	2.9	47
144	Is the hypervalent radical SiH ₅ a stable species? An ab initio study. <i>Chemical Physics Letters</i> , 1990, 166, 49-53.	2.6	15

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145	Hypercoordination in SiH ₅ ⁻ and SiH ₅ .bul.. An electron-count dependence. Inorganic Chemistry, 1990, 29, 3047-3048.	4.0	17