

Wen-Bih Tzeng

List of Publications by Year in descending order

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papers

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186265

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138
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896
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#	ARTICLE	IF	CITATIONS
1	Kinetic energy release measurements of ammonia cluster ions during metastable decomposition and determination of cluster ion binding energies. <i>Journal of Chemical Physics</i> , 1990, 92, 332-339.	3.0	92
2	Ionization Energy of p-Fluoroaniline and Vibrational Levels of p-Fluoroaniline Cation Determined by Mass-Analyzed Threshold Ionization Spectroscopy. <i>Journal of Physical Chemistry A</i> , 1999, 103, 8612-8619.	2.5	83
3	Mass analyzed threshold ionization of deuterium substituted isotopomers of aniline and p-fluoroaniline: Isotope effect and site-specific electronic transition. <i>Journal of Chemical Physics</i> , 2001, 115, 743-751.	3.0	75
4	A 193 nm laser photofragmentation time-of-flight mass spectrometric study of CS ₂ and CS ₂ clusters. <i>Journal of Chemical Physics</i> , 1988, 88, 1658-1669.	3.0	73
5	Dissociation dynamics: Measurements of decay fractions of metastable ammonia cluster ions. <i>Journal of Chemical Physics</i> , 1990, 93, 2506-2512.	3.0	73
6	A Gaussian ab initio study of CH ₂ SH, CH ₂ S ⁺ , CH ₃ S ⁺ , CH ₂ SH ⁺ , CH ₃ SH ⁺ , CH ₃ ⁺ , and CH ₃ SH ⁺ . <i>Journal of Chemical Physics</i> , 1992, 97, 6557-6568.	3.0	66
7	A study of the structures and vibrations of C ₆ H ₅ NH ₂ , C ₆ H ₅ NHD, C ₆ H ₅ ND ₂ , C ₆ D ₅ NH ₂ , C ₆ D ₅ NHD, and C ₆ D ₅ ND ₂ in the S ₁ state by ab initio calculations. <i>Computational and Theoretical Chemistry</i> , 1998, 428, 231-240.	1.5	64
8	Multiphoton ionization of acetone clusters: metastable unimolecular decomposition of acetone cluster ions and the influence of solvation on intracuster ion-molecule reactions. <i>Journal of the American Chemical Society</i> , 1989, 111, 6035-6040.	13.7	55
9	Photophysics of clusters. Intracuster reactions and dynamics of dissociation processes. <i>Journal of the Chemical Society, Faraday Transactions</i> , 1990, 86, 2417.	1.7	52
10	Structures and vibrations of p-methylaniline in the S ₀ and S ₁ states studied by ab initio calculations and resonant two-photon ionization spectroscopy. <i>Journal of Molecular Structure</i> , 1998, 446, 93-102.	3.6	48
11	Mass analyzed threshold ionization spectroscopy of anisole cation and the OCH ₃ substitution effect. <i>Chemical Physics Letters</i> , 2005, 407, 100-104.	2.6	43
12	Spectroscopy of phenylacetylene bound to clusters of ammonia and the surface cluster analogy. <i>Journal of Chemical Physics</i> , 1989, 90, 11-18.	3.0	42
13	Structures and vibrations of ortho-, meta-, and para-fluoroanilines in the S ₀ and S ₁ states by ab initio calculations and resonant two-photon ionization spectroscopy. <i>Journal of the Chemical Society, Faraday Transactions</i> , 1997, 93, 2981-2987.	1.7	42
14	Mass analyzed threshold ionization spectroscopy of p-fluorophenol cation and the p-fluoro substitution effect. <i>Chemical Physics Letters</i> , 2004, 390, 65-70.	2.6	42
15	Molecular beam photoionization study of HgBr ₂ and HgI ₂ . <i>Journal of Chemical Physics</i> , 1983, 78, 50-61.	3.0	39
16	Photodissociation of CH ₂ BrCl at 248 and 193 nm investigated by translational spectroscopy. <i>Chemical Physics Letters</i> , 1994, 227, 467-471.	2.6	39
17	Mass-Analyzed Threshold Ionization Spectroscopy of o-, m-, and p-Methylaniline Cations: Vicinal Substitution Effects on Electronic Transition, Ionization, and Molecular Vibration. <i>Journal of Physical Chemistry A</i> , 2002, 106, 6462-6468.	2.5	38
18	Mass-Analyzed Threshold Ionization Spectroscopy of the Selected Rotamers of Hydroquinone and p-Dimethoxybenzene Cations. <i>Journal of Physical Chemistry A</i> , 2001, 105, 11455-11461.	2.5	36

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19	Ionization energy of o-fluoroaniline and vibrational levels of o-fluoroaniline cation determined by mass-analyzed threshold ionization spectroscopy. <i>Physical Chemistry Chemical Physics</i> , 2000, 2, 3759-3763.	2.8	35
20	Mass analyzed threshold ionization spectroscopy of o-fluorophenol and o-methoxyphenol cations and influence of the nature and relative location of substituents. <i>Chemical Physics</i> , 2006, 323, 429-438.	1.9	35
21	Metastable unimolecular and collision-induced dissociation of hydrogen-bonded clusters: evidence for intracuster molecular rearrangement and the structure of solvated protonated complexes. <i>Journal of the American Chemical Society</i> , 1991, 113, 1960-1969.	13.7	34
22	Mass analyzed threshold ionization spectroscopy of 4-aminobenzonitrile cation. <i>Chemical Physics</i> , 2000, 261, 449-455.	1.9	34
23	Mass analyzed threshold ionization of the ³⁵ Cl and ³⁷ Cl isotopomers of p-chloroaniline. <i>Journal of Chemical Physics</i> , 2000, 113, 4109-4115.	3.0	34
24	Molecular beam photoionization study of H ₂ S. <i>International Journal of Mass Spectrometry and Ion Physics</i> , 1983, 50, 315-329.	1.3	33
25	Mass-analyzed threshold ionization spectroscopy of p-methylphenol and p-ethylphenol cations and the alkyl substitution effect. <i>Journal of Chemical Physics</i> , 2004, 120, 10513-10519.	3.0	32
26	Evaporative dissociation of ammonia cluster ions: quantification of decay fractions and isotope effects. <i>The Journal of Physical Chemistry</i> , 1991, 95, 8306-8309.	2.9	31
27	Species-Selected Mass-Analyzed Threshold Ionization Spectra of m-Fluoroaniline Cation. <i>Applied Spectroscopy</i> , 2001, 55, 120-124.	2.2	31
28	Mass analyzed threshold ionization spectroscopy of 2-aminopyridine cation. <i>Chemical Physics Letters</i> , 2002, 353, 55-62.	2.6	31
29	A study of the unimolecular decomposition of the (C ₂ H ₄) ₂ complex. <i>Journal of Chemical Physics</i> , 1984, 80, 1482-1489.	3.0	29
30	Excited-state structure and vibrations of p-diaminobenzene studied by ab initio calculations. <i>Computational and Theoretical Chemistry</i> , 1998, 434, 247-253.	1.5	29
31	Structure of protonated solvation complexes: ammonia-trimethylamine cluster ions and their metastable decompositions. <i>The Journal of Physical Chemistry</i> , 1991, 95, 585-591.	2.9	28
32	Molecular beam photoionization study of HgCl ₂ . <i>Journal of Chemical Physics</i> , 1983, 78, 37-45.	3.0	27
33	S ₁ → S ₀ transition of phenylacetylene: ab initio and resonant two-photon ionization studies. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 1996, 52, 1703-1716.	3.9	26
34	Intracuster Reaction, Fragmentation, and Structure of Monomethylamine, Dimethylamine, and Trimethylamine Cluster Ions. <i>The Journal of Physical Chemistry</i> , 1996, 100, 15340-15345.	2.9	25
35	Mass-analyzed threshold ionization spectroscopy of the rotamers of p-n-propylphenol cations and configuration effect. <i>Journal of Chemical Physics</i> , 2005, 122, 044311.	3.0	25
36	Intracuster reactions in phenylacetylene ammonia clusters initiated through resonant enhanced ionization. <i>Journal of Chemical Physics</i> , 1989, 90, 19-24.	3.0	24

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37	Protonated acetaldehyde clusters: Stability, structure and metastable unimolecular decomposition. <i>Chemical Physics Letters</i> , 1990, 168, 30-36.	2.6	23
38	A study of the excited state structure and vibrations of hydroquinone by ab initio calculations and resonant two-photon ionization spectroscopy. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 1997, 53, 2595-2604.	3.9	23
39	Simultaneous Detection of C ₆ H ₅ NHD, C ₆ H ₅ ND ₂ , C ₆ D ₅ NH ₂ , C ₆ D ₅ NHD, and C ₆ D ₅ ND ₂ by Resonant Two-Photon Ionization Mass Spectrometry. <i>Applied Spectroscopy</i> , 1998, 52, 890-893.	2.2	23
40	Resonant two-photon ionization and mass-analyzed threshold ionization spectroscopy of the selected rotamers of m-methoxyaniline and o-methoxyaniline. <i>Journal of Molecular Spectroscopy</i> , 2007, 244, 1-8.	1.2	23
41	Vibronic features of p-ethylaniline, p-ethylaniline-NHD, and p-ethylaniline-ND ₂ by resonant two-photon ionization mass spectrometry. <i>Journal of Molecular Structure</i> , 1999, 482-483, 315-322.	3.6	22
42	Cation spectroscopy of 3,4-difluoroaniline by two-color resonant two-photon mass-analyzed threshold ionization. <i>Journal of Molecular Spectroscopy</i> , 2011, 266, 52-56.	1.2	22
43	Vibronic and cation spectroscopy of 2,4-difluoroaniline. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2012, 93, 176-179.	3.9	22
44	Rotamers of p-methoxyphenol cation studied by mass analyzed threshold ionization spectroscopy. <i>Chemical Physics Letters</i> , 2005, 410, 99-103.	2.6	21
45	Cation spectroscopy of o-fluoroanisole and p-fluoroanisole by two-color resonant two-photon mass-analyzed threshold ionization. <i>Chemical Physics Letters</i> , 2012, 524, 38-41.	2.6	21
46	Photoionization study of HgAr. <i>Journal of Chemical Physics</i> , 1985, 82, 648-652.	3.0	20
47	Influence of solvation on dissociation: Metastable unimolecular decomposition of mixed ammonia- ϵ -acetone cluster ions. <i>Chemical Physics Letters</i> , 1990, 166, 343-352.	2.6	20
48	Ammonia-acetone mixed clusters: internal ion-molecule reactions, structure, and bonding. <i>Journal of the American Chemical Society</i> , 1990, 112, 4097-4104.	13.7	20
49	Rotamers of <i>o</i> - and <i>m</i> -Dimethoxybenzenes Studied by Mass-Analyzed Threshold Ionization Spectroscopy and Theoretical Calculations. <i>Journal of Physical Chemistry A</i> , 2010, 114, 11144-11152.	2.5	20
50	Mass analyzed threshold ionization spectroscopy of the ³⁵ Cl and ³⁷ Cl isotopomers of p-chlorophenol and isotope effect. <i>Chemical Physics Letters</i> , 2006, 422, 271-275.	2.6	19
51	Identification of four rotamers of m-methoxystyrene by resonant two-photon ionization and mass analyzed threshold ionization spectroscopy. <i>Journal of Chemical Physics</i> , 2015, 142, 124314.	3.0	19
52	Rotamers of m-fluoroanisole studied by two-color resonant two-photon mass-analyzed threshold ionization spectroscopy. <i>Journal of Molecular Spectroscopy</i> , 2012, 274, 43-47.	1.2	18
53	Mass analyzed threshold ionization of p-bromoaniline: heavy atom effects on electronic transition, ionization, and molecular vibration. <i>Chemical Physics Letters</i> , 2002, 356, 267-276.	2.6	17
54	3-Chloro-4-fluoroaniline studied by resonant two-photon ionization and mass-analyzed threshold ionization spectroscopy. <i>Journal of Molecular Spectroscopy</i> , 2013, 288, 1-6.	1.2	17

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55	Mass analyzed threshold ionization spectroscopy of p-methoxyaniline cation and influence of the OCH ₃ substituent. <i>Chemical Physics Letters</i> , 2003, 370, 44-51.	2.6	16
56	Studies of Structural Isomers <i>o</i> -, <i>m</i> -, and <i>p</i> -Fluorophenylacetylene by Two-Color Resonant Two-Photon Mass-Analyzed Threshold Ionization Spectroscopy. <i>Journal of Physical Chemistry A</i> , 2014, 118, 8277-8286.	2.5	16
57	A study of the unimolecular decomposition of the (C ₂ H ₄) ₃ complex. <i>Journal of Chemical Physics</i> , 1985, 83, 2813-2817.	3.0	15
58	Mass analyzed threshold ionization spectroscopy of 3-aminopyridine cation and vicinal substitution effect. <i>Chemical Physics</i> , 2002, 280, 191-203.	1.9	15
59	Photoionization study of hydrogen sulfide ((H ₂ S) ₂ and (H ₂ S) ₃). <i>Journal of the American Chemical Society</i> , 1983, 105, 7531-7536.	13.7	14
60	Mass analyzed threshold ionization spectroscopy of p-ethylaniline cation: alkyl chain effects on ionization and molecular vibration. <i>Chemical Physics Letters</i> , 2002, 362, 19-25.	2.6	14
61	Mass analyzed threshold ionization spectroscopy of N-methylaniline, N-ethylaniline, and N,N-dimethylaniline cations: Influence of N-alkyl substitution on the ionization energy and molecular vibration. <i>Journal of Chemical Physics</i> , 2003, 118, 4929-4937.	3.0	14
62	Mass analyzed threshold ionization spectroscopy of p-cyanophenol cation and the CN substitution effect. <i>Chemical Physics Letters</i> , 2005, 411, 506-510.	2.6	14
63	Structures and vibrations of p-dimethoxybenzene conformers in the S ₀ and S ₁ states studied by ab initio calculations and resonant two-photon ionization spectroscopy. <i>Journal of Molecular Structure</i> , 1998, 448, 91-100.	3.6	13
64	Identification of Impurities in Phenylacetylene by Species-Selected Mass-Analyzed Threshold Ionization Spectroscopy. <i>Applied Spectroscopy</i> , 2003, 57, 1178-1182.	2.2	13
65	Rotamers of m-aminophenol cation studied by mass analyzed threshold ionization spectroscopy and theoretical calculations. <i>Chemical Physics Letters</i> , 2004, 394, 182-187.	2.6	13
66	Two-color resonant two-photon ionization and mass-analyzed threshold ionization spectroscopy of o-chloroanisole. <i>Journal of Photochemistry and Photobiology A: Chemistry</i> , 2012, 243, 73-79.	3.9	13
67	Resonant two-photon mass-analyzed threshold ionization of 2,5-difluoroaniline. <i>Chemical Physics Letters</i> , 2013, 580, 28-31.	2.6	13
68	Photodissociation of CBrCl ₃ at 248 nm by translational spectroscopy. <i>Chemical Physics Letters</i> , 1994, 222, 141-145.	2.6	12
69	Detection of Styrene Impurities in Phenylacetylene by Resonance-Enhanced Multiphoton Ionization Time-of-Flight Mass Spectrometry. <i>Applied Spectroscopy</i> , 1999, 53, 731-734.	2.2	12
70	Rotamers of 3,4-difluorophenol studied by two-color resonant two-photon mass-analyzed threshold ionization spectroscopy. <i>Journal of Photochemistry and Photobiology A: Chemistry</i> , 2013, 270, 53-59.	3.9	12
71	4-Chloro-3-fluoroaniline studied by resonant two-photon ionization and mass-analyzed threshold ionization spectroscopy. <i>Chemical Physics Letters</i> , 2014, 595-596, 73-76.	2.6	12
72	Reactions of hydrogen halides with clusters of ammonia molecules. <i>The Journal of Physical Chemistry</i> , 1989, 93, 7703-7707.	2.9	11

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73	Observation of the influence of isomeric structures of cluster ions on the dynamics of dissociation: ammonia-triethylamine system. <i>The Journal of Physical Chemistry</i> , 1990, 94, 6927-6930.	2.9	11
74	Multiphoton ionization of ether clusters: intracuster ion-molecule reactions and metastable decompositions. <i>The Journal of Physical Chemistry</i> , 1991, 95, 5080-5085.	2.9	11
75	Evidence of cyclic structures in protonated hydrogen-bonded complexes. <i>Chemical Physics Letters</i> , 1991, 178, 411-418.	2.6	11
76	Mass analyzed threshold ionization spectroscopy of N-deuterium substituted indoline cation: isotope effect on the electronic transition, ionization and molecular vibration. <i>Chemical Physics Letters</i> , 2003, 371, 662-669.	2.6	11
77	Rotamers of m-chloroanisole studied by two-color resonant two-photon mass-analyzed threshold ionization spectroscopy. <i>Chemical Physics</i> , 2013, 425, 114-120.	1.9	11
78	Rotamers of 3,4-difluoroanisole studied by two-color resonant two-photon mass-analyzed threshold ionization spectroscopy. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2013, 102, 365-370.	3.9	11
79	Mass analyzed threshold ionization spectroscopy of 1-methylindole cation. <i>Chemical Physics Letters</i> , 2003, 377, 620-626.	2.6	10
80	Vibronic and cation spectroscopy of m-chloroaniline. <i>Journal of Molecular Spectroscopy</i> , 2011, 269, 248-253.	1.2	10
81	DFT-supported Threshold Ionization Study of Chromium Biphenyl Complexes: Unveiling the Mechanisms of Substituent Influence on Redox Properties of Sandwich Compounds. <i>Chemistry - A European Journal</i> , 2017, 23, 13669-13675.	3.3	10
82	Spectroscopic Investigation of cis-2,4-Difluorophenol Cation by Mass-analyzed Threshold Ionization Spectroscopy. <i>Bulletin of the Korean Chemical Society</i> , 2014, 35, 815-820.	1.9	10
83	A study of the unimolecular decompositions of the $(C_3H_6)_2^+$ and $(C_3H_6)_2^+$ complexes. <i>Journal of Chemical Physics</i> , 1985, 83, 2803-2812.	3.0	9
84	Stability, structure, and binding energies of solvated cluster ions: ammonia-acetonitrile and ammonia-acetaldehyde systems. <i>The Journal of Physical Chemistry</i> , 1991, 95, 5757-5763.	2.9	9
85	Mass analyzed threshold ionization spectroscopy of indoline cation: Cyclization effect and large amplitude vibrations. <i>Journal of Chemical Physics</i> , 2003, 118, 10034-10041.	3.0	9
86	Mass analyzed threshold ionization spectroscopy of p-methylanisole cation and the substitution effect. <i>Chemical Physics Letters</i> , 2005, 414, 276-281.	2.6	9
87	Vibrational spectra and theoretical calculations of p-chlorophenol in the electronically excited S ₁ and ionic ground D ₀ states. <i>Journal of Photochemistry and Photobiology A: Chemistry</i> , 2008, 193, 245-253.	3.9	9
88	Mass-analyzed threshold ionization of deuterium substituted indazole and benzimidazole and site-specific H/D exchange reaction. <i>Chemical Physics Letters</i> , 2010, 501, 6-10.	2.6	9
89	Ionization energy and active cation vibrations of trans-2-fluorostyrene. <i>Journal of Molecular Spectroscopy</i> , 2017, 332, 3-7.	1.2	9
90	Determination of laser beam waist using photoionization time-of-flight mass spectrometer. <i>Review of Scientific Instruments</i> , 1994, 65, 2776-2780.	1.3	8

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91	Vibronic features of 2,6-dimethylaniline, 2,6-dimethylaniline- ¹⁵ NHD, and 2,6-dimethylaniline- ¹⁵ ND ₂ by resonant two-photon ionization mass spectrometry. <i>Journal of the Chemical Society, Faraday Transactions</i> , 1998, 94, 2913-2917.	1.7	8
92	Mass analyzed threshold ionization spectroscopy of deuterium substituted N-methylaniline and N-ethylaniline cations: isotope effect on transition energy and large amplitude vibrations. <i>Chemical Physics</i> , 2003, 295, 97-107.	1.9	8
93	Mass analyzed threshold ionization spectroscopy of 5-methylindole and 3-methylindole cations and the methyl substitution effect. <i>Journal of Chemical Physics</i> , 2004, 120, 5057-5063.	3.0	8
94	Mass analyzed threshold ionization spectroscopy of aza-aromatic bicyclic molecules: Benzimidazole and benzotriazole. <i>Chemical Physics</i> , 2007, 334, 189-195.	1.9	8
95	Molecular structures and vibrations of cis and trans m-cresol in the electronically excited S ₁ and cationic D ₀ states. <i>Journal of Photochemistry and Photobiology A: Chemistry</i> , 2007, 188, 252-259.	3.9	8
96	Active vibrations of indene cation studied by mass-analyzed threshold ionization spectroscopy. <i>Journal of Photochemistry and Photobiology A: Chemistry</i> , 2011, 220, 139-144.	3.9	8
97	Fine Substituent Effects in Sandwich Complexes: A Threshold Ionization Study of Monosubstituted Chromium Bisarene Compounds. <i>Chemistry - A European Journal</i> , 2016, 22, 4690-4694.	3.3	8
98	Site-Specific H/D Exchange of p-Methoxyphenol Studied by Resonant Two-Photon Ionization and Mass-Analyzed Threshold Ionization Spectroscopy. <i>Journal of Physical Chemistry A</i> , 2005, 109, 9481-9487.	2.5	7
99	Vibrational Spectrum of o-Dimethoxybenzene in the S ₁ and D ₀ States. <i>Chinese Journal of Chemistry</i> , 2008, 26, 51-54.	4.9	7
100	Resonant two-photon mass-analyzed threshold ionization spectroscopy of 1-fluoronaphthalene and 2-fluoronaphthalene. <i>Journal of Molecular Spectroscopy</i> , 2012, 281, 40-46.	1.2	7
101	Mass analyzed threshold ionization spectroscopy of 7-azaindole cation. <i>Chemical Physics Letters</i> , 2003, 380, 503-511.	2.6	6
102	Mass analyzed threshold ionization spectroscopy of indazole cation. <i>Chemical Physics Letters</i> , 2005, 411, 86-90.	2.6	6
103	Rotamers of m-cresol cation studied by mass-analyzed threshold ionization spectroscopy. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2007, 67, 989-994.	3.9	6
104	Selected cis- and trans-p-methoxystyrene rotamers studied by mass-analyzed threshold ionization spectroscopy. <i>Chemical Physics Letters</i> , 2011, 503, 25-28.	2.6	6
105	Two-color resonant two-photon ionization and mass-analyzed threshold ionization spectroscopy of 4-chlorostyrene. <i>Chemical Physics Letters</i> , 2017, 682, 34-37.	2.6	6
106	Cation spectra of p-chloroanisole and the heavy atom effect on ionization energy. <i>Chemical Physics Letters</i> , 2019, 731, 136626.	2.6	6
107	Stable shell structures in hydrogen-bonded complexes. <i>Zeitschrift für Physik D-Atoms Molecules and Clusters</i> , 1991, 20, 47-51.	1.0	5
108	Active vibrations of 1-cyanonaphthalene cation studied by mass-analyzed threshold ionization spectroscopy. <i>Chemical Physics Letters</i> , 2013, 558, 20-24.	2.6	5

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109	Laser spectroscopic and computational insights into unexpected structural behaviours of sandwich complexes upon ionization. <i>Dalton Transactions</i> , 2021, 50, 10729-10736.	3.3	5
110	Mass analyzed threshold ionization spectroscopy of p-aminophenol cation and the substitution effect. <i>Chemical Physics</i> , 2004, 305, 285-290.	1.9	4
111	Vibronic and cation spectroscopy of p-ethynylaniline. <i>Chemical Physics Letters</i> , 2012, 543, 19-22.	2.6	4
112	Electronic excited states of chromium and vanadium bisarene complexes revisited: interpretation of the absorption spectra on the basis of TD DFT calculations. <i>Dalton Transactions</i> , 2014, 43, 17703-17711.	3.3	4
113	Rotamers of m-Methylanisole Studied by Mass-Analyzed Threshold Ionization Spectroscopy. <i>Wuli Huaxue Xuebao/ Acta Physico - Chimica Sinica</i> , 2014, 30, 1416-1425.	4.9	4
114	Selected cis- and trans-3-fluorostyrene rotamers studied by two-color resonant two-photon mass-analyzed threshold ionization spectroscopy. <i>Journal of Molecular Spectroscopy</i> , 2015, 316, 72-78.	1.2	4
115	Resonant two-photon ionization and mass-analyzed threshold ionization spectroscopy of 3,5-difluorophenol. <i>Chemical Physics Letters</i> , 2018, 700, 145-148.	2.6	4
116	Cation Vibrations of 1-Methylnaphthalene and 2-Methylnaphthalene through Mass-Analyzed Threshold Ionization Spectroscopy. <i>Journal of Physical Chemistry A</i> , 2019, 123, 5969-5979.	2.5	4
117	Quantum-Chemical Modeling of the Mass-analyzed Threshold Ionization Spectra of Ferrocene and Cobaltocene. <i>High Energy Chemistry</i> , 2020, 54, 414-420.	0.9	4
118	Multiphoton ionization of acetone clusters: metastable unimolecular decomposition of acetone cluster ions and the influence of solvation on intracluster ion-molecule reactions [Erratum to document cited in CA111(9):77364q]. <i>Journal of the American Chemical Society</i> , 1989, 111, 8326-8326.	13.7	3
119	Mass-analyzed Threshold Ionization Spectroscopy of Rotamers of <i>p</i> -ethoxyphenol Cations and Configuration Effect. <i>Chinese Journal of Chemical Physics</i> , 2009, 22, 649-654.	1.3	3
120	Mass-analyzed threshold ionization spectroscopy of deuterium-substituted isotopomers of o-fluoroaniline and m-fluoroaniline cations. <i>Journal of Molecular Spectroscopy</i> , 2011, 269, 49-55.	1.2	3
121	Mass-analyzed threshold ionization spectroscopy of trans-1-methoxynaphthalene cation and the methoxyl substitution effect. <i>Journal of Molecular Spectroscopy</i> , 2013, 284-285, 16-20.	1.2	3
122	Vibronic and cation spectroscopy of selected rotamers of 4-chloro-3-fluorophenol. <i>Molecular Physics</i> , 2014, 112, 2397-2406.	1.7	3
123	TD DFT insights into unusual properties of excited sandwich complexes: structural transformations and vibronic interactions in Rydberg-state bis(l-6-benzene)chromium. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 23988-23997.	2.8	3
124	Mass-analyzed threshold ionization spectroscopy of trans-o-methylanisole. <i>Journal of Molecular Spectroscopy</i> , 2019, 355, 26-31.	1.2	3
125	Multiphoton Ionization of Benzene- ϵ -Ammonia Clusters: Intracluster Reaction and Cluster Ion Stability. <i>Journal of the Chinese Chemical Society</i> , 1994, 41, 505-509.	1.4	2
126	Vibrations and theoretical calculations of p-methylanisole in the first electronically excited S ₁ and ionic ground D ₀ states. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2007, 67, 824-829.	3.9	2

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127	Resonant two-photon ionization and mass-analyzed threshold ionization spectroscopy of p-vinylaniline. <i>Chemical Physics</i> , 2012, 407, 71-75.	1.9	2
128	Mass-analyzed threshold ionization spectroscopy of 2,6-dimethylaniline, 2,6-dimethylaniline-NHD, and 2,6-dimethylaniline-ND ₂ . <i>Chemical Physics Letters</i> , 2012, 551, 50-53.	2.6	2
129	Vibronic and cation spectroscopy of structural isomers p- and m-diaminobenzene and the amino substitution effect. <i>Journal of Photochemistry and Photobiology A: Chemistry</i> , 2013, 251, 94-99.	3.9	2
130	Rydberg state mediated multiphoton ionization of (I ⁷⁺ -C ₇ H ₇)(I ⁵⁺ -C ₅ H ₅)Cr: DFT-supported experimental insights into the molecular and electronic structures of excited sandwich complexes. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 9665-9671.	2.8	2
131	Rotamers of 2, 5-Difluorophenol Studied Using Mass-Analyzed Threshold Ionization Spectroscopy. <i>Wuli Huaxue Xuebao/ Acta Physico - Chimica Sinica</i> , 2016, 32, 893-900.	4.9	2
132	Two-Color Resonant Two-Photon Mass-Analyzed Threshold Ionization of 2,4-Difluoroanisole and the Additivity Relation of Ionization Energy. <i>Journal of Physical Chemistry A</i> , 2020, 124, 10517-10526.	2.5	2
133	Spectroscopy of phenylacetylene-carbon dioxide clusters. <i>Chemical Physics Letters</i> , 1988, 150, 231-234.	2.6	1
134	Investigations on the Photoreactions of Phenothiazine and Phenoxazine in Presence of 9-cyanoanthracene by Using Steady State and Time Resolved Spectroscopic Techniques. <i>Journal of Fluorescence</i> , 2010, 20, 1061-1068.	2.5	1
135	Ultraviolet laser ionization studies of 1-fluoronaphthalene clusters and density functional theory calculations. <i>Chinese Physics B</i> , 2010, 19, 123602.	1.4	1
136	Mass analyzed threshold ionization spectroscopy of methyl-p-aminobenzoate cation. <i>Chemical Physics Letters</i> , 2006, 421, 77-80.	2.6	0
137	LIV Resonant Two-Photon Ionization Spectrum of 1-Naphthol. <i>Wuli Huaxue Xuebao/ Acta Physico - Chimica Sinica</i> , 2009, 25, 2488-2492.	4.9	0
138	Tribute to Cheuk-Yiu Ng. <i>Journal of Physical Chemistry A</i> , 2021, 125, 7353-7355.	2.5	0