Evan Bieske

List of Publications by Year
in descending order

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134	3,547	126907	175258
papers	citations	33	52
		h-index	g-index

135	135	135	2475
all docs	docs citations	times ranked	citing authors

#	Article	IF	Citations
1	Photodissociation dynamics of N3+. Journal of Chemical Physics, 2022, 156, 124307.	3.0	0
2	An ion mobility mass spectrometer coupled with a cryogenic ion trap for recording electronic spectra of charged, isomer-selected clusters. Review of Scientific Instruments, 2022, 93, 043201.	1.3	11
3	Action spectroscopy of deprotomer-selected hydroxycinnamate anions. European Physical Journal D, 2021, 75, 1.	1.3	5
4	Action spectroscopy of isomer-selected luciferin anions. European Physical Journal D, 2021, 75, 1.	1.3	3
5	Electronic spectra of positively charged carbon clustersâ€”C2n+ (n = 6â€“14). Journal of Chemical Physics, 2021, 155, 214302.	3.0	13
6	Photoisomerization of Linear and Stacked Isomers of a Charged Styryl Dye: A Tandem Ion Mobility Study. Journal of the American Society for Mass Spectrometry, 2021, 32, 2842-2851.	2.8	2
7	Actinic Wavelength Action Spectroscopy of the IO^{â€“} Reaction Intermediate. Journal of Physical Chemistry Letters, 2021, 12, 11939-11944.	4.6	1
8	Nonadiabatic Dynamics between Valence, Nonvalence, and Continuum Electronic States in a Heteropolycyclic Aromatic Hydrocarbon. Journal of Physical Chemistry Letters, 2021, 12, 11811-11816.	4.6	4
9	Reversible Photoswitching of Isolated Ionic Hemiindigos with Visible Light. ChemPhysChem, 2020, 21, 680-685.	2.1	14
10	Photophysics of Isolated Rose Bengal Anions. Journal of Physical Chemistry A, 2020, 124, 8429-8438.	2.5	7
11	Electronic Spectrum of the Tropylium Cation in the Gas Phase. Journal of Physical Chemistry Letters, 2020, 11, 8867-8872.	4.6	6
12	N3+: Full-dimensional ground state potential energy surface, vibrational energy levels, and dynamics. Journal of Chemical Physics, 2020, 153, 044302.	3.0	7
13	Photo- and Collision-Induced Isomerization of a Charge-Tagged Norbornadieneâ€“Quadricyclane System. Journal of Physical Chemistry Letters, 2020, 11, 6045-6050.	4.6	15
14	Electronic Spectrum and Photodissociation Chemistry of the 1-Butyn-3-yl Cation, H₃CCHCCH⁺. Journal of Physical Chemistry A, 2020, 124, 2366-2371.	2.5	3
15	Near-infrared reversible photoswitching of an isolated azobenzene-stilbene dye. Chemical Physics Letters, 2020, 741, 137065.	2.6	3
16	Electronic Spectra of Diacetylene Cations (HC₄H⁺) Tagged with Ar and N₂. Journal of Physical Chemistry A, 2019, 123, 7228-7236.	2.5	1
17	Photodetachment and photoreactions of substituted naphthalene anions in a tandem ion mobility spectrometer. Faraday Discussions, 2019, 217, 34-46.	3.2	13
18	Ultrafast photoisomerisation of an isolated retinoid. Physical Chemistry Chemical Physics, 2019, 21, 10567-10579.	2.8	12

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19	Electronic Spectra of the Triacetylene Cation (HC6H+) and Protonated Triacetylene (HC6H2+) Tagged with Ar. Australian Journal of Chemistry, 2019, 72, 260.	0.9	1
20	Photoinitiated Intramolecular Proton Transfer in Deprotonated para-Coumaric Acid. Journal of Physical Chemistry A, 2019, 123, 4419-4430.	2.5	18
21	Structural characterization and gas-phase studies of the [Ag₁₀H₈(L)₆]²⁺ nanocluster dication. Nanoscale, 2019, 11, 22880-22889.	5.6	16
22	Unveiling New Isomers and Rearrangement Routes on the C₇H₈⁺ Potential Energy Surface. Journal of Physical Chemistry A, 2019, 123, 823-830.	2.5	5
23	Linkage Photoisomerization of an Isolated Ruthenium Sulfoxide Complex: Sequential versus Concerted Rearrangement. Inorganic Chemistry, 2018, 57, 5701-5706.	4.0	7
24	Photoswitching an Isolated Donorâ€“Acceptor Stenhouse Adduct. Journal of Physical Chemistry Letters, 2018, 9, 665-671.	4.6	46
25	Reversible Photoisomerization of the Isolated Green Fluorescent Protein Chromophore. Journal of Physical Chemistry Letters, 2018, 9, 2647-2651.	4.6	23
26	Differential-Mobility Spectrometry of 1-Deoxysphingosine Isomers: New Insights into the Gas Phase Structures of Ionized Lipids. Analytical Chemistry, 2018, 90, 5343-5351.	6.5	31
27	FromEtoZand back again: reversible photoisomerisation of an isolated charge-tagged azobenzene. Physical Chemistry Chemical Physics, 2018, 20, 509-513.	2.8	14
28	Double Molecular Photoswitch Driven by Light and Collisions. Physical Review Letters, 2018, 120, 223002.	7.8	24
29	Ion mobility action spectroscopy of flavin dianions reveals deprotomer-dependent photochemistry. Physical Chemistry Chemical Physics, 2018, 20, 19672-19681.	2.8	23
30	Electronic spectrum and photodissociation chemistry of the linear methyl propargyl cation H2C4H3+. Journal of Chemical Physics, 2017, 146, 044307.	3.0	4
31	A Strong cisâ€Effect in an Imidazoleâ€Imidazoliumâ€Substituted Alkene. Angewandte Chemie - International Edition, 2017, 56, 8473-8480.	13.8	3
32	A Strong cis â€Effect in an Imidazoleâ€Imidazoliumâ€Substituted Alkene. Angewandte Chemie, 2017, 129, 8593-8600.	2.0	1
33	Isomerisation of an intramolecular hydrogen-bonded photoswitch: protonated azobis(2-imidazole). Physical Chemistry Chemical Physics, 2017, 19, 12776-12783.	2.8	21
34	Online measurement of photoisomerisation efficiency in solution using ion mobility mass spectrometry. Analyst, The, 2017, 142, 2100-2103.	3.5	5
35	Seleniranium Ions Undergo Ï€-Ligand Exchange via an Associative Mechanism in the Gas Phase. Journal of Organic Chemistry, 2017, 82, 6289-6297.	3.2	10
36	Electronic spectrum of the protonated diacetylene cation (H2C4H+). Journal of Chemical Physics, 2017, 147, 084302.	3.0	7

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37	Photochrome-doped organic films for photonic keypad locks and multi-state fluorescence. Physical Chemistry Chemical Physics, 2017, 19, 19984-19991.	2.8	11
38	Protomer-Specific Photochemistry Investigated Using Ion Mobility Mass Spectrometry. Journal of Physical Chemistry A, 2017, 121, 6021-6027.	2.5	32
39	Photoisomerization of Protonated Azobenzenes in the Gas Phase. Journal of Physical Chemistry A, 2017, 121, 6413-6419.	2.5	23
40	Hydrogen-adduction to open-shell graphene fragments: spectroscopy, thermochemistry and astrochemistry. Chemical Science, 2017, 8, 1186-1194.	7.4	6
41	Monitoring Isomerization of Molecules in Solution Using Ion Mobility Mass Spectrometry. Analytical Chemistry, 2016, 88, 11978-11981.	6.5	24
42	Electronic spectrum of 9-methylanthracenium radical cation. Journal of Chemical Physics, 2016, 144, 154303.	3.0	0
43	Photoisomerization of Î²-Ionone Protonated Schiff Base in the Gas Phase. Journal of Physical Chemistry A, 2016, 120, 6557-6562.	2.5	4
44	Photo and Collision Induced Isomerization of a Cyclic Retinal Derivative: An Ion Mobility Study. Journal of the American Society for Mass Spectrometry, 2016, 27, 1483-1490.	2.8	12
45	Modulating electron injection from an organic dye to a titania nanoparticle with a photochromic energy transfer acceptor. Journal of Materials Chemistry C, 2016, 4, 6215-6219.	5.5	6
46	Ab Initio Characterization of the Electrostatic Complexes Formed by H₂ Molecule and Cr⁺, Mn⁺, Cu⁺, and Zn⁺ Cations. Journal of Physical Chemistry A, 2016, 120, 5006-5015.	2.5	7
47	Blue to near-IR energy transfer cascade within a dye-doped polymer matrix, mediated by a photochromic molecular switch. Physical Chemistry Chemical Physics, 2016, 18, 5095-5098.	2.8	14
48	Electronic spectrum of the propargyl cationâ€‰(H2C3H+)â€‰tagged with Ne and N2. Journal of Chemical Physics, 2015, 143, 184306.	3.0	8
49	Does the triphenylamine-based D35 dye sensitizer form aggregates on metal-oxide surfaces?. Journal of Photochemistry and Photobiology A: Chemistry, 2015, 302, 35-41.	3.9	13
50	Ultraviolet photodissociation action spectroscopy of the N-pyridinium cation. Journal of Chemical Physics, 2015, 142, 014301.	3.0	24
51	Electron Injection and Energy-Transfer Properties of Spiropyranâ€“Cyclodextrin Complexes Coated onto Metal Oxide Nanoparticles: Toward Photochromic Light Harvesting. Journal of Physical Chemistry C, 2015, 119, 14076-14084.	3.1	32
52	Photoisomerization action spectroscopy: flicking the protonated merocyanineâ€“spiropyran switch in the gas phase. Physical Chemistry Chemical Physics, 2015, 17, 25676-25688.	2.8	46
53	Retinal shows its true colours: photoisomerization action spectra of mobility-selected isomers of the retinal protonated Schiff base. Physical Chemistry Chemical Physics, 2015, 17, 22623-22631.	2.8	35
54	An ion mobility mass spectrometer for investigating photoisomerization and photodissociation of molecular ions. Review of Scientific Instruments, 2014, 85, 123109.	1.3	58

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55	Photoisomerization action spectrum of retinal protonated Schiff base in the gas phase. Journal of Chemical Physics, 2014, 140, 164307.	3.0	29
56	Ion Mobility Unlocks the Photofragmentation Mechanism of Retinal Protonated Schiff Base. Journal of Physical Chemistry Letters, 2014, 5, 3195-3199.	4.6	21
57	Suppressing FÃ¶rster Resonance Energy Transfer between Organic Dyes on a Cosensitized Metal Oxide Surface. Journal of Physical Chemistry C, 2014, 118, 19646-19654.	3.1	16
58	Interaction of the Beryllium Cation with Molecular Hydrogen and Deuterium. Journal of Physical Chemistry A, 2014, 118, 6711-6720.	2.5	8
59	Photophysics and aggregation effects of a triphenylamine-based dye sensitizer on metal-oxide nanoparticles suspended in an ion trap. Physical Chemistry Chemical Physics, 2013, 15, 20326.	2.8	13
60	Electronic Spectroscopy of the 1,3-Cyclopentadiene Cation (C5H6+). Journal of Physical Chemistry A, 2013, 117, 11276-11281.	2.5	5
61	Changing the shape of molecular ions: photoisomerization action spectroscopy in the gas phase. Physical Chemistry Chemical Physics, 2013, 15, 9540.	2.8	52
62	Photoisomerization Action Spectroscopy of the Carbocyanine Dye DTC⁺ in the Gas Phase. Journal of Physical Chemistry A, 2013, 117, 13319-13325.	2.5	20
63	Gas-phase electronic spectroscopy of the indene cation (C9H8+). Journal of Chemical Physics, 2013, 138, 224307.	3.0	10
64	Laboratory Spectroscopy of PAHs. Proceedings of the International Astronomical Union, 2013, 9, 247-257.	0.0	2
65	Electronic absorptions of the benzylium cation. Journal of Chemical Physics, 2012, 137, 204304.	3.0	24
66	Electronic Spectra of Gas-Phase Polycyclic Aromatic Nitrogen Heterocycle Cations: Isoquinoline⁺ and Quinoline⁺. Journal of Physical Chemistry A, 2012, 116, 4323-4329.	2.5	32
67	A sting in the tail of flexible molecules: spectroscopic and energetic challenges in the case of p-aminophenethylamine. Physical Chemistry Chemical Physics, 2012, 14, 9219.	2.8	7
68	Effect of multiplicative noise on least-squares parameter estimation with applications to the atomic force microscope. Review of Scientific Instruments, 2012, 83, 055106.	1.3	16
69	Properties of the B+-H2and B+-D2complexes: A theoretical and spectroscopic study. Journal of Chemical Physics, 2012, 137, 124312.	3.0	7
70	Spring constant calibration of atomic force microscope cantilevers of arbitrary shape. Review of Scientific Instruments, 2012, 83, 103705.	1.3	228
71	Attaching molecular hydrogen to metal cations: perspectives from gas-phase infrared spectroscopy. Physical Chemistry Chemical Physics, 2012, 14, 14954.	2.8	36
72	Molecular collision dynamics. Physical Chemistry Chemical Physics, 2011, 13, 8073.	2.8	2

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73	Potential energy surface and rovibrational calculations for the ${m Mg}^+$ Mg +â€“${m H}_2$H2 and ${m Mg}^+$ Mg +â€“${m D}_2$D2 complexes. Journal of Chemical Physics, 2011, 134, 044310.	3.0	14
74	Infrared Spectroscopy of the Ag⁺â^'H₂ Complex: Exploring the Connection Between Vibrational Band-Shifts and Binding Energies. Journal of Physical Chemistry Letters, 2011, 2, 719-724.	4.6	23
75	Rotationally resolved infrared spectrum of the Na+-D2 complex: An experimental and theoretical study. Journal of Chemical Physics, 2011, 134, 214302.	3.0	8
76	Distortion in the thermal noise spectrum and quality factor of nanomechanical devices due to finite frequency resolution with applications to the atomic force microscope. Review of Scientific Instruments, 2011, 82, 095104.	1.3	11
77	Infrared Spectra and ab initio Calculations for Fluoride-acetylene Clusters: F - -(HCCH)n, n=3 - 6. Australian Journal of Chemistry, 2011, 64, 633.	0.9	2
78	Mixing Laser Spectroscopy and Mass Spectrometry-Infrared Spectra of Metal Cationâ€“Hydrogen Complexes. European Journal of Mass Spectrometry, 2010, 16, 415-420.	1.0	2
79	Infrared spectra and density functional theory calculations for Mn+â€“(CH4)n (n=1â€“6) clusters. International Journal of Mass Spectrometry, 2010, 297, 46-54.	1.5	16
80	Infrared Spectra of Mass-Selected Brâ^'â^'(NH3)n and Iâ^'â^'NH3 Clusters. Journal of Physical Chemistry A, 2010, 114, 4762-4769.	2.5	8
81	Structure and properties of the Zn+â€“D2 complex. Journal of Chemical Physics, 2009, 131, 224304.	3.0	15
82	Photoacoustic detection of gases using microcantilevers. Journal of Applied Physics, 2009, 106, .	2.5	21
83	Infrared Spectra of Mass-Selected Mg⁺â^'H₂ and Mg⁺â^'D₂ Complexes. Journal of Physical Chemistry A, 2009, 113, 199-204.	2.5	18
84	Spectroscopic Study of the Benchmark Mn⁺â^'H₂ Complex. Journal of Physical Chemistry A, 2009, 113, 6044-6048.	2.5	24
85	The Cr+â€“D2 cation complex: Accurate experimental dissociation energy, intermolecular bond length, and vibrational parameters. Journal of Chemical Physics, 2009, 131, 164303.	3.0	16
86	Attachment of Molecular Hydrogen to an Isolated Boron Cation: An Infrared and ab initio Study. Journal of the American Chemical Society, 2008, 130, 12986-12991.	13.7	21
87	The Na+â€“H2 cation complex: Rotationally resolved infrared spectrum, potential energy surface, and rovibrational calculations. Journal of Chemical Physics, 2008, 129, 184306.	3.0	28
88	Morphology-Dependent Resonance Emission from Individual Micron-Sized Particles. Springer Series on Fluorescence, 2007, , 415-429.	0.8	2
89	The Al+â€“H2 cation complex: Rotationally resolved infrared spectrum, potential energy surface, and rovibrational calculations. Journal of Chemical Physics, 2007, 127, 164310.	3.0	31
90	Infrared spectra of the Li+â€“(H2)n (n=1â€“3) cation complexes. Journal of Chemical Physics, 2007, 126, 204309.	3.0	39

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91	Interactions between the Chloride Anion and Aromatic Molecules:Â Infrared Spectra of the Cl-â˜C6H5CH3, Cl-â˜C6H5NH2and Cl-â˜C6H5OH Complexes. Journal of Physical Chemistry A, 2007, 111, 7322-7328.	2.5	9
92	Calibration of a quadrupole ion trap for particle mass spectrometry. International Journal of Mass Spectrometry, 2007, 262, 241-246.	1.5	28
93	Infrared Spectra and Ab Initio Calculations for the F-â˜(CH4)n(n= 1â˜8) Anion Clusters. Journal of Physical Chemistry A, 2006, 110, 13736-13743.	2.5	25
94	Observation of nondegenerate cavity modes for a distorted polystyrene microsphere. Optics Letters, 2006, 31, 2211.	3.3	12
95	Infrared spectra of Clâ˜â€“(C6H6)m= 1, 2. Chemical Physics Letters, 2006, 428, 18-22.	2.6	13
96	Apparatus for the study of electronic spectra of collisionally cooled cations: para-dichlorobenzene. Journal of Molecular Structure, 2006, 795, 93-97.	3.6	43
97	Ab initiopotential energy surface, infrared spectra, and dynamics of the ion-molecule complexes between Brâ˜ and H2, D2, and HD. Journal of Chemical Physics, 2006, 125, 114313.	3.0	11
98	Rotationally resolved infrared spectrum of the Li+â€“D2 cation complex. Journal of Chemical Physics, 2006, 125, 044310.	3.0	32
99	Infrared spectra of the Clâ€“â€“C2H4 and Brâ€“â€“C2H4 anion dimers. Physical Chemistry Chemical Physics, 2005, 7, 3419.	2.8	7
100	Infrared Spectra and ab Initio Calculations for the Cl-â˜(CH4)n(n= 1â˜10) Anion Clusters. Journal of Physical Chemistry A, 2005, 109, 8481-8486.	2.5	16
101	Isomeric interconversion in the linear Clâ˜-HD anion complex. Journal of Chemical Physics, 2004, 121, 2085-2093.	3.0	15
102	Infrared spectrum of the I[sup â˜]â€“D[sub 2] anion complex. Journal of Chemical Physics, 2004, 121, 12276.	3.0	14
103	Structures of F - -(CH4)n and Cl - -(CH4)n (n = 1,2) Anion Clusters Elucidated through Ab Initio Calculations and Infrared Spectra. Australian Journal of Chemistry, 2004, 57, 1157.	0.9	18
104	Spectroscopic studies of anion complexes and clusters: A microscopic approach to understanding anion solvation. Chemical Society Reviews, 2003, 32, 231.	38.1	44
105	Ab initio potential energy and dipole moment surfaces, infrared spectra, and vibrational predissociation dynamics of the 35Clâ˜â‹¯H2/D2 complexes. Journal of Chemical Physics, 2003, 119, 12931-12945.	3.0	46
106	Clâ˜â€“C6H6, Brâ˜â€“C6H6, and Iâ˜â€“C6H6 anion complexes: Infrared spectra and ab initio calculations. Journal of Chemical Physics, 2003, 119, 9559-9567.	3.0	49
107	Infrared Investigations of Negatively Charged Complexes and Clusters. International Reviews in Physical Chemistry, 2003, 22, 129-151.	2.3	69
108	Brâ˜-H2 and Iâ˜-H2 anion complexes: Infrared spectra and radial intermolecular potential energy curves. Journal of Chemical Physics, 2002, 117, 3256-3262.	3.0	35

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109	Infrared Spectra of Size Selected Cl-â˜(D2)nand F-â˜(D2)nAnion Clusters. Journal of Physical Chemistry A, 2002, 106, 906-910.	2.5	13
110	The 35Clâ˜â€“H2 and 35Clâ˜â€“D2 anion complexes: Infrared spectra and radial intermolecular potentials. Journal of Chemical Physics, 2001, 115, 824-832.	3.0	54
111	Rotationally resolved infrared spectrum of the Brâ˜â˜D2 anion complex. Journal of Chemical Physics, 2001, 115, 6394-6400.	3.0	36
112	Structural and energetic properties of the Brâ˜â€“C2H2 anion complex from rotationally resolved mid-infrared spectra and ab initio calculations. Journal of Chemical Physics, 2000, 113, 1075-1080.	3.0	21
113	Rotationally resolved infrared spectrum of the Clâ˜â€“H2 anion complex. Journal of Chemical Physics, 2000, 113, 10154-10157.	3.0	36
114	High-Resolution Spectroscopy of Cluster Ions. Chemical Reviews, 2000, 100, 3963-3998.	47.7	468
115	Infrared and ab Initio Study of the Chlorideâ˜Ammonia Anion Complex. Journal of Physical Chemistry A, 2000, 104, 2562-2566.	2.5	23
116	Infrared spectra of Clâ˜â€“(C2H2)nâ€Š(1â©½nâ©½9) anion clusters: Spectroscopic evidence for solvent shell closure. Journal of Chemical Physics, 1999, 110, 9443-9449.	3.0	39
117	Photofragmentation dynamics of the (N2O)2+ and (N2O)3+ clusters: fragment N2O+ A â†•X spectra. Chemical Physics, 1998, 239, 369-378.	1.9	2
118	Infrared predissociation spectra of Nenâ€“HN2+ clusters (n=1â€“5). Journal of Chemical Physics, 1998, 108, 8964-8975.	3.0	47
119	Microsolvation of the ammonium ion in argon: infrared spectra of NH4+î—¸Arn complexes (n = 1â€“7). International Journal of Mass Spectrometry and Ion Processes, 1997, 167-168, 637-647.	1.8	54
120	Dissociation energy of the Arî—¸HN+2 complex. Chemical Physics Letters, 1997, 265, 303-307.	2.6	54
121	Observation of the infrared spectrum of the Î½3 band of the argon-ammonium ionic complex. Chemical Physics Letters, 1996, 250, 266-272.	2.6	28
122	Midâ€infrared spectra of Heâ€“HN+2and He2â€“HN+2. Journal of Chemical Physics, 1996, 104, 3876-3885.	3.0	65
123	Midâ€infrared spectra of the protonâ€bound complexes Nenâ€“HCO+ (n=1,2). Journal of Chemical Physics, 1996, 105, 1770-1777.	3.0	60
124	The infrared spectrum of the H2â€“HCO+ complex. Journal of Chemical Physics, 1995, 102, 5152-5164.	3.0	71
125	Size Effects in Cluster Infrared Spectra: the .nu.1 Band of Arn-HCO+ (n = 1-13). The Journal of Physical Chemistry, 1995, 99, 17118-17129.	2.9	114
126	Electronic spectroscopy of size-selected ionic complexes. Journal of the Chemical Society, Faraday Transactions, 1995, 91, 1.	1.7	56

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127	The infrared spectrum of Heâ€“HCO+. Journal of Chemical Physics, 1995, 103, 1297-1302.	3.0	53
128	A 3.Pl.u .rarw. X 3.SIGMA.g- Electronic Spectrum of N3+. The Journal of Physical Chemistry, 1994, 98, 8896-8902.	2.9	34
129	The van der Waals vibrations of anilineâ€“(argon)2 in the S1 electronic state. Journal of Chemical Physics, 1991, 94, 7019-7028.	3.0	47
130	Mass selected resonance enhanced multiphoton ionization spectroscopy of anilineâ€“Arn (n=3,4,5,â€‰...) van der Waals complexes. Journal of Chemical Physics, 1991, 94, 7029-7037.	3.0	35
131	The Bâ†X electronic spectrum of N+2â€“Ne. Journal of Chemical Physics, 1991, 94, 4749-4755.	3.0	35
132	Photodissociation spectroscopy of aromaticâ€“rare gas cluster ions: Low frequency vibrations in pâ€‹difluorobenzene+â‹…argon. Journal of Chemical Physics, 1990, 92, 4620-4621.	3.0	19
133	The Bâ†X electronic spectrum of N2+â€“He. Journal of Chemical Physics, 1990, 93, 4477-4478.	3.0	36
134	Photo-induced 6Ï€-electrocyclisation and cycloreversion of isolated dithienylethene anions. Physical Chemistry Chemical Physics, 0, , .	2.8	0