

Scott A Reid

List of Publications by Year in descending order

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115
papers

1,852
citations

304368

22
h-index

377514

34
g-index

117
all docs

117
docs citations

117
times ranked

1328
citing authors

#	ARTICLE	IF	CITATIONS
1	Implementation and evaluation of an adaptive online summer preparatory course for general chemistry: Whom does it benefit?. <i>Chemistry Education Research and Practice</i> , 2021, 22, 303-311.	1.4	5
2	Unraveling a trifecta of weak non-covalent interactions: The dissociation energy of the anisole-ammonia 1:1 complex. <i>Chemical Physics Letters</i> , 2021, 762, 138106.	1.2	1
3	Zeroing in on the best early-course metrics to identify at-risk students in general chemistry: an adaptive learning pre-assessment vs. traditional diagnostic exam. <i>International Journal of Science Education</i> , 2021, 43, 552-569.	1.0	10
4	Probing cooperativity in C-H \cdots N and C-H \cdots I interactions: Dissociation energies of aniline \cdots (CH ₄) _n (n = 1, 2) van der Waals complexes from resonant ionization and velocity mapped ion imaging measurements. <i>Journal of Chemical Physics</i> , 2020, 153, 044303.	1.2	6
5	Restructuring a General College Chemistry Sequence Using the ACS Anchoring Concepts Content Map. <i>Journal of Chemical Education</i> , 2020, 97, 651-658.	1.1	8
6	Tribute to Hanna Reisler. <i>Journal of Physical Chemistry A</i> , 2019, 123, 6381-6383.	1.1	0
7	Charge-transfer or excimeric state? Exploring the nature of the excited state in cofacially arrayed polyfluorene derivatives. <i>Journal of Photochemistry and Photobiology A: Chemistry</i> , 2019, 374, 125-130.	2.0	2
8	C-H \cdots I and C-H \cdots O Interactions in Concert: A Study of the Anisole \cdots Methane Complex using Resonant Ionization and Velocity Mapped Ion Imaging. <i>Journal of Physical Chemistry A</i> , 2019, 123, 2874-2880.	1.1	10
9	Strength of I \cdots I Stacking, from Neutral to Cation: Precision Measurement of Binding Energies in an Isolated I \cdots I Stacked Dimer. <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 2058-2061.	2.1	15
10	Reactive pathways in the bromobenzene-ammonia dimer cation radical: Evidence for a roaming halogen radical. <i>Journal of Molecular Structure</i> , 2018, 1172, 113-118.	1.8	2
11	Vertical vs. adiabatic ionization energies in solution and gas-phase: probing ionization-induced reorganization in conformationally-mobile bichromophoric actuators using photoelectron spectroscopy, electrochemistry and theory. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 25615-25622.	1.3	9
12	I \cdots I stacking vs. C-H \cdots I interaction: Excimer formation and charge resonance stabilization in van der Waals clusters of 9,9-dimethylfluorene. <i>Journal of Chemical Physics</i> , 2018, 149, 134314.	1.2	10
13	An Electron-Rich Calix[4]arene-Based Receptor with Unprecedented Binding Affinity for Nitric Oxide. <i>Chemistry - A European Journal</i> , 2018, 24, 17439-17443.	1.7	6
14	Game of Frontier Orbitals: A View on the Rational Design of Novel Charge-Transfer Materials. <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 3978-3986.	2.1	25
15	Molecular Actuators in Action: Electron-Transfer-Induced Conformation Transformation in Cofacially Arrayed Polyfluorenes. <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 4233-4238.	2.1	7
16	Spreading Electron Density Thin: Increasing the Chromophore Size in Polyaromatic Wires Decreases Interchromophoric Electronic Coupling. <i>Journal of Physical Chemistry C</i> , 2018, 122, 17668-17675.	1.5	7
17	The Role of Torsional Dynamics on Hole and Exciton Stabilization in I \cdots I Stacked Assemblies: Design of Rigid Torsionomers of a Cofacial Bifluorene. <i>Angewandte Chemie - International Edition</i> , 2018, 57, 8189-8193.	7.2	16
18	The Role of Torsional Dynamics on Hole and Exciton Stabilization in I \cdots I Stacked Assemblies: Design of Rigid Torsionomers of a Cofacial Bifluorene. <i>Angewandte Chemie</i> , 2018, 130, 8321-8325.	1.6	4

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19	Cofacially Arrayed Polyfluorenes: Spontaneous Formation of π -Stacked Assemblies in the Gas Phase. <i>Journal of Physical Chemistry Letters</i> , 2017, 8, 5272-5276.	2.1	9
20	Effect of Facial Encumbrance on Excimer Formation and Charge Resonance Stabilization in Model Bichromophoric Assemblies. <i>Journal of Physical Chemistry C</i> , 2017, 121, 15580-15588.	1.5	10
21	A flipped classroom redesign in general chemistry. <i>Chemistry Education Research and Practice</i> , 2016, 17, 914-922.	1.4	31
22	First Experimental Evidence for the Diverse Requirements of Excimer vs Hole Stabilization in π -Stacked Assemblies. <i>Journal of Physical Chemistry Letters</i> , 2016, 7, 3042-3045.	2.1	14
23	Two ⁺ 's Company, Three ⁺ 's a Crowd: Exciton Localization in Cofacially Arrayed Polyfluorenes. <i>Journal of Physical Chemistry Letters</i> , 2016, 7, 2915-2920.	2.1	12
24	Impact of the Flipped Classroom on Student Performance and Retention: A Parallel Controlled Study in General Chemistry. <i>Journal of Chemical Education</i> , 2016, 93, 13-23.	1.1	188
25	From Wires to Cables: Attempted Synthesis of 1,3,5-Trifluorenylcyclohexane as a Platform for Molecular Cables. <i>Journal of Organic Chemistry</i> , 2016, 81, 1627-1634.	1.7	0
26	Hydrogen-atom attack on phenol and toluene is ortho-directed. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 8625-8636.	1.3	9
27	Revisiting the Renner-Ellipsoidal effect in the π -stacked state of CCN: Pulsed discharge-supersonic jet single vibronic level emission spectroscopy. <i>Journal of Chemical Physics</i> , 2015, 142, 214304.	0.4	3
28	Towards a global model of spin-orbit coupling in the halocarbenes. <i>Journal of Chemical Physics</i> , 2015, 142, 214304.	1.2	12
29	When isomerisation is electron transfer: the intriguing story of the iso-halocarbons. <i>International Reviews in Physical Chemistry</i> , 2014, 33, 341-370.	0.9	12
30	Photoinduced Electron Transfer in Donor-Acceptor Complexes of Ethylene with Molecular and Atomic Iodine. <i>Journal of Physical Chemistry A</i> , 2014, 118, 6838-6845.	1.1	8
31	Two roaming pathways in the photolysis of CH_3CHO between 328 and 308 nm. <i>Chemical Science</i> , 2014, 5, 4633-4638.	3.7	49
32	Case of the Missing Isomer: Pathways for Molecular Elimination in the Photoinduced Decomposition of 1,1-Dibromoethane. <i>Journal of Physical Chemistry A</i> , 2013, 117, 11915-11923.	1.1	10
33	π -Stacking, C-H \cdots H \cdots , and Halogen Bonding Interactions in Bromobenzene and Mixed Bromobenzene-Benzene Clusters. <i>Journal of Physical Chemistry A</i> , 2013, 117, 13556-13563.	1.1	41
34	Photoisomerization and Photoinduced Reactions in Liquid CCl_4 and CHCl_3 . <i>Journal of Physical Chemistry A</i> , 2013, 117, 13388-13398.	1.1	18
35	Concerted and sequential pathways of proton-coupled electron transfer in hydrogen halide elimination. <i>Chemical Physics Letters</i> , 2013, 556, 35-38.	1.2	8
36	Reactive Pathways in the Chlorobenzene-Ammonia Dimer Cation Radical: New Insights from Experiment and Theory. <i>Journal of Physical Chemistry A</i> , 2013, 117, 12429-12437.	1.1	7

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37	Spectroscopy and dynamics of the predissociated, quasi-linear S ₂ state of chlorocarbene. <i>Journal of Chemical Physics</i> , 2012, 137, 104307.	1.2	11
38	On the electronic spectroscopy of the iso-polyhalomethanes. <i>Chemical Physics Letters</i> , 2012, 551, 64-67.	1.2	1
39	Product state and speed distributions in photochemical triple fragmentations. <i>Faraday Discussions</i> , 2012, 157, 227.	1.6	27
40	Photochemistry of Furyl- and Thienyldiazomethanes: Spectroscopic Characterization of Triplet 3-Thienylcarbene. <i>Journal of the American Chemical Society</i> , 2012, 134, 6443-6454.	6.6	14
41	Excitation Spectra of the Jet-Cooled 4-Phenylbenzyl and 4-(4- ² -Methylphenyl)benzyl Radicals. <i>Journal of Physical Chemistry A</i> , 2012, 116, 10780-10785.	1.1	7
42	Probing radical pathways in electrophilic addition of halogens: Classical vs. bridged intermediates. <i>Chemical Physics Letters</i> , 2012, 554, 86-89.	1.2	7
43	On π -stacking, C-H/ π , and halogen bonding interactions in halobenzene clusters: Resonant two-photon ionization studies of chlorobenzene. <i>Journal of Chemical Physics</i> , 2012, 137, 184307.	1.2	33
44	Spectroscopic and computational studies of matrix-isolated iso-CXBr ₃ (X=F, Cl, Br): Structure, properties, and photochemistry of substituted iso-tribromomethanes. <i>Journal of Molecular Structure</i> , 2012, 1025, 61-68.	1.8	6
45	On the electronic spectroscopy of closed-shell cations derived from resonance-stabilized radicals: Insights from theory and Franck-Condon analysis. <i>Astronomy and Astrophysics</i> , 2012, 541, A8.	2.1	8
46	Pulsed Jet Discharge Matrix Isolation and Computational Study of Bromine Atom Complexes: Br \cdot \cdot BrXCH ₂ (X = H, Cl, Br). <i>Journal of Physical Chemistry A</i> , 2011, 115, 9820-9827.	1.1	7
47	Probing the electronic structure of the nickel monohalides: Spectroscopy of the low-lying electronic states of NiBr and NiCl. <i>Journal of Molecular Spectroscopy</i> , 2011, 269, 36-40.	0.4	8
48	Electronic spectroscopy of the transition of DCO and lifetimes and relative quantum yields of the state. <i>Journal of Molecular Spectroscopy</i> , 2011, 270, 33-39.	0.4	1
49	Spectroscopic and computational studies of matrix-isolated iso-CHBr ₃ : Structure, properties, and photochemistry of iso-bromoform. <i>Journal of Chemical Physics</i> , 2011, 135, 124503.	1.2	24
50	Chemistry at the threshold: Unexpected products, unusual mechanisms, and generally weird things that happen near the energetic threshold for a reaction. , 2011, , .		0
51	Formation and relaxation dynamics of iso-CH ₂ Cl ⁺ I in cryogenic matrices. <i>Journal of Chemical Physics</i> , 2011, 135, 114503.	1.2	19
52	Optical-optical double resonance spectroscopy of the quasi-linear S ₂ state of CHF and CDF. I. Spectroscopic analysis. <i>Journal of Chemical Physics</i> , 2011, 135, 104315.	1.2	9
53	Optical-optical double resonance spectroscopy of the quasi-linear S ₂ state of CHF and CDF. II. Predissociation and mode-specific dynamics. <i>Journal of Chemical Physics</i> , 2011, 135, 104316.	1.2	10
54	Spectroscopy and thermochemistry of a jet-cooled open-shell polyene: 1,4-pentadienyl radical. <i>Journal of Chemical Physics</i> , 2011, 135, 124306.	1.2	12

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55	Single vibronic level emission spectroscopy of the low-lying electronic states of NiI. <i>Chemical Physics Letters</i> , 2010, 497, 168-171.	1.2	6
56	Pulsed-jet discharge matrix isolation and computational study of CX ₂ Br ⁺ (X=H, F). <i>Chemical Physics Letters</i> , 2010, 484, 214-218.	1.2	3
57	Matrix isolation and computational studies of the CF ₂ I radical. <i>Chemical Physics Letters</i> , 2010, 496, 68-73.	1.2	7
58	Matrix isolation and computational study of isodifluorodibromomethane (F ₂ CFBr ⁺ Br): A route to Br ₂ formation in CF ₂ Br ₂ photolysis. <i>Journal of Chemical Physics</i> , 2010, 132, 084503.	1.2	19
59	Isomerization as a Key Path to Molecular Products in the Gas-Phase Decomposition of Halons. <i>Journal of Physical Chemistry Letters</i> , 2010, 1, 3090-3095.	2.1	23
60	Spectroscopic and Computational Studies of the Laser Photolysis of Matrix Isolated 1,2-Dibromoethanes: Formation and Fate of the Bromoethyl Radicals. <i>Journal of Physical Chemistry A</i> , 2010, 114, 9919-9926.	1.1	17
61	Photoinduced Electron Transfer in a Prototypical Mulliken Donor-Acceptor Complex: C ₂ H ₄ -Br ₂ . <i>Journal of Physical Chemistry Letters</i> , 2010, 1, 2618-2621.	2.1	9
62	Direct Observation of Electron-Transfer-Induced Conformational Transformation (Molecular) Tj ETQq0 0 0 rgBT /Overlock 10 Tf 50 467 T 14592-14595.	1.2	12
63	Theoretical and Experimental Spectroscopy of the S ₂ State of CHF and CDF: Dynamically Weighted Multireference Configuration Interaction Calculations for High-Lying Electronic States. <i>Journal of Physical Chemistry Letters</i> , 2010, 1, 641-646.	2.1	57
64	Characterization of iso-CF ₂ I ₂ in frequency and ultrafast time domains. <i>Journal of Chemical Physics</i> , 2010, 132, 124501.	1.2	29
65	Comment on "Facile strategy and mechanism for orthorhombic SnO ₂ thin films" [Appl. Phys. Lett. 89, 231902 (2006)]. <i>Applied Physics Letters</i> , 2009, 94, 186103.	1.5	1
66	The halocarbenes: model systems for understanding the spectroscopy, dynamics and chemistry of carbenes. <i>International Reviews in Physical Chemistry</i> , 2009, 28, 435-480.	0.9	45
67	Fluorescence Excitation and Emission Spectroscopy of the $\tilde{X}^1\Sigma^+$ System of CHI and CDI. <i>Journal of Physical Chemistry A</i> , 2009, 113, 13407-13412.	1.1	6
68	First observation of the elusive iodocarbene: ground state multiplicity and singlet-triplet gap of CHI. <i>Physical Chemistry Chemical Physics</i> , 2008, 10, 6090.	1.3	15
69	Stimulated Emission Pumping Spectroscopy of the $[X^1\Sigma^+]$ State of CHF. <i>Journal of Physical Chemistry A</i> , 2008, 112, 466-471.	1.1	16
70	Unraveling the $\tilde{A}^1\Sigma^+ B^1\Sigma^+ \tilde{X}^1\Sigma^+ A^1\Sigma^+$ Spectrum of CCl ₂ : The Renner-Teller Effect, Barrier to Linearity, and Vibrational Analysis Using an Effective Polyad Hamiltonian. <i>Journal of Physical Chemistry A</i> , 2008, 112, 11355-11362.	1.1	13
71	High resolution probe of spin-orbit coupling and the singlet-triplet gap in chlorocarbene. <i>Journal of Chemical Physics</i> , 2008, 128, 171101.	1.2	23
72	High resolution study of spin-orbit mixing and the singlet-triplet gap in chlorocarbene: Stimulated emission pumping spectroscopy of CH ₃ Cl and CD ₃ Cl. <i>Journal of Chemical Physics</i> , 2008, 129, 104309.	1.2	17

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73	Observation of the predissociated, quasilinear $B\tilde{1}f(A\hat{\epsilon}^21)$ state of CHF by optical-optical double resonance. Journal of Chemical Physics, 2007, 126, 051105.	1.2	19
74	Electronic Spectroscopy of an Isolated Halocarbo-cation: The Iodomethyl Cation CH_2I^+ and Its Deuterated Isotopomers. Journal of Physical Chemistry A, 2007, 111, 10562-10566.	1.1	9
75	Electronic emission spectroscopy and fluorescence lifetime of the $B\tilde{1}f(A\hat{\epsilon}^21)$ state of CHF. Chemical Physics Letters, 2007, 449, 282-285.	1.2	11
76	Single vibronic level emission spectroscopy of the system of dibromocarbene. Journal of Molecular Spectroscopy, 2007, 241, 136-142.	0.4	14
77	Electronic spectroscopy of the system of CDCl. Journal of Molecular Spectroscopy, 2007, 241, 143-150.	0.4	9
78	Electronic spectroscopy, lifetimes, and barrier to linearity in the system of dibromocarbene. Journal of Molecular Spectroscopy, 2007, 241, 180-185.	0.4	10
79	Single vibronic level emission spectroscopy of the system of bromochlorocarbene. Journal of Molecular Spectroscopy, 2007, 246, 113-117.	0.4	5
80	Probing spin-orbit mixing and the singlet-triplet gap in dichloromethylene via $K\alpha$ -sorted emission spectra. Physical Chemistry Chemical Physics, 2006, 8, 4320-4326.	1.3	11
81	Electronic spectroscopy of the $\tilde{A}1A\hat{\epsilon}^31\tilde{1}f(A\hat{\epsilon}^21)$ system of CDF. Physical Chemistry Chemical Physics, 2006, 8, 707.	1.3	17
82	Laser Spectroscopy of a Halocarbo-cation in the Gas Phase: CH_2I^+ . Journal of the American Chemical Society, 2006, 128, 9320-9321.	6.6	9
83	Reassignment of the electronic origin in the system of dibromocarbene. Journal of Molecular Spectroscopy, 2006, 240, 139-140.	0.4	6
84	Electronic spectroscopy of the $\tilde{A}1A\hat{\epsilon}^31\tilde{1}f(A\hat{\epsilon}^21)$ system of CDBr. Journal of Chemical Physics, 2006, 125, 094305.	1.2	6
85	Fluorescence excitation and emission spectroscopy of the $\tilde{A}1A\hat{\epsilon}^31\tilde{1}f(A\hat{\epsilon}^21)$ system of CHBr. Journal of Chemical Physics, 2006, 124, 134302.	1.2	34
86	Fluorescence excitation and single vibronic level emission spectroscopy of the $\tilde{A}1A\hat{\epsilon}^31\tilde{1}f(A\hat{\epsilon}^21)$ system of $CHCl_2$. Journal of Chemical Physics, 2006, 124, 224314.	1.2	32
87	A DFT study of the hyperfine coupling constants of triplet carbenes and biradicals. Computational and Theoretical Chemistry, 2005, 725, 45-53.	1.5	16
88	Dispersed fluorescence spectroscopy of jet-cooled HCF and DCF: Vibrational structure of the $\tilde{X}1A\hat{\epsilon}^21$ state. Journal of Chemical Physics, 2005, 123, 014314.	1.2	21
89	Two-color resonant four-wave mixing spectroscopy of highly predissociated levels in the $\tilde{A}1A12$ state of CH_3S . Journal of Chemical Physics, 2005, 122, 124313.	1.2	12
90	Conformational Changes in the Ligand-binding Domain of a Functional Ionotropic Glutamate Receptor. Journal of Biological Chemistry, 2005, 280, 8633-8636.	1.6	33

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91	Vibrational mode selectivity in hyperfine interactions: Polarization quantum beat spectroscopy of HCF(\tilde{A}^3). Journal of Chemical Physics, 2004, 120, 1164-1167.	1.2	17
92	Polarization quantum beat spectroscopy of HCF(\tilde{A}^3). II. Renner-Teller and spin-orbit mixing in the simplest singlet carbene. Journal of Chemical Physics, 2004, 121, 8874-8879.	1.2	18
93	Fluorescence excitation spectroscopy of the system of jet-cooled HCl in the region 5150-6050 Å. Journal of Molecular Spectroscopy, 2004, 225, 43-47.	0.4	17
94	Polarization quantum beat spectroscopy of HCF(\tilde{A}^3). I. 19F and 1H hyperfine structure and Zeeman effect. Journal of Chemical Physics, 2004, 121, 8869-8873.	1.2	19
95	On the Renner-Teller Effect and Barriers to Linearity and Dissociation in HCF(\tilde{A}^3). Journal of Physical Chemistry A, 2004, 108, 3732-3738.	1.1	30
96	Lifetime lengthening and the Renner-Teller effect in the HCF (\tilde{A}^3 and \tilde{X}^1) system. Chemical Physics Letters, 2003, 378, 548-552.	1.2	32
97	Fluorescence excitation spectroscopy of the system of jet-cooled NH ₂ in the region 2900-4300 Å. Journal of Molecular Spectroscopy, 2003, 219, 37-44.	0.4	16
98	A time-of-flight mass spectrometric study of laser fluence dependencies in SnO ₂ ablation: implications for pulsed laser deposited tin oxide thin films. International Journal of Mass Spectrometry, 2003, 230, 11-17.	0.7	3
99	On the energy dependence of the Zeeman and hyperfine parameters in the state of OH and OD. Chemical Physics, 2003, 291, 61-72.	0.9	18
100	Fluorescence spectra of NH ₂ \tilde{X}^1 and \tilde{B}^1 \tilde{A}^1 bands: Experiment and theory. Journal of Chemical Physics, 2003, 119, 2614-2617.	1.2	4
101	Phase Transformations in Pulsed Laser Deposited Nanocrystalline Tin Oxide Thin Films. Chemistry of Materials, 2003, 15, 564-567.	3.2	35
102	Zeeman quantum-beat spectroscopy of NO ₂ : Eigenstate-resolved Landé gF factors near dissociation threshold. Journal of Chemical Physics, 2002, 116, 525-531.	1.2	11
103	6P7/2-Excited-State Decay Mechanism and Energy-Transfer Processes in KMgF ₃ :Eu ²⁺ and KMgF ₃ :Eu ²⁺ X (X = Tj, ET, Q, 1, 1, 0.784314, 8). Journal of Chemical Physics, 2001, 115, 8868-8875.	3.2	8
104	Study of energy transfer in KMgF ₃ :Eu ²⁺ X (X=Gd, Cr, Ce) by the decay model of 6P7/2 excited state of Eu ²⁺ . Chemical Physics Letters, 2001, 335, 17-22.	1.2	10
105	On the energy dependence of the hyperfine interaction in excited states of NO ₂ . Journal of Chemical Physics, 2001, 115, 8868-8875.	1.2	7
106	On the hyperfine structure of NO ₂ levels near dissociation threshold. Journal of Chemical Physics, 2000, 112, 10067-10069.	1.2	12
107	Pulsed Laser Ablation of Sn and SnO ₂ Targets: A Neutral Composition, Energetics, and Wavelength Dependence. Journal of Physical Chemistry B, 2000, 104, 5324-5330.	1.2	13
108	EXPERIMENTAL STUDIES OF RESONANCES IN UNIMOLECULAR DECOMPOSITION. Annual Review of Physical Chemistry, 1996, 47, 495-525.	4.8	59

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109	Unimolecular Reaction of NO ₂ : Overlapping Resonances, Fluctuations, and the Transition State. The Journal of Physical Chemistry, 1996, 100, 474-487.	2.9	63
110	Resonances and fluctuations in the unimolecular reaction of NO ₂ . Faraday Discussions, 1995, 102, 129.	1.6	14
111	State-specific photofragment yield spectroscopy of jet-cooled methyl nitrite. Chemical Physics Letters, 1993, 209, 22-28.	1.2	15
112	Interaction between overall and internal rotation below, near, and above the summit of a torsional barrier: 1n1€* methylglyoxal. Journal of Chemical Physics, 1992, 97, 2338-2346.	1.2	7
113	A scanning, single mode, LiNbO ₃ , optical parametric oscillator. Optics Communications, 1989, 69, 289-293.	1.0	26
114	Intramolecular vibrational state mixing in glyoxal by stimulated emission pumping spectroscopy. Chemical Physics Letters, 1987, 139, 525-527.	1.2	15
115	ESR spectra of a seven-coordinated pentagonal bipyramidal manganese(II) complex. Inorganica Chimica Acta, 1985, 105, L13-L14.	1.2	3