Terry A Miller

List of Publications by Year in descending order

Source: https://exaly.com/author-pdf/7531095/publications.pdf

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365 papers 10,120 citations

41258 49 h-index 71 g-index

402 all docs 402 docs citations

times ranked

402

3306 citing authors

#	Article	IF	Citations
1	A combined experimental and computational study on the transition of the calcium isopropoxide radical as a candidate for direct laser cooling. Physical Chemistry Chemical Physics, 2022, 24, 8749-8762.	1.3	2
2	Time-resolved measurements of HO2 radical in a heated plasma flow reactor. Combustion and Flame, 2022, 241, 112097.	2.8	6
3	Laser-Induced Fluorescence Spectroscopy of Large Secondary Alkoxy Radicals: Part I. Spectral Overviews and Vibronic Analysis. Journal of Physical Chemistry A, 2021, 125, 1391-1401.	1.1	4
4	Vibronically coupled states: computational considerations and characterisation of vibronic and rovibronic spectroscopic parameters. International Reviews in Physical Chemistry, 2021, 40, 165-298.	0.9	13
5	Electronic spectroscopy of the A1 $if2Aae^2ae^2 A2if2Aae^2a^2A^2if2Aae^2$ transitions of jet-cooled calcium ethoxide ra Vibronic structure of alkaline earth monoalkoxide radicals of $i>C S symmetry. Journal of Chemical Physics, 2021, 155, 024301.$	dicals: 1.2	4
6	Laser-Induced Fluorescence Spectroscopy of Large Secondary Alkoxy Radicals: Part II. Rotational and Fine Structure. Journal of Physical Chemistry A, 2021, 125, 1402-1412.	1.1	3
7	Rotational and fine structure of open-shell molecules in nearly degenerate electronic states. II. Interpretation of experimentally determined interstate coupling parameters of alkoxy radicals. Journal of Chemical Physics, 2020, 153, 174306.	1.2	4
8	Laser-induced fluorescence and dispersed-fluorescence spectroscopy of the \tilde{A}_f 2E \hat{a} 2X \hat{I}_f 2A1 transition of jet-cooled calcium methoxide (CaOCH3) radicals. Journal of Chemical Physics, 2019, 151, 134303.	1.2	18
9	Jon T. Hougen. Journal of Molecular Spectroscopy, 2019, 360, 13-14.	0.4	0
10	First-Principles Calculation of Jahn–Teller Rotational Distortion Parameters. Journal of Physical Chemistry A, 2019, 123, 4990-5004.	1.1	7
11	Quantifying the effects of higher order coupling terms on fits using a second order Jahn-Teller Hamiltonian. Journal of Molecular Spectroscopy, 2018, 343, 102-115.	0.4	9
12	Studies via Near-Infrared Cavity Ringdown Spectroscopy and Electronic Structure Calculations of the Products of the Photolysis of Dihalomethane/N2/O2 Mixtures. Journal of Physical Chemistry A, 2017, 121, 98-112.	1.1	1
13	Modeling the CH Stretch/Torsion/Rotation Couplings in Methyl Peroxy (CH ₃ OO). Journal of Physical Chemistry A, 2017, 121, 9619-9630.	1.1	6
14	Sub-Doppler infrared spectroscopy of resonance-stabilized hydrocarbon intermediates: $\langle i \rangle \hat{1}^{1}/2 \langle i \rangle \langle sub \rangle \langle i \rangle \hat{1}^{1}/2 \langle i \rangle \langle sub \rangle \langle i \rangle \hat{1}^{1}/2 \langle i \rangle \langle sub \rangle \langle sub \rangle \langle i \rangle \langle sub \rangle \langle$	1.3	7
15	Manifestations of Torsion-CH Stretch Coupling in the Infrared Spectrum of CH ₃ OO. Journal of Physical Chemistry A, 2016, 120, 4827-4837.	1.1	9
16	Laser-Induced Fluorescence Spectroscopy of Jet-Cooled <i>t</i> -Butoxy. Journal of Physical Chemistry A, 2015, 119, 11804-11812.	1.1	9
17	Jet cooled cavity ringdown spectroscopy of the AËœ2E″â†XËœ2A2′ transition of the NO3 radical. Journal of Chemical Physics, 2015, 142, 184305.	1.2	29
18	Jet-Cooled Laser-Induced Fluorescence Spectroscopy of Isopropoxy Radical: Vibronic Analysis of $\langle i \rangle B f \langle i \rangle A f \langle i \rangle$	1.1	15

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19	Jet-Cooled Laser-Induced Fluorescence Spectroscopy of Cyclohexoxy: Rotational and Fine Structure of Molecules in Nearly Degenerate Electronic States. Journal of Physical Chemistry A, 2014, 118, 11871-11890.	1.1	17
20	Observation of the electronic transition of C6–C10 peroxy radicals. Chemical Physics Letters, 2014, 601, 149-154.	1.2	5
21	Diffraction using laser-driven broadband electron wave packets. Nature Communications, 2014, 5, 4635.	5.8	68
22	Imaging and Scattering Studies of the Unimolecular Dissociation of the BrCH ₂ CH ₂ ORadical from BrCH ₂ CH ₂ ONO Photolysis at 351 nm. Journal of Physical Chemistry A, 2014, 118, 404-416.	1.1	4
23	Detection and Characterization of Reactive Chemical Intermediates Using Cavity Ringdown Spectroscopy. Springer Series in Optical Sciences, 2014, , 61-91.	0.5	0
24	Laser induced fluorescence study of the - transition of FCH2CH2O. Chemical Physics Letters, 2013, 555, 64-71.	1.2	3
25	Autobiography of Terry A. Miller. Journal of Physical Chemistry A, 2013, 117, 13209-13215.	1.1	0
26	Rotationally resolved $B\hat{f}\hat{a}\dagger X\hat{f}$ electronic spectra of the isopropoxy radical: A comparative study. Journal of Chemical Physics, 2013, 139, 094308.	1.2	20
27	Kinetic measurements of the C2H5O2 radical using time-resolved cavity ring-down spectroscopy with a continuous source. Journal of Chemical Physics, 2013, 139, 094201.	1.2	10
28	Imaging ultrafast molecular dynamics with laser-induced electron diffraction. Nature, 2012, 483, 194-197.	13.7	519
29	Detection and Characterization of Products from Photodissociation of XCH ₂ CH ₂ ONO (X = F, Cl, Br, OH). Journal of Physical Chemistry A, 2012, 116, Aralysis of the <a "="" 1998="" altimg="si61.gif" href="mailto:mml:math xmlns:mml=" http:="" math="" mathml"="" www.w3.org="">http://www.w3.org/1998/Math/MathML altimg="si61.gif"	1.1	12
30	overflow="scroll"> <mml:mrow><mml:mover accent="true"><mml:mrow><mml:mtext>A</mml:mtext></mml:mrow><mml:mrow><mml:mo stretchy="true">â^1/4</mml:mo></mml:mrow></mml:mover><mml:ms>â€"</mml:ms><mml:mover accent="true"><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><mml:mrow><m< td=""><td>1.2</td><td>5</td></m<></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mrow></mml:mover></mml:mrow>	1.2	5
31	stretchy="true">â^1/4electronic fransition of High-Order Harmonic Generation in the Long Wavelength Limit of a Strong Laser Field. IEEE Journal of Selected Topics in Quantum Electronics, 2012, 18, 419-433.	1.9	12
32	Electronic Transition Moment for the $0 < \text{sub} > 0 < / \text{sub} > 0 < / \text{sup} > 0 < / \text{sup} > 0$ Band of the $\tilde{A}f$ $\hat{a}\uparrow \bullet X\hat{l}f$ Transition in the Ethyl Peroxy Radical. Journal of Physical Chemistry A, 2011, 115, 13931-13941.	1.1	6
33	overflow="scroll"> <mml:mrow><mml:mover accent="true"><mml:mrow><mml:mi>A</mml:mi></mml:mrow><mml:mrow><mml:mo stretchy="true">â°¼</mml:mo </mml:mrow></mml:mover </mml:mrow> â€" <mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" altimg="si35.gif"</mml:math 	1.2	2
34	Spectroscopic studies of the \$ilde{A}\$Ö\$ilde{X}\$X̃ electronic spectrum of the β-hydroxyethylperoxy radical: Structure and dynamics. Journal of Chemical Physics, 2011, 135, 184304.	1.2	7
35	The spectroscopic characterization of the methoxy radical. III. Rotationally resolved \$\$skew3ilde{A}^2A_lext{–}skew3ilde{X}^2E\$AÌf2A1â€"XÌf2E electronic and \$\$skew3ilde{X}^2E\$XÌf2E submillimeter wave spectra of partially deuterated CH2DO and CHD2O \$\$skew3ilde{X}^2\$ and \$\$copy of the \$\frac{1}{1} \frac{1}{1} \frac{1}{1	1.2	21
36	xinlhs:mml="http://www.w3.org/1998/Math/MathMt" altimg="si112.gif" display="inline" overflow="scroll"> <mml:mrow><mml:mover accent="true"><mml:mrow><mml:mi>A</mml:mi></mml:mrow><mml:mrow><mml:mo stretchy="true">\a^1/4</mml:mo></mml:mrow><</mml:mover></mml:mrow>	1.2	13

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#	Article	lF	CITATIONS
37	The Ã-X̃ absorption of vinoxy radical revisited: Normal and Herzberg–Teller bands observed via cavity ringdown spectroscopy. Journal of Chemical Physics, 2010, 132, 114302.	1.2	8
38	Observation of the $\tilde{A}f\hat{a}^{\circ}X\hat{I}f$ Electronic Transitions of Cyclopentyl and Cyclohexyl Peroxy Radicals via Cavity Ringdown Spectroscopy. Journal of Physical Chemistry A, 2010, 114, 218-231.	1.1	13
39	Observation of the $\tilde{A}f\hat{a}^{*}X\hat{I}f$ Electronic Transition of the \hat{I}^{2} -Hydroxyethylperoxy Radical. Journal of Physical Chemistry Letters, 2010, 1, 1846-1852.	2.1	16
40	Measurements of the Absolute Absorption Cross Sections of the $\langle i \rangle \tilde{A} f \langle i \rangle \hat{a} \uparrow \langle i \rangle \tilde{A} f \langle i \rangle$ Transition in Organic Peroxy Radicals by Dual-Wavelength Cavity Ring-Down Spectroscopy. Journal of Physical Chemistry A, 2010, 114, 11583-11594.	1.1	12
41	High-resolution cavity ringdown spectroscopy of the jet-cooled propyl peroxy radical C3H7O2. Physical Chemistry Chemical Physics, 2010, 12, 4773.	1.3	16
42	<i>Ãf</i> Âfâ°' <i>XÌf</i> Absorption of Propargyl Peroxy Radical (Hâ°'C≡Câ°'CH ₂ OO ^Â) Cavity Ring-Down Spectroscopic and Computational Study. Journal of Physical Chemistry A, 2010, 114, 12437-12446.): A 1.1	5
43	High-resolution cavity ringdown spectroscopy of the jet-cooled ethyl peroxy radical C2H5O2. Journal of Chemical Physics, 2009, 131, 184303.	1.2	21
44	The spectroscopic characterization of the methoxy radical. II. Rotationally resolved $A \hat{I} = -X \hat{I} = $	1.2	32
45	The spectroscopic characterization of the methoxy radical. I. Rotationally resolved Ã A21–X̃ E2 electronic spectra of CH3O. Journal of Chemical Physics, 2009, 130, 074302.	1.2	41
46	An investigation of harmonic generation in liquid media with a mid-infrared laser. Optics Express, 2009, 17, 20959.	1.7	37
47	High harmonic generation from long wavelength drivers. , 2009, , .		0
48	Computational investigation of the Jahn-Teller effect in the ground and excited electronic states of the tropyl radical. Part I. Theoretical calculation of spectroscopically observable parameters. Journal of Chemical Physics, 2008, 128, 084310.	1.2	19
49	The structure and spectra of organic peroxy radicals. Physical Chemistry Chemical Physics, 2008, 10, 3955.	1.3	47
50	Observation of the $\tilde{A}f\hat{a}^{*}X\hat{I}f$ Electronic Transition of the Isomers and Conformers of Pentyl Peroxy Radical Using Cavity Ringdown Spectroscopy. Journal of Physical Chemistry A, 2008, 112, 1445-1456.	1.1	10
51	The Changing Shapes of Molecules. Science, 2008, 320, 881-882.	6.0	5
52	Experimental investigation of the Jahn-Teller effect in the ground and excited electronic states of the tropyl radical. Part II. Vibrational analysis of the AlfE3â \in 32-XlfE2â \in 32 electronic transition. Journal of Chemical Physics, 2008, 128, 084311.	1.2	22
53	Effect of methyl rotation on the electronic spectrum of the methyl peroxy radical. Journal of Chemical Physics, 2007, 127, 044310.	1.2	24
54	Rovibronic bands of the Alfa†Xlf transition of CH3OO and CD3OO detected with cavity ringdown absorption near 1.2–1.4l¼m. Journal of Chemical Physics, 2007, 127, 044311.	1.2	37

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55	The vibrationless $A \hat{f} \hat{a} + \hat{x} \hat{f}$ transition of the jet-cooled deuterated methyl peroxy radical CD3O2 by cavity ringdown spectroscopy. Journal of Chemical Physics, 2007, 127, 224305.	1.2	25
56	Jahn-Teller and related effects in the silver trimer. II: Vibrational analysis of the AÌfE″2-XÌfE′2 electronic transition. Journal of Chemical Physics, 2007, 126, 124309.	1.2	12
57	Development of the Hamiltonian and matrix elements for partially deuterated methoxy radical. Molecular Physics, 2007, 105, 529-540.	0.8	14
58	Quasi-Fourier-transform limited, scannable, high energy titanium-sapphire laser source for high resolution spectroscopy. Review of Scientific Instruments, 2007, 78, 033102.	0.6	20
59	Jahn-Teller and related effects in the silver trimer. I. The ab initio calculation of spectroscopically observable parameters for the XÌfE′2 and AÌfE″2 electronic states. Journal of Chemical Physics, 2007, 126, 124308.	1.2	20
60	Investigation of Ethyl Peroxy Radical Conformers via Cavity Ringdown Spectroscopy of the $\tilde{A}f$ -Xlf Electronic Transition. Journal of Physical Chemistry A, 2007, 111, 832-840.	1.1	33
61	Electron Resonance of Gaseous Diatomic Molecules. Advances in Chemical Physics, 2007, , 149-248.	0.3	105
62	High-resolution IR cavity ring-down spectroscopy of jet-cooled free radicals and other species. Physical Chemistry Chemical Physics, 2006, 8, 1682.	1.3	27
63	Spectroscopic probing and diagnostics of the geometric structure of the alkoxy and alkyl peroxy radical intermediates. Molecular Physics. 2006, 104, 2581-2593, if display="inline" overflow="scroll" cavity ringdown spectroscopy of the <mm:math <mm:math="" altimg="si16.gif" cavity="" d<="" display="si16.gif" of="" overflow="scroll" ringdown="" spectroscopy="" td="" the=""><td>0.8</td><td>28</td></mm:math>	0.8	28
64	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/ja/dtd" xmlns:tb="http://www.elsevier.com/xml/ja/dtd" xmlns:tb="http://www.elsevier.com/xml/ja/dtd" xmlns:tb="http://www.elsevier.com/xml/ja/dtd" xmlns:tb="http://www.elsevier.com/xml/ja/dtd" xmlns:tb="http://www.w3.org/1998/Math/MathML" xmlns:tb="http	1.2	17
65	xmlns:sb="http://www.elsevier.com/xml/common/struct-bib/dtd" xmlns:ce="ht. Chemical Physics Letter High resolution spectra and conformational analysis of 2-butoxy radical. Journal of Chemical Physics, 2006, 125, 094316.	1.2	14
66	A novel Fourier transform limited, high energy, tunable Ti:Sapphire source. , 2006, , .		0
67	Conformational analysis of the 1- and 2-propyl peroxy radicals. Chemical Physics Letters, 2005, 406, 81-89.	1.2	29
68	Determination of the excited-state structure of 7-azaindole-water cluster using a Franck-Condon analysis. Journal of Chemical Physics, 2005, 123, 224311.	1.2	22
69	Theoretical Determinations of the Ambient Conformational Distribution and Unimolecular Decomposition ofn-Propylperoxy Radical. Journal of Physical Chemistry A, 2005, 109, 3637-3646.	1.1	36
70	Near-IR Cavity Ringdown Spectroscopy and Kinetics of the Isomers and Conformers of the Butyl Peroxy Radical. Journal of Physical Chemistry A, 2005, 109, 11191-11197.	1.1	37
71	Observation of the $\tilde{A}f\hat{a}^*X\hat{I}f$ Electronic Transition of the 1-C3H7O2and 2-C3H7O2Radicals Using Cavity Ringdown Spectroscopy. Journal of Physical Chemistry A, 2005, 109, 1308-1315.	1.1	34
72	Accurate ab initio determination of spectroscopic and thermochemical properties of mono- and dichlorocarbenes. Physical Chemistry Chemical Physics, 2005, 7, 2881.	1.3	43

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73	Jet-cooled laser spectroscopy of the cyclohexoxy radical. Journal of Chemical Physics, 2004, 120, 10579-10593.	1.2	26
74	Dispersed fluorescence spectroscopy of primary and secondary alkoxy radicals. Journal of Chemical Physics, 2004, 121, 11780-11797.	1.2	34
7 5	The rotationally resolved electronic spectra of several conformers of 1-hexoxy and 1-heptoxy. Canadian Journal of Chemistry, 2004, 82, 854-866.	0.6	20
76	Radiative and non-radiative decay of selected vibronic levels of the B state of alkoxy radicals. Chemical Physics Letters, 2003, 380, 749-757.	1.2	23
77	Explorations of Conical Intersections and Their Ramifications for Chemistry Through the Jahn—Teller Effect. ChemInform, 2003, 34, no.	0.1	О
78	Theoretical prediction of spectroscopic constants of 1-alkoxy radicals. Journal of Molecular Spectroscopy, 2003, 220, 276-290.	0.4	34
79	Observation of bands among the four lowest pseudorotational states of 1,3-dioxolane. Journal of Molecular Spectroscopy, 2003, 221, 227-238.	0.4	4
80	The absorption spectroscopy of the lowest pseudorotational states of tetrahydrofuran. Journal of Chemical Physics, 2003, 118, 3589-3599.	1.2	47
81	Rotationally Resolved Electronic Spectra of the $\widehat{B}f\widehat{a}^*X\widehat{I}f$ Transition in Multiple Conformers of 1-Butoxy and 1-Pentoxy Radicals. Journal of Physical Chemistry A, 2003, 107, 5189-5201.	1.1	33
82	Cavity Ringdown Spectroscopy of the $\tilde{A}f$ \hat{a} $\hat{A}f$ Electronic Transition of the CH3C(O)O2Radical. Journal of Physical Chemistry A, 2003, 107, 7704-7712.	1.1	29
83	Dispersed fluorescence spectra of the CCl2 Ãf–XÌf vibronic bands. Physical Chemistry Chemical Physics, 2003, 5, 1352-1358.	1.3	31
84	Explorations of conical intersections and their ramifications for chemistry through the Jahn–Teller effect. Chemical Society Reviews, 2003, 32, 38-49.	18.7	93
85	Rotationally resolved B̃–X̃ electronic spectra of both conformers of the 1-propoxy radical. Journal of Chemical Physics, 2003, 118, 4954-4969.	1.2	38
86	Calculation of the Jahn-Teller effect in benzene cation: Application to spectral analysis. Journal of Chemical Physics, 2002, 117, 10654-10674.	1.2	75
87	Spectroscopy of Σ00→Î10 Vibrational–Tunneling–Rotational Band in Rg · ND3 (Rg = Ne, Ar, Kr). Journal of Molecular Spectroscopy, 2002, 214, 202-215.	0.4	12
88	Laser Excitation Spectra of Large Alkoxy Radicals Containing 5â^12 Carbon Atoms. Journal of Physical Chemistry A, 2001, 105, 2925-2928.	1.1	24
89	Submillimeter wave vibration–rotation spectroscopy of Arâ‹CO and Arâ‹ND3. Journal of Chemical Physics, 2001, 114, 6100-6106.	1.2	28
90	Kinetics of Atomic Nitrogen Photofragment Produced by Laser Photodissociation of N2O. Journal of Physical Chemistry A, 2001, 105, 5977-5983.	1.1	9

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91	Observation of the $\tilde{A}f\hat{a}^{\lambda}\tilde{X}$ electronic transition of the CF3O2 radical. Chemical Physics Letters, 2001, 335, 298-304.	1.2	20
92	The Jahnâ€"Teller and related effects in the cyclopentadienyl radical. I. The ab initio calculation of spectroscopically observable parameters. Journal of Chemical Physics, 2001, 114, 4855-4868.	1.2	71
93	The Jahn–Teller and related effects in the cyclopentadienyl radical. II. Vibrational analysis of the Ãf 2A2″–XÌf 2E1″ electronic transition. Journal of Chemical Physics, 2001, 114, 4869-4882.	1.2	67
94	The structure of floppy molecules: the Rg·XH/D (Rg=Ar, Ne, and Kr, X=O or S) family of complexes. Journal of Molecular Structure, 2000, 525, 1-45.	1.8	40
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