

# Stephen G Kukolich

## List of Publications by Year in descending order

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55  
papers

955  
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430754

18  
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docs citations

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times ranked

593  
citing authors

#	ARTICLE	IF	CITATIONS
1	Calculations and analysis of <sup>55</sup> Mn nuclear quadrupole coupling for asymmetric top acyl methyl manganese pentacarbonyl. <i>Chemical Physics Letters</i> , 2021, 762, 138151.	1.2	2
2	Measurements of microwave and NMR spectra for <sup>15</sup> N substituted formamidinium formate. <i>Journal of Molecular Spectroscopy</i> , 2021, 378, 111478.	0.4	0
3	Calculated molecular properties and microwave spectrum analysis for formamidinium formate. <i>Journal of Molecular Spectroscopy</i> , 2020, 372, 111331.	0.4	2
4	Microwave Spectra and Theoretical Calculations for Two Structural Isomers of Methylmanganese Pentacarbonyl. <i>Inorganic Chemistry</i> , 2020, 59, 6432-6438.	1.9	3
5	Synthesis, microwave spectra, x-ray structure, and high-level theoretical calculations for formamidinium formate. <i>Journal of Chemical Physics</i> , 2019, 150, 094305.	1.2	4
6	Microwave Spectra, Structure, and the Aromatic Character of 1-Chloroborepin. <i>Journal of Physical Chemistry A</i> , 2018, 122, 1542-1549.	1.1	1
7	Microwave measurements of the tropolone- <sup>13</sup> C-formic acid doubly hydrogen bonded dimer. <i>Journal of Chemical Physics</i> , 2016, 144, 044306.	1.2	4
8	Microwave spectra, molecular structure, and aromatic character of 4a,8a-azaboranaphthalene. <i>Journal of Chemical Physics</i> , 2016, 144, 114303.	1.2	7
9	Rotational spectra and gas phase structure of the maleimide - <sup>13</sup> C-Formic acid doubly hydrogen bonded dimer. <i>Journal of Molecular Spectroscopy</i> , 2016, 321, 1-4.	0.4	7
10	Microwave spectra and structure of the cyclopropanecarboxylic acid-formic acid dimer. <i>Journal of Chemical Physics</i> , 2015, 143, 124311.	1.2	11
11	Microwave measurements of cyclopropanecarboxylic acid and - <sup>13</sup> C-OD isotopologue. <i>Journal of Molecular Spectroscopy</i> , 2015, 313, 1-3.	0.4	2
12	Microwave Spectrum for a Second Higher Energy Conformer of Cyclopropanecarboxylic Acid and Determination of the Gas Phase Structure of the Ground State. <i>Journal of Physical Chemistry A</i> , 2015, 119, 10016-10021.	1.1	2
13	Identification and characterization of 1,2-BN cyclohexene using microwave spectroscopy. <i>Chemical Physics Letters</i> , 2015, 639, 88-92.	1.2	4
14	Measurements of deuterium quadrupole coupling in propiolic acid and fluorobenzenes using pulsed-beam Fourier transform microwave spectrometers. <i>Journal of Chemical Physics</i> , 2015, 142, 154306.	1.2	0
15	Gas phase measurements of mono-fluoro-benzoic acids and the dimer of 3-fluoro-benzoic acid. <i>Journal of Chemical Physics</i> , 2015, 142, 144303.	1.2	10
16	Microwave spectrum and molecular structure parameters for the 1,2-cyclohexanedione (mono-enolic)- <sup>13</sup> C-formic acid dimer. <i>Chemical Physics Letters</i> , 2014, 613, 86-89.	1.2	8
17	Calculations and measurements of the deuterium tunneling frequency in the propiolic acid-formic acid dimer and description of a newly constructed Fourier transform microwave spectrometer. <i>Journal of Chemical Physics</i> , 2013, 139, 084316.	1.2	6
18	Microwave Structure for the Propiolic Acid- <sup>13</sup> C-Formic Acid Complex. <i>Journal of Physical Chemistry A</i> , 2013, 117, 9525-9530.	1.1	10

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19	Microwave measurements of proton tunneling and structural parameters for the propiolic acid-formic acid dimer. <i>Journal of Chemical Physics</i> , 2011, 135, 154304.	1.2	49
20	Design, construction, and testing of a large-cavity, 1-10 GHz Flygare-Balle spectrometer. <i>Review of Scientific Instruments</i> , 2011, 82, 094103.	0.6	9
21	Microwave spectrum and structural parameters for the formamide-formic acid dimer. <i>Journal of Chemical Physics</i> , 2010, 133, 174304.	1.2	27
22	Communications: Evidence for proton tunneling from the microwave spectrum of the formic acid-propionic acid dimer. <i>Journal of Chemical Physics</i> , 2010, 132, 201101.	1.2	28
23	Microwave Spectrum, Structural Parameters, and Quadrupole Coupling for 1,2-Dihydro-1,2-azaborine. <i>Journal of the American Chemical Society</i> , 2010, 132, 5501-5506.	6.6	40
24	Microwave measurements and ab initio calculations of structural and electronic properties of N-Et-1,2-azaborine. <i>Journal of Chemical Physics</i> , 2009, 131, 224312.	1.2	24
25	The rotational spectrum and structure for the argon-cyclopentadienyl thallium van der Waals complex: Experimental and computational studies of noncovalent bonding in an organometallic $\eta^5$ -complex. <i>Journal of Chemical Physics</i> , 2008, 129, 054305.	1.2	8
26	Microwave Spectra and Gas-Phase Structural Parameters of Bis( $\eta^5$ -cyclopentadienyl)tungsten Dihydride. <i>Organometallics</i> , 2007, 26, 2070-2076.	1.1	46
27	Microwave Spectroscopy Measurements of the Gas-Phase Structure of Cyclopentadienyltungsten Tricarbonyl Hydride. <i>Organometallics</i> , 2005, 24, 2848-2853.	1.1	4
28	Microwave Measurements of $^{14}\text{N}$ and $\text{D}$ Quadrupole Coupling for (Z)-2-Hydroxypyridine and 2-Pyridone Tautomers. <i>Journal of Physical Chemistry A</i> , 2004, 108, 9531-9539.	1.1	31
29	Microwave Spectroscopy Measurements of Rotational Spectra and DFT Calculations for Two Distinct Structural Isomers of $1,1\text{-}\eta^5$ -Dimethylferrocene. <i>Journal of the American Chemical Society</i> , 2004, 126, 844-850.	6.6	25
30	Molecular Structure of $\eta^5$ -Benzyne from Microwave Measurements. <i>Journal of Physical Chemistry A</i> , 2004, 108, 2645-2651.	1.1	25
31	Measurements of microwave spectra and structural parameters for methylferrocene. <i>Journal of Chemical Physics</i> , 2002, 117, 3741-3747.	1.2	14
32	The gas-phase structure of chloroferrocene from microwave spectra. <i>Journal of Chemical Physics</i> , 2000, 112, 747-751.	1.2	16
33	Structural Characterization of anti- and syn-Allyltricarbonyliron Bromide: Rotational Spectra, Quadrupole Coupling, and Density Functional Calculations. <i>Inorganic Chemistry</i> , 2000, 39, 827-835.	1.9	4
34	Microwave Spectra and the Molecular Structure of Tetracarbonyl ethyleneiron. <i>Journal of the American Chemical Society</i> , 1999, 121, 4023-4030.	6.6	16
35	Microwave measurements and calculations on the molecular structure of tetracarbonyldihydridoruthenium. <i>Journal of Chemical Physics</i> , 1998, 109, 9473-9478.	1.2	8
36	Molecular Structure of Tetracarbonyldihydroiron: Microwave Measurements and Density Functional Theory Calculations. <i>Journal of the American Chemical Society</i> , 1998, 120, 6774-6780.	6.6	21

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37	Measurements of structural and quadrupole coupling parameters for bromoferrocene using microwave spectroscopy. <i>Journal of Chemical Physics</i> , 1997, 107, 6541-6548.	1.2	18
38	Gas-Phase Conformational Analysis of 1,4,7-Trithiacyclononane. <i>Journal of Physical Chemistry A</i> , 1997, 101, 9180-9184.	1.1	8
39	Measurements of the microwave spectrum, Re-H bond length, and Re quadrupole coupling for HRe(CO) <sub>5</sub> . <i>Journal of Chemical Physics</i> , 1993, 99, 6465-6469.	1.2	23
40	Comment on: Structure of H <sub>2</sub> S-SO <sub>2</sub> . <i>Journal of Chemical Physics</i> , 1990, 93, 871-872.	1.2	9
41	Deuterium quadrupole coupling in BD <sub>3</sub> CO. <i>Journal of Chemical Physics</i> , 1982, 77, 4312-4317.	1.2	13
42	Beam maser spectroscopy on J = 1-0, K = 1, and K = 0 transitions in CH <sub>3</sub> CN and CH <sub>3</sub> <sup>13</sup> CN. <i>Journal of Chemical Physics</i> , 1982, 76, 97-101.	1.2	33
43	Beam maser measurements of hyperfine structure in chloroacetylene-d. <i>Journal of Molecular Spectroscopy</i> , 1982, 94, 95-99.	0.4	13
44	Beam maser measurements of distortion effects on quadrupole coupling in NH <sub>3</sub> . <i>Journal of Chemical Physics</i> , 1974, 61, 3780-3784.	1.2	14
45	Microwave Spectroscopy and Molecular Structure: Some Undergraduate Laboratory Projects. <i>American Journal of Physics</i> , 1973, 41, 1084-1086.	0.3	2
46	High-Resolution Molecular Zeeman Measurements in CH <sub>2</sub> F <sub>2</sub> ; Observation of <sup>19</sup> F Chemical Shift Anisotropy. <i>Journal of Chemical Physics</i> , 1972, 56, 4446-4449.	1.2	18
47	Variation of Cl Quadrupole Coupling with Isotopic Substitution in CH <sub>3</sub> Cl. <i>Journal of Chemical Physics</i> , 1972, 57, 4052-4054.	1.2	13
48	High Resolution Measurements of Hyperfine Structure in the Rotational Spectrum of CH <sub>3</sub> NC. <i>Journal of Chemical Physics</i> , 1972, 57, 869-871.	1.2	34
49	Deuterium Quadrupole Coupling in Formyl Fluoride. <i>Journal of Chemical Physics</i> , 1971, 55, 610-612.	1.2	10
50	Molecular Beam Measurement of the Magnetic Susceptibility Anisotropies and Molecular Quadrupole Moment in H <sub>2</sub> CO. <i>Journal of Chemical Physics</i> , 1971, 54, 8-11.	1.2	25
51	High-Resolution Measurements of <sup>35</sup> Cl and D Quadrupole Coupling in CH <sub>2</sub> DCl and CD <sub>3</sub> Cl. <i>Journal of Chemical Physics</i> , 1971, 55, 4488-4493.	1.2	27
52	Measurements of Deuterium Quadrupole Coupling in Formic Acid. <i>Journal of Chemical Physics</i> , 1969, 51, 358-360.	1.2	15
53	Measurements of Hyperfine Structure in NH <sub>2</sub> D. <i>Journal of Chemical Physics</i> , 1968, 49, 5523-5525.	1.2	37
54	Hyperfine Structure of N <sup>15</sup> H <sub>3</sub> . <i>Physical Review</i> , 1968, 172, 59-63.	2.7	28

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55	Measurement of Ammonia Hyperfine Structure with a Two-Cavity Maser. Physical Review, 1967, 156, 83-92.	2.7	123