

# Corey J Evans

## List of Publications by Year in descending order

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47  
papers

1,712  
citations

279798

23  
h-index

276875

41  
g-index

48  
all docs

48  
docs citations

48  
times ranked

898  
citing authors

#	ARTICLE	IF	CITATIONS
1	Noble Gas~Metal Chemical Bonds. Microwave Spectra, Geometries, and Nuclear Quadrupole Coupling Constants of Ar~AuCl and Kr~AuCl. <i>Journal of the American Chemical Society</i> , 2000, 122, 6100-6105.	13.7	182
2	Noble gas~metal chemical bonding? The microwave spectra, structures, and hyperfine constants of Ar~CuX (X=F,~S,~Br). <i>Journal of Chemical Physics</i> , 2000, 112, 9363-9374.	3.0	178
3	The microwave spectra and structures of Ar~AgX (X=F,Cl,Br). <i>Journal of Chemical Physics</i> , 2000, 112, 1321-1329.	3.0	173
4	Noble gas~metal chemical bonding: the microwave spectra, structures and hyperfine constants of Ar~AuF and Ar~AuBr. <i>Physical Chemistry Chemical Physics</i> , 2000, 2, 3943-3948.	2.8	128
5	Confirmation of the Existence of Gold(I) Fluoride, AuF:~ Microwave Spectrum and Structure. <i>Journal of the American Chemical Society</i> , 2000, 122, 1560-1561.	13.7	99
6	Microwave Spectrum, Structure, and Hyperfine Constants of Kr~AgCl: Formation of a Weak Kr~Ag Covalent Bond. <i>Journal of Molecular Spectroscopy</i> , 2001, 206, 33-40.	1.2	74
7	The Pure Rotational Spectra of AuCl and AuBr. <i>Journal of Molecular Spectroscopy</i> , 2000, 203, 105-117.	1.2	69
8	Benzoxazine Atropisomers: Intrinsic Atropisomerization Mechanism and Conversion to High Performance Thermosets. <i>Macromolecules</i> , 2018, 51, 7574-7585.	4.8	61
9	Pure Rotational Spectra, Structures, and Hyperfine Constants of OC~AuX (X = F, Cl, Br). <i>Inorganic Chemistry</i> , 2001, 40, 6123-6131.	4.0	58
10	A three-dimensional multivariate image processing technique for the analysis of FTIR spectroscopic images of multiple tissue sections. <i>BMC Medical Imaging</i> , 2006, 6, 12.	2.7	50
11	The Pure Rotational Spectrum of AuI. <i>Journal of Molecular Spectroscopy</i> , 2001, 205, 344-346.	1.2	44
12	Geometries and Bond Energies of the He~MX, Ne~MX, and Ar~MX (M = Cu, Ag, Au; X = F, Cl) Complexes. <i>Journal of Physical Chemistry A</i> , 2010, 114, 4446-4454.	2.5	43
13	Fourier Transform Microwave Spectroscopy of Cyanides and Isocyanides of Al, Ga, and In. <i>Journal of Molecular Spectroscopy</i> , 2001, 209, 178-191.	1.2	42
14	Pure Rotational Spectrum and Structure of Platinum Monocarbonyl, PtCO. <i>Journal of Physical Chemistry A</i> , 2001, 105, 9659-9663.	2.5	40
15	Parasites under the Spotlight: Applications of Vibrational Spectroscopy to Malaria Research. <i>Chemical Reviews</i> , 2018, 118, 5330-5358.	47.7	40
16	The pure rotational spectrum of ScBr. <i>Physical Chemistry Chemical Physics</i> , 2000, 2, 43-46.	2.8	34
17	The Hyperfine Structures of ScCl and ScF. <i>Journal of Molecular Spectroscopy</i> , 2000, 199, 275-283.	1.2	33
18	Infrared spectroscopy of Li(NH <sub>3</sub> ) <sub>n</sub> clusters for n=4~7. <i>Journal of Chemical Physics</i> , 2006, 125, 034302.	3.0	31

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19	The pure rotational spectra of the lanthanum monohalides, LaF, LaCl, LaBr, LaI. Journal of Molecular Spectroscopy, 2003, 218, 169-179.	1.2	28
20	Easily Processable Thermosets with Outstanding Performance via Smart Twisted Small-Molecule Benzoxazines. Macromolecular Rapid Communications, 2020, 41, e1900625.	3.9	28
21	Elucidating the role of acetylene in <i>ortho</i> -phthalimide functional benzoxazines: design, synthesis, and structure-property investigations. Polymer Chemistry, 2021, 12, 5059-5068.	3.9	27
22	High-resolution IR spectroscopy of jet-cooled CF <sub>3</sub> CH <sub>2</sub> F. Journal of the Chemical Society, Faraday Transactions, 1995, 91, 1723.	1.7	25
23	Hyperfine Structure in the Rotational Spectrum of GaF: A Comparison of Experimental and Calculated Spin-Rotation and Electric Field Gradient Tensors. Journal of Molecular Spectroscopy, 2000, 204, 184-194.	1.2	25
24	High-Resolution FTIR Spectrum of Jet-Cooled CH <sub>3</sub> CHF <sub>2</sub> . The Journal of Physical Chemistry, 1996, 100, 8660-8664.	2.9	22
25	Discovery of the optically forbidden S <sub>1</sub> →S <sub>0</sub> transition of silylidene (H <sub>2</sub> C=Si). Journal of Chemical Physics, 2003, 118, 1642-1648.	3.0	22
26	High resolution FTIR spectroscopy of 1,1,1,2-tetrafluoroethane. Journal of Molecular Spectroscopy, 2003, 218, 48-52.	1.2	15
27	The rotational spectrum and complete heavy atom structure of the chiral molecule verbenone. Journal of Molecular Spectroscopy, 2017, 342, 109-115.	1.2	15
28	The Pure Rotational Spectrum of Yttrium Monoiodide. Journal of Molecular Spectroscopy, 2000, 199, 311-313.	1.2	14
29	The High-Resolution FTIR Spectrum of Jet-Cooled CH <sub>3</sub> CF <sub>2</sub> Cl. Journal of Molecular Spectroscopy, 1997, 182, 342-349.	1.2	12
30	A study of the molecular structure and Renner-Teller effect in the $\tilde{A}^1\Sigma^+ \leftarrow \tilde{X}^1\Sigma^+$ electronic spectrum of jet-cooled boron disulfide, BS <sub>2</sub> . Journal of Chemical Physics, 2003, 119, 2047-2056.	3.0	11
31	The pure rotational spectra of the open-shell diatomic molecules Pbl and Snl. Journal of Chemical Physics, 2015, 143, 244309.	3.0	10
32	Spectroscopic detection of the SiCCl free radical. Journal of Chemical Physics, 2002, 117, 6446-6449.	3.0	9
33	The Fourier transform infrared spectrum of jet-cooled CF <sub>3</sub> CHFCF <sub>3</sub> . Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 1999, 55, 1177-1183.	3.9	8
34	Ab initio predictions of the spectroscopic parameters of the silicon halomethylidyne (Si=C=X; X=F, Cl, Br) free radicals. Journal of Chemical Physics, 2002, 117, 6439-6445.	3.0	8
35	The electronic spectroscopy and molecular structure of the HPCl free radical: A potential III-V semiconductor growth intermediate. Journal of Chemical Physics, 2003, 119, 2037-2046.	3.0	8
36	Effects of End-Caps on the Atropisomerization, Polymerization, and the Thermal Properties of <i>ortho</i> -Imide Functional Benzoxazines. Polymers, 2019, 11, 399.	4.5	8

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37	Synchrotron infrared spectroscopy of the $\hat{1}/24$ , $\hat{1}/28$ , $\hat{1}/210$ , $\hat{1}/211$ and $\hat{1}/214$ fundamental bands of thirane. Journal of Molecular Spectroscopy, 2015, 316, 32-37.	1.2	7
38	CF <sub>3</sub> <sup>+</sup> and CF <sub>2</sub> H <sup>+</sup> : new reagents for n-alkane determination in chemical ionisation reaction mass spectrometry. Analyst, The, 2016, 141, 6564-6570.	3.5	7
39	IR Band Profiling of Dichlorodifluoromethane in the Greenhouse Window: High-Resolution FTIR Spectroscopy of $\hat{1}/2<sub>2</sub>$ and $\hat{1}/2<sub>8</sub>$ . Journal of Physical Chemistry A, 2014, 118, 2480-2487.	2.5	5
40	The millimetre-wave spectrum of estragole. Chemical Physics Letters, 2013, 580, 37-42.	2.6	4
41	Computational Study on the Structures of the [H, Si, N, C, O] Isomers: A Possible Species of Interstellar Interest. Journal of Physical Chemistry A, 2007, 111, 13148-13156.	2.5	3
42	Spectroscopic Investigation of the Electronic $\tilde{A}^1\Sigma^+$ Transition of HSiNC. Journal of Physical Chemistry A, 2009, 113, 8533-8539.	2.5	3
43	An investigation of atropisomerism in ortho $\hat{\epsilon}$ imide substituted 1,3 $\hat{\epsilon}$ benzoxazine by experimental NMR and DFT calculations. Journal of Physical Organic Chemistry, 2019, 32, e3926.	1.9	3
44	Spectroscopic investigation of the $\tilde{A}^1\Sigma^+$ electronic transition of HSiNCO. Journal of Chemical Physics, 2009, 131, 124302.	3.0	2
45	Computational study on the energies and structures of the [H, Si, N, C, S] isomers. Theoretical Chemistry Accounts, 2010, 127, 661-669.	1.4	2
46	Vibrational energies and full analytic potential energy functions of Pbl and Inl from pure microwave data. Journal of Molecular Spectroscopy, 2016, 330, 80-88.	1.2	2
47	Ab initio predictions of the spectroscopic parameters of the germanium halomethylidyne (Ge=C-X; X=F,) Tj ETQq1 1.0.784314 rgBT /Cve	1.5	0