

Wolfgang Jger

List of Publications by Citations

Source: <https://exaly.com/author-pdf/3755897/wolfgang-jager-publications-by-citations.pdf>

Version: 2024-04-27

This document has been generated based on the publications and citations recorded by exaly.com. For the latest version of this publication list, visit the link given above.

The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

82

papers

1,794

citations

26

h-index

38

g-index

88

ext. papers

2,033

ext. citations

4.9

avg, IF

4.94

L-index

#	Paper	IF	Citations
82	Evidence for heavy atom large amplitude motions in RG-cyclopropane van der Waals complexes (RG=Ne, Ar, Kr) from rotation-tunneling spectroscopy. <i>Journal of Chemical Physics</i> , 1997 , 106, 7968-7980	3.9	153
81	Spectroscopic exploration of atomic scale superfluidity in doped helium nanoclusters. <i>Physical Review Letters</i> , 2006 , 97, 183401	7.4	82
80	Rotational spectroscopic investigation of carbonyl sulfide solvated with helium atoms. <i>Journal of Chemical Physics</i> , 2003 , 119, 5457-5466	3.9	76
79	Rotational spectrum, structure, and chlorine nuclear quadrupole coupling constants of the van der Waals complex ArCl ₂ . <i>Journal of Chemical Physics</i> , 1993 , 98, 3726-3731	3.9	69
78	Microwave spectra of metal chlorides produced using laser ablation. <i>Journal of Chemical Physics</i> , 1993 , 99, 3320-3328	3.9	67
77	Pure rotational spectra of the mixed rare gas van der Waals complexes NeXe, ArXe, and KrXe. <i>Journal of Chemical Physics</i> , 1993 , 99, 919-927	3.9	66
76	Microwave spectroscopy of ternary and quaternary van der Waals clusters. <i>International Reviews in Physical Chemistry</i> , 2005 , 24, 301-338	7	56
75	Spectroscopic studies of OCS-doped 4He clusters with 9-72 helium atoms: observation of broad oscillations in the rotational moment of inertia. <i>Journal of Physical Chemistry A</i> , 2007 , 111, 7329-37	2.8	47
74	Microwave-submillimeter wave double-resonance spectrometer for the investigation of van der Waals complexes. <i>Review of Scientific Instruments</i> , 1998 , 69, 4061-4067	1.7	44
73	Direct Spectroscopic detection of the orientation of free OH groups in methyl lactate-(water)(1,2) clusters: hydration of a chiral hydroxy ester. <i>Angewandte Chemie - International Edition</i> , 2014 , 53, 1156-9	16.4	43
72	Investigation of the Ne-NH ₃ van der Waals complex: Rotational spectrum and ab initio calculations. <i>Journal of Chemical Physics</i> , 2001 , 115, 6504-6512	3.9	41
71	A Direct Link from the Gas to the Condensed Phase: A Rotational Spectroscopic Study of 2,2,2-Trifluoroethanol Trimers. <i>Angewandte Chemie - International Edition</i> , 2017 , 56, 6289-6293	16.4	40
70	High resolution spectroscopy of Ne and Ar containing noble gas clusters. <i>Journal of Chemical Physics</i> , 1997 , 107, 4788-4796	3.9	40
69	Rotational spectroscopic study of hydrogen cyanide embedded in small 4He clusters. <i>Journal of Chemical Physics</i> , 2012 , 137, 174303	3.9	39
68	Rotational spectra of the mixed rare gas dimers NeKr and ArKr. <i>Journal of Chemical Physics</i> , 1995 , 103, 2827-2833	3.9	35
67	The microwave rotational spectrum of the van der Waals complex KrCl ₂ . <i>Journal of Chemical Physics</i> , 1993 , 99, 7510-7520	3.9	34
66	The benzoic acid-water complex: a potential atmospheric nucleation precursor studied using microwave spectroscopy and ab initio calculations. <i>Physical Chemistry Chemical Physics</i> , 2014 , 16, 2305-14	2.6	33

65	Microwave spectroscopic investigation of the mixed rare gas van der Waals trimers Ne ₂ Kr and Ne ₂ Ar. <i>Journal of Chemical Physics</i> , 1994 , 100, 4171-4180	3.9	31
64	Rotational spectra of the Kr-H ₂ O van der Waals complex. <i>Molecular Physics</i> , 2000 , 98, 1575-1588	1.7	30
63	The Chiral Trimer and a Metastable Chiral Dimer of Achiral Hexafluoroisopropanol: A Multi-Messenger Study. <i>Angewandte Chemie - International Edition</i> , 2019 , 58, 5080-5084	16.4	30
62	The dynamics of the CO ₂ interaction: Strong Coriolis coupling in CO-para-N ₂ . <i>Journal of Chemical Physics</i> , 2000 , 113, 514-524	3.9	29
61	Rotational spectra and theoretical study of tetramers and trimers of 2-fluoroethanol: dramatic intermolecular compensation for intramolecular instability. <i>Physical Chemistry Chemical Physics</i> , 2018 , 20, 27630-27637	3.6	28
60	Chirped-pulse and cavity-based Fourier transform microwave spectroscopy of a chiral epoxy ester: methyl glycidate. <i>Journal of Physical Chemistry A</i> , 2013 , 117, 13249-54	2.8	27
59	A Direct Link from the Gas to the Condensed Phase: A Rotational Spectroscopic Study of 2,2,2-Trifluoroethanol Trimers. <i>Angewandte Chemie</i> , 2017 , 129, 6386-6390	3.6	26
58	Unusual H-Bond Topology and Bifurcated H-bonds in the 2-Fluoroethanol Trimer. <i>Angewandte Chemie - International Edition</i> , 2015 , 54, 11711-5	16.4	26
57	The rotational spectra of the fluorobenzene-water and p-difluorobenzene-water dimers: Structure and internal dynamics. <i>Journal of Molecular Spectroscopy</i> , 2011 , 268, 47-52	1.3	26
56	Hydration of the simplest keto acid: a rotational spectroscopic and ab initio study of the pyruvic acid-water complex. <i>Physical Chemistry Chemical Physics</i> , 2017 , 19, 4440-4446	3.6	24
55	Rotational Spectra of Two Hydrogen-Bonded Methyl Salicylate Monohydrates: Relative Stability and Tunneling Motions. <i>Journal of Physical Chemistry Letters</i> , 2015 , 6, 3126-3131	6.4	23
54	The rich conformational landscape of perillyl alcohol revealed by broadband rotational spectroscopy and theoretical modelling. <i>Physical Chemistry Chemical Physics</i> , 2019 , 21, 15408-15416	3.6	23
53	Direct observation of rotational transitions of the CO ₂ O dimer. <i>Journal of Chemical Physics</i> , 1999 , 111, 5754-5756	3.9	23
52	Microwave spectra of the Ar ₂ D ₃ van der Waals complex and its partially protonated isotopomers. <i>Journal of Chemical Physics</i> , 2001 , 114, 3968-3976	3.9	21
51	Fourier transform microwave rotational spectra of the Ne ₂ N ₂ O and Ar ₂ N ₂ O van der Waals trimers. <i>Journal of Chemical Physics</i> , 1999 , 111, 3919-3928	3.9	20
50	Discovering the Elusive Global Minimum in a Ternary Chiral Cluster: Rotational Spectra of Propylene Oxide Trimer. <i>Angewandte Chemie - International Edition</i> , 2020 , 59, 22427-22430	16.4	20
49	Decelerating molecules with microwave fields. <i>Physical Review A</i> , 2012 , 85,	2.6	19
48	Microwave rotational spectra of the Kr-NH ₃ van der Waals complex. <i>Molecular Physics</i> , 2001 , 99, 1215-1228		19

47	Chirped-pulse and cavity-based Fourier transform microwave spectra of the methyl lactate-ammonia adduct. <i>Angewandte Chemie - International Edition</i> , 2013 , 52, 4402-5	16.4	18
46	Relative Humidity Dependence of Soot Aggregate Restructuring Induced by Secondary Organic Aerosol: Effects of Water on Coating Viscosity and Surface Tension. <i>Environmental Science and Technology Letters</i> , 2017 , 4, 386-390	11	17
45	Chirality induction and amplification in the 2,2,2-trifluoroethanol-propylene oxide adduct. <i>Angewandte Chemie - International Edition</i> , 2014 , 53, 7277-80	16.4	17
44	Rotational spectra and internal dynamics of NeH ₂ S. <i>Molecular Physics</i> , 2002 , 100, 611-622	1.7	17
43	The Chiral Trimer and a Metastable Chiral Dimer of Achiral Hexafluoroisopropanol: A Multi-Messenger Study. <i>Angewandte Chemie</i> , 2019 , 131, 5134-5138	3.6	15
42	High-resolution rotational spectroscopy of H ₂ -OCS: a study of minor isotopomers of OCS. <i>Molecular Physics</i> , 2008 , 106, 23-31	1.7	15
41	Tunnelling and barrier-less motions in the 2-fluoroethanol-water complex: a rotational spectroscopic and ab initio study. <i>Physical Chemistry Chemical Physics</i> , 2017 , 19, 12221-12228	3.6	14
40	Conformational Panorama and Chirality Controlled Structure-Energy Relationship in a Chiral Carboxylic Acid Dimer. <i>Angewandte Chemie - International Edition</i> , 2020 , 59, 15703-15710	16.4	14
39	Direkter spektroskopischer Nachweis spezifischer Orientierungen freier OH-Gruppen in Methyllactat-(Wasser) _{1,2} -Clustern: Hydratation eines chiralen Hydroxyesters. <i>Angewandte Chemie</i> , 2014 , 126, 1175-1178	3.6	14
38	Rotational spectroscopy of the atmospheric photo-oxidation product o-toluic acid and its monohydrate. <i>Physical Chemistry Chemical Physics</i> , 2016 , 18, 448-57	3.6	13
37	Relationship between Coating-Induced Soot Aggregate Restructuring and Primary Particle Number. <i>Environmental Science & Technology</i> , 2017 , 51, 8376-8383	10.3	13
36	Microwave spectroscopic investigation of the Ne-C ₂ H ₂ van der Waals dimer. <i>Physical Chemistry Chemical Physics</i> , 2003 , 5, 1744-1751	3.6	13
35	Contrasting Effects of Water on the Barriers to Decarboxylation of Two Oxalic Acid Monohydrates: A Combined Rotational Spectroscopic and Ab Initio Study. <i>Journal of Physical Chemistry Letters</i> , 2016 , 7, 1143-7	6.4	13
34	Rotational spectroscopy of the methyl glycidate-water complex: conformation and water and methyl rotor tunnelling motions. <i>Physical Chemistry Chemical Physics</i> , 2017 , 19, 29508-29515	3.6	12
33	Single and double resonance spectroscopy of methanol embedded in superfluid helium nanodroplets. <i>Journal of Chemical Physics</i> , 2014 , 141, 044301	3.9	12
32	Chirality Induction and Amplification in the 2,2,2-Trifluoroethanol-Propylene Oxide Adduct. <i>Angewandte Chemie</i> , 2014 , 126, 7405-7408	3.6	11
31	Enhanced Evaporation of Microscale Droplets With an Infrared Laser. <i>Journal of Heat Transfer</i> , 2017 , 139,	1.8	10
30	Unusual H-Bond Topology and Bifurcated H-bonds in the 2-Fluoroethanol Trimer. <i>Angewandte Chemie</i> , 2015 , 127, 11877-11881	3.6	10

29	Microwave rotational spectra of the Ar ₃ NH ₃ van der Waals tetramer. <i>Journal of Chemical Physics</i> , 2002 , 116, 2379-2387	3.9	10
28	The 1-Naphthol Dimer and Its Surprising Preference for π -Stacking over Hydrogen Bonding. <i>Journal of Physical Chemistry Letters</i> , 2019 , 10, 2836-2841	6.4	9
27	HIGH-RESOLUTION FOURIER-TRANSFORM MICROWAVE SPECTROSCOPY OF METHYL- AND DIMETHYLNAPHTHALENES. <i>Astrophysical Journal</i> , 2015 , 805, 141	4.7	9
26	Microwave rotational spectroscopic investigation of the Ar ₂ NH ₃ van der Waals trimer. <i>Physical Chemistry Chemical Physics</i> , 2002 , 4, 4883-4889	3.6	9
25	Spectroscopic study of the mixed rare-gas molecule van der Waals trimer NeArHCl. <i>Journal of Chemical Physics</i> , 1999 , 110, 4354-4362	3.9	9
24	Rotational spectroscopy of chiral tetrahydro-2-furoic acid: Conformational landscape, conversion, and abundances. <i>Journal of Chemical Physics</i> , 2018 , 149, 224306	3.9	9
23	Microwave and ab initio studies of the internal rotation of ethylene in the Ar-ethylene and Ne-ethylene van der Waals complexes. <i>Journal of Chemical Physics</i> , 2003 , 119, 8449-8463	3.9	8
22	Shapes of rotating normal fluid He ₃ versus superfluid He ₄ droplets in molecular beams. <i>Physical Review B</i> , 2020 , 102,	3.3	8
21	Rotational Spectroscopy of p-Toluic Acid and Its 1:1 Complex with Water. <i>Journal of Physical Chemistry A</i> , 2017 , 121, 8625-8631	2.8	7
20	A microwave spectroscopic and ab initio study of keto-enol tautomerism and isomerism in the cyclohexanone-water complex. <i>Physical Chemistry Chemical Physics</i> , 2019 , 21, 12872-12880	3.6	7
19	Rotational spectroscopic and theoretical study of the perfluorobutyric acid?formic acid complex. <i>Journal of Molecular Spectroscopy</i> , 2017 , 335, 88-92	1.3	5
18	Theoretical study of the microwave spectrum of isotopologues of OCS(He) ₂ . <i>Canadian Journal of Chemistry</i> , 2010 , 88, 1146-1154	0.9	5
17	Discovering the Elusive Global Minimum in a Ternary Chiral Cluster: Rotational Spectra of Propylene Oxide Trimer. <i>Angewandte Chemie</i> , 2020 , 132, 22613-22616	3.6	5
16	Rotational spectroscopy of methyl benzoylformate and methyl mandelate: structure and internal dynamics of a model reactant and product of enantioselective reduction. <i>Physical Chemistry Chemical Physics</i> , 2015 , 17, 21942-9	3.6	4
15	Microwave spectroscopy of the seeded binary and ternary clusters CO-(pH ₂) ₂ , CO-pH ₂ -He, CO-HD, and CO-(oD ₂)(N=1,2). <i>Journal of Chemical Physics</i> , 2015 , 142, 144308	3.9	3
14	Conformational Panorama and Chirality Controlled Structure-Energy Relationship in a Chiral Carboxylic Acid Dimer. <i>Angewandte Chemie</i> , 2020 , 132, 15833-15840	3.6	3
13	The ArNe-N ₂ O van der Waals trimer: a high resolution spectroscopic study of its rotational spectrum, structure and dynamics. <i>Molecular Physics</i> , 2001 , 99, 13-24	1.7	3
12	Microwave spectrum of the complex of 3,3,3-trifluoro-2-(trifluoromethyl)propanoic acid and formic acid. <i>Molecular Physics</i> , 2019 , 117, 1193-1199	1.7	3

11	Conformational Landscape, Chirality Recognition and Chiral Analyses: Rotational Spectroscopy of Tetrahydro-2-Furoic Acid-Propylene Oxide Conformers. <i>ChemPhysChem</i> , 2021 , 22, 455-460	3.2	3
10	Rotational Spectroscopic Study of Quantum Solvation in Isotopologic (pH)-CO Clusters. <i>Journal of Physical Chemistry A</i> , 2017 , 121, 3671-3678	2.8	2
9	A spectroscopic and ab initio study of the hydrogen peroxide-formic acid complex: hindering the internal motion of HO. <i>Physical Chemistry Chemical Physics</i> , 2018 , 20, 21345-21351	3.6	2
8	Aggregation of solutes in bosonic versus fermionic quantum fluids. <i>Science Advances</i> , 2021 , 7, eabk2247	14.3	2
7	Structural and dynamical features of the 2,2,2-trifluoroethanolammonia complex. <i>Physical Chemistry Chemical Physics</i> , 2020 , 22, 23019-23027	3.6	2
6	Higher-Energy Hexafluoroisopropanol-Water Isomer and Its Large Amplitude Motions: Rotational Spectra and DFT Calculations. <i>Journal of Physical Chemistry A</i> , 2021 , 125, 10401-10409	2.8	1
5	Conformational landscape and intricate conformational relaxation paths of 4,4,4-trifluoro-1-butanol: Rotational spectroscopy and quantum chemical calculations. <i>Journal of Molecular Structure</i> , 2020 , 1217, 128359	3.4	1
4	Rotational Spectrum and Molecular Structures of the Binary Aggregates of 1,1,1,3,3,3-Hexafluoro-2-propanol with Ne and Ar. <i>Journal of Physical Chemistry A</i> , 2021 , 125, 5355-5364	2.8	1
3	Unusual binary aggregates of perylene bisimide revealed by their electronic transitions in helium nanodroplets and DFT calculations. <i>Physical Chemistry Chemical Physics</i> , 2021 , 23, 13862-13872	3.6	1
2	A rotational spectroscopic study of (+) and (-)-carveol: further insights into conformational dynamics in monoterpenes and monoterpenoids. <i>Physical Chemistry Chemical Physics</i> , 2021 , 23, 15159-15168	3.6	1
1	Hydrogen bonding interactions in the 1,1,1,3,3,3-hexafluoro-2-propanol-4-dioxane complex: Rotational spectroscopy and density functional theory calculations. <i>Journal of Molecular Spectroscopy</i> , 2021 , 376, 111408	1.3	1