

Andrey A Vigasin

List of Publications by Year in descending order

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68
papers

1,342
citations

361413

20
h-index

361022

35
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68
all docs

68
docs citations

68
times ranked

1180
citing authors

#	ARTICLE	IF	CITATIONS
1	Transient reducing greenhouse warming on early Mars. <i>Geophysical Research Letters</i> , 2017, 44, 665-671.	4.0	178
2	Update of the HITRAN collision-induced absorption section. <i>Icarus</i> , 2019, 328, 160-175.	2.5	105
3	Water vapour self-continuum and water dimers: 1. Analysis of recent work. <i>Journal of Quantitative Spectroscopy and Radiative Transfer</i> , 2011, 112, 1286-1303.	2.3	93
4	Bound, metastable and free states of bimolecular complexes. <i>Infrared Physics</i> , 1991, 32, 461-470.	0.5	63
5	Matrix isolation spectra of the carbon dioxide monomer and dimer revisited. <i>Vibrational Spectroscopy</i> , 2000, 23, 83-94.	2.2	58
6	Explicit correlation treatment of the potential energy surface of CO ₂ dimer. <i>Journal of Chemical Physics</i> , 2014, 140, 234310.	3.0	53
7	Direct absorption spectroscopy of water clusters formed in a continuous slit nozzle expansion. <i>Journal of Chemical Physics</i> , 2009, 131, 204312.	3.0	52
8	Water vapor continuous absorption in various mixtures: possible role of weakly bound complexes. <i>Journal of Quantitative Spectroscopy and Radiative Transfer</i> , 2000, 64, 25-40.	2.3	50
9	Vibrational frequency shifts caused by weak intermolecular interactions. <i>Chemical Physics Letters</i> , 1997, 269, 235-243.	2.6	46
10	On the spectroscopic manifestations of weakly bound complexes in rarefied gases. <i>Chemical Physics Letters</i> , 1985, 117, 85-88.	2.6	33
11	Vibrational frequency shifts and thermodynamic stabilities of (HF) _n isomers (n=4-8). <i>Chemical Physics Letters</i> , 1995, 245, 319-325.	2.6	32
12	Four-photon coherent spectroscopy of orientational motion of H ₂ O molecules in liquid water. <i>Journal of Raman Spectroscopy</i> , 2005, 36, 145-147.	2.5	29
13	H ₂ O-N ₂ collision-induced absorption band intensity in the region of the N ₂ fundamental: <i>ab initio</i> investigation of its temperature dependence and comparison with laboratory data. <i>Philosophical Transactions Series A, Mathematical, Physical, and Engineering Sciences</i> , 2012, 370, 2691-2709.	3.4	28
14	On the possibility to quantify contributions from true bound and metastable pairs to infrared absorption in pressurised water vapour. <i>Molecular Physics</i> , 2010, 108, 2309-2313.	1.7	27
15	Temperature Variations of the Interaction Induced Absorption of CO ₂ in the $\hat{1}\frac{1}{2}1, 2\hat{1}\frac{1}{2}2$ Region: FTIR Measurements and Dimer Contribution. <i>Journal of Molecular Spectroscopy</i> , 2002, 213, 51-56.	1.2	24
16	Contribution of bound, metastable and free states of bimolecular complexes to collision-induced intensity of absorption. <i>Chemical Physics Letters</i> , 1994, 225, 537-541.	2.6	23
17	On the nature of collision-induced absorption in gaseous homonuclear diatomics. <i>Journal of Quantitative Spectroscopy and Radiative Transfer</i> , 1996, 56, 409-422.	2.3	23
18	Intensity and Bandshapes of Collision-Induced Absorption by CO ₂ in the Region of the Fermi Doublet. <i>Journal of Molecular Spectroscopy</i> , 2000, 200, 89-95.	1.2	22

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19	Subdivision of phase space for anisotropically interacting water molecules. <i>Molecular Physics</i> , 1997, 90, 101-106.	1.7	21
20	Statistical physics partitioning and classical trajectory analysis of the phase space in CO ₂ -Ar weakly interacting pairs. <i>Journal of Molecular Structure</i> , 2005, 742, 31-36.	3.6	21
21	Direct FTIR high resolution probe of small and medium size Ar _n (CO ₂) _m van der Waals complexes formed in a slit supersonic expansion. <i>Journal of Molecular Spectroscopy</i> , 2006, 240, 141-152.	1.2	20
22	Collision-Induced Absorption in the Region of the O ₂ Fundamental: Bandshapes and Dimeric Features. <i>Journal of Molecular Spectroscopy</i> , 2000, 202, 59-66.	1.2	18
23	Water vapor continuum: Whether collision-induced absorption is involved?. <i>Journal of Quantitative Spectroscopy and Radiative Transfer</i> , 2014, 148, 58-64.	2.3	18
24	Classical calculation of the equilibrium constants for true bound dimers using complete potential energy surface. <i>Journal of Chemical Physics</i> , 2015, 143, 234304.	3.0	18
25	Nuclear Spin Symmetry Conservation in ¹ H ₂ ¹⁶ O Investigated by Direct Absorption FTIR Spectroscopy of Water Vapor Cooled Down in Supersonic Expansion. <i>Journal of Physical Chemistry A</i> , 2017, 121, 7455-7468.	2.5	18
26	Spectral composition of the water vapour self-continuum absorption within 2.7 and 6.25 μm bands. <i>Journal of Quantitative Spectroscopy and Radiative Transfer</i> , 2019, 228, 97-105.	2.3	18
27	On the water dimer contribution to the OH stretching absorption band profile in pressurized water vapour. <i>Molecular Physics</i> , 2008, 106, 1155-1159.	1.7	16
28	Simulation of collision-induced absorption spectra based on classical trajectories and <i>ab initio</i> potential and induced dipole surfaces. I. Case study of N ₂ -N ₂ rototranslational band. <i>Journal of Chemical Physics</i> , 2019, 151, 194106.	3.0	16
29	On the temperature variations of the integrated absorption intensity in the oxygen fundamental. <i>Journal of Molecular Spectroscopy</i> , 2004, 224, 185-187.	1.2	15
30	Identification of the (CO ₂) ₂ Dimer Vibrations in the $\hat{1}\frac{1}{2}1$, $2\hat{1}\frac{1}{2}2$ Region: Anharmonic Variational Calculations. <i>Journal of Molecular Spectroscopy</i> , 2001, 209, 81-87.	1.2	14
31	Density evolution of absorption bandshapes in the water vapor OH-stretching fundamental and overtone: evidence for molecular aggregation. <i>Journal of Molecular Structure</i> , 2005, 742, 173-181.	3.6	14
32	On the origin of the band structure observed in the collision-induced absorption bands of CO ₂ . <i>Journal of Molecular Spectroscopy</i> , 2003, 218, 260-261.	1.2	12
33	On the influence of van der waals association on IR absorption band shapes of highly compressed carbon dioxide. <i>Infrared Physics</i> , 1989, 29, 575-582.	0.5	11
34	Thermally averaged effective dissociation energy of dimers. <i>Chemical Physics Letters</i> , 1995, 242, 33-38.	2.6	11
35	<i>Ab initio</i> 3D potential energy and dipole moment surfaces for the CH ₄ -Ar complex: Collision-induced intensity and dimer content. <i>Journal of Chemical Physics</i> , 2016, 144, 054304.	3.0	10
36	IR-spectra of (CO ₂) ₂ dimers and collision-induced absorption of carbon dioxide in the region of the fermi doublet ($\hat{1}\frac{1}{2}1$, $2\hat{1}\frac{1}{2}2$). <i>Journal of Quantitative Spectroscopy and Radiative Transfer</i> , 1993, 50, 695-703.	2.3	9

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37	Mass-action law for highly excited dimers. <i>Chemical Physics Letters</i> , 1998, 290, 495-501.	2.6	8
38	Ab initio and multipolar characterisation of the induced dipole surface for CH ₄ –CH ₄ : Application to dipole-forbidden absorption in the Titan’s atmosphere. <i>Journal of Molecular Spectroscopy</i> , 2013, 291, 102-107.	1.2	8
39	Far-infrared collision-induced absorption in rare gas mixtures: Quantum and semi-classical calculations. <i>Journal of Chemical Physics</i> , 2014, 140, .	3.0	8
40	Spin-selective adsorption of water vapor. <i>Doklady Physics</i> , 2002, 47, 842-845.	0.7	7
41	Systematic trends in the vibrational frequency shifts of some molecules trapped in amorphous water ice. <i>Journal of Molecular Structure</i> , 2003, 658, 101-113.	3.6	7
42	Extensive ab initio study of the integrated IR intensity in the N ₂ fundamental collision-induced band. <i>Molecular Physics</i> , 2008, 106, 1227-1231.	1.7	7
43	The nature of the absorption bandshape density evolution for the first overtone of CO compressed by N ₂ . <i>Infrared Physics</i> , 1993, 34, 289-298.	0.5	6
44	Simultaneous infrared absorption in a mixture of CO ₂ and H ₂ O: The role of hydrogen-bonded aggregates. <i>Journal of Quantitative Spectroscopy and Radiative Transfer</i> , 1994, 52, 295-301.	2.3	6
45	High-resolution CARS spectroscopy of small carbon dioxide clusters: investigation of the CO ₂ dimer in the Fermi dyad. <i>Journal of Molecular Structure</i> , 1997, 410-411, 47-50.	3.6	6
46	Simulation of collision-induced absorption spectra based on classical trajectories and ab initio potential and induced dipole surfaces. II. CO ₂ –Ar rototranslational band including true dimer contribution. <i>Journal of Chemical Physics</i> , 2021, 155, 064301.	3.0	6
47	Thermally Averaged Spectroscopic Parameters of the Weakly Bound Dimers. <i>Journal of Molecular Spectroscopy</i> , 2001, 205, 9-15.	1.2	5
48	Breakdown of conventional rovibrational selection rules for field- or collision-induced absorption in symmetric linear molecules. <i>Physical Review A</i> , 2020, 102, .	2.5	5
49	Spatial distribution of CO ₂ dimers in axisymmetric gas jets expanding in a vacuum. <i>Journal of Applied Mechanics and Technical Physics</i> , 1989, 30, 52-57.	0.5	4
50	CCSD(T) potential energy and induced dipole surfaces for N ₂ –H ₂ (D ₂): Retrieval of the collision-induced absorption integrated intensities in the regions of the fundamental and first overtone vibrational transitions. <i>Journal of Chemical Physics</i> , 2012, 137, 114308.	3.0	4
51	Greenhouse effect in planetary atmospheres caused by molecular symmetry breaking in intermolecular interactions. <i>Izvestiya - Atmospheric and Oceanic Physics</i> , 2017, 53, 164-173.	0.9	4
52	Trajectory-based Simulation of Far-infrared Collision-induced Absorption Profiles of CH ₄ –N ₂ for Modeling Titan’s Atmosphere. <i>Astrophysical Journal, Supplement Series</i> , 2022, 258, 33.	7.7	4
53	Herman-Wallis factor for a molecule of type HCN. <i>Journal of Quantitative Spectroscopy and Radiative Transfer</i> , 2022, 288, 108274.	2.3	4
54	Structure and molecular spectroscopy of gas-phase complexes. <i>Journal of Structural Chemistry</i> , 1988, 28, 735-764.	1.0	3

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55	Comprehensive classical analysis of partition function and some observables for weakly interacting polyatomic dimers. <i>Journal of Chemical Physics</i> , 2018, 149, 194304.	3.0	3
56	Theory of rovibrational line intensities in allowed and collision-induced absorption spectra of linear molecules. <i>Physical Review A</i> , 2021, 104, .	2.5	3
57	Systematization of published research graphics characterizing weakly bound molecular complexes with carbon dioxide. , 2017, , .		3
58	Kinetics of dimer formation in rarefied water vapor streams. <i>Journal of Applied Mechanics and Technical Physics</i> , 1981, 22, 66-71.	0.5	2
59	Structure and properties of associates of water. <i>Journal of Structural Chemistry</i> , 1983, 24, 102-131.	1.0	2
60	ON THE MODELLING OF ABSORPTION IN THE $\hat{v}_{1/2}^3$ BAND FAR WING OF CO ₂ PERTURBED BY ARGON. <i>Journal of Quantitative Spectroscopy and Radiative Transfer</i> , 1999, 61, 743-749.	2.3	2
61	On the concept of excluded volume for weakly associating gas. <i>Molecular Physics</i> , 2012, 110, 2957-2961.	1.7	2
62	Diffraction of light by absorbing inclusions in solids. <i>Soviet Journal of Quantum Electronics</i> , 1977, 7, 370-372.	0.1	1
63	Application of spontaneous Raman spectroscopy to the study of molecular association in a freely expanding gas jet. <i>Journal of Applied Spectroscopy</i> , 1991, 55, 777-780.	0.7	1
64	The effect of perturbing gas density on the intensity distribution in rovibrational absorption bands. <i>Chemical Physics</i> , 2006, 325, 404-410.	1.9	1
65	Systematization of published research plots in spectroscopy of weakly bounded complexes of molecular oxygen and nitrogen. , 2018, , .		1
66	Thermal stresses in laser materials containing inclusions. <i>Soviet Journal of Quantum Electronics</i> , 1976, 6, 522-524.	0.1	0
67	<title>Partially deuterated water dimer: are there perspectives for its spectroscopic detection in the atmosphere?</title>. , 2006, 6580, 246.		0
68	<title>Ab initio simulation of collision-induced intensity in the N$\langle inf \rangle \langle roman \rangle 2 \langle /roman \rangle \langle /inf \rangle \langle /math \rangle$ fundamental</title>. , 2006, 6580, 90.		0