

Sm Melikova

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/1950361/publications.pdf>

Version: 2024-02-01

65
papers

994
citations

430874

18
h-index

477307

29
g-index

65
all docs

65
docs citations

65
times ranked

629
citing authors

#	ARTICLE	IF	CITATIONS
1	Unusual spectroscopic properties of CF ₃ H dissolved in liquified Ar, N ₂ , CO, and CO ₂ . <i>Chemical Physics Letters</i> , 2002, 352, 301-310.	2.6	69
2	Proton Transfer Equilibria in Schiff Bases with Steric Repulsion. <i>Journal of Physical Chemistry A</i> , 2004, 108, 2131-2138.	2.5	66
3	A Cryosolution Infrared Study of the Complexes of Fluoroform with Ammonia and Pyridine: Evidence for a C ^{δ+} -H ^{δ-} ⋯N Pseudo Blue-Shifting Hydrogen Bond. <i>Journal of Physical Chemistry A</i> , 2005, 109, 3038-3044.	2.5	63
4	Blue shifted F ₃ CH⋯FCD ₃ and Cl ₃ CH⋯FCD ₃ weakly H-bound complexes. <i>Cryospectroscopic and ab initio study</i> . <i>Chemical Physics</i> , 2005, 313, 225-243.	1.9	59
5	Ab Initio Studies of Electron Acceptor-Donor Interactions with Blue- and Red-Shifted Hydrogen Bonds. <i>ChemPhysChem</i> , 2005, 6, 1282-1292.	2.1	59
6	Interpretation of Hydrogen/Deuterium Isotope Effects on NMR Chemical Shifts of [FHF] ⁺ Ion Based on Calculations of Nuclear Magnetic Shielding Tensor Surface. <i>Zeitschrift Fur Physikalische Chemie</i> , 2003, 217, 1549-1564.	2.8	41
7	Spectroscopic and structural consequences of intramolecular hydrogen bond formation in ortho-dimethylaminomethylphenol. <i>Journal of Molecular Structure</i> , 2001, 559, 127-145.	3.6	40
8	Strong and weak effects caused by non covalent interactions between chloroform and selected electron donor molecules. <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 14223.	2.8	32
9	Cryospectroscopic and ab initio studies of haloform⋯trimethylamine H-bonded complexes. <i>Physical Chemistry Chemical Physics</i> , 2009, 11, 1551.	2.8	30
10	Comparative studies of blue shifting and red shifting effects in fluoroform and acetylene cryogenic solutions. <i>Journal of Molecular Structure</i> , 2004, 705, 49-61.	3.6	29
11	Gas phase FT-IR spectra and structure of aminoalcohols with intramolecular hydrogen bonds. <i>Vibrational Spectroscopy</i> , 1999, 20, 69-83.	2.2	28
12	FTIR study of the hydrogen bond symmetry in protonated homodimers of pyridine and collidine in solution. <i>Journal of Molecular Structure</i> , 2012, 1018, 39-44.	3.6	27
13	Infrared spectra and relative stability of the F ₃ CH/NH ₃ H-bonded complex in liquefied Xe. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2005, 61, 1595-1602.	3.9	24
14	Solvent effect on the blue shifted weakly H-bound F ₃ CH⋯FCD ₃ complex. <i>Journal of Molecular Structure</i> , 2008, 880, 64-68.	3.6	24
15	Dimerization and solvent-assisted proton dislocation in the low-barrier hydrogen bond of a Mannich base: a low-temperature NMR study. <i>Magnetic Resonance in Chemistry</i> , 2001, 39, S81-S90.	1.9	23
16	New look at the Badger-Bauer rule: Correlations of spectroscopic IR and NMR parameters with hydrogen bond energy and geometry. FHF complexes. <i>Journal of Molecular Structure</i> , 2018, 1164, 129-136.	3.6	22
17	Interplay of π -stacking and inter-stacking interactions in two-component crystals of neutral closed-shell aromatic compounds: periodic DFT study. <i>RSC Advances</i> , 2020, 10, 27899-27910.	3.6	20
18	Infrared studies of acetylene dissolved in liquefied Ar, Kr, N ₂ , CO, and CO ₂ . <i>Journal of Molecular Structure</i> , 2002, 614, 305-313.	3.6	18

#	ARTICLE	IF	CITATIONS
19	A cryosolution FTIR and ab initio study of the blue shifting C δ -H δ -F hydrogen bonded complexes F2ClCH δ -FCD3 and Cl2FCH δ -FCD3. Chemical Physics, 2008, 354, 71-79.	1.9	18
20	Effect of electrical and mechanical anharmonicity on vibrational spectra of H-bonded complexes: phenol-tctdot;B (B = acetonitrile, pyridine) systems. Journal of Molecular Structure, 1998, 448, 239-246.	3.6	17
21	Influence of hydrogen bonding on the conformation of ortho -aminomethylphenol. Journal of Molecular Structure, 2000, 523, 223-239.	3.6	15
22	Infrared studies of weak B δ -HCl (B = HCl, CO, Xe, N2) complex formation in solid Kr solutions. Journal of Molecular Structure, 1998, 448, 231-237.	3.6	14
23	Car-Parrinello Molecular Dynamics Study of a Blue-Shifted Intermolecular Weak-Hydrogen-Bond System. ChemPhysChem, 2005, 6, 1719-1724.	2.1	14
24	The model analysis of the effects of electrical and mechanical anharmonicity on vibrational spectra of H-bonded complexes. Journal of Molecular Structure, 2000, 552, 273-282.	3.6	13
25	Towards understanding the spectroscopic features of enflurane. The fundamental and overtone bands of CH stretching vibrations. Chemical Physics Letters, 2014, 604, 68-71.	2.6	13
26	Temperature dependence studies and model calculations of $\hat{\nu}_{1/2}(\text{OH})$ and $\hat{\nu}_{1/2}(\text{OD})$ band shapes of salicylaldehyde. Journal of Molecular Structure, 1996, 381, 141-148.	3.6	12
27	Higher-order transitions in the IR spectrum of the weak OC δ -HCl complex dissolved in liquid CO. Chemical Physics Letters, 2000, 325, 425-432.	2.6	12
28	Car-Parrinello Molecular Dynamics Study of the Blue-Shifted F3CH δ -... δ -... δ -FCD3 System in Liquid N2. ChemPhysChem, 2006, 7, 1221-1228.	2.1	12
29	The gas phase FTIR studies of chloroform+B and halothane+B (B=TMA, FCD3) mixtures. Journal of Molecular Structure, 2014, 1072, 32-37.	3.6	12
30	Theoretical vibrational study of hydrogen-bonded complexes : a simple anharmonic model. Journal De Chimie Physique Et De Physico-Chimie Biologique, 1992, 89, 607-613.	0.2	12
31	FTIR and Raman spectra of CH(D)FCICF2OCHF derivatives of enflurane. Experimental and ab initio study. Chemical Physics, 2015, 453-454, 26-34.	1.9	11
32	Temperature dependence of the $\hat{\nu}_{1/2s}(\text{OH})$ band shape of ortho-Mannich bases in gas phase and liquid xenon solution. Vibrational Spectroscopy, 1994, 7, 265-274.	2.2	10
33	Theoretical study of Hal3CH/F2CD2 (Hal=F,Cl) and F3CH/FH heterodimers with blue shifted hydrogen bonds. Chemical Physics, 2006, 327, 193-201.	1.9	10
34	Cooperative effects in blue-shifted hydrogen bonded cluster of from first principles simulations. Chemical Physics, 2009, 361, 129-136.	1.9	10
35	Cryospectroscopic and ab initio anharmonic studies of acetylene δ -trimethylamine H-bonded complex. Chemical Physics, 2010, 375, 92-100.	1.9	8
36	Infrared spectroscopy and ab initio study of hydrogen bonded Cl3CD δ -N(CH3)3 complex in the gas phase. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2015, 136, 95-99.	3.9	8

#	ARTICLE	IF	CITATIONS
37	The solubility of light olefins in liquid nitrogen: Part 2. <i>Cryogenics</i> , 1979, 19, 649-651.	1.7	7
38	Study on the solubility light alkanes in liquid nitrogen. <i>Cryogenics</i> , 1980, 20, 48-51.	1.7	7
39	Vibrational spectra of $\text{O}^{\delta-}\text{H}^{\delta+}\text{Cl}$ complex in Kr solutions at liquid to solid phase transition. <i>Journal of Molecular Structure</i> , 1999, 511-512, 233-240.	3.6	7
40	Degeneracy Lifting Effect in the FTIR Spectrum of Fluoroform Trapped in a Nitrogen Matrix. An Experimental and Carê“Parrinello Molecular Dynamics Study. <i>Journal of Physical Chemistry A</i> , 2016, 120, 3497-3503.	2.5	7
41	FTIR cryospectroscopic and ab initio studies of desfluraneâ€“dimethyl ether Hâ€“bonded complexes. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2017, 184, 163-168.	3.9	7
42	Stretching mode interaction of equal and unequal pairs of $\text{SeH}^{\delta+}$ and $\text{SeD}^{\delta-}$ defects in alkali halides. <i>Radiation Effects and Defects in Solids</i> , 1999, 151, 311-315.	1.2	6
43	$\text{CH}^{\delta+}\text{B}$ interactions in acetylene containing solutions: experimental and theoretical DFT studies. <i>Journal of Molecular Structure</i> , 2003, 645, 295-302.	3.6	6
44	The infrared study of fluoroformâ€“methyl fluoride mixtures in argon and nitrogen matrices. Evidence of nonlinear blue-shifting complex formation. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2018, 203, 185-194.	3.9	6
45	Effect of liquid to solid phase transition on rotational and vibrational broadening of vibrational bands of CF_3Cl in Xe solutions. <i>Vibrational Spectroscopy</i> , 2000, 24, 277-286.	2.2	5
46	Vibrational spectra and conformational analysis of desflurane. A cryosolution and ab initio study. <i>Chemical Physics Letters</i> , 2015, 637, 77-82.	2.6	5
47	Investigation of the IR spectra of weakly hydrogen-bonded complex $\text{Cl}_3\text{CH}^{\delta+}\text{O}(\text{CD}_3)_2$ in a cryosolution in liquid krypton. <i>Optics and Spectroscopy (English Translation of Optika I Spektroskopiya)</i> , 2016, 120, 242-249.	0.6	5
48	Conformational origin of temperature changes in the IR spectrum of isoflurane. A cryosolution and ab initio study. <i>Journal of Molecular Structure</i> , 2017, 1134, 374-380.	3.6	5
49	IR spectra of halothaneâ€“acetone complex in liquefied noble gases (Kr and Xe). <i>Optics and Spectroscopy (English Translation of Optika I Spektroskopiya)</i> , 2017, 123, 30-37.	0.6	5
50	Cryosolution infrared study of hydrogen bonded halothane acetylene complex. <i>Journal of Molecular Structure</i> , 2018, 1160, 434-439.	3.6	3
51	Noncovalent interactions between isoflurane and dimethyl ether. Spectroscopic evidence of trimer formation. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2019, 223, 117363.	3.9	3
52	The analysis of the band structure in $\hat{1}/2 ₁$ of $\text{HCN}\dots\text{HF}$ complex. <i>Journal De Chimie Physique Et De Physico-Chimie Biologique</i> , 1994, 91, 163-171.	0.2	3
53	Vibrational spectra of Freons dissolved in liquid argon. <i>Journal of Applied Spectroscopy</i> , 1985, 43, 999-1005.	0.7	2
54	An optical cryostat with a variable layer for examining the spectra of solutions in liquefied gases. <i>Journal of Applied Spectroscopy</i> , 1986, 45, 1000-1001.	0.7	2

#	ARTICLE	IF	CITATIONS
55	Infrared studies of CO ₂ doped Xe solutions in gas, liquid and solid phases. The fundamental $\hat{\nu}_3$ band and the Coriolis perturbed Fermi doublet ($\hat{\nu}_{21}+\hat{\nu}_{21}$, $\hat{\nu}_{21}+\hat{\nu}_{211}$). Journal of Molecular Structure, 2001, 596, 179-183.	3.6	2
56	IR spectra of CH ₃ F in liquid and solid noble gas solutions. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2000, 56, 1813-1818.	3.9	1
57	FTIR studies of HCl dissolved in liquid CO: Anharmonic effects in the weak O=C-HCl complex. Journal of Molecular Structure, 2007, 844-845, 64-69.	3.6	1
58	Spectroscopic evidence of reversible micro-crystallization of carbon dioxide dissolved in solid xenon. Chemical Physics Letters, 2013, 583, 54-59.	2.6	1
59	Evidence of noncovalent interactions between sevoflurane and dimethyl ether. FTIR cryospectroscopic and ab initio studies. Journal of Molecular Structure, 2020, 1221, 128852.	3.6	1
60	Spectral Diagnostics of the Dynamics of the Formation of a Homoconjugated Complex [HCN.H.NCH] ⁺ . Optics and Spectroscopy (English Translation of Optika i Spektroskopiya), 2020, 128, 467-469.	0.6	1
61	IR cryospectroscopic manifestation of complex formation between methoxyflurane and dimethyl ether in liquid Xe. Chemical Physics Letters, 2020, 742, 137134.	2.6	1
62	Application of the method of cryospectroscopy to the study of the molecular composition of gases. Journal of Applied Spectroscopy, 1992, 56, 224-230.	0.7	0
63	Quantitative analysis of sulfur hexafluoride for molecular impurities using IR absorption spectra of solutions in liquid argon. Journal of Applied Spectroscopy, 1992, 56, 360-364.	0.7	0
64	Shape of the IR bands of CH ₄ : The CH ₄ -Kr system in different phase states. Optics and Spectroscopy (English Translation of Optika i Spektroskopiya), 2000, 88, 169-175.	0.6	0
65	Cryospectroscopic evidence of trimer formation between halothane and trimethylamine, stabilized by hydrogen and halogen bonds. Journal of Molecular Structure, 2021, 1243, 130766.	3.6	0