

Konstantin S Rutkowski

List of Publications by Year in descending order

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#	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	A Cryosolution Infrared Study of the Complexes of Fluoroform with Ammonia and Pyridine: Evidence for a C-H...N Pseudo Blue-Shifting Hydrogen Bond. <i>Journal of Physical Chemistry A</i> , 2005, 109, 3038-3044.	2.5	63
3	Blue shifted F ₃ CH...FCD ₃ and Cl ₃ CH...FCD ₃ weakly H-bound complexes. <i>Cryospectroscopic and ab initio study. Chemical Physics</i> , 2005, 313, 225-243.	1.9	59
4	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
5	Gas phase IR spectra of systems with intramolecular hydrogen bonds. <i>Journal of Molecular Structure</i> , 1994, 322, 195-203.	3.6	34
6	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
7	Cryospectroscopic and ab initio studies of haloform...trimethylamine H-bonded complexes. <i>Physical Chemistry Chemical Physics</i> , 2009, 11, 1551.	2.8	30
8	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
9	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
10	The C-H...N versus O...N hydrogen bonding in the ammonia complexes with CH ₃ OH and CH ₃ SH. <i>Journal of Molecular Structure</i> , 2012, 1009, 96-102.	3.6	26
11	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
12	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
13	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
14	A cryosolution FTIR and ab initio study of the blue shifting C-H...F hydrogen bonded complexes F ₂ ClCH...FCD ₃ and Cl ₂ FCH...FCD ₃ . <i>Chemical Physics</i> , 2008, 354, 71-79.	1.9	18
15	Infrared studies of weak B...HCl (B = HCl, CO, Xe, N ₂) complex formation in solid Kr solutions. <i>Journal of Molecular Structure</i> , 1998, 448, 231-237.	3.6	14
16	Car-Parrinello Molecular Dynamics Study of a Blue-Shifted Intermolecular Weak-Hydrogen-Bond System. <i>ChemPhysChem</i> , 2005, 6, 1719-1724.	2.1	14
17	Towards understanding the spectroscopic features of enflurane. The fundamental and overtone bands of CH stretching vibrations. <i>Chemical Physics Letters</i> , 2014, 604, 68-71.	2.6	13
18	Temperature dependence studies and model calculations of $\hat{\nu}_{1/2}(\text{OH})$ and $\hat{\nu}_{1/2}(\text{OD})$ band shapes of salicylaldehyde. <i>Journal of Molecular Structure</i> , 1996, 381, 141-148.	3.6	12

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19	Higher-order transitions in the IR spectrum of the weak $\text{OC}\ddot{\text{C}}\text{HCl}$ complex dissolved in liquid CO . <i>Chemical Physics Letters</i> , 2000, 325, 425-432.	2.6	12
20	Carà€Parrinello Molecular Dynamics Study of the Blue-Shifted $\text{F}_3\text{CH}\ddot{\text{C}}\text{H}\ddot{\text{C}}\text{H}\ddot{\text{C}}\text{H}\text{F}$ System in Liquid N_2 . <i>ChemPhysChem</i> , 2006, 7, 1221-1228.	2.1	12
21	First-principles study of fluoroform adsorption on a hexagonal ice (0001) surface: weak hydrogen bondsâ€”strong structural effects. <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 14101.	2.8	12
22	The gas phase FTIR studies of chloroform+B and halothane+B (B=TMA, FCD3) mixtures. <i>Journal of Molecular Structure</i> , 2014, 1072, 32-37.	3.6	12
23	FTIR and Raman spectra of $\text{CH(D)FCICF}_2\text{OCHF}$ derivatives of enflurane. Experimental and ab initio study. <i>Chemical Physics</i> , 2015, 453-454, 26-34.	1.9	11
24	Temperature dependence of the $\hat{1}/_2\text{s(OH)}$ band shape of ortho-Mannich bases in gas phase and liquid xenon solution. <i>Vibrational Spectroscopy</i> , 1994, 7, 265-274.	2.2	10
25	Theoretical study of $\text{Hal}_3\text{CH/F}_2\text{CD}_2$ (Hal=F,Cl) and $\text{F}_3\text{CH/FH}$ heterodimers with blue shifted hydrogen bonds. <i>Chemical Physics</i> , 2006, 327, 193-201.	1.9	10
26	Cooperative effects in blue-shifted hydrogen bonded cluster of from first principles simulations. <i>Chemical Physics</i> , 2009, 361, 129-136.	1.9	10
27	Cryospectroscopic and ab initio anharmonic studies of acetyleneâ€”trimethylamine H-bonded complex. <i>Chemical Physics</i> , 2010, 375, 92-100.	1.9	8
28	Infrared spectroscopy and ab initio study of hydrogen bonded $\text{Cl}_3\text{CD}\ddot{\text{A}}\text{N(CH}_3)_3$ complex in the gas phase. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2015, 136, 95-99.	3.9	8
29	Vibrational spectra of $\text{OC}\ddot{\text{C}}\text{HCl}$ complex in Kr solutions at liquid to solid phase transition. <i>Journal of Molecular Structure</i> , 1999, 511-512, 233-240.	3.6	7
30	Degeneracy Lifting Effect in the FTIR Spectrum of Fluoroform Trapped in a Nitrogen Matrix. An Experimental and Carà€Parrinello Molecular Dynamics Study. <i>Journal of Physical Chemistry A</i> , 2016, 120, 3497-3503.	2.5	7
31	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
32	$\text{CH}\ddot{\text{C}}\text{H}\ddot{\text{B}}$ interactions in acetylene containing solutions: experimental and theoretical DFT studies. <i>Journal of Molecular Structure</i> , 2003, 645, 295-302.	3.6	6
33	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
34	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
35	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
36	Investigation of the IR spectra of weakly hydrogen-bonded complex $\text{Cl}_3\text{CH}\ddot{\text{C}}\text{H}\ddot{\text{O}}(\text{CD}_3)_2$ in a cryosolution in liquid krypton. <i>Optics and Spectroscopy (English Translation of Optika i Spektroskopiya)</i> , 2016, 120, 242-249.	0.6	5

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37	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
38	IR spectra of halothane-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
39	Spectroscopic manifestation of molecular rotation dynamics in dense media: CO fundamental band in liquid and solid CO-Kr and CO-Xe solutions. <i>Journal of Molecular Liquids</i> , 2001, 92, 251-261.	4.9	4
40	Spectroscopic evidence of weak complex formation between N,N-dimethylformamide and fluoroform in the gas phase. <i>Journal of Molecular Structure</i> , 2018, 1160, 328-332.	3.6	4
41	Evolution of IR spectra of a weakly-bound OCO^-HCl complex with increasing CO_2 density from the gas to liquid phase. <i>Journal of Molecular Structure</i> , 2001, 598, 205-211.	3.6	3
42	Cryosolution infrared study of hydrogen bonded halothane acetylene complex. <i>Journal of Molecular Structure</i> , 2018, 1160, 434-439.	3.6	3
43	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
44	Infrared studies of CO_2 doped Xe solutions in gas, liquid and solid phases. The fundamental ν_3 band and the Coriolis perturbed Fermi doublet ($\nu_1 + \nu_2$, $\nu_1 + \nu_2$). <i>Journal of Molecular Structure</i> , 2001, 596, 179-183.	3.6	2
45	Identification of H-Bonds in a Calcified Aortic Valve. <i>Technical Physics Letters</i> , 2019, 45, 927-929.	0.7	2
46	IR-IR double-resonance studies of vibrational relaxation of CD_3F in solid and liquid Xe, Kr, Ar solutions near the melting point. <i>Chemical Physics</i> , 1998, 237, 403-411.	1.9	1
47	IR spectra of CH_3F in liquid and solid noble gas solutions. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2000, 56, 1813-1818.	3.9	1
48	FTIR studies of HCl dissolved in liquid CO: Anharmonic effects in the weak OCO^-HCl complex. <i>Journal of Molecular Structure</i> , 2007, 844-845, 64-69.	3.6	1
49	Spectroscopic evidence of reversible micro-crystallization of carbon dioxide dissolved in solid xenon. <i>Chemical Physics Letters</i> , 2013, 583, 54-59.	2.6	1
50	Evidence of noncovalent interactions between sevoflurane and dimethyl ether. FTIR cryospectroscopic and ab initio studies. <i>Journal of Molecular Structure</i> , 2020, 1221, 128852.	3.6	1
51	Spectral Diagnostics of the Dynamics of the Formation of a Homoconjugated Complex $[\text{HCN} \cdot \text{H} \cdot \text{NCH}]^+$. <i>Optics and Spectroscopy (English Translation of Optika i Spektroskopiya)</i> , 2020, 128, 467-469.	0.6	1
52	IR cryospectroscopic manifestation of complex formation between methoxyflurane and dimethyl ether in liquid Xe. <i>Chemical Physics Letters</i> , 2020, 742, 137134.	2.6	1
53	Vibrational relaxation studies of systems with strong intermolecular interactions at low temperatures in the gas phase and in cryogenic solutions. <i>Advances in Molecular Relaxation and Interaction Processes</i> , 1982, 22, 223-231.	0.5	0
54	Shape of the IR bands of CH_4 : The CH_4 -Kr system in different phase states. <i>Optics and Spectroscopy (English Translation of Optika i Spektroskopiya)</i> , 2000, 88, 169-175.	0.6	0

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55	Cryospectroscopic evidence of trimer formation between halothane and trimethylamine, stabilized by hydrogen and halogen bonds. <i>Journal of Molecular Structure</i> , 2021, 1243, 130766.	3.6	0