

Mustafa Senyel

List of Publications by Year in descending order

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papers

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#	ARTICLE	IF	CITATIONS
1	FT-IR and NMR investigation of 1-phenylpiperazine: A combined experimental and theoretical study. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2007, 67, 793-801.	3.9	52
2	Computational study on favipiravir adsorption onto undoped- and silicon-decorated C60 fullerenes. <i>Journal of Theoretical and Computational Chemistry</i> , 2017, 16, 1750011.	1.8	25
3	Vibrational spectroscopic investigation and conformational analysis of 1-heptylamine: A comparative density functional study. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2013, 114, 668-680.	3.9	22
4	Theoretical and experimental vibrational spectroscopic study of 3-piperidino-propylamine. <i>Journal of Molecular Structure</i> , 2009, 923, 120-126.	3.6	19
5	Vibrational spectroscopic study on some Hofmann type clathrates: $M(1\text{-Phenylpiperazine})_2Ni(CN)_4 \cdot 2G$ (M=Ni, Co and Cd; G=aniline). <i>Journal of Molecular Structure</i> , 2009, 919, 41-46.	3.6	19
6	4-Mercaptophenylboronic acid: Conformation, FT-IR, Raman, OH stretching and theoretical studies. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2015, 144, 131-138.	3.9	19
7	FT-IR spectroscopic investigation of some Hofmann type complexes: $M(1\text{-phenylpiperazine})_2Ni(CN)_4$ (M=Ni, Co, Cd, Pd or Mn). <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2008, 70, 367-375.	3.9	16
8	Conformational analysis and vibrational spectroscopic investigation of 3-phenylpropylamine. <i>Vibrational Spectroscopy</i> , 2009, 50, 277-284.	2.2	15
9	1H , ^{13}C , ^{15}N NMR and $nJ(C, H)$ coupling constants investigation of 3-piperidino-propylamine: A combined experimental and theoretical study. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2008, 71, 830-834.	3.9	14
10	FT-IR and Raman spectroscopic and quantum chemical investigations of some metal halide complexes of 1-phenylpiperazine. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2012, 88, 144-155.	3.9	13
11	Title is missing!. <i>Journal of Inclusion Phenomena and Macrocyclic Chemistry</i> , 2001, 39, 175-180.	1.6	12
12	Title is missing!. <i>Journal of Inclusion Phenomena and Macrocyclic Chemistry</i> , 2003, 45, 129-137.	1.6	11
13	Vibrational spectra, DFT calculations, conformational stabilities and assignments of the fundamentals of the 1-butylpiperazine. <i>Journal of Molecular Structure</i> , 2016, 1122, 324-330.	3.6	10
14	FT-IR Spectroscopic Study on Some 4-(3-Cyclohexen-1-yl)pyridine Metal(II) Tetracyanonickelate Complexes. <i>Zeitschrift Fur Naturforschung - Section A Journal of Physical Sciences</i> , 2005, 60, 532-536.	1.5	9
15	Vibrational spectroscopic and thermal studies of some 3-phenylpropylamine complexes. <i>Vibrational Spectroscopy</i> , 2009, 51, 299-307.	2.2	9
16	Vibrational investigation of 1-cyclopentylpiperazine: A combined experimental and theoretical study. <i>Science China: Physics, Mechanics and Astronomy</i> , 2014, 57, 1654-1661.	5.1	9
17	One dimensional coordination polymers: Synthesis, crystal structures and spectroscopic properties. <i>Journal of Molecular Structure</i> , 2016, 1123, 327-334.	3.6	9
18	Syntheses, structural characterization and spectroscopic studies of cadmium(II)-metal(II) cyanide complexes with 4-(2-aminoethyl)pyridine. <i>Journal of Molecular Structure</i> , 2017, 1130, 80-88.	3.6	9

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19	Title is missing!. Journal of Inclusion Phenomena and Macrocyclic Chemistry, 2001, 39, 169-174.	1.6	8
20	Interaction between doped C ₆₀ fullerenes and piperazine-2,3,5,6-tetraone: DFT simulation. Main Group Metal Chemistry, 2018, 41, 63-66.	1.6	8
21	Experimental and theoretical NMR study of 4-(3-cyclohexen-1-yl)pyridine. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2008, 69, 1252-1256.	3.9	7
22	Molecular structure, NMR analyses, density functional theory and ab initio Hartree-Fock calculations of 3-phenylpropylamine. Comptes Rendus Chimie, 2009, 12, 808-815.	0.5	7
23	Synthesis, spectroscopic, thermal and structural properties of 4-(2-aminoethyl)pyridinium tetracyanometallate(II) complexes. Journal of Molecular Structure, 2017, 1136, 281-287.	3.6	7
24	One-dimensional coordination polymers with 4-(2-aminoethyl)pyridine: Synthesis, crystal structures and spectroscopic properties. Polyhedron, 2017, 123, 56-61.	2.2	6
25	FT-IR Spectroscopic Study on Some Hofmann-Td Type Complexes: Ni(4-Phenylpyridine) ₂ M(CN) ₄ (M = Cd) Tj ETQq _{1,1.5} 0.7843 ₁₄ rgBT	1.5	5
26	FT-Raman spectroscopic spectra on 3-phenylpropylamine inclusion compounds. Journal of Raman Spectroscopy, 2011, 42, 558-562.	2.5	5
27	DFT, FT-Raman and FT-IR investigations of 1-cyclobutylpiperazine. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2012, 90, 109-117.	3.9	5
28	Density functional theory study on the adsorption of valproic acid to doped fullerenes. Main Group Metal Chemistry, 2018, 41, 67-71.	1.6	5
29	Syntheses, crystal structures, spectroscopic properties and thermal decompositions of one dimensional coordination polymers with 4-(2-aminoethyl)pyridine and cyanide ligands: [M(1/4-4aepy) ₂ (H ₂ O) ₂][M' ₂ (CN) ₄] (M = Cu or Zn, M' = Ni or Pd). Journal of Molecular Structure, 2019, 1176, 641-649.	3.6	5
30	¹ H, ¹³ C NMR and nJCH coupling constants investigation of 4-Phenylpyridine: A combined experimental and theoretical study. Physics Letters, Section A: General, Atomic and Solid State Physics, 2007, 371, 300-306.	2.1	4
31	The synthesis, molecular structure, FT-IR and XRD spectroscopic investigation of 4-[(2-[(2-furylmethyl)imino]methyl)-4-methoxyphenoxy]methyl]benzonitrile: A comparative DFT study. Journal of Molecular Structure, 2011, 991, 12-17.	3.6	4
32	Conformation and NH stretching of 1,1-dihalogenoheptan-1-amines [CH ₃ (CH ₂) ₅ CX ₂ NH ₂ ; X=F, Cl or Br]: Halogen and solvent effects. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2015, 139, 171-178.	3.9	4
33	FT-IR spectroscopic investigation of some Hofmann type complexes: M(2-(1-Cyclohexenyl)ethylamine) ₂ Ni(CN) ₄ (M=Ni, Co). Journal of Molecular Structure, 2008, 892, 311-315.	3.6	3
34	Molecular structure, NMR analyses, density functional theory and ab initio Hartree-Fock calculations of 4,4'-diaminoctafluorobiphenyl. Journal of Molecular Structure, 2008, 891, 151-156.	3.6	3
35	Vibrational spectroscopic and conformational studies of 1-(4-pyridyl)piperazine. Chemical Papers, 2010, 64, .	2.2	3
36	Vibrational spectroscopic investigations of some Hofmann-Td-type complexes: Ni(2-(1-cyclohexenyl)ethylamine) ₂ M(CN) ₄ (M=Cd or Hg). Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2010, 76, 435-438.	3.9	3

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37	Vibrational spectroscopic study on some Hofmann-Td type clathrates: Ni(4-phenylpyridine) ₂ M(CN) ₄ ·2G (M=Cd or Hg, G=1,4-dioxane). <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2011, 78, 1487-1491.	3.9	3
38	Sensor application of doped C60 fullerenes in detection of 1-(3-trifluoromethylphenyl)piperazine as an alternative to ecstasy. <i>Main Group Metal Chemistry</i> , 2019, 42, 23-27.	1.6	3
39	Synthesis and vibrational spectroscopic analysis of some Hofmann type clathrates. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2011, 82, 383-388.	3.9	2
40	Vibrational spectroscopic study on some Hofmann type clathrates: M(2-(1-cyclohexenyl)ethylamine) ₂ Ni(CN) ₄ ·2benzene (M=Ni and Cd). <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2011, 79, 308-311.	3.9	1
41	Vibrational spectroscopic analysis of some Hofmann-Td type complexes. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2012, 97, 678-682.	3.9	1
42	Crystallographic, vibrational and DFT studies of 1-(2-hydroxy-4,5-dimethylphenyl)ethanone. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2015, 149, 762-770.	3.9	1
43	Molecular, vibrational and electronic structure of 4-bromo-2-halogenobenzaldehydes: Halogen and solvent effects. <i>Physical Sciences Reviews</i> , 2017, 2, .	0.8	1