

Mustafa Senyel

List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	Sensor application of doped C ₆₀ fullerenes in detection of 1-(3-trifluoromethylphenyl)piperazine as an alternative to ecstasy. Main Group Metal Chemistry, 2019, 42, 23-27.	1.6	3
2	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'CN ₄] (M= Cu or Zn, M'A= Ni or Pd). Journal of Molecular Structure, 2019, 1176, 641-649.	3.6	5
3	Density functional theory study on the adsorption of valproic acid to doped fullerenes. Main Group Metal Chemistry, 2018, 41, 67-71.	1.6	5
4	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
5	Computational study on favipiravir adsorption onto undoped- and silicon-decorated C ₆₀ fullerenes. Journal of Theoretical and Computational Chemistry, 2017, 16, 1750011.	1.8	25
6	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
7	One-dimensional coordination polymers with 4-(2-aminoethyl)pyridine: Synthesis, crystal structures and spectroscopic properties. Polyhedron, 2017, 123, 56-61.	2.2	6
8	Syntheses, structural characterization and spectroscopic studies of cadmium(II)-metal(II) cyanide complexes with 4-(2-aminoethyl)pyridine. Journal of Molecular Structure, 2017, 1130, 80-88.	3.6	9
9	Molecular, vibrational and electronic structure of 4-bromo-2-halogenobenzaldehydes: Halogen and solvent effects. Physical Sciences Reviews, 2017, 2, .	0.8	1
10	Vibrational spectra, DFT calculations, conformational stabilities and assignments of the fundamentals of the 1-butylpiperazine. Journal of Molecular Structure, 2016, 1122, 324-330.	3.6	10
11	One dimensional coordination polymers: Synthesis, crystal structures and spectroscopic properties. Journal of Molecular Structure, 2016, 1123, 327-334.	3.6	9
12	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
13	4-Mercaptophenylboronic acid: Conformation, FT-IR, Raman, OH stretching and theoretical studies. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2015, 144, 131-138.	3.9	19
14	Crystallographic, vibrational and DFT studies of 1-(2-hydroxy-4,5-dimethylphenyl)ethanone. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2015, 149, 762-770.	3.9	1
15	Vibrational investigation of 1-cyclopentylpiperazine: A combined experimental and theoretical study. Science China: Physics, Mechanics and Astronomy, 2014, 57, 1654-1661.	5.1	9
16	Vibrational spectroscopic investigation and conformational analysis of 1-heptylamine: A comparative density functional study. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2013, 114, 668-680.	3.9	22
17	Vibrational spectroscopic analysis of some Hofmann-T _d type complexes. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2012, 97, 678-682.	3.9	1
18	FT-IR and Raman spectroscopic and quantum chemical investigations of some metal halide complexes of 1-phenylpiperazine. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2012, 88, 144-155.	3.9	13

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19	DFT, FT-Raman and FT-IR investigations of 1-cyclobutylpiperazine. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2012, 90, 109-117.	3.9	5
20	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
21	Vibrational spectroscopic study on some Hofmann-Td type clathrates: Ni(4-phenylpyridine)2M(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
22	FT-Raman spectroscopic spectra on 3-phenylpropylamine inclusion compounds. <i>Journal of Raman Spectroscopy</i> , 2011, 42, 558-562.	2.5	5
23	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. <i>Journal of Molecular Structure</i> , 2011, 991, 12-17.	3.6	4
24	Vibrational spectroscopic study on some Hofmann type clathrates: M(2-(1-cyclohexenyl)ethylamine)2Ni(CN)4·2benzene (M=Ni and Cd). <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2011, 79, 308-311.	3.9	1
25	Vibrational spectroscopic and conformational studies of 1-(4-pyridyl)piperazine. <i>Chemical Papers</i> , 2010, 64, .	2.2	3
26	Vibrational spectroscopic investigations of some Hofmann-Td-type complexes: Ni(2-(1-cyclohexenyl)ethylamine)2M(CN)4 (M=Cd or Hg). <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2010, 76, 435-438.	3.9	3
27	Theoretical and experimental vibrational spectroscopic study of 3-piperidino-propylamine. <i>Journal of Molecular Structure</i> , 2009, 923, 120-126.	3.6	19
28	Conformational analysis and vibrational spectroscopic investigation of 3-phenylpropylamine. <i>Vibrational Spectroscopy</i> , 2009, 50, 277-284.	2.2	15
29	Vibrational spectroscopic study on some Hofmann type clathrates: M(1-Phenylpiperazine)2Ni(CN)4·2G (M=Ni, Co and Cd; G=aniline). <i>Journal of Molecular Structure</i> , 2009, 919, 41-46.	3.6	19
30	Vibrational spectroscopic and thermal studies of some 3-phenylpropylamine complexes. <i>Vibrational Spectroscopy</i> , 2009, 51, 299-307.	2.2	9
31	Molecular structure, NMR analyses, density functional theory and ab initio Hartree-Fock calculations of 3-phenylpropylamine. <i>Comptes Rendus Chimie</i> , 2009, 12, 808-815.	0.5	7
32	FT-IR spectroscopic investigation of some Hofmann type complexes: M(2-(1-Cyclohexenyl)ethylamine)2Ni(CN)4 (M=Ni, Co). <i>Journal of Molecular Structure</i> , 2008, 892, 311-315.	3.6	3
33	Experimental and theoretical NMR study of 4-(3-cyclohexen-1-yl)pyridine. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2008, 69, 1252-1256.	3.9	7
34	FT-IR spectroscopic investigation of some Hofmann type complexes: M(1-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
35	1H, 13C, 15N 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
36	Molecular structure, NMR analyses, density functional theory and ab initio Hartree-Fock calculations of 4,4-diaminoctafluorobiphenyl. <i>Journal of Molecular Structure</i> , 2008, 891, 151-156.	3.6	3

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37	1H, 13C NMR and η JCH 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
38	FT-IR and NMR investigation of 1-phenylpiperazine: A combined experimental and theoretical study. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2007, 67, 793-801.	3.9	52
39	FT-IR Spectroscopic Study on Some 4-(3-Cyclohexen-1-yl)pyridine Metal(II) Tetracyanonickelate Complexes. Zeitschrift Fur Naturforschung - Section A Journal of Physical Sciences, 2005, 60, 532-536.	1.5	9
40	FT-IR Spectroscopic Study on Some Hofmann-Td Type Complexes: Ni(4-Phenylpyridine)2M(CN)4 (M = Cd) Tj ETQq0,0,0 rgBT _{1.5} /Overlock 1		
41	Title is missing!. Journal of Inclusion Phenomena and Macrocyclic Chemistry, 2003, 45, 129-137.	1.6	11
42	Title is missing!. Journal of Inclusion Phenomena and Macrocyclic Chemistry, 2001, 39, 169-174.	1.6	8
43	Title is missing!. Journal of Inclusion Phenomena and Macrocyclic Chemistry, 2001, 39, 175-180.	1.6	12