

Poul Erik Hansen

List of Publications by Year in descending order

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160
papers

4,434
citations

117625

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166
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166
docs citations

166
times ranked

3860
citing authors

#	ARTICLE	IF	CITATIONS
1	Isotope effects in nuclear shielding. <i>Progress in Nuclear Magnetic Resonance Spectroscopy</i> , 1988, 20, 207-255.	7.5	250
2	Bromophenols in Marine Algae and Their Bioactivities. <i>Marine Drugs</i> , 2011, 9, 1273-1292.	4.6	208
3	Pharmacological Profile of Xanthohumol, a Prenylated Flavonoid from Hops (<i>Humulus lupulus</i>). <i>Molecules</i> , 2015, 20, 754-779.	3.8	174
4	¹³ C NMR of polycyclic aromatic compounds. A review. <i>Magnetic Resonance in Chemistry</i> , 1979, 12, 109-142.	0.7	156
5	Isotope Effects on Nuclear Shielding. <i>Annual Reports on NMR Spectroscopy</i> , 1984, 15, 105-234.	1.5	151
6	The influence of organic matter on sorption and fate of glyphosate in soil – Comparing different soils and humic substances. <i>Environmental Pollution</i> , 2009, 157, 2865-2870.	7.5	104
7	NMR and IR Investigations of Strong Intramolecular Hydrogen Bonds. <i>Molecules</i> , 2017, 22, 552.	3.8	95
8	Magnetic susceptibility: Solutions, emulsions, and cells. <i>Concepts in Magnetic Resonance</i> , 2003, 18A, 56-71.	1.3	87
9	Unraveling the Electronic and Vibrational Contributions to Deuterium Isotope Effects on ¹³ C Chemical Shifts Using ab Initio Model Calculations. Analysis of the Observed Isotope Effects on Sterically Perturbed Intramolecular Hydrogen-Bonded-Hydroxy Acyl Aromatics. <i>Journal of the American Chemical Society</i> , 1998, 120, 9063-9069.	13.7	84
10	Deuterium-Induced Isotope Effects on ¹³ C Chemical Shifts as a Probe for Tautomerism in Enolic ^{1,2} -Diketones. <i>Magnetic Resonance in Chemistry</i> , 1996, 34, 467-478.	1.9	77
11	Isotope Effects on Chemical Shifts as an Analytical Tool in Structural Studies of Intramolecular Hydrogen Bonded Compounds. <i>Current Organic Chemistry</i> , 2000, 4, 19-54.	1.6	73
12	Deuterium isotope effects on ¹³ C chemical shifts of intramolecularly hydrogen-bonded Schiff bases. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1999, , 2809-2817.	0.9	70
13	Sorption of Polycyclic Aromatic Compounds to Humic Acid As Studied by High-Performance Liquid Chromatography. <i>Environmental Science & Technology</i> , 1997, 31, 1102-1108.	10.0	66
14	Phosphate pool dynamics in the arbuscular mycorrhizal fungus <i>Glomus intraradices</i> studied by in vivo ³¹ P NMR spectroscopy. <i>New Phytologist</i> , 2004, 162, 783-794.	7.3	66
15	Deuterium and ¹⁸ O isotope effects on ¹³ C chemical shifts of sterically hindered and/or intramolecularly hydrogen-bonded-hydroxy acyl aromatics. <i>Magnetic Resonance in Chemistry</i> , 1994, 32, 399-408.	1.9	61
16	Characterization of Salt Bridges to Lysines in the Protein G B1 Domain. <i>Journal of the American Chemical Society</i> , 2009, 131, 4674-4684.	13.7	61
17	Bilirubin Acidity. Titrimetric and ¹³ C NMR Studies.. <i>Acta Chemica Scandinavica</i> , 1979, 33b, 281-293.	0.7	59
18	Deuterium isotope effects on the ¹³ C nuclear shielding of intramolecularly hydrogen-bonded systems. <i>Magnetic Resonance in Chemistry</i> , 1986, 24, 903-910.	1.9	58

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19	Substituent effects on deuterium isotope effects on nuclear shielding of intramolecularly hydrogen-bonded aromatic ketones, aldehydes and esters. <i>Magnetic Resonance in Chemistry</i> , 1993, 31, 23-37.	1.9	58
20	Studies based on deuterium isotope effect on ¹³ C chemical shifts. <i>Progress in Nuclear Magnetic Resonance Spectroscopy</i> , 2004, 45, 1-29.	7.5	55
21	Deuterium isotope effects on ¹³ C nuclear shielding as a measure of tautomeric equilibria. <i>Magnetic Resonance in Chemistry</i> , 1982, 18, 58-61.	0.7	50
22	³¹ P NMR for the study of P metabolism and translocation in arbuscular mycorrhizal fungi. <i>Plant and Soil</i> , 2000, 226, 245-253.	3.7	50
23	Oneâ€bond deuterium isotope effects on ¹⁵ N chemical shifts in Schiff bases. <i>Zeitschrift Fur Elektrotechnik Und Elektrochemie</i> , 1998, 102, 410-413.	0.9	49
24	A ¹⁹ F NMR study of Câ€“â€“ halogen bonding. <i>Chemical Physics</i> , 2011, 381, 5-10.	1.9	47
25	Isotope effects on chemical shifts of proteins and peptides. <i>Magnetic Resonance in Chemistry</i> , 2000, 38, 1-10.	1.9	45
26	Deuterium Isotope Effects on ¹³ C and ¹⁵ N Nuclear Shielding in Intramolecularly Hydrogen-Bonded Compounds. Investigation of Enamine Derivatives.. <i>Acta Chemica Scandinavica</i> , 1990, 44, 826-832.	0.7	44
27	Tautomerism of enolic triacetylmethane, 2-acyl-1,3-cycloalkanediones, 5-acyl Meldrumâ€™s acids and 5-acyl-1,3-dimethylbarbituric acids studied by means of deuterium isotope effects on ¹³ C chemical shifts. <i>Magnetic Resonance in Chemistry</i> , 1998, 36, 315-324.	1.9	42
28	Deuterium isotope effects on ¹³ C nuclear shielding of amino and acetamido compounds. Tautomerism and intramolecular hydrogen bonding. <i>Magnetic Resonance in Chemistry</i> , 1992, 30, 786-795.	1.9	40
29	Deuterium Isotope Effects on ¹³ C Chemical Shifts of o-Hydroxyacyl Aromatics. Intramolecular Hydrogen Bonding. <i>Magnetic Resonance in Chemistry</i> , 1997, 35, 520-528.	1.9	40
30	Hydrogen bonding and tautomerism studied by isotope effects on chemical shifts. <i>Journal of Molecular Structure</i> , 1994, 321, 79-87.	3.6	39
31	Schiff bases of gossypol: an NMR and DFT study. <i>Magnetic Resonance in Chemistry</i> , 2005, 43, 302-308.	1.9	39
32	Deuterium isotope effects on ¹³ C and ¹⁵ N nuclear shielding in o-hydroxyazo dyes. <i>Magnetic Resonance in Chemistry</i> , 1984, 22, 569-572.	0.7	35
33	Variable temperature ¹ H and ¹³ C NMR spectroscopic investigation of the enolâ€“enethiol tautomerism of ^{Î²} -thioxoketones. Isotope effects due to deuterium chelation. <i>Journal of Molecular Structure</i> , 2000, 552, 45-62.	3.6	35
34	Steric compression and twist in o-hydroxy acyl aromatics with intramolecular hydrogen bonding. <i>Journal of Molecular Structure</i> , 2005, 749, 155-168.	3.6	35
35	Photovoltaic Performance and Characteristics of Dyeâ€Sensitized Solar Cells Prepared with the N719 Thermal Degradation Products [Ru(LH) ₂ (NCS)(4-tert-butylpyridine)][N(Bu) ₄] and [Ru(LH) ₂ (NCS)(1-methylbenzimidazole)][N(Bu) ₄]. <i>European Journal of Inorganic Chemistry</i> , 2011, 2011, 2533-2539.	2.0	35
36	Hydrogen bonding monitored by deuterium isotope effects on carbonyl ¹³ C chemical shift in BPTI: intra-residue hydrogen bonds in antiparallel ^{Î²} -sheet. <i>International Journal of Biological Macromolecules</i> , 1991, 13, 2-8.	7.5	34

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37	Variable-temperature NMR studies of 2,6-dihydroxy acylaromatic compounds. Deuterium isotope effects on chemical shifts, isotopic perturbation of equilibrium and barriers to rotation. <i>Magnetic Resonance in Chemistry</i> , 1993, 31, 893-902.	1.9	34
38	Ab Initio Calculations of Deuterium Isotope Effects on Hydrogen and Nitrogen Nuclear Magnetic Shielding in the Hydrated Ammonium Ion.. <i>Acta Chemica Scandinavica</i> , 1992, 46, 1065-1071.	0.7	34
39	¹⁷ O chemical shifts and deuterium isotope effects on ¹³ C chemical shifts of intramolecularly hydrogen-bonded compounds. <i>Magnetic Resonance in Chemistry</i> , 1998, 36, 921-928.	1.9	32
40	Strong intramolecular hydrogen bonding involving nitro- and acetyl groups. Deuterium isotope effects on chemical shifts. <i>Journal of Molecular Structure</i> , 2006, 789, 81-91.	3.6	32
41	Effect of Different Humic Substances on the Fate of Diuron and Its Main Metabolite 3,4-Dichloroaniline in Soil. <i>Environmental Science & Technology</i> , 2008, 42, 8687-8691.	10.0	32
42	Cytotoxic Geranylated Xanthenes and O-Alkylated Derivatives of .ALPHA.-Mangostin. <i>Chemical and Pharmaceutical Bulletin</i> , 2009, 57, 830-834.	1.3	32
43	Isotope effect on chemical shifts in hydrogen-bonded systems. <i>Journal of Labelled Compounds and Radiopharmaceuticals</i> , 2007, 50, 967-981.	1.0	30
44	OH stretching frequencies in systems with intramolecular hydrogen bonds: Harmonic and anharmonic analyses. <i>Chemical Physics</i> , 2011, 389, 107-115.	1.9	30
45	Deuterium isotope effects on ¹⁵ N, ¹³ C and ¹ H chemical shifts of proton sponges. <i>Journal of Molecular Structure</i> , 2002, 615, 121-140.	3.6	29
46	Binding of Topotecan to a Nicked DNA Oligomer in Solution. <i>Chemistry - A European Journal</i> , 2008, 14, 2788-2794.	3.3	29
47	Intramolecular hydrogen bonding of the enol forms of ¹² -ketoamides and ¹² -kethioamides. Deuterium isotope effects on ¹³ C chemical shifts. <i>Journal of Molecular Structure</i> , 1996, 378, 45-59.	3.6	28
48	Density Functional Theory Study of Intramolecular Hydrogen Bonding and Proton Transfer in <i>o</i> -Hydroxyaryl Ketimines. <i>Journal of Physical Chemistry A</i> , 2008, 112, 3478-3485.	2.5	28
49	CH, CD, CC and HH coupling constants in isotopically enriched cyclobutene. <i>Magnetic Resonance in Chemistry</i> , 1981, 15, 288-293.	0.7	27
50	Primary tritium and deuterium isotope effects on chemical shifts of compounds having an intramolecular hydrogen bond. <i>Magnetic Resonance in Chemistry</i> , 2000, 38, 525-535.	1.9	27
51	Progress of Bromophenols in Marine Algae from 2011 to 2020: Structure, Bioactivities, and Applications. <i>Marine Drugs</i> , 2020, 18, 411.	4.6	27
52	Temperature coefficient of NH chemical shifts of thioamides and amides in relation to structure. <i>Journal of Molecular Structure</i> , 2004, 700, 91-103.	3.6	26
53	Deuterium isotope effects on ¹⁵ N backbone chemical shifts in proteins. <i>Journal of Biomolecular NMR</i> , 2009, 44, 119-126.	2.8	26
54	Theoretical and NMR Studies of Deuterium Isotopic Perturbation of Hydrogen Bonding in Symmetrical Dihydroxy Compounds. <i>Journal of Organic Chemistry</i> , 2010, 75, 1331-1342.	3.2	26

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55	Isotope Effects on Chemical Shifts in the Study of Intramolecular Hydrogen Bonds. <i>Molecules</i> , 2015, 20, 2405-2424.	3.8	26
56	Intramolecular hydrogen bonding and tautomerism of acylpyran-2,4-diones, -2,4,6-triones and acylpyridinediones and benzannelated derivatives. Deuterium isotope effects on ¹³ C NMR chemical shifts. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1995, , 1901.	0.9	25
57	Sorption of Polycyclic Aromatic Compounds to Humic and Fulvic Acid HPLC Column Materials. <i>Journal of Environmental Quality</i> , 2001, 30, 526-537.	2.0	25
58	Multiple Binding Modes of the Camptothecin Family to DNA Oligomers. <i>Chemistry - A European Journal</i> , 2004, 10, 5776-5787.	3.3	25
59	Variable-temperature NMR study of the enol forms of benzoylacetones. <i>Magnetic Resonance in Chemistry</i> , 2005, 43, 992-998.	1.9	25
60	The structure of the phototransformation product of monothiodibenzoylmethane. <i>Chemical Physics Letters</i> , 2001, 350, 502-508.	2.6	24
61	Solution and solid state ¹³ C NMR and X-ray studies of genistein complexes with amines. Potential biological function of the C-7, C-5, and C-4-OH groups. <i>Organic and Biomolecular Chemistry</i> , 2003, 1, 3578-3585.	2.8	24
62	Long-Range Intrinsic and Equilibrium Deuterium Isotope Effects on ¹⁹ F Chemical Shifts.. <i>Acta Chemica Scandinavica</i> , 1997, 51, 881-888.	0.7	24
63	Substituent Conformational effects in vicinal ¹³ C- ¹³ C spin-spin coupling constants. <i>Magnetic Resonance in Chemistry</i> , 1979, 12, 169-173.	0.7	23
64	Deuterium Isotope Effects on ¹³ C Chemical Shifts of Enaminones.. <i>Acta Chemica Scandinavica</i> , 1997, 51, 1016-1023.	0.7	23
65	On prediction of OH stretching frequencies in intramolecularly hydrogen bonded systems. <i>Journal of Molecular Structure</i> , 2012, 1018, 8-13.	3.6	22
66	Structural Studies of ¹² C-Diketones and Their Implications on Biological Effects. <i>Pharmaceuticals</i> , 2021, 14, 1189.	3.8	22
67	Carbon-carbon coupling constants of 1-phenylazo-2-naphthol and 2-phenylazo-1-naphthol obtained by the SEMINA-1 technique. <i>Magnetic Resonance in Chemistry</i> , 1986, 24, 772-776.	1.9	21
68	Deuterium isotope effects on ¹⁷ O chemical shifts of intramolecularly hydrogen bonded systems. <i>Journal of Molecular Structure</i> , 1999, 509, 171-181.	3.6	21
69	Thioacetylacetone: Structural and Vibrational Assignments. <i>ChemPhysChem</i> , 2004, 5, 495-502.	2.1	21
70	Characterisation of the PT-form of o-hydroxy acylaromatic Schiff bases by NMR spectroscopy and DFT calculations. <i>Journal of Molecular Structure</i> , 2004, 707, 69-75.	3.6	21
71	Detection of salt bridges to lysines in solution in barnase. <i>Chemical Communications</i> , 2013, 49, 9824-9826.	4.1	21
72	Deuterium isotope effects on ¹⁴ N and ¹⁵ N nuclear shielding in simple nitrogen-containing compounds. <i>Magnetic Resonance in Chemistry</i> , 1985, 23, 973-976.	1.9	20

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73	Trichloromethyl compounds – Natural background concentrations and fates within and below coniferous forests. <i>Science of the Total Environment</i> , 2010, 408, 6223-6234.	8.0	20
74	Ab Initio Calculations of Deuterium Isotope Effects on Chemical Shifts of Salt-Bridged Lysines. <i>Journal of Physical Chemistry B</i> , 2011, 115, 3208-3215.	2.6	20
75	A Reinvestigation of One-Bond Deuterium Isotope Effects on Nitrogen and on Proton Nuclear Shielding for the Ammonium Ion.. <i>Acta Chemica Scandinavica</i> , 1989, 43, 222-232.	0.7	20
76	Reverse quantitative structure–activity relationship for modelling the sorption of esfenvalerate to dissolved organic matter. <i>Chemosphere</i> , 2002, 49, 1317-1325.	8.2	19
77	Preparation and structural characterization of a new class of stable thioketones: ortho-hydroxythioacetophenones. <i>Tetrahedron Letters</i> , 2006, 47, 8433-8435.	1.4	19
78	Intramolecular hydrogen bonding of novelo-hydroxythioacetophenones and related compounds evaluated by deuterium isotope effects on ¹³ C chemical shifts. <i>Magnetic Resonance in Chemistry</i> , 2007, 45, 245-252.	1.9	19
79	Long-range deuterium isotope effects in tautomeric α -thioxoketones. ¹ H and ¹³ C NMR study. <i>Journal of Physical Organic Chemistry</i> , 1991, 4, 225-232.	1.9	18
80	Intramolecular hydrogen bonding in 8-quinolinol N-oxides, quinaldinic acid N-oxides and quinoline-2-carboxamide N-oxide. Deuterium isotope effects on ¹³ C chemical shifts. <i>Journal of Molecular Structure</i> , 1997, 436-437, 189-199.	3.6	18
81	A spectrochemometric approach to tautomerism and hydrogen-bonding in 3-acyltetronic acids. <i>Journal of Molecular Structure</i> , 2006, 790, 80-88.	3.6	18
82	Long-Range Deuterium Isotope Effects on ¹³ C Chemical Shifts of Intramolecularly Hydrogen-Bonded N-Substituted 3-(Cycloamine)thiopropionamides or Amides: – A Case of Electric Field Effects. <i>Journal of Organic Chemistry</i> , 2007, 72, 4108-4116.	3.2	18
83	Factor Analysis of Deuterium Isotope Effects on ¹³ C NMR Chemical Shifts in Schiff Bases. <i>Chemistry - A European Journal</i> , 2005, 11, 4758-4766.	3.3	17
84	Conformational and tautomeric eccentricities of 2-acetyl-1,8-dihydroxynaphthalenes. <i>Magnetic Resonance in Chemistry</i> , 2007, 45, 106-117.	1.9	17
85	Long-range deuterium isotope effects on ¹³ C chemical shifts of intramolecularly hydrogen-bonded compounds. <i>Purpurogallins</i> . <i>Magnetic Resonance in Chemistry</i> , 1993, 31, 71-74.	1.9	16
86	NMR of a series of novel hydroxyflavothiones. <i>Magnetic Resonance in Chemistry</i> , 2009, 47, 1043-1054.	1.9	16
87	Deuterium Isotope Effects on ¹³ C-NMR Chemical Shifts of 10-Hydroxybenzo[h]quinolines. <i>Molecules</i> , 2013, 18, 4544-4560.	3.8	15
88	Structural studies on Mannich bases of 2-Hydroxy-3,4,5,6-tetrachlorobenzene. An UV, IR, NMR and DFT study. A mini-review. <i>Journal of Molecular Structure</i> , 2016, 1119, 235-239.	3.6	15
89	Computational Prediction of ¹ H and ¹³ C NMR Chemical Shifts for Protonated Alkylpyrroles: Electron Correlation and Not Solvation is the Salvation. <i>ChemPhysChem</i> , 2019, 20, 78-91.	2.1	15
90	A Spectroscopic Overview of Intramolecular Hydrogen Bonds of NH– O,S,N Type. <i>Molecules</i> , 2021, 26, 2409.	3.8	15

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91	Deuterium isotope effects on ¹³ C chemical shifts of nitromalonamide. <i>Magnetic Resonance in Chemistry</i> , 2008, 46, 726-729.	1.9	14
92	The effect of 4-tert-butylpyridine and Li ⁺ on the thermal degradation of TiO ₂ -bound ruthenium dye N719. <i>Solar Energy</i> , 2013, 88, 23-30.	6.1	14
93	Intramolecular Hydrogen Bonds in Normal and Sterically Compressed o-Hydroxy Aromatic Aldehydes. Isotope Effects on Chemical Shifts and Hydrogen Bond Strength. <i>Molecules</i> , 2019, 24, 4533.	3.8	14
94	A Simple, Effective, Green Method for the Regioselective 3-Acylation of Unprotected Indoles. <i>Molecules</i> , 2015, 20, 19605-19619.	3.8	13
95	A concept for stimulated proton transfer in 1-(phenyldiazenyl)naphthalen-2-ols. <i>Dyes and Pigments</i> , 2018, 156, 91-99.	3.7	13
96	Isotopic Perturbation of Equilibrium in 2,6-Dihydroxybenzoyl Compounds. A ¹³ C and ¹ H NMR Investigation.. <i>Acta Chemica Scandinavica</i> , 1988, 42b, 423-432.	0.7	13
97	Deuterium Isotope Effects on Carbonyl Carbon Chemical Shifts of BPTI. Hydrogen Bonding and Structure Determination in Proteins.. <i>Acta Chemica Scandinavica</i> , 1989, 43, 710-712.	0.7	13
98	Reactivity and diastereoselectivity of Michael additions of amines to achiral $\hat{1}\pm, \hat{1}^2$ -unsaturated thioamides. <i>Tetrahedron</i> , 2001, 57, 8705-8718.	1.9	12
99	1,1- $\hat{1}\pm, \hat{1}\pm$ - $\hat{1}\pm$ - $\hat{1}\pm$ -(2,4,6-Trihydroxybenzene-1,3,5-triyl)triethanone tautomerism revisited. <i>Tetrahedron Letters</i> , 2014, 55, 354-357.	1.4	12
100	The possible tautomerism of the potential rotary switch 2-(2-(2-Hydroxy-4-nitrophenyl)hydrazono)-1-phenylbutane-1,3-dione. <i>Dyes and Pigments</i> , 2017, 144, 249-261.	3.7	12
101	Cellular Uptake and Intracellular Phosphorylation of GS-441524: Implications for Its Effectiveness against COVID-19. <i>Viruses</i> , 2021, 13, 1369.	3.3	12
102	NH Stretching Frequencies of Intramolecularly Hydrogen-Bonded Systems: An Experimental and Theoretical Study. <i>Molecules</i> , 2021, 26, 7651.	3.8	12
103	Ab initio calculations of external charge effects on the isotropic ¹³ C, ¹⁵ N and ¹⁷ O nuclear shieldings of amides. <i>Chemical Physics Letters</i> , 1994, 224, 275-282.	2.6	11
104	Deuterium isotope effects on ¹³ C and ¹⁵ N chemical shifts of intramolecularly hydrogen-bonded enamino-carbonyl derivatives of Meldrum's and Tetronic acid. <i>Journal of Molecular Structure</i> , 2010, 976, 377-391.	3.6	11
105	Dye-sensitized solar cells and complexes between pyridines and iodines. A NMR, IR and DFT study. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2012, 98, 247-251.	3.9	11
106	Arresting consecutive steps of a photochromic reaction: studies of $\hat{1}^2$ -thioxoketones combining laser photolysis with NMR detection. <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 9128-9137.	2.8	11
107	A Reinvestigation of the Ionic Liquid Diisopropylethylammonium Formate by NMR and DFT Methods. <i>Journal of Physical Chemistry B</i> , 2016, 120, 11279-11286.	2.6	11
108	Molecular structure and intramolecular hydrogen bond strength of 3-methyl-4-amino-3-penten-2-one and its N Me and N-Ph substitutions by experimental and theoretical methods. <i>Journal of Molecular Structure</i> , 2019, 1184, 233-245.	3.6	11

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109	2DeltaH(D) and 1DeltaN(D) Isotope Effects on Nuclear Shielding of Ammonium Ions in Complexes with Crown Ethers and Cryptands.. Acta Chemica Scandinavica, 1993, 47, 777-788.	0.7	11
110	The role of ring substituents on hydrogen bonding of 5-cyano-2-hydroxyacetophenone and 2-hydroxy-4-methoxy-5-nitroacetophenone in the ground and excited states. Journal of Molecular Structure, 2007, 844-845, 77-82.	3.6	10
111	Photochromism in p-methylbenzoylthioacetone and related \hat{I}^2 -thioxoketones. Chemical Physics, 2007, 338, 11-22.	1.9	10
112	Methodological problems in determining TCAA in soils—the discovery of novel natural trichloroacetyl containing compounds and their interference with a common method for determining TCAA in soil and vegetation. Journal of Environmental Monitoring, 2010, 12, 672-680.	2.1	10
113	A new chromanone acid from the bark of Calophyllum dryobalanoides. Phytochemistry Letters, 2012, 5, 287-291.	1.2	10
114	Determination of the tautomeric equilibria of pyridoyl benzoyl \hat{I}^2 -diketones in the liquid and solid state through the use of deuterium isotope effects on ^1H and ^{13}C NMR chemical shifts and spin coupling constants. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2015, 136, 107-112.	3.9	10
115	Ring current and anisotropy effects on OH chemical shifts in resonance-assisted intramolecular H-bonds. Tetrahedron Letters, 2018, 59, 2288-2292.	1.4	10
116	NMR of Natural Products as Potential Drugs. Molecules, 2021, 26, 3763.	3.8	10
117	^{13}C -NMR Chemical Shift Databases as a Quick Tool to Evaluate Structural Models of Humic Substances—!2010-01-15—!2010-02-02—!2010-06-18—!. The Open Magnetic Resonance Journal, 2010, 3, 96-105.	0.5	10
118	New deuterium isotope effects on ^{13}C and ^{19}F chemical shifts across intramolecular hydrogen bonds of non-resonance assisted systems. Tetrahedron Letters, 2005, 46, 839-842.	1.4	9
119	Photochromism and polarization spectroscopy of p-methyl(thiobenzoyl)acetone. Chemical Physics, 2006, 328, 205-215.	1.9	9
120	Oliveridepsidones A— $^{\text{D}}$, antioxidant depsidones from <i>Garcinia oliveri</i> . Magnetic Resonance in Chemistry, 2012, 50, 242-245.	1.9	9
121	Letter to the editor: Sequence-specific resonance assignments of the potent cytotoxin equinatoxin II. Journal of Biomolecular NMR, 2000, 18, 281-282.	2.8	8
122	Microwave-Assisted Facile and Rapid Friedel—Crafts Benzoylation of Arenes Catalyzed by Bismuth Trifluoromethanesulfonate. Synthetic Communications, 2014, 44, 2921-2929.	2.1	8
123	NMR, MP2, and DFT study of thiophenoxyketenimines (<i>o</i> -thio—Schiff bases): Determination of the preferred form. Magnetic Resonance in Chemistry, 2018, 56, 172-182.	1.9	8
124	^1H NMR Spin-echo Spectroscopy of Human Erythrocytes. Transformation of Exogenous Compounds. NMR in Biomedicine, 1990, 3, 248-258.	2.8	7
125	Greener Friedel—Crafts Acylation Using Microwave—Enhanced Reactivity of Bismuth Triflate in the Friedel—Crafts Benzoylation of Aromatic Compounds with Benzoic Anhydride. ChemistrySelect, 2017, 2, 571-575.	1.5	7
126	Application of the HECADÉ method to the measurement of long-range heteronuclear ^{13}C , ^1H spin-spin coupling constants in tautomeric \hat{I}^2 -sulfonylenamines. Magnetic Resonance in Chemistry, 2000, 38, 839-844.	1.9	6

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127	The tautomeric equilibrium and stereochemistry of β^2 -sulfonyl enamines. <i>New Journal of Chemistry</i> , 2002, 26, 1060-1069.	2.8	6
128	The application of high performance liquid chromatography humic acid columns in determination of K^+ of polycyclic aromatic compounds. <i>Environmental Toxicology and Chemistry</i> , 2003, 22, 741-745.	4.3	6
129	Molecular structure, intramolecular hydrogen bond strength, vibrational assignment, and spectroscopic insight of 4-phenylamino-3-penten-2-one and its derivatives: A theoretical and experimental study. <i>Journal of Molecular Liquids</i> , 2021, 334, 116035.	4.9	6
130	^1H NMR of compounds with low water solubility in the presence of erythrocytes: effects of emulsion phase separation. <i>European Biophysics Journal</i> , 2001, 30, 69-74.	2.2	5
131	Secondary thioamide group deformations in different surroundings: The case of intramolecular $\text{NH}\cdots\text{N}$ hydrogen bond – An X-ray study combined with theoretical calculations. <i>Journal of Molecular Structure</i> , 2008, 892, 438-445.	3.6	5
132	Isotope effects on chemical shifts in the study of hydrogen bonded biological systems. <i>Progress in Nuclear Magnetic Resonance Spectroscopy</i> , 2020, 120-121, 109-117.	7.5	5
133	Derivatives of usnic acid cause cytostatic effect in Caco-2 cells. <i>Natural Product Research</i> , 2020, 35, 1-7.	1.8	5
134	Azo-hydrazone molecular switches: Synthesis and NMR conformational investigation. <i>Magnetic Resonance in Chemistry</i> , 2021, 59, 1116-1125.	1.9	5
135	The best density functional theory functional for the prediction of ^1H and ^{13}C chemical shifts of protonated alkylpyrroles. <i>Journal of Computational Chemistry</i> , 2021, 42, 1248-1262.	3.3	5
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