Wiktor Kozminski

List of Publications by Year in Descending Order

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

46 29 139 2,944 g-index h-index citations papers 3,321 5.12 141 3.7 L-index ext. citations avg, IF ext. papers

#	Paper	IF	Citations
139	Mechanism of Atg9 recruitment by Atg11 in the cytoplasm-to-vacuole targeting pathway <i>Journal of Biological Chemistry</i> , 2022 , 101573	5.4	2
138	X-ray wavefunction refinement and comprehensive structural studies on bromo-substituted analogues of 2-deoxy-d-glucose in solid state and solution <i>RSC Advances</i> , 2022 , 12, 8345-8360	3.7	
137	Structure, dynamics, and function of SrnR, a transcription factor for nickel-dependent gene expression. <i>Metallomics</i> , 2021 ,	4.5	1
136	Experimental and Computational Studies on Structure and Energetic Properties of Halogen Derivatives of 2-Deoxy-D-Glucose. <i>International Journal of Molecular Sciences</i> , 2021 , 22,	6.3	2
135	Hyperphosphorylation of Human Osteopontin and Its Impact on Structural Dynamics and Molecular Recognition. <i>Biochemistry</i> , 2021 , 60, 1347-1355	3.2	2
134	Order from disorder in the sarcomere: FATZ forms a fuzzy but tight complex and phase-separated condensates with a ctinin. <i>Science Advances</i> , 2021 , 7,	14.3	1
133	Metal Exchange in the Interprotein Zn -Binding Site of the Rad50 Hook Domain: Structural Insights into Cd -Induced DNA-Repair Inhibition. <i>Chemistry - A European Journal</i> , 2020 , 26, 3297-3313	4.8	7
132	H, C and N backbone resonance assignment of BRCA1 fragment 219-504. <i>Biomolecular NMR Assignments</i> , 2020 , 14, 289-293	0.7	О
131	Novel Cyclic Biphalin Analogues by Ruthenium-Catalyzed Ring Closing Metathesis: and Biological Profile. <i>ACS Medicinal Chemistry Letters</i> , 2019 , 10, 450-456	4.3	4
130	Structural analysis of 25-hydroxycholesterol stereoisomers differing in configuration in position 17 and 20, by three-dimensional NMR spectra. <i>Steroids</i> , 2019 , 143, 49-52	2.8	
129	High-dimensional NMR methods for intrinsically disordered proteins studies. <i>Methods</i> , 2018 , 148, 81-87	4.6	9
128	Insight into human insulin aggregation revisited using NMR derived translational diffusion parameters. <i>Journal of Biomolecular NMR</i> , 2018 , 71, 101-114	3	1
127	The Two Isoforms of Lyn Display Different Intramolecular Fuzzy Complexes with the SH3 Domain. <i>Molecules</i> , 2018 , 23,	4.8	8
126	Structure and dynamics of Helicobacter pylori nickel-chaperone HypA: an integrated approach using NMR spectroscopy, functional assays and computational tools. <i>Journal of Biological Inorganic Chemistry</i> , 2018 , 23, 1309-1330	3.7	12
125	Conformational Equilibrium of Cinchonidine in CD Solution. Alternative NMR/DFT Approach. <i>Journal of Physical Chemistry A</i> , 2018 , 122, 7832-7841	2.8	4
124	Structure and Dynamics of the Huntingtin Exon-1 N-Terminus: Alsolution NMR Perspective. <i>Journal of the American Chemical Society</i> , 2017 , 139, 1168-1176	16.4	40
123	H, N, C resonance assignment of plant dehydrin early response to dehydration 10 (ERD10). <i>Biomolecular NMR Assignments</i> , 2017 , 11, 127-131	0.7	3

122	The RxLR Motif of the Host Targeting Effector AVR3a of Is Cleaved before Secretion. <i>Plant Cell</i> , 2017 , 29, 1184-1195	11.6	55
121	Nonuniform Sampling Methods in NMR Data Acquisition 2017 , 418-422		
120	Reconnaissance of reactivity of an Ag(II)SO4 one-electron oxidizer towards naphthalene derivatives. <i>New Journal of Chemistry</i> , 2017 , 41, 10742-10749	3.6	10
119	Reconstruction of non-uniformly sampled five-dimensional NMR spectra by signal separation algorithm. <i>Journal of Biomolecular NMR</i> , 2017 , 68, 129-138	3	15
118	Joint non-uniform sampling of all incremented time delays for quicker acquisition in protein relaxation studies. <i>Journal of Biomolecular NMR</i> , 2017 , 68, 155-161	3	11
117	Five and four dimensional experiments for robust backbone resonance assignment of large intrinsically disordered proteins: application to Tau3x protein. <i>Journal of Biomolecular NMR</i> , 2016 , 65, 193-203	3	7
116	Nuclear overhauser spectroscopy of chiral CHD methylene groups. <i>Journal of Biomolecular NMR</i> , 2016 , 64, 27-37	3	
115	1H, 15N, 13C resonance assignment of human GAP-43. <i>Biomolecular NMR Assignments</i> , 2016 , 10, 171-4	0.7	7
114	Biochemical and Structural Characterization of the Interaction between the Siderocalin NGAL/LCN2 (Neutrophil Gelatinase-associated Lipocalin/Lipocalin 2) and the N-terminal Domain of Its Endocytic Receptor SLC22A17. <i>Journal of Biological Chemistry</i> , 2016 , 291, 2917-30	5.4	25
113	Artifacts in time-resolved NUS: A case study of NOE build-up curves from 2D NOESY. <i>Journal of Magnetic Resonance</i> , 2016 , 265, 108-16	3	13
112	Metal-coupled folding as the driving force for the extreme stability of Rad50 zinc hook dimer assembly. <i>Scientific Reports</i> , 2016 , 6, 36346	4.9	22
111	Amino acid recognition for automatic resonance assignment of intrinsically disordered proteins. Journal of Biomolecular NMR, 2016 , 64, 239-53	3	7
110	The solution structure of the MANEC-type domain from hepatocyte growth factor activator inhibitor-1 reveals an unexpected PAN/apple domain-type fold. <i>Biochemical Journal</i> , 2015 , 466, 299-309	3.8	14
109	Applications of high dimensionality experiments to biomolecular NMR. <i>Progress in Nuclear Magnetic Resonance Spectroscopy</i> , 2015 , 90-91, 49-73	10.4	27
108	(1)H, (15)N, (13)C resonance assignment of human osteopontin. <i>Biomolecular NMR Assignments</i> , 2015 , 9, 289-92	0.7	5
107	Analysis of complex reacting mixtures by time-resolved 2D NMR. <i>Analytical Chemistry</i> , 2015 , 87, 1337-4.	3 7.8	28
106	Six- and seven-dimensional experiments by combination of sparse random sampling and projection spectroscopy dedicated for backbone resonance assignment of intrinsically disordered proteins. Journal of Biomolecular NMR, 2015 , 63, 283-90	3	14
105	High resolution 4D HPCH experiment for sequential assignment of (13)C-labeled RNAs via phosphodiester backbone. <i>Journal of Biomolecular NMR</i> , 2015 , 63, 291-8	3	О

104	Synthesis of rigid tryptophan mimetics by the diastereoselective Pictet-Spengler reaction of Phomo-tryptophan derivatives with chiral Amino aldehydes. <i>Journal of Peptide Science</i> , 2015 , 21, 893-9	0 ² ;1	2
103	(13)C-detected NMR experiments for automatic resonance assignment of IDPs and multiple-fixing SMFT processing. <i>Journal of Biomolecular NMR</i> , 2015 , 62, 179-90	3	6
102	Accelerating diffusion-ordered NMR spectroscopy by joint sparse sampling of diffusion and time dimensions. <i>Angewandte Chemie - International Edition</i> , 2014 , 53, 6464-7	16.4	21
101	Comparison of electrochemical- and nuclear magnetic resonance spectroscopy methods for determination of diffusion coefficients in gel environment. <i>Electrochimica Acta</i> , 2014 , 144, 228-234	6.7	1
100	Accelerating Diffusion-Ordered NMR Spectroscopy by Joint Sparse Sampling of Diffusion and Time Dimensions. <i>Angewandte Chemie</i> , 2014 , 126, 6582-6585	3.6	5
99	C4f/H4Pselective, non-uniformly sampled 4D HC(P)CH experiment for sequential assignments of (13)C-labeled RNAs. <i>Journal of Biomolecular NMR</i> , 2014 , 60, 91-8	3	8
98	"CON-CON" assignment strategy for highly flexible intrinsically disordered proteins. <i>Journal of Biomolecular NMR</i> , 2014 , 60, 209-18	3	28
97	Backbone and partial side chain assignment of the microtubule binding domain of the MAP1B light chain. <i>Biomolecular NMR Assignments</i> , 2014 , 8, 123-7	0.7	3
96	Probing local backbone geometries in intrinsically disordered proteins by cross-correlated NMR relaxation. <i>Angewandte Chemie - International Edition</i> , 2013 , 52, 4604-6	16.4	13
95	Selective diagonal-free (13)C, (13)C-edited aliphatic-aromatic NOESY experiment with non-uniform sampling. <i>Journal of Biomolecular NMR</i> , 2013 , 56, 217-26	3	7
94	Biosynthetic engineered B28(K)-B29(P) human insulin monomer structure in water and in water/acetonitrile solutions. <i>Journal of Biomolecular NMR</i> , 2013 , 55, 303-9	3	6
93	Protonation-dependent conformational variability of intrinsically disordered proteins. <i>Protein Science</i> , 2013 , 22, 1196-205	6.3	27
92	☐H, ☐C and ☐N resonance assignments of human BASP1. <i>Biomolecular NMR Assignments</i> , 2013 , 7, 315-9	0.7	9
91	4D Non-uniformly sampled C,C-NOESY experiment for sequential assignment of 13C, 15N-labeled RNAs. <i>Journal of Biomolecular NMR</i> , 2013 , 57, 1-9	3	8
90	High-dimensionality 13C direct-detected NMR experiments for the automatic assignment of intrinsically disordered proteins. <i>Journal of Biomolecular NMR</i> , 2013 , 57, 353-61	3	39
89	☐H, ☐C, and ☑N backbone and side chain resonance assignments of the C-terminal DNA binding and dimerization domain of v-Myc. <i>Biomolecular NMR Assignments</i> , 2013 , 7, 321-4	0.7	4
88	Study of near-symmetric cyclodextrins by compressed sensing 2D NMR. <i>Magnetic Resonance in Chemistry</i> , 2013 , 51, 110-5	2.1	13
87	(1)H, (13)C, and (15)N chemical shifts assignments for human endothelial monocyte-activating polypeptide EMAP II. <i>Biomolecular NMR Assignments</i> , 2013 , 7, 25-9	0.7	4

86	High-dimensional NMR spectra for structural studies of biomolecules. ChemPhysChem, 2013, 14, 3015-	253.2	27	
85	Iterative thresholding algorithm for multiexponential decay applied to PGSE NMR data. <i>Analytical Chemistry</i> , 2013 , 85, 1828-33	7.8	53	
84	Peptides and peptidoaldehydes as substrates for the Pictet-Spengler reaction. <i>Journal of Peptide Science</i> , 2013 , 19, 433-40	2.1	4	
83	Probing Local Backbone Geometries in Intrinsically Disordered Proteins by Cross-Correlated NMR Relaxation. <i>Angewandte Chemie</i> , 2013 , 125, 4702-4704	3.6		
82	Suppression of sampling artefacts in high-resolution four-dimensional NMR spectra using signal separation algorithm. <i>Journal of Magnetic Resonance</i> , 2012 , 214, 91-102	3	43	
81	Speeding up sequence specific assignment of IDPs. <i>Journal of Biomolecular NMR</i> , 2012 , 53, 293-301	3	63	
80	Crystal and electronic structure, lattice dynamics and thermal properties of Ag(I)(SO3)R (R = F, CF3) Lewis acids in the solid state. <i>Dalton Transactions</i> , 2012 , 41, 2034-47	4.3	23	
79	TSAR: a program for automatic resonance assignment using 2D cross-sections of high dimensionality, high-resolution spectra. <i>Journal of Biomolecular NMR</i> , 2012 , 54, 81-95	3	20	
78	Insights from impedance spectroscopy into the mechanism of thermal decomposition of M(NH2BH3), M = H, Li, Na, Li(0.5)Na(0.5), hydrogen stores. <i>Physical Chemistry Chemical Physics</i> , 2012 , 14, 5778-84	3.6	23	
77	High dimensional and high resolution pulse sequences for backbone resonance assignment of intrinsically disordered proteins. <i>Journal of Biomolecular NMR</i> , 2012 , 52, 329-37	3	48	
76	Generalized Fourier transform for non-uniform sampled data. <i>Topics in Current Chemistry</i> , 2012 , 316, 79-124		26	
75	Polymorphism of a Model Arylboronic Azaester: Combined Experimental and Computational Studies. <i>Crystal Growth and Design</i> , 2011 , 11, 1835-1845	3.5	25	
74	5D 13C-detected experiments for backbone assignment of unstructured proteins with a very low signal dispersion. <i>Journal of Biomolecular NMR</i> , 2011 , 50, 1-11	3	68	
73	Phase transition induced improvement in H2 desorption kinetics: the case of the high-temperature form of Y(BH4)3. <i>Physical Chemistry Chemical Physics</i> , 2011 , 13, 8847-51	3.6	28	
72	A Traceless, Solid-Supported Synthesis of 町urn Mimetics Based on the Hexahydropyrazino[1,2-a]pyrazine-1,2-dione Scaffold. <i>Synthesis</i> , 2010 , 2010, 221-232	2.9	4	
71	Non-uniform frequency domain for optimal exploitation of non-uniform sampling. <i>Journal of Magnetic Resonance</i> , 2010 , 205, 286-92	3	61	
70	Iterative algorithm of discrete Fourier transform for processing randomly sampled NMR data sets. Journal of Biomolecular NMR, 2010 , 47, 65-77	3	74	
69	Strategy for complete NMR assignment of disordered proteins with highly repetitive sequences based on resolution-enhanced 5D experiments. <i>Journal of Biomolecular NMR</i> , 2010 , 48, 169-77	3	77	

Polymorphism of Fluoroargentates(II): Facile Collapse of a Layered Network of EK2AgF4 Due to 68 the Insufficient Size of the Potassium Cation. European Journal of Inorganic Chemistry, 2010, 2010, 2919- $\frac{20}{20}$ 25 $\frac{13}{20}$ Random sampling in multidimensional NMR spectroscopy. Progress in Nuclear Magnetic Resonance 67 10.4 92 Spectroscopy, **2010**, 57, 420-34 A set of 4D NMR experiments of enhanced resolution for easy resonance assignment in proteins. 66 3 30 Journal of Magnetic Resonance, **2010**, 202, 109-16 Determination of heteronuclear coupling constants from 3D HSQC-TOCSY experiment with 65 19 2.1 optimized random sampling of evolution time space. Magnetic Resonance in Chemistry, 2009, 47, 205-9 Spatial structure and NMR spectra of strained [2.2.2] cyclophanes. Magnetic Resonance in Chemistry, 64 2.1 5 2009. 47. 407-14 Complete (1)H and (13)C signal assignment of prenol-10 with 3D NMR spectroscopy. Magnetic 63 2.1 14 Resonance in Chemistry, 2009, 47, 825-9 Narrow peaks and high dimensionalities: exploiting the advantages of random sampling. Journal of 62 66 3 Magnetic Resonance, **2009**, 197, 219-28 KAgF3, K2AgF4 and K3Ag2F7: important steps towards a layered antiferromagnetic 61 3.3 32 fluoroargentate(II),. CrystEngComm, 2009, 11, 1702 Determination of spin-spin couplings from ultrahigh resolution 3D NMR spectra obtained by optimized random sampling and multidimensional Fourier transformation. Journal of the American 60 16.4 23 Chemical Society, 2008, 130, 5404-5 Optimization of random time domain sampling in multidimensional NMR. Journal of Magnetic 85 59 Resonance, 2008, 192, 123-30 Direct insight into insulin aggregation by 2D NMR complemented by PFGSE NMR. Proteins: 58 4.2 19 Structure, Function and Bioinformatics, 2008, 71, 1057-65 Diastereoselective PictetBpengler condensation of tryptophan with &mino aldehydes as chiral 2.4 57 carbonyl components. Tetrahedron, 2008, 64, 1506-1514 Three-dimensional NMR Spectroscopy of organic molecules by random sampling of evolution time 56 2.1 12 space and multidimensional Fourier transformation. Magnetic Resonance in Chemistry, 2007, 45, 171-4 J(F,H), J(C,H) and J(H,H) couplings involving the individual methyl group protons in 1,2,3,4-tetrachloro-5,6,7,8-tetrafluoro-9-methyltriptycene. Evidence of blue-shifting hydrogen 2.1 55 bond. Magnetic Resonance in Chemistry, 2007, 45, 1040-4 Lineshapes and artifacts in Multidimensional Fourier Transform of arbitrary sampled NMR data 67 54 3 sets. Journal of Magnetic Resonance, 2007, 188, 344-56 NMR of Cyclodextrins and Their Complexes **2006**, 231-254 53 12 Two-dimensional Fourier transform of arbitrarily sampled NMR data sets. Journal of Magnetic 52 3 114 Resonance, 2006, 179, 323-8 The studies of tautomerism in 6-mercaptopurine derivatives by 1Ha3C, 1Ha5N NMR and 13C, 15N 21 CPMAS-experimental and quantum chemical approach. Journal of Molecular Structure, **2006**, 785, 205-21 $^{\circ}5^{\circ}4$

(2000-2006)

50	Random sampling of evolution time space and Fourier transform processing. <i>Journal of Biomolecular NMR</i> , 2006 , 36, 157-68	3	94	
49	Efficient compensation of low-frequency magnetic field disturbances in NMR with fluxgate sensors. Journal of Magnetic Resonance, 2005 , 174, 287-91	3	14	
48	A protein backbone psi and phi angle dependence of 2J(N(i),C alpha(i-1)): the new NMR experiment and quantum chemical calculations. <i>Journal of Biomolecular NMR</i> , 2005 , 31, 87-95	3	9	
47	The DQ-HN[CACB] and DQ-HN(CO)[CACB] sequences with evolution of double quantum Calpha-Cbeta coherences. <i>Journal of Magnetic Resonance</i> , 2004 , 171, 186-91	3	3	
46	The set of triple-resonance sequences with a multiple quantum coherence evolution period. Journal of Magnetic Resonance, 2004 , 171, 338-44	3	3	
45	NMR studies of chiral recognition by cyclodextrins. <i>Chirality</i> , 2004 , 16, 90-105	2.1	95	
44	15N, 13C and 1H nuclear magnetic shielding and spin⊞pin coupling in gaseous 15N-enriched methylamine. <i>Journal of Molecular Structure</i> , 2004 , 704, 305-309	3.4	16	
43	15NH4+ ion movement inside d(G4T4G4)2 G-quadruplex is accelerated in the presence of smaller Na+ ions. <i>Organic and Biomolecular Chemistry</i> , 2004 , 2, 1970-3	3.9	32	
42	A Study of Multiple Complexation of 日 和 DCyclodextrins: Surprisingly Differing Stoichiometries of 和 DCyclodextrin Complexes. <i>Supramolecular Chemistry</i> , 2004 , 16, 287-292	1.8	10	
41	Multiple quadrature detection in reduced dimensionality experiments. <i>Journal of Biomolecular NMR</i> , 2003 , 26, 157-66	3	64	
40	The new HMQC-based technique for the quantitative determination of heteronuclear coupling constants. Application for the measurement of 3J(HR),P(i+1)) in DNA oligomers. <i>Journal of Magnetic Resonance</i> , 2003 , 160, 120-5	3	11	
39	On the impossibility of determination of stepwise binding constants for the 1 : 2 complex of (+)-camphor with alpha-cyclodextrin. <i>Organic and Biomolecular Chemistry</i> , 2003 , 1, 581-4	3.9	16	
38	An improved 33S nuclear magnetic shielding scale from the gas-phase study of COS. <i>Magnetic Resonance in Chemistry</i> , 2002 , 40, 563-565	2.1	23	
37	15N, 13C and 1H nuclear magnetic shielding and spin pin coupling constants of 1-13C, 15N-enriched acetonitrile in gaseous mixtures with SF6 and CO2. <i>Chemical Physics Letters</i> , 2002 , 358, 263-270	2.5	28	
36	An experimental test of C-N bond twisting in the TICT state: syn-anti photoisomerization in 2-(N-methyl-N-isopropylamino)-5-cyanopyridine. <i>Journal of the American Chemical Society</i> , 2002 , 124, 2406-7	16.4	72	
35	Effects of Intermolecular Interactions on 33S Magnetic Shielding in Gaseous SF6. <i>Journal of Physical Chemistry A</i> , 2002 , 106, 2829-2832	2.8	22	
34	The tautomeric equilibrium and stereochemistry of	3.6	6	
33	Application of adiabatic inversion pulses for elimination of baseline distortions in Fourier transform NMR. A natural abundance 170 NMR spectrum for gaseous acetone. <i>Magnetic Resonance in Chemistry</i> , 2000 , 38, 459-462	2.1	29	

32	Application of the HECADE method to the measurement of long-range heteronuclear 13C,1H spinBpin coupling constants in tautomeric 最ulfonylenamines. <i>Magnetic Resonance in Chemistry</i> , 2000 , 38, 839-844	2.1	5
31	Sensitivity improvement and new acquisition scheme of heteronuclear active-coupling-pattern-tilting spectroscopy. <i>Journal of Magnetic Resonance</i> , 2000 , 142, 294-9	3	64
30	A dynamic NMR study of self-inclusion of a pendant group in amphiphilic 6-thiophenyl-6-deoxycyclodextrins. <i>Journal of Molecular Structure</i> , 2000 , 519, 33-36	3.4	13
29	NMR manifestations and molecular dynamics modeling of chiral recognition of pinenes by Eyclodextrin. <i>Journal of Molecular Structure</i> , 2000 , 523, 205-212	3.4	22
28	1H, 13C, 15N NMR and X-Ray Diffractometry in Structural Studies of Macrocyclic Lactams Containing Pyridine Moiety. <i>Supramolecular Chemistry</i> , 2000 , 12, 229-235	1.8	12
27	Application of adiabatic inversion pulses for elimination of baseline distortions in Fourier transform NMR. A natural abundance 170 NMR spectrum for gaseous acetone 2000 , 38, 459		1
26	Rosavin as a product of glycosylation by Rhodiola rosea (roseroot) cell cultures. <i>Plant Cell, Tissue and Organ Culture</i> , 1999 , 56, 105-110	2.7	27
25	Simplified multiplet pattern HSQC-TOCSY experiment for accurate determination of long-range heteronuclear coupling constants. <i>Journal of Magnetic Resonance</i> , 1999 , 137, 408-12	3	45
24	A pure-phase homonuclear J-modulated HMQC experiment with tilted cross-peak patterns for an accurate determination of homonuclear coupling constants. <i>Journal of Magnetic Resonance</i> , 1999 , 141, 185-90	3	11
23	The new active-coupling-pattern tilting experiment for an efficient and accurate determination of homonuclear coupling constants. <i>Journal of Magnetic Resonance</i> , 1998 , 134, 189-93	3	13
22	Synthesis and Characterization of 1,2-Disubstituted Vinylsilanes and Their Geometric Differentiation with 3J(29Si,1H)-Coupling Constants. Application of a Novel Heteronuclear J-Resolved NMR Experiment. <i>Organometallics</i> , 1997 , 16, 3128-3134	3.8	9
21	HECADE: HMQC- and HSQC-Based 2D NMR Experiments for Accurate and Sensitive Determination of Heteronuclear Coupling Constants from E.COSY-Type Cross Peaks. <i>Journal of Magnetic Resonance</i> , 1997 , 124, 383-392	3	83
20	Pure-Phase Homo- and Heteronuclear J Spectra with Tilted Cross Peaks for an Accurate Determination of Coupling Constants. <i>Journal of Magnetic Resonance</i> , 1997 , 125, 193-6	3	7
19	187Os NMR Study of (B-Arene)osmium(II) Complexes: Separation of Electronic and Steric Ligand Effects Organometallics, 1996 , 15, 3124-3135	3.8	31
18	57Fe NMR Study of Ligand Effects in Cyclopentadienyliron Complexes Organometallics, 1996 , 15, 2469)- <u>3</u> . \$ 77	29
17	An Analysis of the Bonding Properties of Benz[a]azulene by X-Ray, NMR, and Computational Studies. <i>Helvetica Chimica Acta</i> , 1996 , 79, 837-854	2	9
16	57Fe, 13C Coupling Constants from Inverse Detection Experiments at Natural Isotope Abundance. <i>Magnetic Resonance in Chemistry</i> , 1996 , 34, 89-92	2.1	10
15	Sensitive Measurement of One-Bond Carbontarbon Spin Coupling Constants at Natural Isotope Abundance. <i>Magnetic Resonance in Chemistry</i> , 1996 , 34, 311-315	2.1	14

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14	Sensitive Measurement and Unambiguous Assignment of Long-Range13C,13C Coupling Constants at Natural Isotope Abundance. <i>Journal of Magnetic Resonance Series A</i> , 1996 , 122, 245-247		15
13	An X-ray study of some 2,3-diphenyltetrazolium salts. <i>Journal of Chemical Crystallography</i> , 1995 , 25, 29	-35 .5	6
12	15N and 13C solid-state nuclear magnetic resonance study of 5-thiomethyltetrazole. <i>Solid State Nuclear Magnetic Resonance</i> , 1995 , 4, 121-4	3.1	1
11	Spin l attice Relaxation Times of Transition-Metal Nuclei from Inverse-Detection Experiments*. <i>Journal of Magnetic Resonance Series A</i> , 1995 , 116, 262-265		8
10	Characteristic coupling constants1J(13C-13C)of some mesoionic methylides containing a diphenyltetrazolium ring. <i>Spectroscopy</i> , 1994 , 12, 21-23		6
9	Multinuclear magnetic resonance study of some mesoionic 1,3-diphenyltetrazoles with various exocyclic groups. <i>Magnetic Resonance in Chemistry</i> , 1994 , 32, 284-287	2.1	10
8	13C and 15N NMR study of mesoionic type A and type B tetrazoles with four nitrogen atoms in the exocyclic group. <i>Magnetic Resonance in Chemistry</i> , 1994 , 32, 746-748	2.1	5
7	14N and 15N NMR study of 2,1,3-thiadiazolium-5-olate and related compounds. <i>Journal of Molecular Structure</i> , 1994 , 323, 177-179	3.4	3
6	A multinuclear NMR study of some mesoionic 1,3-dimethyltetrazoles, 1- and 2-methyltetrazoles and related compounds. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1994 , 1327-1332		20
5	An X-ray diffraction study of some mesoionic 2,3-diphenyltetrazoles. <i>Journal of Crystallographic and Spectroscopic Research</i> , 1993 , 23, 133-141		8
4	15N, 14N, 13C and 1H NMR study of mesoionic methylides and thiocarbonyl ylides with a 2,3-diphenyltetrazolium ring. <i>Journal of Molecular Structure</i> , 1993 , 295, 15-18	3.4	5
3	A multinuclear NMR study on some cyclic aminimides and related compounds. <i>Journal of Molecular Structure</i> , 1991 , 243, 365-368	3.4	3
2	A 13C and 15N NMR study of some mesoionic 4-hydroxy pyrazole derivatives in various solvents. Journal of Molecular Structure, 1991 , 243, 369-372	3.4	6
1	13C and 15N NMR study of 2,3-diphenyltetrazolium-5-olate and 5-thiolate. <i>Magnetic Resonance in Chemistry</i> , 1990 , 28, 1027-1029	2.1	16