

Gareth A Morris

List of Publications by Citations

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252
papers

10,023
citations

48
h-index

91
g-index

279
ext. papers

10,755
ext. citations

5.5
avg, IF

6.2
L-index

| # | Paper | IF | Citations |
|-----|---|------|-----------|
| 252 | Enhancement of nuclear magnetic resonance signals by polarization transfer. <i>Journal of the American Chemical Society</i> , 1979 , 101, 760-762 | 16.4 | 1689 |
| 251 | Sensitivity enhancement in nitrogen-15 NMR: polarization transfer using the INEPT pulse sequence. <i>Journal of the American Chemical Society</i> , 1980 , 102, 428-429 | 16.4 | 320 |
| 250 | Pure shift 1H NMR: a resolution of the resolution problem?. <i>Angewandte Chemie - International Edition</i> , 2010 , 49, 3901-3 | 16.4 | 208 |
| 249 | A one-shot sequence for high-resolution diffusion-ordered spectroscopy. <i>Magnetic Resonance in Chemistry</i> , 2002 , 40, S147-S152 | 2.1 | 207 |
| 248 | Ultrahigh-resolution NMR spectroscopy. <i>Angewandte Chemie - International Edition</i> , 2014 , 53, 6990-2 | 16.4 | 201 |
| 247 | An improved method for heteronuclear chemical shift correlation by two-dimensional NMR. <i>Journal of Magnetic Resonance</i> , 1981 , 42, 501-505 | | 192 |
| 246 | High-Resolution Diffusion-Ordered 2D Spectroscopy (HR-DOSY) - A New Tool for the Analysis of Complex Mixtures. <i>Journal of Magnetic Resonance Series B</i> , 1995 , 108, 170-172 | | 186 |
| 245 | Spin echo NMR spectra without J modulation. <i>Chemical Communications</i> , 2012 , 48, 811-3 | 5.8 | 175 |
| 244 | Pulse sequences for high-resolution diffusion-ordered spectroscopy (HR-DOSY). <i>Magnetic Resonance in Chemistry</i> , 1998 , 36, 706-714 | 2.1 | 165 |
| 243 | Experimental chemical shift correlation maps in nuclear magnetic resonance spectroscopy. <i>Journal of the Chemical Society Chemical Communications</i> , 1978 , 684 | | 162 |
| 242 | Pure shift proton DOSY: diffusion-ordered 1H spectra without multiplet structure. <i>Chemical Communications</i> , 2007 , 933-5 | 5.8 | 150 |
| 241 | Modern NMR techniques for structure elucidation. <i>Magnetic Resonance in Chemistry</i> , 1986 , 24, 371-403 | 2.1 | 150 |
| 240 | Selective excitation in Fourier transform nuclear magnetic resonance. <i>Journal of Magnetic Resonance</i> , 1978 , 29, 433-462 | | 149 |
| 239 | Reference deconvolution methods. <i>Progress in Nuclear Magnetic Resonance Spectroscopy</i> , 1997 , 31, 197-257 | 15.4 | 146 |
| 238 | Simultaneously enhancing spectral resolution and sensitivity in heteronuclear correlation NMR spectroscopy. <i>Angewandte Chemie - International Edition</i> , 2013 , 52, 11616-9 | 16.4 | 141 |
| 237 | Quantitative interpretation of diffusion-ordered NMR spectra: can we rationalize small molecule diffusion coefficients?. <i>Angewandte Chemie - International Edition</i> , 2013 , 52, 3199-202 | 16.4 | 140 |
| 236 | Experimental chemical shift correlation maps from heteronuclear two-dimensional NMR spectroscopy. 1. Carbon-13 and proton chemical shifts of raffinose and its subunits. <i>Journal of the American Chemical Society</i> , 1981 , 103, 4703-4711 | 16.4 | 111 |

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| 235 | Correlation of proton chemical shifts by two-dimensional Fourier transform NMR. <i>Journal of Magnetic Resonance</i> , 1981 , 42, 164-168 | | 106 |
| 234 | Improving the accuracy of pulsed field gradient NMR diffusion experiments: Correction for gradient non-uniformity. <i>Journal of Magnetic Resonance</i> , 2009 , 198, 121-31 | 3 | 105 |
| 233 | A three-dimensional DOSY-HMQC experiment for the high-resolution analysis of complex mixtures. <i>Journal of Magnetic Resonance</i> , 1998 , 131, 131-8 | 3 | 105 |
| 232 | Simple proton spectra from complex spin systems: pure shift NMR spectroscopy using BIRD. <i>Angewandte Chemie - International Edition</i> , 2011 , 50, 9716-7 | 16.4 | 101 |
| 231 | High-resolution NMR and diffusion-ordered spectroscopy of port wine. <i>Journal of Agricultural and Food Chemistry</i> , 2004 , 52, 3736-43 | 5.7 | 101 |
| 230 | True chemical shift correlation maps: a TOCSY experiment with pure shifts in both dimensions. <i>Journal of the American Chemical Society</i> , 2010 , 132, 12770-2 | 16.4 | 98 |
| 229 | Biexponential fitting of diffusion-ordered NMR data: practicalities and limitations. <i>Analytical Chemistry</i> , 2006 , 78, 3040-5 | 7.8 | 97 |
| 228 | Ultrahigh-resolution total correlation NMR spectroscopy. <i>Journal of the American Chemical Society</i> , 2014 , 136, 11867-9 | 16.4 | 96 |
| 227 | Quantifying end-to-end conformational communication of chirality through an achiral peptide chain. <i>Angewandte Chemie - International Edition</i> , 2009 , 48, 5962-5 | 16.4 | 92 |
| 226 | Decoupling two-dimensional NMR spectroscopy in both dimensions: pure shift NOESY and COSY. <i>Angewandte Chemie - International Edition</i> , 2012 , 51, 6460-3 | 16.4 | 90 |
| 225 | Foldamer-mediated remote stereocontrol: >1,60 asymmetric induction. <i>Angewandte Chemie - International Edition</i> , 2014 , 53, 151-5 | 16.4 | 87 |
| 224 | Speedy component resolution: an improved tool for processing diffusion-ordered spectroscopy data. <i>Analytical Chemistry</i> , 2008 , 80, 3777-82 | 7.8 | 83 |
| 223 | "Perfecting" WATERGATE: clean proton NMR spectra from aqueous solution. <i>Chemical Communications</i> , 2013 , 49, 358-60 | 5.8 | 79 |
| 222 | Helix persistence and breakdown in oligoureas of metaphenylenediamine: apparent diastereotopicity as a spectroscopic marker of helix length in solution. <i>Journal of the American Chemical Society</i> , 2008 , 130, 15193-202 | 16.4 | 74 |
| 221 | One-dimensional DOSY. <i>Journal of Magnetic Resonance</i> , 2001 , 153, 103-12 | 3 | 74 |
| 220 | Matrix-assisted diffusion-ordered spectroscopy: mixture resolution by NMR using SDS micelles. <i>Magnetic Resonance in Chemistry</i> , 2010 , 48, 550-3 | 2.1 | 69 |
| 219 | Measuring couplings in crowded NMR spectra: pure shift NMR with multiplet analysis. <i>Chemical Communications</i> , 2015 , 51, 15410-3 | 5.8 | 67 |
| 218 | Measuring screw-sense preference in a helical oligomer by comparison of ¹³ C NMR signal separation at slow and fast exchange. <i>Journal of the American Chemical Society</i> , 2011 , 133, 3712-5 | 16.4 | 67 |

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| 217 | Resolution and assignment of the 270-MHz proton spectrum of cellobiose by homo- and heteronuclear two-dimensional NMR. <i>Journal of the American Chemical Society</i> , 1980 , 102, 1745-1747 | 16.4 | 66 |
| 216 | Sample convection in liquid-state NMR: why it is always with us, and what we can do about it. <i>Journal of Magnetic Resonance</i> , 2015 , 252, 120-9 | 3 | 64 |
| 215 | Improving the Interpretation of Small Molecule Diffusion Coefficients. <i>Analytical Chemistry</i> , 2018 , 90, 3987-3994 | 7.8 | 63 |
| 214 | Isomer resolution by micelle-assisted diffusion-ordered spectroscopy. <i>Analytical Chemistry</i> , 2009 , 81, 4548-50 | 7.8 | 63 |
| 213 | Improving pulse sequences for 3D diffusion-ordered NMR spectroscopy: 2DJ-IDOSY. <i>Analytical Chemistry</i> , 2004 , 76, 5418-22 | 7.8 | 63 |
| 212 | Diastereomeric ratio determination by high sensitivity band-selective pure shift NMR spectroscopy. <i>Chemical Communications</i> , 2014 , 50, 2512-4 | 5.8 | 60 |
| 211 | Local covariance order diffusion-ordered spectroscopy: a powerful tool for mixture analysis. <i>Journal of the American Chemical Society</i> , 2011 , 133, 7640-3 | 16.4 | 59 |
| 210 | A General Method for Extracting Individual Coupling Constants from Crowded (1)H NMR Spectra. <i>Angewandte Chemie - International Edition</i> , 2016 , 55, 1090-3 | 16.4 | 59 |
| 209 | J-modulation effects in DOSY experiments and their suppression: the Oneshot45 experiment. <i>Journal of Magnetic Resonance</i> , 2011 , 208, 270-8 | 3 | 54 |
| 208 | Improving pulse sequences for 3D DOSY: COSY-IDOSY. <i>Chemical Communications</i> , 2005 , 1737-9 | 5.8 | 54 |
| 207 | Simultaneous enhancement of chemical shift dispersion and diffusion resolution in mixture analysis by diffusion-ordered NMR spectroscopy. <i>Chemical Communications</i> , 2011 , 47, 7063-4 | 5.8 | 52 |
| 206 | 2D and 3D DOSY methods for studying mixtures of oligomeric dimethylsiloxanes. <i>Physical Chemistry Chemical Physics</i> , 2004 , 6, 3221 | 3.6 | 52 |
| 205 | A simple pulse sequence for selective excitation in Fourier transform NMR. <i>Journal of Magnetic Resonance</i> , 1976 , 23, 171-175 | | 51 |
| 204 | Accurate determination of one-bond heteronuclear coupling constants with "pure shift" broadband proton-decoupled CLIP/CLAP-HSQC experiments. <i>Journal of Magnetic Resonance</i> , 2014 , 239, 130-8 | 3 | 46 |
| 203 | Conformational Switching of a Foldamer in a Multicomponent System by pH-Filtered Selection between Competing Noncovalent Interactions. <i>Journal of the American Chemical Society</i> , 2015 , 137, 6680-91 | 16.4 | 44 |
| 202 | Ultrahigh-Resolution Diffusion-Ordered Spectroscopy. <i>Angewandte Chemie - International Edition</i> , 2016 , 55, 15579-15582 | 16.4 | 44 |
| 201 | "Perfecting" pure shift HSQC: full homodecoupling for accurate and precise determination of heteronuclear couplings. <i>Chemical Communications</i> , 2014 , 50, 15702-5 | 5.8 | 44 |
| 200 | Residue-specific NH exchange rates studied by NMR diffusion experiments. <i>Journal of Magnetic Resonance</i> , 2007 , 187, 97-104 | 3 | 42 |

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| 199 | Indirect measurement of proton relaxation rates by [INEPT] polarization transfer to carbon-13: Proton spin-lattice relaxation in cholesteryl acetate solutions. <i>Journal of Magnetic Resonance</i> , 1980 , 41, 185-188 | | 42 |
| 198 | Acid-catalyzed degradation of clarithromycin and erythromycin B: a comparative study using NMR spectroscopy. <i>Journal of Medicinal Chemistry</i> , 2000 , 43, 467-74 | 8.3 | 41 |
| 197 | The catalytic resting state of asymmetric homogeneous hydrogenation. Exchange processes delineated by nuclear magnetic resonance saturation-transfer (DANTE) techniques. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1987 , 1583 | | 41 |
| 196 | Measurement of carbon-13-proton coupling-constants in oligosaccharides by two-dimensional carbon-13 N.M.R. spectroscopy. <i>Carbohydrate Research</i> , 1980 , 82, 175-184 | 2.9 | 41 |
| 195 | Resolving natural product epimer spectra by matrix-assisted DOSY. <i>Organic and Biomolecular Chemistry</i> , 2011 , 9, 7062-4 | 3.9 | 40 |
| 194 | Improving pulse sequences for 3D DOSY: convection compensation. <i>Journal of Magnetic Resonance</i> , 2005 , 177, 203-11 | 3 | 39 |
| 193 | Experimental chemical shift correlation maps from heteronuclear two-dimensional nuclear magnetic resonance spectroscopy. II: Carbon-13 and proton chemical shifts of α -D-glucopyranose oligomers. <i>Canadian Journal of Chemistry</i> , 1982 , 60, 2431-2441 | 0.9 | 39 |
| 192 | The GNAT: A new tool for processing NMR data. <i>Magnetic Resonance in Chemistry</i> , 2018 , 56, 546-558 | 2.1 | 38 |
| 191 | Simple Proton Spectra from Complex Spin Systems: Pure Shift NMR Spectroscopy Using BIRD. <i>Angewandte Chemie</i> , 2011 , 123, 9890-9891 | 3.6 | 38 |
| 190 | PSYCHE Pure Shift NMR Spectroscopy. <i>Chemistry - A European Journal</i> , 2018 , 24, 13988-14000 | 4.8 | 36 |
| 189 | Flavonoid mixture analysis by matrix-assisted diffusion-ordered spectroscopy. <i>Journal of Natural Products</i> , 2012 , 75, 131-4 | 4.9 | 36 |
| 188 | Unmixing the NMR spectra of similar species - vive la difference. <i>Chemical Communications</i> , 2013 , 49, 10510-2 | 5.8 | 35 |
| 187 | Left-handed helical preference in an achiral peptide chain is induced by an L-amino acid in an N-terminal type II turn. <i>Journal of Organic Chemistry</i> , 2013 , 78, 2248-55 | 4.2 | 35 |
| 186 | Suppression of t1 Noise in 2D NMR Spectroscopy by Reference Deconvolution. <i>Journal of Magnetic Resonance Series A</i> , 1993 , 101, 351-356 | | 35 |
| 185 | Increasing the quantitative bandwidth of NMR measurements. <i>Chemical Communications</i> , 2016 , 52, 29165-8 | 5.9 | 34 |
| 184 | Ultraclean pure shift NMR. <i>Chemical Communications</i> , 2017 , 53, 10188-10191 | 5.8 | 34 |
| 183 | Homoleptic Trigonal Planar Lanthanide Complexes Stabilized by Superbulky Silylamide Ligands. <i>Organometallics</i> , 2015 , 34, 2314-2325 | 3.8 | 34 |
| 182 | Ultrahigh-Resolution NMR Spectroscopy. <i>Angewandte Chemie</i> , 2014 , 126, 7110-7112 | 3.6 | 34 |

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| 181 | Mechanism for the degradation of erythromycin A and erythromycin A 2'-ethyl succinate in acidic aqueous solution. <i>Journal of Physical Chemistry A</i> , 2007 , 111, 10098-104 | 2.8 | 34 |
| 180 | Diffusion NMR and trilinear analysis in the study of reaction kinetics. <i>Chemical Communications</i> , 2009 , 1252-4 | 5.8 | 33 |
| 179 | A novel NMR method for screening soluble compound libraries. <i>Chemical Communications</i> , 2001 , 239-240 | 0.8 | 33 |
| 178 | Direct observation of the magnetization exchange dynamics responsible for magnetization transfer contrast in human cartilage in vitro. <i>Magnetic Resonance in Medicine</i> , 1992 , 28, 97-104 | 4.4 | 33 |
| 177 | Flaws in foldamers: conformational uniformity and signal decay in achiral helical peptide oligomers. <i>Chemical Science</i> , 2015 , 6, 2313-2322 | 9.4 | 32 |
| 176 | Novel artemisinin and curcumin micellar formulations: drug solubility studies by NMR spectroscopy. <i>Journal of Pharmaceutical Sciences</i> , 2009 , 98, 3666-75 | 3.9 | 32 |
| 175 | A simple flowcell for reaction monitoring by NMR. <i>Magnetic Resonance in Chemistry</i> , 2010 , 48, 516-22 | 2.1 | 32 |
| 174 | Application of ³¹ P-NMR saturation transfer techniques to investigate phospholipid motion and organization in model and biological membranes. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 1980 , 598, 206-11 | 3.8 | 32 |
| 173 | Convection in liquid-state NMR: expect the unexpected. <i>RSC Advances</i> , 2016 , 6, 95173-95176 | 3.7 | 32 |
| 172 | High resolution ¹³ C DOSY: the DEPTSE experiment. <i>Journal of Magnetic Resonance</i> , 2011 , 211, 25-9 | 3 | 30 |
| 171 | A Practical Method for Automated Shimming with Normal Spectrometer Hardware. <i>Journal of Magnetic Resonance</i> , 1997 , 125, 197-201 | 3 | 30 |
| 170 | Suppressing exchange effects in diffusion-ordered NMR spectroscopy. <i>Journal of Magnetic Resonance</i> , 2014 , 238, 16-9 | 3 | 29 |
| 169 | Quantitative Interpretation of Diffusion-Ordered NMR Spectra: Can We Rationalize Small Molecule Diffusion Coefficients?. <i>Angewandte Chemie</i> , 2013 , 125, 3281-3284 | 3.6 | 29 |
| 168 | A diffusion-ordered NMR spectroscopy study of the solubilization of artemisinin by octanoyl-6-O-ascorbic acid micelles. <i>Journal of Pharmaceutical Sciences</i> , 2002 , 91, 2265-70 | 3.9 | 29 |
| 167 | Reaction kinetics studied using diffusion-ordered spectroscopy and multiway chemometrics. <i>Analytical Chemistry</i> , 2010 , 82, 2102-8 | 7.8 | 28 |
| 166 | Z-spectroscopy with Alternating-Phase Irradiation. <i>Journal of Magnetic Resonance</i> , 2010 , 207, 242-50 | 3 | 28 |
| 165 | Correction of systematic errors in CORE processing of DOSY data. <i>Magnetic Resonance in Chemistry</i> , 2006 , 44, 655-60 | 2.1 | 28 |
| 164 | Foldamer-Mediated Remote Stereocontrol: >1,60 Asymmetric Induction. <i>Angewandte Chemie</i> , 2014 , 126, 155-159 | 3.6 | 27 |

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| 163 | Detection of potential TNA and RNA nucleoside precursors in a prebiotic mixture by pure shift diffusion-ordered NMR spectroscopy. <i>Chemistry - A European Journal</i> , 2013 , 19, 4586-95 | 4.8 | 27 |
| 162 | Concerted use of homo- and hetero-nuclear 2D NMR: ¹³ C and ¹ H assignment of sucrose octaacetate. <i>Magnetic Resonance in Chemistry</i> , 1986 , 24, 179-182 | 2.1 | 27 |
| 161 | Constant time gradient HSQC-iDOSY: practical aspects. <i>Magnetic Resonance in Chemistry</i> , 2009 , 47, 1081-5 | 2.1 | 26 |
| 160 | Complete Assignment of the ¹ H and ¹³ C NMR Spectra of Steroidal Sapogenins: Smilagenin and Sarsapogenin. <i>Magnetic Resonance in Chemistry</i> , 1997 , 35, 441-446 | 2.1 | 26 |
| 159 | The Behavior of Multiplet Signals under Radiation Damping Conditions. I. Classical Effects. <i>Journal of Magnetic Resonance Series A</i> , 1995 , 117, 109-112 | | 26 |
| 158 | Real-time pure shift ¹ H HSQC of proteins: a real improvement in resolution and sensitivity. <i>Journal of Biomolecular NMR</i> , 2015 , 62, 43-52 | 3 | 25 |
| 157 | T1-diffusion-ordered spectroscopy: nuclear magnetic resonance mixture analysis using parallel factor analysis. <i>Analytical Chemistry</i> , 2009 , 81, 8119-25 | 7.8 | 25 |
| 156 | Dependence of the ¹ H NMR chemical shifts of ring F resonances on the orientation of the 27-methyl group of spirostane-type steroidal sapogenins. <i>Phytochemistry</i> , 1998 , 47, 255-7 | 4 | 25 |
| 155 | Carbon-13 and proton two-dimensional NMR study of the Ormosia alkaloids panamine, ormosanine, and ormosinine. <i>Journal of the American Chemical Society</i> , 1983 , 105, 2538-2544 | 16.4 | 25 |
| 154 | ¹⁹ F DOSY NMR analysis for spin systems with nJFF couplings. <i>Magnetic Resonance in Chemistry</i> , 2014 , 52, 172-7 | 2.1 | 24 |
| 153 | Simultaneously Enhancing Spectral Resolution and Sensitivity in Heteronuclear Correlation NMR Spectroscopy. <i>Angewandte Chemie</i> , 2013 , 125, 11830-11833 | 3.6 | 24 |
| 152 | Identification of a further transient species relating to rhodium-complex catalysed asymmetric hydrogenation. <i>Journal of the Chemical Society Chemical Communications</i> , 1983 , 664 | | 24 |
| 151 | Particle size measurement of lipoprotein fractions using diffusion-ordered NMR spectroscopy. <i>Analytical and Bioanalytical Chemistry</i> , 2012 , 402, 2407-15 | 4.4 | 23 |
| 150 | Decoupling Two-Dimensional NMR Spectroscopy in Both Dimensions: Pure Shift NOESY and COSY. <i>Angewandte Chemie</i> , 2012 , 124, 6566-6569 | 3.6 | 23 |
| 149 | NMR measurements of diffusion in concentrated samples: avoiding problems with radiation damping. <i>Analytical and Bioanalytical Chemistry</i> , 2004 , 378, 1568-73 | 4.4 | 23 |
| 148 | General Analytical Solutions of the Bloch Equations. <i>Journal of Magnetic Resonance Series A</i> , 1994 , 107, 236-238 | | 23 |
| 147 | Molecular characterisation of oxymethylene-linked poly(oxyethylene). <i>British Polymer Journal</i> , 1987 , 19, 509-516 | | 23 |
| 146 | Minimising research bottlenecks by decluttering NMR spectra. <i>Chemistry - A European Journal</i> , 2015 , 21, 6623-30 | 4.8 | 22 |

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| 145 | A simple method for suppressing dispersion-mode contributions in NMR spectra: The β pseudo echo <i>Journal of Magnetic Resonance</i> , 1981 , 43, 333-338 | | 22 |
| 144 | High resolution, high field magnetic resonance imaging of joints: unexpected features in proton images of cartilage. <i>British Journal of Radiology</i> , 1990 , 63, 907-9 | 3.4 | 21 |
| 143 | Weak satellite signals in high-resolution NMR spectra: Separating the wheat from the chaff. <i>Journal of Magnetic Resonance</i> , 1981 , 42, 341-345 | | 21 |
| 142 | A new tool for NMR analysis of complex systems: selective pure shift TOCSY. <i>RSC Advances</i> , 2016 , 6, 10063-10066 | 3.7 | 20 |
| 141 | Improving accuracy in DOSY and diffusion measurements using triaxial field gradients. <i>Journal of Magnetic Resonance</i> , 2016 , 270, 24-30 | 3 | 20 |
| 140 | Probing the Anions Mediated Associative Behavior of Tin-12 Oxo-Macrocations by Pulsed Field Gradient NMR Spectroscopy. <i>Journal of Physical Chemistry C</i> , 2010 , 114, 16087-16091 | 3.8 | 20 |
| 139 | Suppression of artefacts in nuclear overhauser effect difference spectroscopy by reference deconvolution. <i>Magnetic Resonance in Chemistry</i> , 1989 , 27, 1085-1089 | 2.1 | 20 |
| 138 | Very broadband diffusion-ordered NMR spectroscopy: (19)F DOSY. <i>Chemical Communications</i> , 2016 , 52, 6892-4 | 5.8 | 20 |
| 137 | Anatomising proton NMR spectra with pure shift 2D J-spectroscopy: A cautionary tale. <i>Chemical Physics Letters</i> , 2017 , 683, 398-403 | 2.5 | 19 |
| 136 | Filter diagonalization method for processing PFG NMR data. <i>Journal of Magnetic Resonance</i> , 2013 , 234, 125-34 | 3 | 19 |
| 135 | Diffusion-Ordered Spectroscopy 2009 , | | 19 |
| 134 | Improved DECRA processing of DOSY data: correcting for non-uniform field gradients. <i>Magnetic Resonance in Chemistry</i> , 2007 , 45, 656-60 | 2.1 | 19 |
| 133 | Reference Deconvolution Using Multiplet Reference Signals. <i>Journal of Magnetic Resonance Series A</i> , 1995 , 116, 206-214 | | 19 |
| 132 | Matrix-assisted diffusion-ordered spectroscopy: application of surfactant solutions to the resolution of isomer spectra. <i>Magnetic Resonance in Chemistry</i> , 2012 , 50, 458-65 | 2.1 | 18 |
| 131 | Effects of radiation damping on Z-spectra. <i>Journal of Magnetic Resonance</i> , 2006 , 183, 203-12 | 3 | 18 |
| 130 | Concerted use of two-dimensional NMR techniques in the ab initio assignment of complex spectra: Complete proton and carbon-13 assignment of oligomycin A. <i>Magnetic Resonance in Chemistry</i> , 1985 , 23, 676-683 | 2.1 | 18 |
| 129 | Carbon-13 nuclear magnetic resonance spectra with coherent proton decoupling: peak-height distortions within spin multiplets. <i>Journal of the American Chemical Society</i> , 1978 , 100, 5637-5640 | 16.4 | 18 |
| 128 | Relaxation-encoded NMR experiments for mixture analysis: REST and beer. <i>Chemical Communications</i> , 2017 , 53, 7461-7464 | 5.8 | 17 |

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| 127 | Cleaning up NMR spectra with reference deconvolution for improving multivariate analysis of complex mixture spectra. <i>Journal of Chemometrics</i> , 2014 , 28, 656-662 | 1.6 | 17 |
| 126 | Difluorinated analogues of shikimic acid. <i>Tetrahedron</i> , 2003 , 59, 4827-4841 | 2.4 | 17 |
| 125 | Hoechst 33258 and its complex with the oligonucleotide d(CGCGAATTCGCG) 2: ¹ H NMR assignments and dynamics. <i>Magnetic Resonance in Chemistry</i> , 1992 , 30, 1064-1069 | 2.1 | 17 |
| 124 | Synthesis of (1,2-benzenediolato(2-)-O,O')oxobis(1-piperidinolato(1-)-O,N)molybdenum(VI), [MoO(C ₅ H ₁₀ NO) ₂ (C ₆ H ₄ O ₂)], and structure determination by correlated proton-carbon-13 two-dimensional NMR spectroscopy and x-ray crystallography. <i>Inorganic Chemistry</i> , 1985 , 24, 4070-4077 | 5.1 | 17 |
| 123 | Systematic comparison of sets of (¹³ C NMR spectra that are potentially identical. Confirmation of the configuration of a cuticular hydrocarbon from the cane beetle <i>Antitrogonus parvulus</i> . <i>Journal of Organic Chemistry</i> , 2014 , 79, 7477-90 | 4.2 | 16 |
| 122 | Pure shift ¹ H NMR, a robust method for revealing heteronuclear couplings in complex spectra. <i>RSC Advances</i> , 2014 , 4, 8278-8282 | 3.7 | 16 |
| 121 | Is nevirapine atropisomeric? Experimental and computational evidence for rapid conformational inversion. <i>Organic and Biomolecular Chemistry</i> , 2012 , 10, 716-9 | 3.9 | 16 |
| 120 | Design, synthesis, and evaluation of stable and taste-free erythromycin proprodrugs. <i>Journal of Medicinal Chemistry</i> , 2005 , 48, 3878-84 | 8.3 | 16 |
| 119 | Silicon-29 diffusion-ordered NMR spectroscopy (DOSY) as a tool for studying aqueous silicates. <i>Chemical Communications</i> , 2001 , 2422-3 | 5.8 | 16 |
| 118 | Spirostanol glycoside from fruits of <i>Asparagus officinalis</i> . <i>Phytochemistry</i> , 1988 , 27, 3324-3325 | 4 | 16 |
| 117 | Precise measurement of long-range heteronuclear coupling constants by a novel broadband proton-proton-decoupled CPMG-HSQC method. <i>Chemistry - A European Journal</i> , 2015 , 21, 3472-9 | 4.8 | 15 |
| 116 | Analyzing and correcting spectrometer temperature sensitivity. <i>Journal of Magnetic Resonance</i> , 2001 , 152, 234-46 | 3 | 15 |
| 115 | Combined Use of Gradient-Enhanced Techniques and Reference Deconvolution for Ultralow T ₁ Noise in 2D NMR Spectroscopy. <i>Journal of Magnetic Resonance Series A</i> , 1996 , 123, 246-252 | | 15 |
| 114 | Analysis of virtual one-bond coupling effects in heteronuclear chemical shift correlation 2D N.M.R. spectra. <i>Molecular Physics</i> , 1987 , 61, 467-483 | 1.7 | 15 |
| 113 | Real-time broadband proton-homodecoupled CLIP/CLAP-HSQC for automated measurement of heteronuclear one-bond coupling constants. <i>RSC Advances</i> , 2016 , 6, 87848-87855 | 3.7 | 15 |
| 112 | The synthesis of 2-oxyalkyl-cyclohex-2-enones, related to the bioactive natural products COTC and antheminone A, which possess anti-tumour properties. <i>Tetrahedron</i> , 2010 , 66, 9049-9060 | 2.4 | 14 |
| 111 | Dehydration of Quinate Derivatives: Synthesis of a Difluoromethylene Homologue of Shikimic Acid. <i>Synlett</i> , 2002 , 2002, 0358-0360 | 2.2 | 14 |
| 110 | Determination of the barrier to C-N bond rotation in captopril: Application of reference deconvolution to line-shape analysis. <i>Magnetic Resonance in Chemistry</i> , 1990 , 28, 820-823 | 2.1 | 14 |

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| 109 | The structure of erythromycin A in [2H6]DMSO and buffered D2O: full assignments of the 1H and 13C NMR spectra. <i>Journal of the Chemical Society Perkin Transactions II</i> , 1991 , 1489 | | 14 |
| 108 | Compensation of instrumental imperfections by deconvolution using an internal reference signal. <i>Journal of Magnetic Resonance</i> , 1988 , 80, 547-552 | | 14 |
| 107 | "Tailored detection" of nuclear magnetic resonance signals: application to the assignment of carbon-13 spectra. <i>Journal of the American Chemical Society</i> , 1978 , 100, 6763-6764 | 16.4 | 14 |
| 106 | Dissect and Divide: Putting NMR Spectra of Mixtures under the Knife. <i>Journal of the American Chemical Society</i> , 2019 , 141, 5766-5771 | 16.4 | 13 |
| 105 | Practical aspects of real-time pure shift HSQC experiments. <i>Magnetic Resonance in Chemistry</i> , 2018 , 56, 993-1005 | 2.1 | 13 |
| 104 | Matrix-assisted diffusion-ordered spectroscopy: choosing a matrix. <i>Magnetic Resonance in Chemistry</i> , 2016 , 54, 815-820 | 2.1 | 13 |
| 103 | Resolving complex mixtures: trilinear diffusion data. <i>Journal of Biomolecular NMR</i> , 2014 , 58, 251-7 | 3 | 13 |
| 102 | The acyl nitroso Diels-Alder (ANDA) reaction of sorbate derivatives: an X-ray and 15N NMR study with an application to amino-acid synthesis. <i>Organic and Biomolecular Chemistry</i> , 2009 , 7, 4531-8 | 3.9 | 13 |
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