

Alexandar L Hansen

List of Publications by Year in descending order

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41
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394421

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docs citations

42
times ranked

1494
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|---|------|-----------|
| 1 | Fundamental and practical aspects of machine learning for the peak picking of biomolecular NMR spectra. <i>Journal of Biomolecular NMR</i> , 2022, 76, 49-57. | 2.8 | 5 |
| 2 | 2D NMR-Based Metabolomics with HSQC/TOCSY NOAH Supersequences. <i>Analytical Chemistry</i> , 2021, 93, 6112-6119. | 6.5 | 28 |
| 3 | From Selection to Instruction and Back: Competing Conformational Selection and Induced Fit Pathways in Abiotic Hosts. <i>Angewandte Chemie - International Edition</i> , 2021, 60, 19942-19948. | 13.8 | 18 |
| 4 | Increasing sensitivity and versatility in NMR supersequences with new HSQC-based modules. <i>Journal of Magnetic Resonance</i> , 2021, 329, 107027. | 2.1 | 12 |
| 5 | Observation of Sub-Microsecond Protein Methyl-Side Chain Dynamics by Nanoparticle-Assisted NMR Spin Relaxation. <i>Journal of the American Chemical Society</i> , 2021, 143, 13593-13604. | 13.7 | 10 |
| 6 | DEEP picker is a deep neural network for accurate deconvolution of complex two-dimensional NMR spectra. <i>Nature Communications</i> , 2021, 12, 5229. | 12.8 | 55 |
| 7 | 4,15-Dimethyl-7,12-diazoniatriacyclo[10.4.0.0 ^{2,7}]hexadeca-1(12),2,4,6,13,15-hexaene dibromide monohydrate. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2020, 76, 1467-1471. | 0.5 | 0 |
| 8 | Functional protein dynamics on uncharted time scales detected by nanoparticle-assisted NMR spin relaxation. <i>Science Advances</i> , 2019, 5, eaax5560. | 10.3 | 32 |
| 9 | Extreme Nonuniform Sampling for Protein NMR Dynamics Studies in Minimal Time. <i>Journal of the American Chemical Society</i> , 2019, 141, 16829-16838. | 13.7 | 12 |
| 10 | Real-Time Pure Shift HSQC NMR for Untargeted Metabolomics. <i>Analytical Chemistry</i> , 2019, 91, 2304-2311. | 6.5 | 25 |
| 11 | Resonance assignments of wild-type and two cysteine-free variants of the four-helix bundle protein, Rop. <i>Biomolecular NMR Assignments</i> , 2018, 12, 345-350. | 0.8 | 1 |
| 12 | Quantitative Binding Behavior of Intrinsically Disordered Proteins to Nanoparticle Surfaces at Individual Residue Level. <i>Chemistry - A European Journal</i> , 2018, 24, 16997-17001. | 3.3 | 21 |
| 13 | Non-Uniform and Absolute Minimal Sampling for High-Throughput Multidimensional NMR Applications. <i>Chemistry - A European Journal</i> , 2018, 24, 11535-11544. | 3.3 | 14 |
| 14 | Differential Conformational Dynamics Encoded by the Linker between Quasi RNA Recognition Motifs of Heterogeneous Nuclear Ribonucleoprotein H. <i>Journal of the American Chemical Society</i> , 2018, 140, 11661-11673. | 13.7 | 11 |
| 15 | ¹ H, ¹³ C, ¹⁵ N resonance assignment of recombinant <i>Euplotes raikovi</i> protein Er-23. <i>Biomolecular NMR Assignments</i> , 2018, 12, 291-295. | 0.8 | 0 |
| 16 | Absolute Minimal Sampling of Homonuclear 2D NMR TOCSY Spectra for High-Throughput Applications of Complex Mixtures. <i>Angewandte Chemie - International Edition</i> , 2017, 56, 8149-8152. | 13.8 | 16 |
| 17 | The Michaelis Complex of Arginine Kinase Samples the Transition State at a Frequency That Matches the Catalytic Rate. <i>Journal of the American Chemical Society</i> , 2017, 139, 4846-4853. | 13.7 | 14 |
| 18 | Absolute Minimal Sampling of Homonuclear 2D NMR TOCSY Spectra for High-Throughput Applications of Complex Mixtures. <i>Angewandte Chemie</i> , 2017, 129, 8261-8264. | 2.0 | 8 |

| # | ARTICLE | IF | CITATIONS |
|----|---|------|-----------|
| 19 | Synthesis of 6-phosphofructose aspartic acid and some related Amadori compounds. Carbohydrate Research, 2016, 431, 1-5. | 2.3 | 16 |
| 20 | Residue-Specific Interactions of an Intrinsically Disordered Protein with Silica Nanoparticles and Their Quantitative Prediction. Journal of Physical Chemistry C, 2016, 120, 24463-24468. | 3.1 | 28 |
| 21 | Absolute Minimal Sampling in High-Dimensional NMR Spectroscopy. Angewandte Chemie - International Edition, 2016, 55, 14169-14172. | 13.8 | 19 |
| 22 | Absolut minimales Sampling in der hochdimensionalen NMR-Spektroskopie. Angewandte Chemie, 2016, 128, 14376-14379. | 2.0 | 5 |
| 23 | Rapid Determination of Fast Protein Dynamics from NMR Chemical Exchange Saturation Transfer Data. Angewandte Chemie, 2016, 128, 3169-3171. | 2.0 | 1 |
| 24 | Rapid Determination of Fast Protein Dynamics from NMR Chemical Exchange Saturation Transfer Data. Angewandte Chemie - International Edition, 2016, 55, 3117-3119. | 13.8 | 15 |
| 25 | Kinetic Cooperativity in Human Pancreatic Glucokinase Originates from Millisecond Dynamics of the Small Domain. Angewandte Chemie, 2015, 127, 8247-8250. | 2.0 | 7 |
| 26 | Kinetic Cooperativity in Human Pancreatic Glucokinase Originates from Millisecond Dynamics of the Small Domain. Angewandte Chemie - International Edition, 2015, 54, 8129-8132. | 13.8 | 29 |
| 27 | Characterizing Slow Chemical Exchange in Nucleic Acids by Carbon CEST and Low Spin-Lock Field ^{13}C NMR Spectroscopy. Journal of the American Chemical Society, 2014, 136, 20-23. | 13.7 | 82 |
| 28 | Measurement of histidine pK _a values and tautomer populations in invisible protein states. Proceedings of the National Academy of Sciences of the United States of America, 2014, 111, E1705-12. | 7.1 | 111 |
| 29 | Probing slowly exchanging protein systems via ^{13}C -CEST: monitoring folding of the Im7 protein. Journal of Biomolecular NMR, 2013, 55, 279-289. | 2.8 | 24 |
| 30 | The Role of Ligands on the Equilibria Between Functional States of a G Protein-Coupled Receptor. Journal of the American Chemical Society, 2013, 135, 9465-9474. | 13.7 | 156 |
| 31 | Quantifying Millisecond Exchange Dynamics in Proteins by CPMG Relaxation Dispersion NMR Using Side-Chain ^1H Probes. Journal of the American Chemical Society, 2012, 134, 3178-3189. | 13.7 | 55 |
| 32 | Nonnative Interactions in the FF Domain Folding Pathway from an Atomic Resolution Structure of a Sparsely Populated Intermediate: An NMR Relaxation Dispersion Study. Journal of the American Chemical Society, 2011, 133, 10974-10982. | 13.7 | 37 |
| 33 | Characterizing RNA dynamics at atomic resolution using solution-state NMR spectroscopy. Nature Methods, 2011, 8, 919-931. | 19.0 | 131 |
| 34 | Quantifying millisecond time-scale exchange in proteins by CPMG relaxation dispersion NMR spectroscopy of side-chain carbonyl groups. Journal of Biomolecular NMR, 2011, 50, 347-355. | 2.8 | 28 |
| 35 | Variable helix elongation as a tool to modulate RNA alignment and motional couplings. Journal of Magnetic Resonance, 2010, 202, 117-121. | 2.1 | 19 |
| 36 | Extending the Range of Microsecond-to-Millisecond Chemical Exchange Detected in Labeled and Unlabeled Nucleic Acids by Selective Carbon ^{13}C NMR Spectroscopy. Journal of the American Chemical Society, 2009, 131, 3818-3819. | 13.7 | 109 |

| # | ARTICLE | IF | CITATIONS |
|----|--|------|-----------|
| 37 | Characterizing Complex Dynamics in the Transactivation Response Element Apical Loop and Motional Correlations with the Bulge by NMR, Molecular Dynamics, and Mutagenesis. <i>Biophysical Journal</i> , 2008, 95, 3906-3915. | 0.5 | 65 |
| 38 | Dynamics of Large Elongated RNA by NMR Carbon Relaxation. <i>Journal of the American Chemical Society</i> , 2007, 129, 16072-16082. | 13.7 | 85 |
| 39 | Characterizing the relative orientation and dynamics of RNA A-form helices using NMR residual dipolar couplings. <i>Nature Protocols</i> , 2007, 2, 1536-1546. | 12.0 | 56 |
| 40 | Insight into the CSA tensors of nucleobase carbons in RNA polynucleotides from solution measurements of residual CSA: Towards new long-range orientational constraints. <i>Journal of Magnetic Resonance</i> , 2006, 179, 299-307. | 2.1 | 56 |
| 41 | Pulsed field gradient NMR investigation of solubilization equilibria in amino acid and dipeptide terminated micellar and polymeric surfactant solutions. <i>Magnetic Resonance in Chemistry</i> , 2002, 40, 755-761. | 1.9 | 13 |