

# Raiker Witter

## List of Publications by Year in descending order

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Version: 2024-02-01

47  
papers

1,610  
citations

331670

21  
h-index

289244

40  
g-index

48  
all docs

48  
docs citations

48  
times ranked

1805  
citing authors

| #  | ARTICLE   | IF   | CITATIONS |
|----|---|------|-----------|
| 1  | Synthesis and characterization of $\text{Ca}_{1-x}\text{Sm}_x\text{F}_{2+x}$ (2+) Tj ETQq1 1 0.784314 rgBT<br>Processing Communications, 2021, 3, e226.   | 0.9  | 3         |
| 2  | $\text{CaF}_2$ solid-state electrolytes prepared by vapor pressure exposure and solid synthesis for defect and ionic conductivity tuning. Material Design and Processing Communications, 2020, 2, e76.                                | 0.9  | 5         |
| 3  | Screening of Nutraceuticals and Plant Extracts for Inhibition of Amyloid- $\beta^2$ Fibrillation. Journal of Alzheimer's Disease, 2020, 73, 1003-1012.  | 2.6  | 5         |
| 4  | Compound's Pre-Screening of <i>Withania somnifera</i> , <i>Bacopa monnieri</i> , and <i>Centella asiatica</i> Extracts. Journal of Biosciences and Medicines, 2020, 08, 80-98.  | 0.2  | 0         |
| 5  | Investigation of backbone dynamics and local geometry of bio-molecules using calculated NMR chemical shifts and anisotropies. Journal of Biomolecular NMR, 2019, 73, 727-741.   | 2.8  | 4         |
| 6  | Introducing Interlayer Electrolytes: Toward Room-Temperature High-Potential Solid-State Rechargeable Fluoride Ion Batteries. ACS Applied Energy Materials, 2019, 2, 1553-1562.  | 5.1  | 38        |
| 7  | Crystal phase and surface defect driven synthesis of $\text{Pb}_{1-x}\text{Sn}_x\text{F}_2$ solid solution electrolyte for fluoride ion batteries. Journal of Electroanalytical Chemistry, 2019, 845, 154-159.                        | 3.8  | 17        |
| 8  | Testing Mg as an anode against $\text{BiF}_3$ and $\text{SnF}_2$ cathodes for room temperature rechargeable fluoride ion batteries. Materials Letters, 2019, 244, 159-162.  | 2.6  | 22        |
| 9  | Mechanochemical synthesis of solid-state electrolyte $\text{Sm}_{1-x}\text{Ca}_x\text{F}_3$ for batteries and other electrochemical devices. Materials Letters, 2019, 244, 22-26.   | 2.6  | 13        |
| 10 | Surface defect-enhanced conductivity of calcium fluoride for electrochemical applications. Material Design and Processing Communications, 2019, 1, e44.   | 0.9  | 13        |
| 11 | $^1\text{H}$ line width dependence on MAS speed in solid state NMR – Comparison of experiment and simulation. Journal of Magnetic Resonance, 2018, 291, 32-39.  | 2.1  | 80        |
| 12 | Structure and electrochemical properties of $\text{Na}_{2-x}\text{V}_3\text{P}_2\text{O}_{13}$ ( $x = 0$ and 1): a promising cathode material for sodium-ion batteries. Journal of Materials Chemistry A, 2018, 6, 6947-6958.         | 10.3 | 9         |
| 13 | Medical Plants and Nutraceuticals for Amyloid- $\beta^2$ Fibrillation Inhibition. Journal of Alzheimer's Disease Reports, 2018, 2, 239-252.   | 2.2  | 9         |
| 14 | Synthesis of Fast Fluoride-Ion-Conductive Fluorite-Type $\text{Ba}_{1-x}\text{Sb}_x\text{F}_{2+x}$ (0.1 to 0.4): A Potential Solid Electrolyte for Fluoride-Ion Batteries. ACS Applied Materials & Interfaces, 2018, 10, 17249-17256. | 8.0  | 37        |
| 15 | Room-Temperature, Rechargeable Solid-State Fluoride-Ion Batteries. ACS Applied Energy Materials, 2018, 1, 4766-4775.  | 5.1  | 80        |
| 16 | Lithiation-driven structural transition of $\text{VO}_2\text{F}$ into disordered rock-salt $\text{Li}_x\text{VO}_2\text{F}$ . RSC Advances, 2016, 6, 65112-65118.   | 3.6  | 19        |
| 17 | Molecular dynamics simulations on PGLa using NMR orientational constraints. Journal of Biomolecular NMR, 2015, 63, 265-274.   | 2.8  | 8         |
| 18 | Fast Atomic Charge Calculation for Implementation into a Polarizable Force Field and Application to an Ion Channel Protein. Journal of Chemistry, 2015, 2015, 1-14.   | 1.9  | 6         |

| #  | ARTICLE   | IF   | CITATIONS |
|----|---|------|-----------|
| 19 | Li <sup>+</sup> intercalation in isostructural Li <sub>2</sub> VO <sub>3</sub> and Li <sub>2</sub> VO <sub>2</sub> F with O <sup>2-</sup> and mixed O <sup>2-</sup> /F <sup>-</sup> anions. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 17288-17295.   | 2.8  | 67        |
| 20 | Disordered Lithium-Rich Oxyfluoride as a Stable Host for Enhanced Li <sup>+</sup> Intercalation Storage. <i>Advanced Energy Materials</i> , 2015, 5, 1401814.   | 19.5 | 162       |
| 21 | Beneficial effects of stoichiometry and nanostructure for a LiBH <sub>4</sub> -MgH <sub>2</sub> hydrogen storage system. <i>Journal of Materials Chemistry A</i> , 2014, 2, 66-72.  | 10.3 | 18        |
| 22 | Solid Electrolytes for Fluoride Ion Batteries: Ionic Conductivity in Polycrystalline Tysonite-Type Fluorides. <i>ACS Applied Materials &amp; Interfaces</i> , 2014, 6, 2103-2110.   | 8.0  | 131       |
| 23 | Altered reaction pathways of eutectic LiBH <sub>4</sub> -Mg(BH <sub>4</sub> ) <sub>2</sub> by nanoconfinement. <i>Journal of Materials Chemistry A</i> , 2013, 1, 3379.   | 10.3 | 52        |
| 24 | Nanostructured Fluorite-Type Fluorides As Electrolytes for Fluoride Ion Batteries. <i>Journal of Physical Chemistry C</i> , 2013, 117, 4943-4950.   | 3.1  | 145       |
| 25 | Multilayered core-shell structure of polyol-stabilized calcium fluoride nanoparticles characterized by NMR. <i>Journal of Colloid and Interface Science</i> , 2013, 390, 250-257.   | 9.4  | 20        |
| 26 | Influence of Nanoconfinement on Reaction Pathways of Complex Metal Hydrides. <i>Energy Procedia</i> , 2012, 29, 731-737.  | 1.8  | 11        |
| 27 | <sup>29</sup> Si NMR Shielding Tensors in Triphenylsilanes – <sup>29</sup> Si Solid State NMR Experiments and DFT-GLO Calculations. <i>Zeitschrift Fur Anorganische Und Allgemeine Chemie</i> , 2012, 638, 935-944.   | 1.2  | 22        |
| 28 | Catalytic Influence of Various Cerium Precursors on the Hydrogen Sorption Properties of NaAlH <sub>4</sub> . <i>Advanced Energy Materials</i> , 2012, 2, 560-568.   | 19.5 | 38        |
| 29 | Calculation of fluorine chemical shift tensors for the interpretation of oriented <sup>19</sup> F-NMR spectra of gramicidin A in membranes. <i>Physical Chemistry Chemical Physics</i> , 2009, 11, 7048.  | 2.8  | 30        |
| 30 | Spectral assignments and anisotropy data of cellulose <sup>13</sup> C-NMR chemical shift data of cellulose <sup>13</sup> C determined by INADEQUATE and RAI techniques applied to uniformly <sup>13</sup> C-labeled bacterial celluloses of different <i>Gluconacetobacter xylinus</i> strains. <i>Magnetic Resonance in Chemistry</i> , 2008, 46, 1030-1036. | 1.9  | 16        |
| 31 | Solid state <sup>19</sup> F NMR parameters of fluorine-labeled amino acids. Part I: Aromatic substituents. <i>Journal of Magnetic Resonance</i> , 2008, 191, 7-15.  | 2.1  | 57        |
| 32 | Solid-State <sup>19</sup> F NMR Spectroscopy Reveals That Trp <sub>41</sub> Participates in the Gating Mechanism of the M2 Proton Channel of Influenza A Virus. <i>Journal of the American Chemical Society</i> , 2008, 130, 918-924.   | 13.7 | 47        |
| 33 | Low-E probe for <sup>19</sup> F- <sup>1</sup> H NMR of dilute biological solids. <i>Journal of Magnetic Resonance</i> , 2007, 189, 182-189.   | 2.1  | 39        |
| 34 | All-atom molecular dynamics simulations using orientational constraints from anisotropic NMR samples. <i>Journal of Biomolecular NMR</i> , 2007, 38, 23-39.   | 2.8  | 27        |
| 35 | <sup>13</sup> C Chemical Shift Constrained Crystal Structure Refinement of Cellulose I <sub>II</sub> and Its Verification by NMR Anisotropy Experiments. <i>Macromolecules</i> , 2006, 39, 6125-6132.   | 4.8  | 74        |
| 36 | NMR Chemical Shift Powder Pattern Recoupling at High Spinning Speed and Theoretical Tensor Evaluation Applied to Silk Fibroin. <i>Journal of the American Chemical Society</i> , 2006, 128, 2236-2243.  | 13.7 | 22        |

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|----|---|-----|-----------|
| 37 | 3D Structure Elucidation Using NMR Chemical Shifts. ChemInform, 2005, 36, no.   | 0.0 | 0         |
| 38 | 3D Structure Elucidation Using NMR Chemical Shifts. Annual Reports on NMR Spectroscopy, 2004, , 53-104.   | 1.5 | 17        |
| 39 | Crystal Structure Refinements of Cellulose Polymorphs using Solid State <sup>13</sup> C Chemical Shifts. Cellulose, 2003, 10, 189-199.  | 4.9 | 39        |
| 40 | Powder pattern recoupling at 10kHz spinning speed applied to cellulose. Journal of Magnetic Resonance, 2003, 161, 35-42.  | 2.1 | 23        |
| 41 | Chemical shift driven geometry optimization. Journal of Computational Chemistry, 2002, 23, 298-305.   | 3.3 | 26        |
| 42 | Structure determination of a pseudotriptide zinc complex with the COSMOS-NMR force field and DFT methods. Journal of Biomolecular NMR, 2002, 24, 277-289.   | 2.8 | 17        |
| 43 | Calculation of solid-state <sup>13</sup> C NMR spectra of cellulose I <sup>±</sup> , I <sup>±2</sup> and II using a semi-empirical approach and molecular dynamics. Macromolecular Chemistry and Physics, 2000, 201, 1930-1939. | 2.2 | 19        |
| 44 | Complexation of metal ions by pseudotriptides with different functionalized N-alkyl residues. International Journal of Peptide Research and Therapeutics, 2000, 7, 133-141.   | 0.1 | 5         |
| 45 | Complexation of metal ions by pseudotriptides with different functionalized N-alkyl residues. International Journal of Peptide Research and Therapeutics, 2000, 7, 133-141.   | 0.1 | 1         |
| 46 | New 2D NMR experiments for determining the structure of phosphate glasses: a review. Journal of Non-Crystalline Solids, 2000, 263-264, 61-72.   | 3.1 | 42        |
| 47 | Measurements of chain length distributions in calcium phosphate glasses using 2D double quantum NMR. Solid State Nuclear Magnetic Resonance, 1998, 13, 189-200.   | 2.3 | 75        |