

# Kaan Atak

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/9475453/publications.pdf>

Version: 2024-02-01

28  
papers

528  
citations

687363

13  
h-index

642732

23  
g-index

29  
all docs

29  
docs citations

29  
times ranked

950  
citing authors

| #  | ARTICLE  | IF  | CITATIONS |
|----|--|-----|-----------|
| 1  | The electronic structure and deexcitation pathways of an isolated metalloporphyrin ion resolved by metal L-edge spectroscopy. <i>Chemical Science</i> , 2021, 12, 3966-3976.   | 7.4 | 3         |
| 2  | Probing Structural Information of Gas-Phase Peptides by Near-Edge X-ray Absorption Mass Spectrometry. <i>Journal of the American Society for Mass Spectrometry</i> , 2021, 32, 670-684.  | 2.8 | 5         |
| 3  | Site-Selective Dissociation upon Sulfur L-Edge X-ray Absorption in a Gas-Phase Protonated Peptide. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 1215-1221.   | 4.6 | 20        |
| 4  | Electronic Structure of Aqueous [Co(bpy) <sub>3</sub> ] <sup>2+/3+</sup> Electron Mediators. <i>Inorganic Chemistry</i> , 2019, 58, 4731-4740.   | 4.0 | 17        |
| 5  | Mimicking cellular phospholipid bilayer packing creates predictable crystalline molecular metal-organophosphonate macrocycles and cages. <i>CrystEngComm</i> , 2018, 20, 2152-2158.  | 2.6 | 6         |
| 6  | Chemical bonding in aqueous hexacyano cobaltate from photon- and electron-detection perspectives. <i>Scientific Reports</i> , 2017, 7, 40811.  | 3.3 | 14        |
| 7  | Undistorted X-ray Absorption Spectroscopy Using s-Core-Orbital Emissions. <i>Journal of Physical Chemistry A</i> , 2016, 120, 2808-2814.   | 2.5 | 21        |
| 8  | Analysis of the Electronic Structure of Aqueous Urea and Its Derivatives: A Systematic Soft X-ray DFT Approach. <i>Chemistry - A European Journal</i> , 2016, 22, 12040-12049.   | 3.3 | 6         |
| 9  | Joint Analysis of Radiative and Non-Radiative Electronic Relaxation Upon X-ray Irradiation of Transition Metal Aqueous Solutions. <i>Scientific Reports</i> , 2016, 6, 24659.  | 3.3 | 38        |
| 10 | Impacts of Conformational Geometries in Fluorinated Alkanes. <i>Scientific Reports</i> , 2016, 6, 31382.   | 3.3 | 4         |
| 11 | Influence of the Outer Ligands on Metal-to-Ligand Charge Transfer in Solvated Manganese Porphyrins. <i>Inorganic Chemistry</i> , 2016, 55, 22-28.  | 4.0 | 10        |
| 12 | Fluorination-dependent molecular orbital occupancy in ring-shaped perfluorocarbons. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 18337-18343.  | 2.8 | 6         |
| 13 | Local Energy Gap Opening Induced by Hemin Dimerization in Aqueous Solution. <i>Journal of Physical Chemistry B</i> , 2015, 119, 3058-3062.   | 2.6 | 11        |
| 14 | Ti <sup>3+</sup> Aqueous Solution: Hybridization and Electronic Relaxation Probed by State-Dependent Electron Spectroscopy. <i>Journal of Physical Chemistry B</i> , 2015, 119, 10607-10615.   | 2.6 | 14        |
| 15 | Intermolecular bonding of hemin in solution and in solid state probed by N K-edge X-ray spectroscopies. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 29000-29006.  | 2.8 | 9         |
| 16 | Co(III) protoporphyrin IX chloride in solution: spin-state and metal coordination revealed from resonant inelastic X-ray scattering and electronic structure calculations. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 3409-3414. | 2.8 | 12        |
| 17 | Bokarev <i>et al.</i> Reply. <i>Physical Review Letters</i> , 2014, 112, 129303.   | 7.8 | 5         |
| 18 | Chemical Bonding in Aqueous Ferrocyanide: Experimental and Theoretical X-ray Spectroscopic Study. <i>Journal of Physical Chemistry B</i> , 2014, 118, 1555-1563.   | 2.6 | 61        |

| #  | ARTICLE   | IF   | CITATIONS |
|----|---|------|-----------|
| 19 | Electronic structural insights into efficient MnO <sub>x</sub> catalysts. Journal of Materials Chemistry A, 2014, 2, 18199-18203.   | 10.3 | 40        |
| 20 | Electronic Structure of Hemin in Solution Studied by Resonant X-ray Emission Spectroscopy and Electronic Structure Calculations. Journal of Physical Chemistry B, 2014, 118, 9938-9943.               | 2.6  | 16        |
| 21 | Assistance of the Iron Porphyrin Ligands to the Binding Interaction between the Fe Center and Small Molecules in Solution. Journal of Physical Chemistry B, 2014, 118, 9371-9377.                     | 2.6  | 7         |
| 22 | Direct Observation of Molecular Orbital Mixing in a Solvated Organometallic Complex. Angewandte Chemie - International Edition, 2013, 52, 9841-9844.  | 13.8 | 60        |
| 23 | Local electronic structure of aqueous zinc acetate: oxygen K-edge X-ray absorption and emission spectroscopy on micro-jets. Physical Chemistry Chemical Physics, 2013, 15, 8046.                      | 2.8  | 14        |
| 24 | Nature of the Chemical Bond of Aqueous Fe <sup>2+</sup> Probed by Soft X-ray Spectroscopies and ab Initio Calculations. Journal of Physical Chemistry B, 2013, 117, 12613-12618.                      | 2.6  | 44        |
| 25 | Probing Coster-Kronig Transitions in Aqueous Fe <sup>2+</sup> Solution Using Inverse Partial and Partial Fluorescence Yield at the L-Edge. Journal of Physical Chemistry Letters, 2012, 3, 1619-1623. | 4.6  | 36        |
| 26 | DMSO-Water Clustering in Solution Observed in Soft X-ray Spectra. Journal of Physical Chemistry Letters, 2012, 3, 3697-3701.  | 4.6  | 30        |
| 27 | The Chemical Bond in Carbonyl and Sulfinyl Groups Studied by Soft X-ray Spectroscopy and ab Initio Calculations. ChemPhysChem, 2012, 13, 3106-3111.   | 2.1  | 12        |
| 28 | Chaoticity analysis of the current through pure, hydrogenated and hydrophobically modified PEG-Si thin films under varying relative humidity. Open Physics, 2009, 7, .                                | 1.7  | 3         |