

# Mihails Arhangeliskis

## List of Publications by Year in descending order

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Version: 2024-02-01

52  
papers

1,188  
citations

331259  
21  
h-index

395343  
33  
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61  
all docs

61  
docs citations

61  
times ranked

1513  
citing authors

#	ARTICLE	IF	CITATIONS
1	Metal-organic frameworks as hypergolic additives for hybrid rockets. <i>Chemical Science</i> , 2022, 13, 3424-3436.	3.7	14
2	Mechanochemical Synthesis of Polymorphic Urea-Adipic Acid Cocrystal as a Sustained-Release Nitrogen Source. <i>ChemSusChem</i> , 2022, 15, e202102445.	3.6	9
3	Open versus Interpenetrated: Switchable Supramolecular Trajectories in Mechanosynthesis of a Halogen-Bonded Borromean Network. <i>CheM</i> , 2021, 7, 146-154.	5.8	17
4	Mechanochemical reactivity inhibited, prohibited and reversed by liquid additives: examples from crystal-form screens. <i>Chemical Science</i> , 2021, 12, 3264-3269.	3.7	25
5	Influence of ligand composition on crystal structure formation - isostructurality and morphotropism. <i>CrystEngComm</i> , 2021, 23, 317-323.	1.3	6
6	Mechanosynthesis of Eutectics of Anti-inflammatory Drug Ethenzamide - A Comparison with Analogous Cocrystals. <i>Chemistry Methods</i> , 2021, 1, 408-414.	1.8	3
7	Real-Time Observation of "Soft-Magic-Size Clusters during Hydrolysis of the Model Metallodrug Bismuth Disalicylate. <i>Journal of the American Chemical Society</i> , 2021, 143, 16332-16336.	6.6	5
8	Polymorphism and surface diversity arising from stress-induced transformations - the case of multicomponent forms of carbamazepine. <i>Acta Crystallographica Section B: Structural Science, Crystal Engineering and Materials</i> , 2021, 77, 54-67.	0.5	8
9	Simplifying and expanding the scope of boron imidazolate framework (BIF) synthesis using mechanochemistry. <i>Chemical Science</i> , 2021, 12, 14499-14506.	3.7	7
10	Understanding stress-induced disorder and breakage in organic crystals: beyond crystal structure anisotropy. <i>Chemical Science</i> , 2021, 12, 14270-14280.	3.7	5
11	Mechanochemical Synthesis and Physicochemical Characterization of Previously Unreported Praziquantel Solvates with 2-Pyrrolidone and Acetic Acid. <i>Pharmaceutics</i> , 2021, 13, 1606.	2.0	10
12	Disappearing Polymorphs in Metal-Organic Framework Chemistry: Unexpected Stabilization of a Layered Polymorph over an Interpenetrated Three-Dimensional Structure in Mercury Imidazolate. <i>Chemistry - A European Journal</i> , 2020, 26, 1811-1818.	1.7	25
13	Challenging the Ostwald rule of stages in mechanochemical cocrystallisation. <i>Chemical Science</i> , 2020, 11, 10092-10100.	3.7	49
14	Linker Substituents Control the Thermodynamic Stability in Metal-Organic Frameworks. <i>Journal of the American Chemical Society</i> , 2020, 142, 21720-21729.	6.6	36
15	Solvent-free ageing reactions of rare earth element oxides: from geomimetic synthesis of new metal-organic materials towards a simple, environmentally friendly separation of scandium. <i>Green Chemistry</i> , 2020, 22, 4364-4375.	4.6	8
16	<i>Ab Initio</i> Prediction of Metal-Organic Framework Structures. <i>Chemistry of Materials</i> , 2020, 32, 5835-5844.	3.2	11
17	Mechanochemical reactions of cocrystals: comparing theory with experiment in the making and breaking of halogen bonds in the solid state. <i>Chemical Communications</i> , 2020, 56, 8293-8296.	2.2	18
18	Mechanochemical Formation of Racemic Praziquantel Hemihydrate with Improved Biopharmaceutical Properties. <i>Pharmaceutics</i> , 2020, 12, 289.	2.0	21

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19	On the kinetics of solvate formation through mechanochemistry. <i>CrystEngComm</i> , 2019, 21, 2097-2104.	1.3	14
20	Hypergolic Triggers as Co-crystal Formers: Co-crystallization for Creating New Hypergolic Materials with Tunable Energy Content. <i>Angewandte Chemie - International Edition</i> , 2019, 58, 18399-18404.	7.2	25
21	Hypergolic Triggers as Co-crystal Formers: Co-crystallization for Creating New Hypergolic Materials with Tunable Energy Content. <i>Angewandte Chemie</i> , 2019, 131, 18570-18575.	1.6	7
22	Metal-Organic Frameworks as Fuels for Advanced Applications: Evaluating and Modifying the Combustion Energy of Popular MOFs. <i>Chemistry of Materials</i> , 2019, 31, 4882-4888.	3.2	21
23	Theoretical Prediction and Experimental Evaluation of Topological Landscape and Thermodynamic Stability of a Fluorinated Zeolitic Imidazolate Framework. <i>Chemistry of Materials</i> , 2019, 31, 3777-3783.	3.2	31
24	Hypergolic zeolitic imidazolate frameworks (ZIFs) as next-generation solid fuels: Unlocking the latent energetic behavior of ZIFs. <i>Science Advances</i> , 2019, 5, eaav9044.	4.7	52
25	Cocrystal Dissociation under Controlled Humidity: A Case Study of Caffeine-Glutaric Acid Cocrystal Polymorphs. <i>Organic Process Research and Development</i> , 2019, 23, 845-851.	1.3	34
26	Cocrystal trimorphism as a consequence of the orthogonality of halogen- and hydrogen-bonds synthons. <i>Chemical Communications</i> , 2019, 55, 14066-14069.	2.2	13
27	Air oxidation of sulfur mustard gas simulants using a pyrene-based metal-organic framework photocatalyst. <i>Beilstein Journal of Nanotechnology</i> , 2019, 10, 2422-2427.	1.5	14
28	Halogen-bonded cocrystallization with phosphorus, arsenic and antimony acceptors. <i>Nature Communications</i> , 2019, 10, 61.	5.8	78
29	Experimental and Theoretical Investigation of Structures, Stoichiometric Diversity, and Bench Stability of Cocrystals with a Volatile Halogen Bond Donor. <i>Crystal Growth and Design</i> , 2018, 18, 2387-2396.	1.4	19
30	Mechanochemical Synthesis of Olanzapine Salts and Their Hydration Stability Study Using Powder X-ray Diffraction. <i>Crystal Growth and Design</i> , 2018, 18, 2138-2150.	1.4	27
31	Computational evaluation of metal pentazolate frameworks: inorganic analogues of azolate metal-organic frameworks. <i>Chemical Science</i> , 2018, 9, 3367-3375.	3.7	39
32	Crystallization at Solvent Interfaces Enables Access to a Variety of Cocrystal Polymorphs and Hydrates. <i>Crystal Growth and Design</i> , 2018, 18, 3263-3268.	1.4	15
33	Towards the systematic crystallisation of molecular ionic cocrystals: insights from computed crystal form landscapes. <i>Faraday Discussions</i> , 2018, 211, 401-424.	1.6	20
34	Structure searching methods: general discussion. <i>Faraday Discussions</i> , 2018, 211, 133-180.	1.6	3
35	Crystal structure evaluation: calculating relative stabilities and other criteria: general discussion. <i>Faraday Discussions</i> , 2018, 211, 325-381.	1.6	7
36	Time-Dependent Density-Functional Theory for Modeling Solid-State Fluorescence Emission of Organic Multicomponent Crystals. <i>Journal of Physical Chemistry A</i> , 2018, 122, 7514-7521.	1.1	9

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37	Welcoming Gallium- and Indium-Fumarate MOFs to the Family: Synthesis, Comprehensive Characterization, Observation of Porous Hydrophobicity, and CO <sub>2</sub> Dynamics. ACS Applied Materials & Interfaces, 2018, 10, 28582-28596.	4.0	30
38	Real-time monitoring of mechanochemical formation of pharmaceutical cocrystals using synchrotron X-ray diffraction. Acta Crystallographica Section A: Foundations and Advances, 2018, 74, a277-a277.	0.0	0
39	Understanding the stepwise mechanism in the formation of halogen-bonded organic cocrystals by mechanochemistry. Acta Crystallographica Section A: Foundations and Advances, 2018, 74, a372-a372.	0.0	0
40	Experimental and Theoretical Evaluation of the Stability of True MOF Polymorphs Explains Their Mechanochemical Interconversions. Journal of the American Chemical Society, 2017, 139, 7952-7957.	6.6	93
41	Crystal structure landscape of ethenzamide: a physicochemical property study. CrystEngComm, 2017, 19, 826-833.	1.3	37
42	Combined use of solid-state NMR spectroscopy and theoretical modelling as a method of structure determination. Acta Crystallographica Section A: Foundations and Advances, 2017, 73, a421-a421.	0.0	0
43	Rationalization of the Color Properties of Fluorescein in the Solid State: A Combined Computational and Experimental Study. Chemistry - A European Journal, 2016, 22, 10065-10073.	1.7	24
44	On the predictability of supramolecular interactions in molecular cocrystals – the view from the bench. CrystEngComm, 2016, 18, 5434-5439.	1.3	47
45	Investigation of an Amide-Pseudo Amide Hydrogen Bonding Motif within a Series of Theophylline:Amide Cocrystals. Crystal Growth and Design, 2016, 16, 51-58.	1.4	30
46	Solid-state photoreactivity of 9-substituted acridizinium bromide salts. CrystEngComm, 2014, 16, 10830-10836.	1.3	12
47	Polymorphs, hydrates and solvates of a co-crystal of caffeine with anthranilic acid. Acta Crystallographica Section B: Structural Science, Crystal Engineering and Materials, 2014, 70, 72-80.	0.5	62
48	Advantages of mechanochemical cocrystallisation in the solid-state chemistry of pigments: colour-tuned fluorescein cocrystals. CrystEngComm, 2013, 15, 6289.	1.3	67
49	Solid state grinding as a tool to aid enantiomeric resolution by cocrystallisation. Chemical Communications, 2012, 48, 11340.	2.2	46
50	Mechanochemical synthesis of pyrazine:dicarboxylic acid cocrystals and a study of dissociation by quantitative phase analysis. CrystEngComm, 2012, 14, 5203.	1.3	34
51	Crystal structure of 1-(N-morpholinomethyl)spirobi(4-methyl-3-oxo-2,5-dioxo-1-silacyclopentane) dihydrate, C <sub>11</sub> H <sub>19</sub> NO <sub>7</sub> Si · 2H <sub>2</sub> O. Zeitschrift Fur Kristallographie - New Crystal Structures, 2011, 226, .	0.1	0
52	Effect of Vehicle Composition on the Preparation of Different Types of Dapsone Crystals for Topical Drug Delivery. Molecular Pharmaceutics, 0, , .	2.3	1