

# Jan ÅEejka

## List of Publications by Year in descending order

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docs citations

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#	ARTICLE	IF	CITATIONS
1	Intriguing Cytotoxicity of the Street Dissociative Anesthetic Methoxphenidine: Unexpected Impurities Spotted. <i>International Journal of Molecular Sciences</i> , 2022, 23, 2083.	1.8	2
2	Low-temperature polymorphs of lacosamide. <i>Journal of Crystal Growth</i> , 2021, 562, 126085.	0.7	3
3	Chemistry of 2,14-Dithiacalix[4]arene: Searching for the Missing Fifth Conformer. <i>Journal of Organic Chemistry</i> , 2021, 86, 9788-9801.	1.7	2
4	Nucleophile-induced transformation of phenoxathiin-based thiacalixarenes. <i>Organic and Biomolecular Chemistry</i> , 2021, 19, 8075-8085.	1.5	2
5	Synthesis and identification of deschloroketamine metabolites in rats' urine and a quantification method for deschloroketamine and metabolites in rats' serum and brain tissue using liquid chromatography tandem mass spectrometry. <i>Drug Testing and Analysis</i> , 2020, 12, 343-360.	1.6	9
6	Inherent chirality through a simple dialkylation of 2,14-dithiacalix[4]arene. <i>New Journal of Chemistry</i> , 2020, 44, 14496-14504.	1.4	4
7	Regioselective formation of the quinazoline moiety on the upper rim of calix[4]arene as a route to inherently chiral systems. <i>New Journal of Chemistry</i> , 2020, 44, 6490-6500.	1.4	6
8	Formal Transition-Metal-Catalyzed Phosphole C-H Activation for the Synthesis of Pentasubstituted Phospholes. <i>Organic Letters</i> , 2020, 22, 2187-2190.	2.4	10
9	Dihydrogen Bond in the Aminoborane Complex of a Nicergoline Intermediate. <i>Molecules</i> , 2019, 24, 2548.	1.7	1
10	Epoxidation is the preferred pathway of first-stage metabolism of abiraterone acetate in brown bullhead ( <i>Ameiurus nebulosus</i> ). <i>Environmental Science and Pollution Research</i> , 2019, 26, 34896-34904.	2.7	1
11	Microbatch under-oil salt screening of organic cations: single-crystal growth of active pharmaceutical ingredients. <i>IUCr</i> , 2019, 6, 145-151.	1.0	10
12	Template synthesis and structure of Co(II), Ni(II), and Cu(II) complexes with pyridoxilydenetaurinate Schiff base ligand. <i>Inorganica Chimica Acta</i> , 2018, 477, 248-256.	1.2	12
13	A high throughput screening method for the nano-crystallization of salts of organic cations. <i>Chemical Science</i> , 2018, 9, 3716-3722.	3.7	23
14	Synthesis, absolute configuration and <i>in vitro</i> cytotoxicity of deschloroketamine enantiomers: rediscovered and abused dissociative anaesthetic. <i>New Journal of Chemistry</i> , 2018, 42, 19360-19368.	1.4	14
15	Sodium Aspirin Salts: Crystallization and Characterization. <i>Crystal Growth and Design</i> , 2018, 18, 5287-5294.	1.4	6
16	Novel screening method for the crystallization of salts of organic cations. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2018, 74, e143-e143.	0.0	0
17	Preparation, characterization, and the selective antimicrobial activity of N-alkylammonium 8-diethyleneglycol cobalt bis-dicarbollide derivatives. <i>Journal of Organometallic Chemistry</i> , 2017, 827, 23-31.	0.8	28
18	Structural characterization of various salts of trospium: from small change of anion to huge unit cell. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2016, 72, s351-s351.	0.0	0

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19	Structural characterization of tenofovir disoproxil fumarate Form I using X-ray and electron diffraction and a study of its conversion to related solid forms. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2016, 72, s362-s362.	0.0	0
20	Bowl-shaped TrĀĳger's bases and their recognition properties. <i>Chemical Communications</i> , 2016, 52, 10664-10667.	2.2	13
21	<i>CrystalCMP</i>: an easy-to-use tool for fast comparison of molecular packing. <i>Journal of Applied Crystallography</i> , 2016, 49, 2172-2183.	1.9	72
22	Crystalline forms of dihydroergocornine salts. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2016, 72, s364-s364.	0.0	0
23	X-ray structural analysis of sodium salt of acetylsalicylic acid. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2016, 72, s357-s357.	0.0	0
24	Growing cocrystals by stoichiometric cosublimation. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2015, 71, s457-s457.	0.0	0
25	Systematic solvate screening of trospium chloride: discovering hydrates of a long-established pharmaceutical. <i>CrystEngComm</i> , 2015, 17, 4712-4721.	1.3	20
26	Synthesis of Heavy Fluorous Ruthenium Metathesis Catalysts Using the Stereoselective Addition of Polyfluoroalkyllithium to Sterically Hindered Diimines. <i>Organometallics</i> , 2015, 34, 3327-3334.	1.1	20
27	Increasing dissolution of trospium chloride by co-crystallization with urea. <i>Journal of Crystal Growth</i> , 2014, 399, 19-26.	0.7	13
28	Structural flexibility of 2-hetaryl chromium aminocarbene complexes: Experimental and theoretical evidence. <i>Inorganica Chimica Acta</i> , 2014, 421, 439-445.	1.2	4
29	Evidence for the Cyclic CN<sub>2</sub> Carbene in Solution. <i>Organic Letters</i> , 2014, 16, 852-855.	2.4	7
30	Direct CĀĤH sulfonylation of purines and deazapurines. <i>Organic and Biomolecular Chemistry</i> , 2013, 11, 5189.	1.5	57
31	Synthesis and characterisation of a new naphthalene tris-TrĀĳgerĀ™s base derivativeĀ™ a chiral molecular clip. <i>Tetrahedron Letters</i> , 2013, 54, 308-311.	0.7	12
32	Characterizing Crystal Disorder of Trospium Chloride: A Comprehensive, 13C CP/MAS NMR, DSC, FTIR, and XRPD Study. <i>Journal of Pharmaceutical Sciences</i> , 2013, 102, 1235-1248.	1.6	15
33	Trospium Chloride: Unusual Example of Polymorphism Based on Structure Disorder. <i>Crystal Growth and Design</i> , 2013, 13, 5193-5203.	1.4	17
34	Modification of pharmaceutical substances for structure analysis. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2013, 69, s680-s680.	0.3	0
35	Trospium chloride: polymorphism versus molecular disorder. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2013, 69, s678-s678.	0.3	1
36	Stereochemistry of disordered fluorinated precursors of NHC ligands. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2013, 69, s554-s554.	0.3	0

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37	Synthesis and structural studies of flavin and alloxazine adducts with O-nucleophiles. Journal of Molecular Structure, 2011, 1004, 178-187.	1.8	14
38	Polyfluoroalkylated tripyrazolymethane ligands: Synthesis and complexes. Journal of Fluorine Chemistry, 2011, 132, 434-440.	0.9	5
39	( <i>R</i> )-2,2-Bis[ <i>N</i> -(3,5-dichlorophenyl)ureido]-1,1-binaphthalene chloroform disolvate. Acta Crystallographica Section E: Structure Reports Online, 2011, 67, o384-o385.	0.2	3
40	Trifluoromethylated (tetrahydropyrrolo) quinazolinones by a new three-component reaction and facile assignment of the regio- and stereoisomers formed by NMR spectroscopy. Magnetic Resonance in Chemistry, 2010, 48, 375-385.	1.1	8
41	Nostotrebins 6, a bis(cyclopentenedione) with cholinesterase inhibitory activity isolated from Nostoc sp. str. Lukejovj 27/97. Journal of Enzyme Inhibition and Medicinal Chemistry, 2010, 25, 414-420.	2.5	12
42	Spiro Tröger's Base Derivatives: Another Structural Phoenix?. Organic Letters, 2010, 12, 1872-1875.	2.4	20
43	Multi-component crystals of dihydroergocornine. Acta Crystallographica Section A: Foundations and Advances, 2010, 66, s223-s223.	0.3	0
44	Enantioseparations of non-benzenoid and oligo-Tröger's bases by HPLC on Whelk O1 column. Tetrahedron: Asymmetry, 2009, 20, 1918-1923.	1.8	16
45	Synthesis of bis(polyfluoroalkylated)imidazolium salts as key intermediates for fluororous NHC ligands. Journal of Fluorine Chemistry, 2009, 130, 966-973.	0.9	43
46	5-Ethyl-4a-methoxy-1,3-dimethyl-4a,5-dihydrobenzo[ <i>g</i> ]pteridine-2,4(1 <i>H</i> ,3 <i>H</i> )dione. Acta Crystallographica Section E: Structure Reports Online, 2009, 65, o1536-o1537.	0.2	1
47	Crystal forms of trospium chloride. Acta Crystallographica Section A: Foundations and Advances, 2009, 65, s304-s304.	0.3	1
48	Molecular and solid-state structure of methyl [2-(acridin-9-ylimino)-3-( <i>tert</i> -butylamino)-4-oxothiazolidin-5-ylidene]acetate. Journal of Molecular Structure, 2008, 875, 419-426.	1.8	9
49	calix-Tris-Tröger's bases – a new cavitand family. Chemical Communications, 2007, , 3835.	2.2	24
50	Synthesis of 3-Fluorofuran-2(5 <i>H</i> )-ones Based on <i>Z</i> / <i>E</i> Photoisomerisation and Cyclisation of 2-Fluoro-4-hydroxybut-2-en-1-ylideneacetates. European Journal of Organic Chemistry, 2007, 2007, 5917-5925.	1.2	13
51	Limitations of the Wittig-Horner-type annulation of fluorobutenolide moiety to 3-hydroxyquinoline-2,4(2 <i>H</i> ,3 <i>H</i> )-diones. Novel modifications of the Perkow reaction including fluorinated acyloxy leaving groups. Tetrahedron, 2007, 63, 10549-10561.	1.0	8
52	Structures of cabergoline anhydrate form II and novel cabergoline solvates. Structural Chemistry, 2006, 17, 131-137.	1.0	5
53	Crystal structure of taxuspinnanane A acetonitrile solvate, C <sub>47</sub> H <sub>59</sub> NO <sub>14</sub> · CH <sub>3</sub> CN. Zeitschrift Fur Kristallographie - New Crystal Structures, 2006, 221, 97-100.	0.1	1
54	Rearrangement of Substituted 2,4,4,6-Tetraaryl-4 <i>H</i> -thiopyrans to Triaryl-3 <i>H</i> -benzo[3,4]cyclopenta[1,2- <i>b</i> ]thiophene.. ChemInform, 2005, 36, no.	0.1	0

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55	Norleucine, a natural occurrence in a novel ergot alkaloid $\hat{3}$ -ergokryptinine. <i>Amino Acids</i> , 2005, 29, 145-150.	1.2	5
56	Crystal structures of mollugin and lucidin. <i>Journal of Chemical Crystallography</i> , 2005, 35, 621-627.	0.5	12
57	(5R,6R,7R,9R,13R,14S)-21-Cyclopropylmethyl-6,14-endo-ethano-2 $\hat{2}$ ,3 $\hat{2}$ ,4 $\hat{2}$ ,5 $\hat{2}$ ,7,8-hexahydro-4 $\hat{2}$ ,4 $\hat{2}$ ,5 $\hat{2}$ ,5 $\hat{2}$ -tetramethylohydrochloride methanol solvate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2005, 61, o2274-o2276.	0.2	0
58	Crystal structure of fluticasone propionate, C <sub>25</sub> H <sub>31</sub> F <sub>3</sub> O <sub>5</sub> S. <i>Zeitschrift Fur Kristallographie - New Crystal Structures</i> , 2005, 220, 153-154.	0.1	1
59	Interesting Solvent Area in Crystal Structures of Two Natural Ergot Alkaloids. <i>Collection of Czechoslovak Chemical Communications</i> , 2005, 70, 41-50.	1.0	1
60	Crystal structure of 1-methyl-10 $\hat{1}$ -methoxy-9,10-dihydrolysergol, C <sub>18</sub> H <sub>24</sub> N <sub>2</sub> O <sub>2</sub> . <i>Zeitschrift Fur Kristallographie - New Crystal Structures</i> , 2005, 220, 217-218.	0.1	1
61	Crystal structure of 1-hydroxymethyl-10 $\hat{1}$ -methoxy-9,10-dihydrolysergol, C <sub>18</sub> H <sub>24</sub> N <sub>2</sub> O <sub>3</sub> . <i>Zeitschrift Fur Kristallographie - New Crystal Structures</i> , 2005, 220, .	0.1	1
62	Derivatives of (Phenylsulfanyl)benzoic Acids with Multiple Antileukotrienic Activity. <i>Collection of Czechoslovak Chemical Communications</i> , 2004, 69, 2098-2120.	1.0	7
63	Cabergoline, form VII. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2004, 60, o1167-o1169.	0.2	2
64	Restricted Conformational Flexibility of Furanose Derivatives: $\hat{2}$ Ab Initio Interpretation of Their Nuclear Spin $\hat{2}$ Spin Coupling Constants. <i>Journal of Physical Chemistry A</i> , 2004, 108, 6365-6372.	1.1	17
65	Rearrangement of Substituted 2,4,4,6-Tetraaryl-4H-thiopyrans to Triaryl-3aH-benzo[3,4]cyclopenta[1,2-b]thiophene. <i>Collection of Czechoslovak Chemical Communications</i> , 2004, 69, 1631-1642.	1.0	2
66	Simvastatin. <i>Acta Crystallographica Section C: Crystal Structure Communications</i> , 2003, 59, o428-o430.	0.4	20
67	Pergolide mesylate form II. <i>Acta Crystallographica Section C: Crystal Structure Communications</i> , 2003, 59, o575-o576.	0.4	1
68	Structural Study of Dopamine Agonist Lisuride. <i>Collection of Czechoslovak Chemical Communications</i> , 2003, 68, 2150-2158.	1.0	0
69	Substituted 2,4,4,6-tetraaryl-4H-selenopyrans: preparation, photocolouration and 4H-selenopyran ring geometry; an X-ray and DFT calculation study. <i>Perkin Transactions II RSC</i> , 2002, , 1909-1916.	1.1	3
70	Structure Types of Dihydroergotoxine Mesylates. <i>Collection of Czechoslovak Chemical Communications</i> , 2002, 67, 490-501.	1.0	1
71	X-ray study on polymorphism of some ergopeptine drugs. <i>Acta Crystallographica Section A: Foundations and Advances</i> , 2002, 58, c126-c126.	0.3	0
72	Photoisomerization of 2,4,4,6-tetraaryl-4H-selenopyrans: a new heterocyclic ring contraction. <i>Mendeleev Communications</i> , 2001, 11, 90-91.	0.6	2

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73	Selective Reduction of Peptidic Ergot Alkaloids. Collection of Czechoslovak Chemical Communications, 2000, 65, 1762-1776.	1.0	4
74	Sterically Crowded Heterocycles. XII. Atropisomerism of (1-Aryl-3,5-diphenyl-1H-pyrrol-2-yl)(phenyl)methanones. Collection of Czechoslovak Chemical Communications, 2000, 65, 651-666.	1.0	3
75	Crystal Structures of Dihydro-1 $\pm$ -ergokryptine and Dihydro-1 $^2$ -ergokryptine Mesylates. Collection of Czechoslovak Chemical Communications, 2000, 65, 1329-1338.	1.0	2
76	MarchingCubeELD - electron density visualization program for small molecules and its application on cyclosporins structures investigation. Acta Crystallographica Section A: Foundations and Advances, 2000, 56, s196-s196.	0.3	0
77	Comprehensive Comparison of the Ergot Alkaloids Family with 5-HT1A Receptor Pharmacophore Models. Acta Crystallographica Section A: Foundations and Advances, 2000, 56, s282-s282.	0.3	0
78	Synthesis and Crystal Structure of Pergolide Sulfoxide. Collection of Czechoslovak Chemical Communications, 1998, 63, 803-812.	1.0	3
79	Preparation of New Optically Active Furanoids Containing Various Aza-Heterocyclic Moieties. Collection of Czechoslovak Chemical Communications, 1997, 62, 894-912.	1.0	1
80	Preparation of new trifluoromethyl substituted tri- and tetracyclic heterocycles with Peganin skeleton from a methyl 3,3,3-trifluoropyruvate / 2-aminobenzylamine adduct. Tetrahedron Letters, 1996, 37, 6939-6942.	0.7	13
81	Crystal Structures of Ergot Alkaloid Derivatives. Ergometrine Maleate and Methylegometrine Maleate. Collection of Czechoslovak Chemical Communications, 1996, 61, 1396-1404.	1.0	4
82	Sterically Crowded Heterocycles. VIII. Preparative Photoisomerization of Some Imidazo[1,2-a]pyridines. Collection of Czechoslovak Chemical Communications, 1996, 61, 1473-1488.	1.0	1
83	Sterically Crowded Heterocycles. VI. Relative Configuration of Diastereoisomeric (Z)-1,3-Diphenyl-3-(5-methyl-2-phenylimidazo[1,2-a]pyridin-3-yl)prop-2-en-1-ols. Collection of Czechoslovak Chemical Communications, 1996, 61, 1380-1385.	1.0	1
84	Absolute Crystal Structure Determination of Ergot Alkaloid - Dihydroergocristine Methanesulfonate Monohydrate. Collection of Czechoslovak Chemical Communications, 1995, 60, 1333-1342.	1.0	6