

Joel T Mague

List of Publications by Year in descending order

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481
papers

3,433
citations

185998

28
h-index

276539

41
g-index

485
all docs

485
docs citations

485
times ranked

2665
citing authors

#	ARTICLE	IF	CITATIONS
1	Diastereoselective Synthesis of Polysubstituted Tetrahydropyrans and Thiacyclohexanes via Indium Trichloride Mediated Cyclizations. <i>Journal of Organic Chemistry</i> , 2001, 66, 739-747.	1.7	114
2	Rationalization of Noncovalent Interactions within Six New M ^{II} /8-Aminoquinoline Supramolecular Complexes (M ^{II} = Mn, Cu, and Cd): A Combined Experimental and Theoretical DFT Study. <i>Crystal Growth and Design</i> , 2015, 15, 1351-1361.	1.4	97
3	Novel Carbocycle Enlargement in Aqueous Medium. <i>Journal of the American Chemical Society</i> , 1996, 118, 4216-4217.	6.6	78
4	On the importance of tetrel bonding interactions in lead(II) complexes with (iso)nicotinohydrazide based ligands and several anions. <i>Dalton Transactions</i> , 2016, 45, 10708-10716.	1.6	78
5	Crystal engineering with coordination compounds of NiII, CoII, and CrIII bearing dipicolinic acid driven by the nature of the noncovalent interactions. <i>CrystEngComm</i> , 2014, 16, 5352.	1.3	73
6	Tetranuclear Rhodium(I) Macrocycle Containing Cyclodiphosphazane [Rh ₂ (μ ₄ -Cl) ₂ (CO) ₂ {(tBuNP(OC ₆ H ₄ OMe-o)) ₂ -P ₂ } ₂] and Its Reversible Conversion into trans-[Rh(CO)Cl{(tBuNP(OC ₆ H ₄ OMe-o)) ₂ -P ₂ } ₂]. <i>Organometallics</i> , 2005, 24, 3780-3783.	1.1	68
7	Synthesis, structural and molecular characterization of 2,2-diphenyl-2H,3H,5H,6H,7H-imidazo[2,1-b][1,3]thiazin-3-one. <i>Journal of Molecular Structure</i> , 2019, 1197, 369-376.	1.8	60
8	On the importance of non covalent interactions in the structure of coordination Cu(II) and Co(II) complexes of pyrazine- and pyridine-dicarboxylic acid derivatives: experimental and theoretical views. <i>CrystEngComm</i> , 2014, 16, 6149-6158.	1.3	57
9	Crystal engineering with coordination compounds of 2,6-dicarboxy-4-hydroxypyridine and 9-aminoacridine fragments driven by different nature of the face-to-face π-π stacking. <i>CrystEngComm</i> , 2014, 16, 1359-1377.	1.3	56
10	Synthesis of a novel phenytoin derivative: Crystal structure, Hirshfeld surface analysis and DFT calculations. <i>Journal of Molecular Structure</i> , 2020, 1205, 127630.	1.8	56
11	Importance of polarization assisted/resonance assisted hydrogen bonding interactions and unconventional interactions in crystal formations of five new complexes bearing chelidamic acid through a proton transfer mechanism. <i>RSC Advances</i> , 2015, 5, 72923-72936.	1.7	50
12	Facile Reductive Elimination of Ethane from Strained Dimethylpalladium(II) Complexes. <i>Journal of the American Chemical Society</i> , 2001, 123, 4081-4082.	6.6	47
13	Dodecaphenyltetracene. <i>Angewandte Chemie - International Edition</i> , 2019, 58, 2831-2833.	7.2	45
14	Thioether-Functionalized Ferrocenyl-bis(phosphonite), Fe{(C ₅ H ₄)P(OC ₁₀ H ₆ (μ ₄ -S)C ₁₀ H ₆ O)} ₂ : Synthesis, Coordination Behavior, and Application in Suzuki-Miyaura Cross-Coupling Reactions. <i>Inorganic Chemistry</i> , 2007, 46, 10268-10275.	1.9	43
15	Structural scope of six new layered to pillar-layered hybrid inorganic-organic networks bearing [BW ₁₂ O ₄₀] ⁵⁺ and lanthanoid-cluster; database study toward ligand role in assemblies. <i>CrystEngComm</i> , 2016, 18, 6724-6737.	1.3	41
16	An inorganic-organic hybrid material based on a Keggin-type polyoxometalate@Dysprosium as an effective and green catalyst in the synthesis of 4-aminochromenes via multicomponent reactions. <i>Applied Organometallic Chemistry</i> , 2020, 34, e5793.	1.7	41
17	Synthesis, spectroscopic characterization, crystal structure, DFT, molecular docking and <i>in vitro</i> antibacterial potential of novel quinoline derivatives. <i>Journal of Molecular Structure</i> , 2020, 1209, 127940.	1.8	40
18	Synthesis, crystal structure, DFT calculations, Hirshfeld surface analysis, energy frameworks, molecular dynamics and docking studies of novel isoxazolequinoxaline derivative (IZQ) as anti-cancer drug. <i>Journal of Molecular Structure</i> , 2021, 1232, 130004.	1.8	40

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19	Two Triazole-Based Phosphine Ligands Prepared via Temperature-Mediated Li/H Exchange: Cu ^I and Au ^I Complexes and Structural Studies. <i>Inorganic Chemistry</i> , 2016, 55, 8514-8526.	1.9	39
20	Novel Co(II) and Cu(II) coordination complexes constructed from pyrazole-acetamide: Effect of hydrogen bonding on the self assembly process and antioxidant activity. <i>Journal of Inorganic Biochemistry</i> , 2019, 191, 21-28.	1.5	39
21	Synthetic and Structural Studies of the Coordination Behavior of 2-Pyridylbis(diphenylphosphino)methane. <i>Inorganic Chemistry</i> , 2001, 40, 1962-1971.	1.9	38
22	Synthesis and derivatization, structures and transition metal chemistry of a new large bite bis(phosphinite) derived from bis(2-hydroxy-1-naphthyl)methane. <i>Dalton Transactions RSC</i> , 2002, , 4617-4621.	2.3	36
23	Hydrothermal synthesis, X-ray structure and DFT and magnetic studies of a (H ₂ SiW ₁₂ O ₄₀) ²⁻ based one-dimensional linear coordination polymer. <i>Dalton Transactions</i> , 2015, 44, 8824-8832.	1.6	34
24	Novel antioxidant quinoxaline derivative: Synthesis, crystal structure, theoretical studies, antidiabetic activity and molecular docking study. <i>Journal of Molecular Structure</i> , 2021, 1239, 130484.	1.8	34
25	Sterically Demanding Phosphines with 2,6-Dibenzhydryl-4-methylphenyl Core: Synthesis of Ru ^{II} , Pd ^{II} , and Pt ^{II} Complexes, and Structural and Catalytic Studies. <i>Inorganic Chemistry</i> , 2018, 57, 7468-7480.	1.9	33
26	Synthesis, crystal structure, spectroscopic characterization, Hirshfeld surface analysis, molecular docking studies and DFT calculations, and antioxidant activity of 2-oxo-1,2-dihydroquinoline-4-carboxylate derivatives. <i>Journal of Molecular Structure</i> , 2019, 1188, 255-268.	1.8	32
27	Curcumin and derivatives. <i>Acta Crystallographica Section C: Crystal Structure Communications</i> , 2004, 60, o608-o610.	0.4	31
28	Supramolecular network of a framework material supported by the anion-π linkage of Keggin-type heteropolyoxotungstates: experimental and theoretical insights. <i>Dalton Transactions</i> , 2021, 50, 1895-1900.	1.6	31
29	Hierarchical Assembly of Organic/Inorganic Building Molecules with π-π Interactions. <i>Advanced Functional Materials</i> , 2008, 18, 1526-1535.	7.8	29
30	Two polyoxometalate-based hybrids constructed from trinuclear lanthanoid clusters with single-molecule magnet behavior. <i>Polyhedron</i> , 2021, 194, 114903.	1.0	29
31	Potential antidiabetic activity and molecular docking studies of novel synthesized 3,6-dimethyl-5-oxo-pyrido[3,4-f][1,2,4]triazepino[2,3-a]benzimidazole and 10-amino-2-methyl-4-oxo pyrimido[1,2-a]benzimidazole derivatives. <i>Journal of Molecular Modeling</i> , 2018, 24, 179.	0.8	28
32	Tribenzodecacyclene and Hexabenzodecacyclene. <i>Journal of Organic Chemistry</i> , 2015, 80, 4824-4827.	1.7	27
33	Synthesis, crystal structure, DFT, molecular dynamics simulation and evaluation of the anticorrosion performance of a new pyrazolotriazole derivative. <i>Journal of Molecular Structure</i> , 2019, 1176, 290-297.	1.8	27
34	Buffer and Salt Effects in Aqueous Host-Guest Systems: Screening, Competitive Binding, or Both?. <i>Journal of the American Chemical Society</i> , 2021, 143, 18605-18616.	6.6	27
35	The Hairpin Furans: Easily Prepared Hybrids of Helicenes and Twisted Acenes. <i>Angewandte Chemie - International Edition</i> , 2015, 54, 13957-13960.	7.2	25
36	Synthesis, transition metal chemistry and catalytic reactions of ferrocenylbis(phosphonite), [Fe{C ₅ H ₄ P(OC ₆ H ₃ (OMe-o)(C ₃ H ₅ -p)) ₂ } ₂]. <i>Dalton Transactions</i> , 2013, 42, 11695.	1.6	24

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37	Diamondoid-Type Copper Coordination Polymers Containing Soft Cyclodiphosphazane Ligands. <i>Inorganic Chemistry</i> , 2015, 54, 6063-6065.	1.9	24
38	Self-Assembled Cyclophane-Type Copper(I) Complexes of 2,4,6-Tris(diphenylphosphino)-1,3,5-triazine and Their Catalytic Application. <i>Inorganic Chemistry</i> , 2015, 54, 10985-10992.	1.9	24
39	Synthesis, biological activity and molecular modeling of a new series of condensed 1,2,4-triazoles. <i>Bioorganic Chemistry</i> , 2019, 92, 103193.	2.0	23
40	Docking of disordered independent molecules of novel crystal structure of		

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55	A possible potential COVID-19 drug candidate: Diethyl 2-(2-(2-(3-methyl-2-oxoquinoxalin-1(2H)-yl)acetyl)hydrazono)malonate: Docking of disordered independent molecules of a novel crystal structure, HSA/DFT/XRD and cytotoxicity. <i>Arabian Journal of Chemistry</i> , 2022, 15, 103595.	2.3	17
56	Mono-, Bi-, Tri- and Tetranuclear Palladium(II), Copper(I), and Gold(I) Complexes of Morpholine- and N-Methylpiperazine-Functionalized Cyclodiphosph(III)azans, cis-[(tBuN- $\frac{1}{4}$) ₂ (PNC4H8X) ₂] (X = O, NMe). <i>European Journal of Inorganic Chemistry</i> , 2010, 2010, 4201-4210.	1.0	16
57	Co(III) and Fe(III) complexes of Schiff bases derived from 2,4-dihydroxybenzaldehyde <i>S</i>-allyl-isothiosemicarbazonehydrobromide. <i>Journal of Coordination Chemistry</i> , 2013, 66, 3915-3925.	0.8	16
58	Two new copper and nickel complexes of pyridine-2,6-dicarboxylic acid N-oxide and their proton transferred salts: Solid state and DFT insights. <i>Inorganica Chimica Acta</i> , 2015, 438, 135-145.	1.2	16
59	Synthesis, X-ray characterization and DFT study of a novel Fe(III)â€pyridine-2,6-dicarboxylic acid N-oxide complex with unusual coordination mode. <i>Inorganica Chimica Acta</i> , 2016, 449, 44-51.	1.2	16
60	Synthesis, crystal structure, spectroscopic characterization, Hirshfeld surface analysis, and DFT calculations of 1,4-dimethyl-2-oxo-pyrimido[1,2-a]benzimidazole hydrate. <i>Journal of Molecular Structure</i> , 2018, 1152, 154-162.	1.8	16
61	HSA-interaction studies of uranyl complexes of alkyl substituted isothiosemicarbazone. <i>Journal of Molecular Structure</i> , 2019, 1193, 53-61.	1.8	16
62	Synthesis and Spectroscopic and Thermal Decomposition Studies of Alkali Metal Salts of 2-Oximidopropionate. <i>Inorganic Chemistry</i> , 1997, 36, 2656-2661.	1.9	15
63	Surface-grafted lanthanoid complexes of the tungstosilicate polyanion [SiW ₁₂ O ₄₀] ⁴⁻ : a synthetic, structural and computational investigation. <i>Acta Crystallographica Section C, Structural Chemistry</i> , 2018, 74, 1300-1309.	0.2	15
64	Solvent induced supramolecular polymorphism in Cu(II) coordination complex built from 1,2,4-triazolo[1,5-a]pyrimidine: Crystal structures and anti-oxidant activity. <i>Journal of Inorganic Biochemistry</i> , 2020, 208, 111092.	1.5	15
65	2-pyridylbis(diphenylphosphino)methane chemistry. Synthesis and structures of [Cu($\frac{1}{4}$ ·2·1(Ph ₂ P) ₂ ·CHC ₅ H ₄ N)(THF)] ₂ (BF ₄) ₂ and [Ni(Ph ₂ PCH ₂ C ₅ H ₄ N) ₂]-[NiCl ₄] ^{0.85} CH ₂ Cl ₂ and [Ni(Ph ₂ PCH ₂ C ₅ H ₄ N) ₂]-[NiCl ₄] ^{0.85} CH ₂ Cl ₂ . <i>Journal of Chemical Crystallography</i> , 1997, 27, 603-608.	0.5	14
66	Facile and convenient synthesis of 2,4-disubstituted and 2,3,4-trisubstituted 1,3-thiazoles. <i>Journal of Sulfur Chemistry</i> , 2016, 37, 162-175.	1.0	14
67	Redox-Active Metallodithiolene Groups Separated by Insulating Tetraphosphinobenzene Spacers. <i>Inorganic Chemistry</i> , 2018, 57, 4023-4038.	1.9	14
68	Title is missing!. <i>Journal of Chemical Crystallography</i> , 2001, 31, 295-300.	0.5	13
69	1,1- ² -Bis(dipyrrolylphosphino)ferrocene: Synthesis, coordination chemistry and structural studies. <i>Journal of Organometallic Chemistry</i> , 2016, 824, 15-24.	0.8	13
70	A Straightforward Approach for the Synthesis of Novel Derivatives of Benzo[b]pyrazolo[5- ² ,1- ² :2,3]pyrimido[4,5- ²][1,4]thiazine. <i>Journal of Heterocyclic Chemistry</i> , 2016, 53, 1231-1235.	1.4	13
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73	Ethyl 2-(2,5-dioxo-4,4-diphenylimidazolidin-1-yl)acetate. IUCrData, 2017, 2, .	0.1	13
74	1-Benzyl-3-methylquinoxalin-2(1 <i>H</i>)-one. IUCrData, 2018, 3, .	0.1	13
75	Ring Opening of Epoxides by Pendant Silanols. Organic Letters, 2022, 24, 939-943.	2.4	13
76	Synthesis of a tert-butyl substituted bis(silirane) and comparison with its methyl and phenyl analogs. Journal of Organometallic Chemistry, 2011, 696, 1957-1963.	0.8	12
77	Synthesis and Structures of [LCu(I)(SSi ⁺ Pr ₃)] (L = triphos, carbene) and Related Compounds. Inorganic Chemistry, 2016, 55, 9173-9177.	1.9	12
78	Synthesis, single crystal X-ray characterization, and solution studies of Zn(II)-, Cu(II)-, Ag(I)- and Ni(II)-pyridine-2,6-dipicolinate N-oxide complexes with different topologies and coordination modes. Inorganica Chimica Acta, 2017, 458, 84-96.	1.2	12
79	An inorganic-organic hybrid supramolecular framework based on the $\text{[Mo}_8\text{O}_{26}]^{4-}$ cluster and cobalt complex of aspartic acid: X-ray structure and DFT study. Acta Crystallographica Section C, Structural Chemistry, 2019, 75, 469-477.	0.2	12
80	Syntheses of N-substituted benzimidazolone derivatives: DFT calculations, Hirshfeld surface analysis, molecular docking studies and antibacterial activities. Journal of Molecular Structure, 2020, 1200, 127174.	1.8	12
81	Syntheses of novel 1,5-benzodiazepine derivatives: Crystal structures, spectroscopic characterizations, Hirshfeld surface analyses, molecular docking studies, DFT calculations, corrosion inhibition anticipation, and antibacterial activities. Journal of Heterocyclic Chemistry, 2021, 58, 270-289.	1.4	12
82	Crystal structure, Hirshfeld surface analysis and interaction energy and DFT studies of 3-[(2 <i>Z</i>)-2-[(2,4-dichlorophenyl)methylidene]-3-oxo-3,4-dihydro-2 <i>H</i> -1,4-benzothiazin-4-yl]propanenitrile. Acta Crystallographica Section E: Crystallographic Communications, 2019, 75, 721-727.	0.2	12
83	Title is missing!. Journal of Chemical Crystallography, 2000, 30, 311-320.	0.5	11
84	Crystal structure of 2-(2,3-dimethylanilino)- <i>N</i> -[(1 <i>E</i>)-2-hydroxybenzylidene]benzohydrazide. Acta Crystallographica Section E: Crystallographic Communications, 2015, 71, o957-o958.	0.2	11
85	A Novel Method for the Synthesis of Furo-imidazo[3.3.3]propellanes from Thiocarbonohydrazides. Synlett, 2016, 27, 412-416.	1.0	11
86	Structure and characterization of charge transfer complexes of benzo[1,2- <i>b</i> :3,4- <i>b'</i> :5,6- <i>b''</i>]trithiophene [C ₃ h-BTT]. CrystEngComm, 2017, 19, 6355-6364.	1.3	11
87	Photocatalytic H ₂ -Evolution by Homogeneous Molybdenum Sulfide Clusters Supported by Dithiocarbamate Ligands. Inorganic Chemistry, 2019, 58, 16458-16474.	1.9	11
88	New alkyl (cyclohexyl) 2-oxo-1-(prop-2-yn-1-yl)-1, 2-dihydroquinoline-4-carboxylates: Synthesis, crystal structure, spectroscopic characterization, hirshfeld surface analysis, molecular docking studies and DFT calculations. Journal of Molecular Structure, 2021, 1227, 129520.	1.8	11
89	Synthesis, structural characterisation and theoretical studies of a novel pyridazine derivative: Investigations of anti-inflammatory activity and inhibition of β -glucosidase. Journal of Molecular Structure, 2021, 1234, 130177.	1.8	11
90	Highly Regio- and Diastereoselective Tethered Aza-Wacker Cyclizations of Alkenyl Phosphoramidates. Journal of Organic Chemistry, 2021, 86, 14732-14758.	1.7	11

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91	Aryl acetylene inhibitors for cytochrome P450-based monooxygenase isozymes. Journal of Chemical Crystallography, 1997, 27, 183-189.	0.5	10
92	Title is missing!. Journal of Chemical Crystallography, 2003, 33, 391-402.	0.5	10
93	Ni(II), Co(II), and Cu(II) complexes incorporating 2-pyrazinecarboxylic acid: Synthesis, characterization, electrochemical evaluation, and catalytic activity for the synthesis of 2H-indazolo[2,1-b]phthalazine-triones. Chinese Journal of Catalysis, 2015, 36, 1101-1108.	6.9	10
94	Hairpin Furans and Giant Biaryls. Journal of Organic Chemistry, 2016, 81, 3838-3847.	1.7	10
95	Palladium-catalyzed regioselective direct CH arylation of 2-pyrazolo[3,4-d]pyrimidines. Comptes Rendus Chimie, 2017, 20, 927-933.	0.2	10
96	Synthesis, crystal structure, Hirshfeld surface analysis, and DFT calculations of new 1-[(1-benzyl-1H-1,2,3-triazol-4-yl)methyl]-6-methoxy-1H-benzimidazol-2(3H)-one. Chemical Data Collections, 2018, 17-18, 472-482.	1.1	10
97	Synthesis, anticancer evaluation in vitro, DFT, Hirshfeld surface analysis of some new 4-(1,3-benzothiazol-2-yl)-3-methyl-1-phenyl-4,5-dihydro-1H-pyrazol-5-one derivatives. Journal of Molecular Structure, 2019, 1198, 126910.	1.8	10
98	Synthesis, spectroscopic characterization, thermal, XRD crystal structure, the PLATON structural analysis, and theoretical studies of a new 1,2,4-triazolo-[1,5-a]pyrimidines derivatives. Journal of Molecular Structure, 2019, 1184, 12-24.	1.8	10
99	Synthesis, characterization and bioactivity studies of new dithiocarbamate complexes. New Journal of Chemistry, 2020, 44, 8878-8889.	1.4	10
100	Tethered Silanoxyiodination of Alkenes. Journal of Organic Chemistry, 2021, 86, 9233-9243.	1.7	10
101	Synthesis, Crystal structure, Hirshfeld surface Analysis and computational approach of new 2-methylbenzimidazo[1,2-a]pyrimidin-4(1H)-one. Journal of Molecular Structure, 2021, 1239, 130497.	1.8	10
102	Coordination complexes constructed from pyrazole-acetamide and pyrazole-quinoxaline: effect of hydrogen bonding on the self-assembly process and antibacterial activity. RSC Advances, 2022, 12, 5324-5339.	1.7	10
103	Structural characterization and excited-state properties of luminescent Tris-(2-(3-methyl-5-yl)pyrazol-5-yl)ethane-1,1-dithiolane-2,2-dione. Journal of Molecular Structure, 2022, 1250, 126910.	0.5	9
104	Synthesis and Identification of Novel Potential Molecules Against COVID-19 Main Protease Through Structure-Guided Virtual Screening Approach. Applied Biochemistry and Biotechnology, 2021, 193, 3602-3623.	1.4	9
105	3-Ethyl-5,5-diphenylimidazolidine-2,4-dione. IUCrData, 2017, 2, .	0.1	9
106	Synthesis and crystal structure of 2-azido-N-phenylacetamide, C ₈ H ₈ N ₂ O. Zeitschrift Fur Kristallographie - New Crystal Structures, 2021, 236, 133-134.	0.1	9
107	Greener pastures in evaluating antidiabetic drug for a quinoxaline Derivative: Synthesis, Characterization, Molecular Docking, in vitro and HSA/DFT/XRD studies. Arabian Journal of Chemistry, 2022, 15, 103851.	2.3	9
108	A co-crystal of 2-(1-pyrenyl)-4,4,5,5-tetramethyl-4,5-dihydro-1H-imidazole-3-oxide-1-oxyl with octafluoronaphthalene. CrystEngComm, 2013, 15, 831-835.	1.3	8

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109	Square planar nickel(II) complexes derived from 5-bromo-2-hydroxybenzaldehyde S-ethylisothiosemicarbazone: Preparation, characterization and structural studies. <i>Polyhedron</i> , 2014, 80, 243-249.	1.0	8
110	The first mixed-ligand coordination compound involving 8-aminoquinoline and pyridine-2,6-dicarboxylate: synthesis, X-ray crystal structure, and DFT studies. <i>Journal of Coordination Chemistry</i> , 2015, 68, 3599-3610.	0.8	8
111	An Unexpected Diastereoselective Synthesis of Novel Substituted Pyridines via One-Pot, Four-Component Reaction. <i>Synlett</i> , 2016, 27, 1669-1673.	1.0	8
112	Synthesis and HSA-interaction of a new mixed ligand Cu-isothiosemicarbazonato complex with adenine nucleobase. <i>Polyhedron</i> , 2020, 179, 114357.	1.0	8
113	Group 10 Metal Dithiolene Bis(isonitrile) Complexes: Synthesis, Structures, Properties, and Reactivity. <i>Organometallics</i> , 2020, 39, 2854-2870.	1.1	8
114	Synthesis and Structures of Polyphenylphenanthrenes. <i>Chemistry - A European Journal</i> , 2020, 26, 8458-8464.	1.7	8
115	Unusual rearrangementâ€‘remercuration reactions of allylic silanols. <i>Organic Chemistry Frontiers</i> , 2021, 8, 5361-5368.	2.3	8
116	Dimroth rearrangement-based synthesis of novel derivatives of [1,3]selenazolo[5,4-e][1,2,4]triazolo[1,5-c]pyrimidine as a new class of selenium-containing heterocyclic architecture. <i>Molecular Diversity</i> , 2022, 26, 923-937.	2.1	8
117	Crystal structure and Hirshfeld surface analysis of N-[(E)-(4-methylbenzylidene)amino]phenyl)-2-(5-methyl-1H-pyrazol-3-yl)acetamide hemihydrate. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2019, 75, 154-158.	0.2	8
118	Ethyl 2-[(3-methylquinoxalin-2-yl)sulfanyl]acetate. <i>IUCrData</i> , 2017, 2, .	0.1	8
119	Synthesis of New Pyrimido[4,5â€‘e][1,2,4]triazolo[3,4â€‘b][1,3,4]thiadiazine Derivatives via S/N Smiles Rearrangement. <i>Journal of Heterocyclic Chemistry</i> , 2017, 54, 235-241.	1.4	7
120	Unexpected synthesis of novel 2-pyrone derivatives: crystal structures, Hirshfeld surface analysis and computational studies. <i>Journal of Biomolecular Structure and Dynamics</i> , 2021, 39, 4859-4877.	2.0	7
121	Synthesis, Crystal Structure and Computational Investigation of New 4-Phenylâ€‘decahydroâ€‘1Hâ€‘1,5â€‘benzodiazepinâ€‘2-one as Potent Inhibitor of Muâ€‘opioid Receptor. <i>ChemistrySelect</i> , 2020, 5, 4601-4607.	0.7	7
122	Synthesis and Characterization of Novel Functionally Substituted Planar Pyrimidothienoisquinolines and Nonplanar (3aR, 4S)-Tj ETQq0 0 0 rgBT /Overlock 10 Tf 50 217 Td (2aS)-pyrazolo	0.2	7
123	Crystal structure and Hirshfeld surface analysis of (4Z)-1-butyl-4-(2-oxopropylidene)-2,3,4,5-tetrahydro-1H-1,5-benzodiazepin-2-one. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2018, 74, 1669-1673.	0.2	7
124	Crystal structure and Hirshfeld surface analysis of a new benzodiazepine derivative: 4-dichloromethyl-2,3-dihydro-1H-1,5-benzodiazepin-2-one. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2019, 75, 33-37.	0.2	7
125	3-[2-(5-Oxo-4,4-diphenyl-2-sulfanylideneimidazolidin-1-yl)ethyl]-1,3-oxazolidin-2-one. <i>IUCrData</i> , 2017, 2, .	0.1	7
126	2-{3-[2-(2-Chlorophenyl)ethyl]-2-oxo-1,2-dihydroquinoxalin-1-yl}acetohydrazide. <i>IUCrData</i> , 2017, 2, .	0.1	7

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127	3-Methyl-5,5-diphenylimidazolidine-2,4-dione. IUCrData, 2017, 2, .	0.1	7
128	Coordination complexes of zinc and manganese based on pyridine-2,5-dicarboxylic acid <i>N</i> -oxide: DFT studies and antiproliferative activities consideration. RSC Advances, 2021, 11, 37403-37412.	1.7	7
129	Nitrophenyl-Group-Containing Heterocycles. I. Synthesis, Characterization, Crystal Structure, Anticancer Activity, and Antioxidant Properties of Some New 5,6,7,8-Tetrahydroisoquinolines Bearing 3(4)-Nitrophenyl Group. ACS Omega, 2022, 7, 8767-8776.	1.6	7
130	Lower rim modification of calix[4]arenes to incorporate a single group functionality. Single crystal X-ray structures of 5-(3-bromopropyl)-25,26,27,28-tetrahydroxycalix[4]arene and 25,27-diallyloxy-26,28-dibenzoyloxycalix[4]arene. Supramolecular Chemistry, 1993, 2, 53-60.	1.5	6
131	2-Cyano-1-methylpyridinium nitrate. Acta Crystallographica Section E: Structure Reports Online, 2012, 68, o1653-o1653.	0.2	6
132	Effect of solvents in mixed-ligand supramolecular self-assembly architectures. Polyhedron, 2017, 134, 41-49.	1.0	6
133	Synthesis, X-ray and Fluorescence Characteristics of Pyrimido[5,4- <i>e</i>]thiazolo[3,2- <i>a</i>]pyrimidine as a Novel Heterocyclic System. Journal of Fluorescence, 2017, 27, 1183-1190.	1.3	6
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241	Synthesis, structural elucidation, and antioxidant activity of new phenolic derivatives containing piperidine moiety: Experimental and theoretical investigations. <i>Journal of Heterocyclic Chemistry</i> , 2021, 58, 1268-1277.	1.4	2
242	Crystal structure, Hirshfeld surface analysis and interaction energy calculation of 1-decyl-2,3-dihydro-1-benzimidazol-2-one. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2021, 77, 559-563.	0.2	2
243	Triclinic form of 1,2,4,5-tetracyclohexylbenzene. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2008, 64, o376-o376.	0.2	2
244	Methyl (2Z)-((2Z)-2-((2E)-[1-(4-methylphenyl)ethylidene]hydrazinylidene)-4-oxo-3-phenyl-1,3-thiazolidin-5-ylidene)ethanoate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2013, 69, o1401-o1402.	0.2	2
245	Dimethyl (2Z)-2-[4-((1Z)-1-{2-[(2Z,5Z)-5-(2-methoxy-2-oxoethylidene)-4-oxo-3-phenyl-1,3-thiazolidin-2-ylidene]hydrazin-1-ylidene}ethyl)anilino]butanoate. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2013, 69, o1844-o1845.	0.2	2
246	Crystal structure and Hirshfeld surface analysis of ethyl 2-[4-[(3-methyl-2-oxo-1,2-dihydroquinoxalin-1-yl)methyl]-1H-1,2,3-triazol-1-yl]acetate. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2018, 74, 1648-1652.	0.2	2
247	Crystal structure, DFT calculations and Hirshfeld surface analysis of 3-(4-methylphenyl)-6-nitro-1H-indazole. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2018, 74, 1857-1861.	0.2	2
248	Crystal structure and Hirshfeld surface analysis of 3,4-dihydro-2-(2,4-dioxo-6-methylpyran-3-ylidene)-4-(4-pyridin-4-yl)-1,5-benzodiazepine. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2019, 75, 94-98.	0.2	2
249	Crystal structure, Hirshfeld surface analysis and interaction energy and DFT studies of 2-chloroethyl 2-oxo-1-(prop-2-yn-1-yl)-1,2-dihydroquinoline-4-carboxylate. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2019, 75, 1411-1417.	0.2	2
250	Crystal structure and Hirshfeld surface analysis of tert-butyl-2-(phenylethynyl)imidazo[1,2-a]pyridin-3-amine. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2019, 75, 1564-1567.	0.2	2
251	Crystal structure, Hirshfeld surface analysis and DFT studies of 1,3-bis[2-methoxy-4-(prop-2-en-1-yl)phenoxy]propane. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2020, 76, 344-348.	0.2	2
252	4-Benzyl-2-(4-chlorobenzylidene)-3,4-dihydro-2H-1,4-benzothiazin-3(4H)-one. <i>IUCrData</i> , 2016, 1, .	0.1	2

#	ARTICLE	IF	CITATIONS
253	Dichlorido{2-[(5-methyl-1 <i>H</i> -pyrazol-3-yl) <i>N</i> (<i>N</i>) ² methyl]-1 <i>H</i> -1,3-benzimidazole- <i>N</i> (<i>N</i>) ³]} IUCrData, 2017, 2, .	0.1	2
254	trans-Bis(N-{2-[2-(3-methyl-1 <i>H</i> -pyrazol-5-yl) <i>N</i> (<i>N</i>)acetamido- <i>O</i>]phenyl}benzamide)bis(perchlorato- <i>O</i>)copper(II). IUCrData, 2017, 2, .	0.1	2
255	1-[[3-(Thiophen-2-yl)-4,5-dihydro-1,2-oxazol-5-yl]methyl]-2,3-dihydro-1 <i>H</i> -indole-2,3-dione. IUCrData, 2017, 2, .	0.1	2
256	7-Acetyl-8-(4-chlorophenyl)-3-ethylsulfanyl-6-hydroxy-1,6-dimethyl-5,6,7,8-tetrahydroisoquinoline-4-carbonitrile. IUCrData, 2017, 2, .	0.1	2
257	(<i>E</i>)- <i>N</i> -[(Anthracen-9-yl)methylidene]hydroxylamine. IUCrData, 2017, 2, .	0.1	2
258	Methyl 2-[[[(6 <i>S</i> *,7 <i>R</i> *,8 <i>S</i> *)-7-acetyl-8-(4-chlorophenyl)-4-cyano-6-hydroxy-1,6-dimethyl-5,6,7,8-tetrahydroisoquinolin-3-yl]sulfanyl]acetate. IUCrData, 2017, 2, .	0.1	2
259	Ethyl 2-allyl-2-(5-methyl-1,2,4-triazolo[1,5- <i>a</i>]pyrimidin-7-yl)pent-4-enoate. IUCrData, 2017, 2, .	0.1	2
260	3-Benzyl-5,5-diphenylimidazolidine-2,4-dione. IUCrData, 2018, 3, .	0.1	2
261	(<i>Z</i>)-1-(1,3-Diphenyl-1 <i>H</i> -pyrazol-4-yl)- <i>N</i> -phenylmethanimine <i>N</i> -oxide. IUCrData, 2018, 3, .	0.1	2
262	Ethyl 2-(3-oxo-1,2,3,4-tetrahydroquinoxalin-2-yl)acetate. IUCrData, 2018, 3, .	0.1	2
263	Ethyl 2-[(2- <i>E</i> -4-decyl-3-oxo-1,2,3,4-tetrahydroquinoxalin-2-ylidene]acetate. IUCrData, 2018, 3, .	0.1	2
264	Ethyl 2-[2-(4-oxo-4 <i>H</i> -chromen-2-yl)phenoxy]acetate. IUCrData, 2018, 3, .	0.1	2
265	5-[[2-Hydroxyethyl]sulfanyl]methyl]quinolin-8-ol. IUCrData, 2019, 4, .	0.1	2
266	1-Bromo-2,4,6-tricyclohexylbenzene. Acta Crystallographica Section E: Structure Reports Online, 2008, 64, o335-o335.	0.2	2
267	2-(5-Methoxy-2-methyl-1 <i>H</i> -indol-3-yl)- <i>N</i> (<i>N</i>)-[1- <i>E</i> ,2- <i>E</i> -3-phenylprop-2-en-1-ylidene]acetohydrazide. Acta Crystallographica Section E: Structure Reports Online, 2013, 69, o1493-o1493.	0.2	2
268	<i>N</i> (<i>N</i>)-[1- <i>E</i> ,2- <i>E</i> -4-Methoxybenzylidene]-2-(5-methoxy-2-methyl-1 <i>H</i> -indol-3-yl)acetohydrazide. Acta Crystallographica Section E: Structure Reports Online, 2013, 69, o1660-o1661.	0.2	2
269	Crystal structure of 2-cyano-1-methylpyridinium tetrafluoroborate. Acta Crystallographica Section E: Crystallographic Communications, 2015, 71, o697-o698.	0.2	2
270	1-(5-Nitro-1 <i>H</i> -indazol-1-yl)ethanone. IUCrData, 2016, 1, .	0.1	2

#	ARTICLE	IF	CITATIONS
271	Cyclohexane-1,4-diammonium dithiocyanate. IUCrData, 2016, 1, .	0.1	2
272	4-Methyl-3,4-dihydro-2H-1,4-benzothiazin-3-one. IUCrData, 2017, 2, .	0.1	2
273	Ethyl 2-(6-nitro-1 <i>H</i> -indazol-1-yl)acetate. IUCrData, 2017, 2, .	0.1	2
274	1-(6-Nitro-1 <i>H</i> -indazol-1-yl)ethanone. IUCrData, 2017, 2, .	0.1	2
275	(4 <i>Z</i>)-4-(2-Oxopropylidene)-2,3,4,5-tetrahydro-1 <i>H</i> -1,5-benzodiazepin-2-one. IUCrData, 2017, 2, .	0.1	2
276	5-Nitro-1-(prop-2-en-1-yl)-1 <i>H</i> -indazole. IUCrData, 2017, 2, .	0.1	2
277	Ethyl 2-(4-benzyl-3-methyl-6-oxo-1,6-dihydropyridazin-1-yl)acetate: crystal structure and Hirshfeld surface analysis. Acta Crystallographica Section E: Crystallographic Communications, 2019, 75, 392-396.	0.2	2
278	Crystal structure, DFT study and Hirshfeld surface analysis of 1-nonyl-2,3-dihydro-1 <i>H</i> -indole-2,3-dione. Acta Crystallographica Section E: Crystallographic Communications, 2019, 75, 1140-1144.	0.2	2
279	Crystal structure, Hirshfeld surface analysis and DFT studies of 6-bromo-3-(12-bromododecyl)-2-(4-nitrophenyl)-4 <i>H</i> -imidazo[4,5- <i>b</i>]pyridine. Acta Crystallographica Section E: Crystallographic Communications, 2020, 76, 677-682.	0.2	2
280	A rare example of a double metamagnetic transition leading to 2D and 3D long-range order in the two-dimensional pyrazine- and azido-bridged cobalt(<i>II</i>) compound [Co(py ₂ N ₃) ₂]. Dalton Transactions, 2022, 51, 5617-5623.	1.6	2
281	Crystal structure, Hirshfeld surface analysis, interaction energy and DFT calculations and energy frameworks of methyl 6-chloro-1-methyl-2-oxo-1,2-dihydroquinoline-4-carboxylate. Acta Crystallographica Section E: Crystallographic Communications, 2022, 78, 425-432.	0.2	2
282	Synthesis, virtual screening and computational approach of a quinoxaline derivative as potent anti-HIV agent targeting the reverse transcriptase enzyme. Journal of Biomolecular Structure and Dynamics, 0, , 1-14.	2.0	2
283	Crystal structure and Hirshfeld surface analysis of 2-chloro- <i>N</i> -(4-methoxyphenyl)acetamide. Acta Crystallographica Section E: Crystallographic Communications, 2022, 78, 687-690.	0.2	2
284	Title is missing!. Journal of Chemical Crystallography, 1998, 28, 925-929.	0.5	1
285	Title is missing!. Journal of Chemical Crystallography, 2003, 33, 497-501.	0.5	1
286	Silylated gallium and indium chalcogenide ring systems as potential precursors to ME (E=O, S) materials. Open Chemistry, 2013, 11, 1225-1238.	1.0	1
287	2,2-[(1,3,4-Thiadiazole-2,5-diyl)bis(sulfaneyl)]diacetonitrile. Acta Crystallographica Section E: Structure Reports Online, 2013, 69, o1855-o1855.	0.2	1
288	N-[(<i>E</i>)-Benzylidene]-2-(6-methoxynaphthalen-2-yl)propanohydrazide. Acta Crystallographica Section E: Structure Reports Online, 2013, 69, o1614-o1614.	0.2	1

#	ARTICLE	IF	CITATIONS
289	Crystal structure of (E)-4-{2-(2,4-dinitrophenyl)hydrazin-1-ylidene}methyl}-3-methyl-1-phenyl-5-(1H-pyrrol-1-yl)-1H-pyrazole. Acta Crystallographica Section E: Structure Reports Online, 2014, 70, o1246-o1247.	0.2	1
290	3-[2-(4-Fluorophenyl)-2-oxoethyl]-5,5-diphenylimidazolidine-2,4-dione. Acta Crystallographica Section E: Structure Reports Online, 2014, 70, o226-o227.	0.2	1
291	4-Phenyl-1,2,4-triazaspiro[4.4]non-1-ene-3-thione. Acta Crystallographica Section E: Structure Reports Online, 2014, 70, o433-o434.	0.2	1
292	4-Chloro-N-[(3Z)-2-oxo-2,3-dihydro-1H-indol-3-ylidene]benzohydrazide. Acta Crystallographica Section E: Structure Reports Online, 2014, 70, o47-o47.	0.2	1
293	N-(12-Amino-9,10-dihydro-9,10-ethanoanthracen-11-yl)-4-methylbenzenesulfonamide. Acta Crystallographica Section E: Structure Reports Online, 2014, 70, o248-o249.	0.2	1
294	3-Amino-5,5-diphenylimidazolidine-2,4-dione. Acta Crystallographica Section E: Structure Reports Online, 2014, 70, o262-o263.	0.2	1
295	4-[(1,3-Dioxoisindolin-2-yl)methyl]benzenesulfonamide. Acta Crystallographica Section E: Structure Reports Online, 2014, 70, o291-o292.	0.2	1
296	Methyl 2-[(2Z,5Z)-4-oxo-3-phenyl-2-{2-[(1E)-1,2,3,4-tetrahydronaphthalen-1-ylidene]hydrazin-1-ylidene}-1,3-thiazolidin-5-ylidene)acetate. Acta Crystallographica Section E: Structure Reports Online, 2014, 70, o436-o437.	0.2	1
297	N-[(E)-2-Chlorobenzylidene]-2-(6-methoxynaphthalen-2-yl)propanohydrazide. Acta Crystallographica Section E: Structure Reports Online, 2014, 70, o631-o632.	0.2	1
298	4-Phenyl-1,2,4-triazaspiro[4.6]undec-1-ene-3-thione. Acta Crystallographica Section E: Structure Reports Online, 2014, 70, o640-o640.	0.2	1
299	(Z)-1-[(2E)-3,4-Diphenyl-2,3-dihydro-1,3-thiazol-2-ylidene]-2-[1-(4-hydroxyphenyl)ethylidene]hydrazinium bromide including an unknown solvate. Acta Crystallographica Section E: Structure Reports Online, 2014, 70, o647-o648.	0.2	1
300	(2E)-4-(4-Bromophenyl)-2-{2-[(1E)-cyclopentylidene]hydrazin-1-ylidene}-3-phenyl-2,3-dihydro-1,3-thiazole. Acta Crystallographica Section E: Structure Reports Online, 2014, 70, o669-o669.	0.2	1
301	Ethyl 2-[(4-amino-5-cyano-6-(methylsulfanyl)pyridin-2-yl)carbamoyl]methylsulfanyl]acetate monohydrate. Acta Crystallographica Section E: Structure Reports Online, 2014, 70, o745-o746.	0.2	1
302	Crystal structure of ethyl 2-[(4Z)-3,5-dioxo-1-phenylpyrazolidin-4-ylidene]methyl]amino)acetate. Acta Crystallographica Section E: Structure Reports Online, 2014, 70, o938-o939.	0.2	1
303	Crystal structure of 3-methyl-1-phenyl-5-(1H-pyrrol-1-yl)-1H-pyrazole-4-carbaldehyde. Acta Crystallographica Section E: Structure Reports Online, 2014, 70, o1131-o1132.	0.2	1
304	Crystal structure of 2-[12-methyl-14-phenyl-10,13,14,16-tetraazatetracyclo[7.7.0.0.2,7.0]hexadeca-1(16),2,4,6,9,11(15),12-hepta-8-ylidene]propanoate. Acta Crystallographica Section E: Structure Reports Online, 2014, 70, o1244-o1245.	0.2	1
305	Crystal structure of 5-(4,5-dihydro-1H-imidazol-2-yl)-3-methyl-1-phenyl-1H-pyrazolo[3,4-b]pyrazin-6-amine. Acta Crystallographica Section E: Structure Reports Online, 2014, 70, o1212-o1213.	0.2	1
306	Crystal structure of ethyl 2-{2-[(1Z)-1-hydroxy-3-(4-nitrophenyl)-3-oxoprop-1-en-1-yl]phenoxy}acetate. Acta Crystallographica Section E: Crystallographic Communications, 2015, 71, o917-o918.	0.2	1

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307	Crystal structure of [(E)-{2-[3-(2-[(1E)-[(carbamothioylamino)imino]methyl]phenoxy)propoxy]phenyl}methylidene)amino]thiourea with an unknown solvate. Acta Crystallographica Section E: Crystallographic Communications, 2015, 71, o530-o531.	0.2	1
308	Crystal structure of poly[di-μ ₄ -aqua-μ ₂ -[(1Z)-2-(4-chlorophenyl)-1-cyanoethenyl]-1,2,3,4-tetrazol-1-ido-μ ₂ -N1]sodium]. Acta Crystallographica Section E: Crystallographic Communications, 2015, 71, m102-m103.	0.2	1
309	Non-covalent interactions in tungsten-doped sodium ammonium decavanadate decahydrate. Journal of the Iranian Chemical Society, 2016, 13, 773-777.	1.2	1
310	Convenient one-pot access to novel densely functionalized pyrano[2,3-d][1,3,4]thiadiazolo[3,2-a]pyrimidines via three component reaction. Research on Chemical Intermediates, 2017, 43, 4683-4696.	1.3	1
311	Synthesis, resolution and crystal structures of two enantiomeric rhodamine derivatives. Acta Crystallographica Section E: Crystallographic Communications, 2017, 73, 327-333.	0.2	1
312	Crystal structure of unsymmetrical μ ₂ -diimine palladium(II) complex cis-[[ArN=C(Me)μ ^κ (Et)C=NAr]PdCl ₂] [Ar = 2,6-(iPr) ₂ C ₆ H ₃]. Acta Crystallographica Section E: Crystallographic Communications, 2017, 73, 1148-1150.	0.2	1
313	Synthesis, Crystal Structure, DFT Calculations and Hirshfeld Surface Analysis of 5-Bromo-1-decyl-2,3-dihydro-1H-indolin-2-one. Journal of Chemical Crystallography, 2020, 50, 330-337.	0.5	1
314	Crystal structure and Hirshfeld surface analysis of 2-[[7-acetyl-8-(4-chlorophenyl)-4-cyano-6-hydroxy-1,6-dimethyl-5,6,7,8-tetrahydroisoquinolin-3-yl]sulfanyl]-N,N'-bis(4-chlorophenyl)acetamide. Acta Crystallographica Section E: Crystallographic Communications, 2021, 77, 527-531.	0.2	1
315	Crystal structure and Hirshfeld surface analysis of 2-[[7-acetyl-4-cyano-6-hydroxy-8-(4-methoxyphenyl)-1,6-dimethyl-5,6,7,8-tetrahydroisoquinolin-3-yl]sulfanyl]-N,N'-bis(2-phenyl)acetamide. Acta Crystallographica Section E: Crystallographic Communications, 2021, 77, 663-667.	0.2	1
316	Crystal structure, Hirshfeld surface analysis and density functional theory study of benzyl 2-oxo-1-(prop-2-yn-1-yl)-1,2-dihydroquinoline-4-carboxylate. Acta Crystallographica Section E: Crystallographic Communications, 2021, 77, 824-828.	0.2	1
317	Synthesis and characterization of Mo(0) and W(0) complexes of bis(azol-1-yl)methane based bisphosphines. Journal of Coordination Chemistry, 2021, 74, 2253-2262.	0.8	1
318	Crystal structures of three sterically congested disilanes. Acta Crystallographica Section E: Crystallographic Communications, 2017, 73, 448-452.	0.2	1
319	Crystal structure and Hirshfeld surface analysis of 1-[(1-butyl-1 <i>H</i> -1,2,3-triazol-4-yl)methyl]-3-methylquinoxalin-2(1 <i>H</i>)-one. Acta Crystallographica Section E: Crystallographic Communications, 2018, 74, 1815-1820.	0.2	1
320	Crystal structure, Hirshfeld surface analysis and DFT studies of 1-benzyl-3-[(1-benzyl-1 <i>H</i> -1,2,3-triazol-5-yl)methyl]-2,3-dihydro-1 <i>H</i> -1,3-benzodiazol-2-one monohydrate. Acta Crystallographica Section E: Crystallographic Communications, 2020, 76, 95-101.	0.2	1
321	Crystal structure, Hirshfeld surface analysis and interaction energy and DFT studies of 1-(1,3-benzothiazol-2-yl)-3-(2-hydroxyethyl)imidazolidin-2-one. Acta Crystallographica Section E: Crystallographic Communications, 2020, 76, 370-376.	0.2	1
322	1-Ethyl-4-phenyl-1,5-benzodiazepine-2-thione. IUCrData, 2017, 2, .	0.1	1
323	3-Phenylisoxazolin-5-one: a redetermination. IUCrData, 2017, 2, .	0.1	1
324	3-Hydroxy-3-(2-oxo-2,3-dihydro-1H-indol-3-yl)-2,3-dihydro-1H-indol-2-one. IUCrData, 2017, 2, .	0.1	1

#	ARTICLE	IF	CITATIONS
325	1-[(1-[[[(1S,2R,6R,8R,9S)-4,4,11,11-Tetramethyl-3,5,7,10,12-pentaoxatricyclo[7.3.0.0.2,6]dodecan-8-yl)methyl]-1H-1,2,3-triazol-4-yl)methyl]methyl]pyridin-2-yl]methane. IUCrData, 2017, 2, .	0.1	1
326	2-[(Prop-2-yn-1-yl)amino]anilinium chloride. IUCrData, 2017, 2, .	0.1	1
327	Bis(2-formylphenyl) benzene-1,2-dicarboxylate. IUCrData, 2018, 3, .	0.1	1
328	(Pyridin-2-yl)methyl 6-bromo-2-oxo-1-[(pyridin-2-yl)methyl]-1,2-dihydroquinoline-4-carboxylate. IUCrData, 2018, 3, .	0.1	1
329	(3 <i>R</i> ,4 <i>Z</i>)-1,3-Diethyl-4-(2-oxopropylidene)-2,3,4,5-tetrahydro-1 <i>H</i> -1,5-benzodiazepin-2-one. IUCrData, 2018, 3, .	0.1	1
330	2-Methyl-4-(pyridin-2-yl)-3 <i>H</i> -1,5-benzodiazepine. IUCrData, 2018, 3, .	0.1	1
331	Crystal structure of (E)-N-[[3-methyl-1-phenyl-5-(1 <i>H</i> -pyrrol-1-yl)-1 <i>H</i> -pyrazol-4-yl]methylidene]hydroxylamine. Acta Crystallographica Section E: Structure Reports Online, 2014, 70, o1216-o1217.	0.2	1
332	Crystal structure of (4-methoxyphenyl)[(4-methoxyphenyl)phosphonato]dioxidophosphate(1 ⁻) 2-amino-6-benzyl-3-ethoxycarbonyl-4,5,6,7-tetrahydrothieno[2,3-c]pyridin-6-ium. Acta Crystallographica Section E: Crystallographic Communications, 2015, 71, o997-o998.	0.2	1
333	Crystal structure of 3-benzyl-1-[(1,2,3,4-tetrahydronaphthalen-1-ylidene)amino]thiourea. Acta Crystallographica Section E: Crystallographic Communications, 2015, 71, o974-o975.	0.2	1
334	5-Nitro-1-(prop-2-yn-1-yl)-1 <i>H</i> -indazole. IUCrData, 2016, 1, .	0.1	1
335	2-(2-Amino-5-methyl-1,2,4-triazolo[1,5-a]pyrimidin-7-yl)acetohydrazide monohydrate. IUCrData, 2016, 1, .	0.1	1
336	2-[(2 <i>Z</i>)-2-Benzylidene-3-oxo-3,4-dihydro-2 <i>H</i> -1,4-benzothiazin-4-yl]acetic acid. IUCrData, 2016, 1, .	0.1	1
337	N,N,N-Triethylethanaminium 5,11,17,23-tetra-tert-butyl-25-[(ethoxycarbonyl)methoxy]-26,28-dihydroxy-27-oxido-2,8,14,20-tetrathiacalix[4]arene 1 a molecular salt. IUCrData, 2016, 1, .	0.1	1
338	(E)-N ² -Benzylidene-2-phenylquinoline-4-carbohydrazide. IUCrData, 2016, 1, .	0.1	1
339	IUCrData, 2016, 1, .	0.1	1
340	3-[2-(9 <i>H</i> -Carbazol-9-yl)ethyl]-4-phenyl-1 <i>H</i> -1,2,4-triazole-5(4 <i>H</i>)-thione dimethyl sulfoxide monosolvate. IUCrData, 2016, 1, .	0.1	1
341	(4 <i>Z</i>)-1-Dodecyl-4-(2-oxopropylidene)-2,3,4,5-tetrahydro-1 <i>H</i> -1,5-benzodiazepin-2-one. IUCrData, 2016, 1, .	0.1	1
342	(2 <i>Z</i>)-2-Benzylidene-4-[(1-benzyl-1 <i>H</i> -1,2,3-triazol-4-yl)methyl]-3,4-dihydro-2 <i>H</i> -1,4-benzothiazin-3-one. IUCrData, 2016, 1, .	0.1	1

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343	2-Methyl-3-(3-methylisoxazol-5-yl)-4-oxo-4 <i>H</i> -pyrido[1,2- <i>a</i>]pyrimidin-1-ium chloride. IUCrData, 2016, 1, .	0.1	1
344	(5 <i>Z</i>)-3-(2-Oxopropyl)-5-(3,4,5-trimethoxybenzylidene)-1,3-thiazolidine-2,4-dione. IUCrData, 2016, 1, .	0.1	1
345	(<i>E</i>)-4-Methoxy-3,5-dimethyl-2-[(3-nitrophenyl)ethenyl]pyridine. IUCrData, 2016, 1, .	0.1	1
346	2,3-Dihydrobenz[4,5]imidazo[2,1- <i>b</i>]thiazole. IUCrData, 2016, 1, .	0.1	1
347	N-{2-[2-(5-Methyl-1 <i>H</i> -pyrazol-3-yl)acetamido]phenyl}benzamide monohydrate. IUCrData, 2017, 2, .	0.1	1
348	3-Chloro-6-nitro-1-[(1-octyl-1 <i>H</i> -1,2,3-triazol-4-yl)methyl]-1 <i>H</i> -indazole. IUCrData, 2017, 2, .	0.1	1
349	(4 <i>Z</i>)-4-(2-Oxopropylidene)-1,3-bis(prop-2-en-1-yl)-2,3,4,5-tetrahydro-1 <i>H</i> -1,5-benzodiazepin-2-one. IUCrData, 2017, 2, .	0.1	1
350	(2 <i>Z</i>)-2-Benzylidene-4-octadecyl-3,4-dihydro-2 <i>H</i> -1,4-benzothiazin-3-one. IUCrData, 2017, 2, .	0.1	1
351	(2 <i>Z</i>)-2-(4-Chlorobenzylidene)-4-[2-(2-oxooxazoliden-3-yl)ethyl]-3,4-dihydro-2 <i>H</i> -1,4-benzothiazin-3-one. IUCrData, 2017, 2, .	0.1	1
352	1-(3-Bromo-6-nitro-1 <i>H</i> -indazol-1-yl)ethan-1-one. IUCrData, 2017, 2, .	0.1	1
353	1-Chloro-3-(6-nitro-1 <i>H</i> -indazol-1-yl)propan-2-ol. IUCrData, 2017, 2, .	0.1	1
354	3-Acetyl-2-methyl-4 <i>H</i> -pyrido[1,2- <i>a</i>]pyrimidin-4-one. IUCrData, 2017, 2, .	0.1	1
355	4-[(<i>E</i>)-(2-Hydroxynaphthalen-1-yl)methylidene]amino}-1,5-dimethyl-2-phenyl-2,3-dihydro-1 <i>H</i> -pyrazol-3-one: a new polymorph (<i>i</i> ² -phase). IUCrData, 2017, 2, .	0.1	1
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357	A monoclinic modification of (4 <i>Z</i>)-1-benzyl-4-(2-oxopropylidene)-2,3,4,5-tetrahydro-1 <i>H</i> -1,5-benzodiazepin-2-one. IUCrData, 2018, 3, .	0.1	1
358	1-Methyl-4-phenyl-3-[4-(trifluoromethyl)phenyl]-1 <i>H</i> -pyrazolo[3,4- <i>d</i>]pyrimidine. IUCrData, 2018, 3, .	0.1	1
359	4-Phenyl-5 <i>a</i> ,6,7,8,9,9 <i>a</i> -hexahydro-1 <i>H</i> -1,5-benzodiazepin-2(5 <i>H</i>)-one. IUCrData, 2018, 3, .	0.1	1
360	Diethyl 4-(4-chloro-2-propyl-1 <i>H</i> -imidazol-5-yl)-2,6-dimethyl-1,4-dihydropyridine-3,5-dicarboxylate monohydrate. IUCrData, 2018, 3, .	0.1	1

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363	Crystal structure, computational study and Hirshfeld surface analysis of ethyl (2 <i>S</i> ,3 <i>R</i>)-3-(3-amino-1 <i>H</i> -1,2,4-triazol-1-yl)-2-hydroxy-3-phenylpropanoate. Acta Crystallographica Section E: Crystallographic Communications, 2019, 75, 1919-1924.	0.2	1
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374	3,6-Dichloro-9-(prop-2-yn-1-yl)-9 <i>H</i> -carbazole. Acta Crystallographica Section E: Structure Reports Online, 2014, 70, o27-o27.	0.2	0
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390	Crystal structure of 4-[(E)-(4-hydroxybenzylidene)amino]-1,5-dimethyl-2-phenyl-1H-pyrazol-3(2H)-one. Acta Crystallographica Section E: Crystallographic Communications, 2015, 71, o947-o948.	0.2	0
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400	Crystal structures of two hydrazide derivatives of mefenamic acid, 3-(2,3-dimethylanilino)- <i>N</i> -[β -(<i>E</i>)-(furan-2-yl)methylidene]benzohydrazide and <i>N</i> -[β -(<i>E</i>)-benzylidene]-2-(2,3-dimethylanilino)benzohydrazide. Acta Crystallographica Section E: Crystallographic Communications, 2021, 77, 242-246.	0.2	0
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402	Crystal structure and Hirshfeld surface analysis of (3 <i>S</i> ,3 <i>R</i> ,6 <i>S</i>)-3-(1,3-diphenyl-1 <i>H</i> -pyrazol-4-yl)-5-(4-methoxyphenyl)-2-phenyl-3,3a,4,5,6,6a-hexahydro-2 <i>H</i> -indolizino[1,2- <i>b</i>]pyridine. Acta Crystallographica Section E: Crystallographic Communications, 2021, 77, 356-359.	0.2	0
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416	Crystal structure of (5Z)-5-(5-bromo-2-hydroxybenzylidene)-1,3-thiazolidine-2,4-dione. Acta Crystallographica Section E: Crystallographic Communications, 2015, 71, o919-o920.	0.2	0
417	Crystal structure of 3-(prop-2-en-1-yl)-1-[[[(1E)-1,2,3,4-tetrahydronaphthalen-1-ylidene]amino]thiourea. Acta Crystallographica Section E: Crystallographic Communications, 2015, 71, o976-o977.	0.2	0
418	(4Z)-1-Benzyl-4-(2-oxopropylidene)-2,3,4,5-tetrahydro-1H-1,5-benzodiazepin-2-one. IUCrData, 2016, 1, .	0.1	0
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420	Methyl 2-phenylquinoline-4-carboxylate. IUCrData, 2016, 1, .	0.1	0
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425	Ethyl 2-(6-bromo-2-phenyl-3H-imidazo[4,5-b]pyridin-3-yl)acetate. IUCrData, 2016, 1, .	0.1	0
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428	5-Fluoro-1-(prop-2-en-1-yl)-2,3-dihydro-1H-indole-2,3-dione. IUCrData, 2017, 2, .	0.1	0
429	2-(2-Oxo-4-phenyl-2,3-dihydro-1H-1,5-benzodiazepin-1-yl)acetic acid. IUCrData, 2017, 2, .	0.1	0
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434	5-[[5-(4-Chlorophenyl)-3-methyl-1H-pyrazol-1-yl]methyl]-1,3,4-oxadiazole-2(3H)-thione. IUCrData, 2017, 2, .	0.1	0
435	1-(3-Phenyl-1H-pyrazol-5-yl)-1,2,3,4-tetrahydroquinoxaline-2,3-dione. IUCrData, 2017, 2, .	0.1	0
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438	1,4-Di-n-octyl-1,2,3,4-tetrahydroquinoxaline-2,3-dione. IUCrData, 2017, 2, .	0.1	0
439	5-Fluoro-1-[(1-[(1 <i>S</i> ,2 <i>R</i> ,6 <i>R</i> ,8 <i>S</i> ,9 <i>R</i>)-4,4,11,11-tetramethyl-3,5,7,10,12-pentaoxatricyclo[7.3.0.0 ^{2,6}]dodecan-8-yl)methyl]-1H-1,2,3-triazol-5-yl)methyl]-1H-1,2,3-triazole. IUCrData, 2017, 2, .	0.1	0
440	1-(Prop-2-en-1-yl)-3-[(prop-2-en-1-yl)oxy]quinoxalin-2(1H)-one. IUCrData, 2017, 2, .	0.1	0
441	1-[(1-Benzyl-1H-1,2,3-triazol-4-yl)methyl]-1H-1,3-benzodiazole. IUCrData, 2017, 2, .	0.1	0
442	5-Methyl-4-(3-methyl-1-phenyl-1H-pyrazol-4-yl)-2-phenyl-2,3-dihydro-1H-pyrazol-3-one. IUCrData, 2017, 2, .	0.1	0
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445	1,5-Dimethyl-3,5-diphenyl-1,5-dihydro-3 <i>H</i> -spiro[pyrazolo[3,4- <i>d</i>]pyrimidine-4,2- ² [1,3,4]-thiadiazole]. IUCrData, 2017, 2, .	0.1	0
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447	(Pyridin-2-yl)methyl 2-oxo-1-[(pyridin-2-yl)methyl]-1,2-dihydroquinoline-4-carboxylate hemihydrate. IUCrData, 2017, 2, .	0.1	0
448	1-(3-Chloro-6-nitro-1H-indazol-1-yl)ethan-1-one. IUCrData, 2017, 2, .	0.1	0
449	1-Methyl-4-phenyl-1H-pyrazolo[3,4- <i>d</i>]pyrimidine. IUCrData, 2017, 2, .	0.1	0
450	(3 <i>E</i>)-4-(3,4,5-Trimethoxyphenyl)but-3-en-2-one. IUCrData, 2017, 2, .	0.1	0

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451	4-(Prop-2-yn-1-ylsulfanyl)-1H-pyrazolo[3,4-d]pyrimidine. IUCrData, 2017, 2, .	0.1	0
452	Nâ€²-[(1Z)-1-(3-Methyl-5-oxo-1-phenyl-4,5-dihydro-1H-pyrazol-4-ylidene)ethyl]-2-[(4-methylphenyl)sulfanyl]acetohydrazide. IUCrData, 2017, 2, .	0.1	0
453	Ethyl 3-amino-2-cyano-4-(4-methoxyphenyl)-6-methylthieno[2,3-b]pyridine-5-carboxylate. IUCrData, 2017, 2, .	0.1	0
454	5-Acetyl-3-amino-4-(4-methoxyphenyl)-6-methylthieno[2,3-b]pyridine-2-carbonitrile. IUCrData, 2017, 2, .	0.1	0
455	3-Bromo-6-nitro-1-(prop-2-en-1-yl)-1H-indazole. IUCrData, 2018, 3, .	0.1	0
456	Nâ€²-[(1E)-2,5-Dimethoxybenzylidene]pyridine-2-carbohydrazide. IUCrData, 2018, 3, .	0.1	0
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459	1-[(Oxiran-2-yl)methyl]-3-phenyl-1,2-dihydroquinoxalin-2-one. IUCrData, 2018, 3, .	0.1	0
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