

El Mokhtar Essassi

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/3109156/publications.pdf>

Version: 2024-02-01

23
papers

199
citations

1163117

8
h-index

1058476

14
g-index

23
all docs

23
docs citations

23
times ranked

188
citing authors

#	ARTICLE	IF	CITATIONS
1	Co(II) and Zn(II) pyrazolyl-benzimidazole complexes with remarkable antibacterial activity. <i>New Journal of Chemistry</i> , 2020, 44, 2210-2221.	2.8	54
2	Catalyst- and Base-Controlled Site-Selective sp^2 and sp^3 Direct Arylation of 5,7-Dimethyl-2-phenylpyrazolo[1,5-a]pyrimidine Using Aryl Bromides. <i>European Journal of Organic Chemistry</i> , 2012, 2012, 2572-2578.	2.4	34
3	A one-pot process for the microwave-assisted synthesis of 7-substituted pyrazolo[1,5-a]pyrimidine. <i>RSC Advances</i> , 2016, 6, 3301-3306.	3.6	15
4	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. <i>Journal of Heterocyclic Chemistry</i> , 2021, 58, 270-289.	2.6	12
5	Coordination complexes constructed from pyrazole-acetamide and pyrazole-quinoxaline: effect of hydrogen bonding on the self-assembly process and antibacterial activity. <i>RSC Advances</i> , 2022, 12, 5324-5339.	3.6	10
6	One-Pot S_NAr /Direct Pd-Catalyzed CH Arylation Functionalization of Pyrazolo[1,5-a]pyrimidine at the C3 and C7 Positions. <i>European Journal of Organic Chemistry</i> , 2018, 2018, 3936-3942.	2.4	9
7	In Vitro Evaluation of the Multidrug Resistance Reversing Activity of Novel Imidazo[4,5-b]pyridine Derivatives. <i>Anticancer Research</i> , 2018, 38, 3999-4003.	1.1	9
8	Crystal structure and Hirshfeld surface analysis of N -{2-[(<i>E</i>)-(4-methylbenzylidene)amino]phenyl}-2-(5-methyl-1 <i>H</i> -pyrazol-3-yl)acetamide hemihydrate. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2019, 75, 154-158.	0.5	8
9	Crystal structure and Hirshfeld surface analysis of a new benzodiazepine derivative: 4-dichloromethyl-2,3-dihydro-1 <i>H</i> -1,5-benzodiazepin-2-one. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2019, 75, 33-37.	0.5	7
10	Crystal structure, DFT study and Hirshfeld surface analysis of ethyl 6-chloro-2-ethoxyquinoline-4-carboxylate. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2019, 75, 912-916.	0.5	7
11	Synthesis and crystal structure of 1-octyl-3-phenylquinoxalin-2(1 <i>H</i>)-one, $C_{22}H_{26}N_2O$. <i>Zeitschrift Fur Kristallographie - New Crystal Structures</i> , 2021, 236, 173-175.	0.3	5
12	Synthesis and crystal structure of (<i>E</i>)-1-benzyl-3-(4-methoxystyryl)quinoxalin-2(1 <i>H</i>)-one, $C_{24}H_{20}N_2O_2$. <i>Zeitschrift Fur Kristallographie - New Crystal Structures</i> , 2020, 235, 1323-1325.	0.3	5
13	7-Chloro-5-methyl-2-phenylpyrazolo[1,5-a]pyrimidine. <i>Acta Crystallographica Section E: Structure Reports Online</i> , 2013, 69, o749-o749.	0.2	4
14	Crystal structure, Hirshfeld surface analysis and DFT study of 1-ethyl-3-phenyl-1,2-dihydroquinoxalin-2-one. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2021, 77, 18-22.	0.5	3
15	Crystal structure, Hirshfeld surface analysis and density functional theory study of 6-methyl-2-[(5-methylisoxazol-3-yl)methyl]-1 <i>H</i> -benzimidazole. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2021, 77, 396-401.	0.5	3
16	Crystal structure, Hirshfeld surface analysis and DFT study of N -{2-(2-amino-5-methylphenyl)-2-(5-methyl-1 <i>H</i> -pyrazol-3-yl)acetamide}. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2021, 77, 638-642.	0.5	3
17	One-Step Synthesis of novel $N1$ -substituted benzimidazole derivatives: Experimental and theoretical investigations. <i>Journal of Heterocyclic Chemistry</i> , 0, , .	2.6	3
18	A new synthetic route for the preparation of 2,2,5-trimethyl-7,7-dihydro-6,7-bipyrazolo[1,5-a]pyrimidine-3,3-dicarbonitrile, structural elucidation, Hirshfeld surface analysis, energy framework, density functional theory and molecular docking investigations. <i>Journal of the Chinese Chemical Society</i> , 2022, 69, 717-730.	1.4	3

#	ARTICLE	IF	CITATIONS
19	Crystal structure, Hirshfeld surface analysis and density functional theory study of 1-nonyl-3-phenylquinoxalin-2-one. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2021, 77, 1037-1042.	0.5	2
20	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. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2022, 78, 425-432.	0.5	2
21	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. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2021, 77, 824-828.	0.5	1
22	Crystal structure, Hirshfeld surface analysis, DFT and molecular docking investigation of 2-(2-oxo-1,3-oxazolidin-3-yl)ethyl 2-[2-(2-oxo-1,3-oxazolidin-3-yl)ethoxy]quinoline-4-carboxylate. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2021, 77, 28-33.	0.5	0
23	Crystal structure and Hirshfeld surface analysis of hexyl 1-hexyl-2-oxo-1,2-dihydroquinoline-4-carboxylate. <i>Acta Crystallographica Section E: Crystallographic Communications</i> , 2020, 76, 642-645.	0.5	0