Tahar Lakhlifi

List of Publications by Year
in descending order

4 papers	55 citations	1937685 4 h-index	2272923 4 g-index
4 all docs	4 docs citations	4 times ranked	74 citing authors

#	Article	IF	Citations
1	3D-QSAR modeling and molecular docking studies on a series of 2,5 disubstituted 1,3,4-oxadiazoles. Journal of Molecular Structure, 2017, 1145, 278-284.	3.6	22
2	In Silico Exploration of Aryl Halides Analogues as CheckpointKinase 1 Inhibitors by Using 3D QSAR, Molecular Docking Study,and ADMET Screening. Advanced Pharmaceutical Bulletin, 2019, 9, 84-92.	1.4	17
3	Integrated 3D-QSAR, molecular docking, and molecular dynamics simulation studies on 1,2,3-triazole based derivatives for designing new acetylcholinesterase inhibitors. Turkish Journal of Chemistry, 2021, 45, 647-660.	1.2	10
4	Rational design of novel potential EGFR inhibitors by 3D-QSAR, molecular docking, molecular dynamics simulation, and pharmacokinetics studies. Chemical Data Collections, 2022, 39, 100851.	2.3	6