

# Tahar Lakhli

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/817149/publications.pdf>

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