

Giulia Culetta

List of Publications by Year in descending order

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8
papers

113
citations

1478505

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h-index

1588992

8
g-index

8
all docs

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docs citations

8
times ranked

139
citing authors

#	ARTICLE	IF	CITATIONS
1	In Silico Design, Synthesis, and Biological Evaluation of Anticancer Arylsulfonamide Endowed with Anti-Telomerase Activity. <i>Pharmaceuticals</i> , 2022, 15, 82.	3.8	11
2	Immunoproteasome and Non-Covalent Inhibition: Exploration by Advanced Molecular Dynamics and Docking Methods. <i>Molecules</i> , 2021, 26, 4046.	3.8	3
3	Evaluation of the IKK β Binding of Indicaxanthin by Induced-Fit Docking, Binding Pose Metadynamics, and Molecular Dynamics. <i>Frontiers in Pharmacology</i> , 2021, 12, 701568.	3.5	24
4	Exploring the SARS-CoV-2 Proteome in the Search of Potential Inhibitors via Structure-Based Pharmacophore Modeling/Docking Approach. <i>Computation</i> , 2020, 8, 77.	2.0	30
5	Comparing molecular dynamics-derived pharmacophore models with docking: A study on CDK-2 inhibitors. <i>Chemical Data Collections</i> , 2020, 28, 100485.	2.3	8
6	Pharmacophore-Based Design of New Chemical Scaffolds as Translational Readthrough-Inducing Drugs (TRIDs). <i>ACS Medicinal Chemistry Letters</i> , 2020, 11, 747-753.	2.8	13
7	A Definitive Pharmacophore Modelling Study on CDK2 ATP Pocket Binders: Tracing the Path of New Virtual High-Throughput Screenings. <i>Current Drug Discovery Technologies</i> , 2020, 17, 740-747.	1.2	2
8	Indicaxanthin, a multi-target natural compound from <i>Opuntia ficus-indica</i> fruit: From its poly-pharmacological effects to biochemical mechanisms and molecular modelling studies. <i>European Journal of Medicinal Chemistry</i> , 2019, 179, 753-764.	5.5	22