Sehan Lee

List of Publications by Year in descending order

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1163117 1474206 9 155 8 9 citations h-index g-index papers 10 10 10 239 citing authors docs citations times ranked all docs

#	Article	IF	CITATION
1	A mechanism-based 3D-QSAR approach for classification and prediction of acetylcholinesterase inhibitory potency of organophosphate and carbamate analogs. Journal of Computer-Aided Molecular Design, 2016, 30, 347-363.	2.9	34
2	Structure-Based Understanding of Binding Affinity and Mode of Estrogen Receptor \hat{l}_{\pm} Agonists and Antagonists. PLoS ONE, 2017, 12, e0169607.	2.5	27
3	Development of 3D-QSAR Model for Acetylcholinesterase Inhibitors Using a Combination of Fingerprint, Molecular Docking, and Structure-Based Pharmacophore Approaches. Toxicological Sciences, 2015, 148, 60-70.	3.1	25
4	Calculation of the Solvation Free Energy of Neutral and Ionic Molecules in Diverse Solvents. Journal of Chemical Information and Modeling, 2011, 51, 105-114.	5.4	17
5	In Silico Site-Directed Mutagenesis Informs Species-Specific Predictions of Chemical Susceptibility Derived From the Sequence Alignment to Predict Across Species Susceptibility (SeqAPASS) Tool. Toxicological Sciences, 2018, 166, 131-145.	3.1	17
6	3D-QSAR study of steroidal and azaheterocyclic human aromatase inhibitors using quantitative profile of protein–ligand interactions. Journal of Cheminformatics, 2018, 10, 2.	6.1	14
7	Development of Surface-SFED Models for Polar Solvents. Journal of Chemical Information and Modeling, 2012, 52, 440-448.	5.4	9
8	A generalized G-SFED continuum solvation free energy calculation model. Proceedings of the National Academy of Sciences of the United States of America, 2013, 110, E662-7.	7.1	9
9	Incorporation of Hydrogen Bond Angle Dependency into the Generalized Solvation Free Energy Density Model. Journal of Chemical Information and Modeling, 2018, 58, 761-772.	5.4	3