

Yuan Xiaohui

List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	Structure Based Affinity Maturation and Characterizing of SARS-CoV Antibody CR3022 against SARS-CoV-2 by Computational and Experimental Approaches. <i>Viruses</i> , 2022, 14, 186.	3.3	6
2	Interaction Analysis of Odorant-Binding Protein 12 from <i>Sirex noctilio</i> and Volatiles from Host Plants and Symbiotic Fungi Based on Molecule Dynamics Simulation. <i>Agronomy</i> , 2022, 12, 861.	3.0	2
3	Mechanistic Insights to the Binding of Antibody CR3022 Against RBD from SARS-CoV and HCoV-19/SARS-CoV-2: A Computational Study. <i>Combinatorial Chemistry and High Throughput Screening</i> , 2021, 24, 1069-1082.	1.1	10
4	Computational identification of potential chemoprophylactic agents according to dynamic behavior of peroxisome proliferator-activated receptor gamma. <i>RSC Advances</i> , 2021, 11, 147-159.	3.6	5
5	Computational and Experimental Approaches to Decipher the Binding Mechanism of General Odorant-Binding Protein 2 from <i>Athetis lepigone</i> to Chlorpyrifos and Phoxim. <i>Journal of Agricultural and Food Chemistry</i> , 2021, 69, 88-100.	5.2	18
6	Computational Interaction Analysis of <i>Sirex noctilio</i> Odorant-Binding Protein (SnocOBP7) Combined with Female Sex Pheromones and Symbiotic Fungal Volatiles. <i>Agronomy</i> , 2021, 11, 2461.	3.0	6
7	Computational Simulation of HIV Protease Inhibitors to the Main Protease (Mpro) of SARS-CoV-2: Implications for COVID-19 Drugs Design. <i>Molecules</i> , 2021, 26, 7385.	3.8	6
8	Assessing the performance of the g_mmpbsa tools to simulate the inhibition of oseltamivir to influenza virus neuraminidase by molecular mechanics Poisson-Boltzmann surface area methods. <i>Journal of the Chinese Chemical Society</i> , 2020, 67, 46-53.	1.4	43
9	Rapid Structure-Based Screening Informs Potential Agents for Coronavirus Disease (COVID-19) Outbreak*. <i>Chinese Physics Letters</i> , 2020, 37, 058701.	3.3	11
10	Key Amino Acid Residues Influencing Binding Affinities of Pheromone-Binding Protein from <i>Athetis lepigone</i> to Two Sex Pheromones. <i>Journal of Agricultural and Food Chemistry</i> , 2020, 68, 6092-6103.	5.2	17
11	Organophosphorus insecticide interacts with the pheromone-binding proteins of <i>Athetis lepigone</i> : Implication for olfactory dysfunction. <i>Journal of Hazardous Materials</i> , 2020, 397, 122777.	12.4	30
12	Molecular Docking and Molecular Dynamics (MD) Simulation of Human Anti-Complement Factor H (CFH) Antibody Ab42 and CFH Polypeptide. <i>International Journal of Molecular Sciences</i> , 2019, 20, 2568.	4.1	27
13	Novel binding patterns between ganoderic acids and neuraminidase: Insights from docking, molecular dynamics and MM/PBSA studies. <i>Journal of Molecular Graphics and Modelling</i> , 2016, 65, 27-34.	2.4	24
14	Phylogenetic and structural analysis of major surface proteins hemagglutinin and neuraminidase of novel avian influenza virus A H7N9 from chinese patient. <i>Chemical Research in Chinese Universities</i> , 2013, 29, 934-940.	2.6	5
15	A Novel Antibody Humanization Method Based on Epitopes Scanning and Molecular Dynamics Simulation. <i>PLoS ONE</i> , 2013, 8, e80636.	2.5	24
16	Structure-Based High-Throughput Epitope Analysis of Hexon Proteins in B and C Species Human Adenoviruses (HAdVs). <i>PLoS ONE</i> , 2012, 7, e32938.	2.5	19
17	The Antiadenovirus Activities of Cinnamaldehyde In Vitro. <i>Laboratory Medicine</i> , 2009, 40, 669-674.	1.2	14
18	Molecular modeling and epitopes mapping of human adenovirus type 3 hexon protein. <i>Vaccine</i> , 2009, 27, 5103-5110.	3.8	53