

# Yutaka Akiyama

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/8552353/publications.pdf>

Version: 2024-02-01

17  
papers

193  
citations

1306789

7  
h-index

1125271

13  
g-index

18  
all docs

18  
docs citations

18  
times ranked

241  
citing authors

#	ARTICLE	IF	CITATIONS
1	Large-Scale Membrane Permeability Prediction of Cyclic Peptides Crossing a Lipid Bilayer Based on Enhanced Sampling Molecular Dynamics Simulations. <i>Journal of Chemical Information and Modeling</i> , 2021, 61, 3681-3695.	2.5	50
2	Multiple grid arrangement improves ligand docking with unknown binding sites: Application to the inverse docking problem. <i>Computational Biology and Chemistry</i> , 2018, 73, 139-146.	1.1	30
3	MEGADOCK-Web: an integrated database of high-throughput structure-based protein-protein interaction predictions. <i>BMC Bioinformatics</i> , 2018, 19, 62.	1.2	25
4	A prospective compound screening contest identified broader inhibitors for Sirtuin 1. <i>Scientific Reports</i> , 2019, 9, 19585.	1.6	15
5	Computational prediction of plasma protein binding of cyclic peptides from small molecule experimental data using sparse modeling techniques. <i>BMC Bioinformatics</i> , 2018, 19, 527.	1.2	9
6	QEX: target-specific druglikeness filter enhances ligand-based virtual screening. <i>Molecular Diversity</i> , 2019, 23, 11-18.	2.1	9
7	GHOSTX: A Fast Sequence Homology Search Tool for Functional Annotation of Metagenomic Data. <i>Methods in Molecular Biology</i> , 2017, 1611, 15-25.	0.4	8
8	Comprehensive Fungal Community Analysis of House Dust Using Next-Generation Sequencing. <i>International Journal of Environmental Research and Public Health</i> , 2020, 17, 5842.	1.2	8
9	PKRank: a novel learning-to-rank method for ligand-based virtual screening using pairwise kernel and RankSVM. <i>Artificial Life and Robotics</i> , 2018, 23, 205-212.	0.7	6
10	Solving Generalized Polyomino Puzzles Using the Ising Model. <i>Entropy</i> , 2022, 24, 354.	1.1	6
11	Plasma protein binding prediction focusing on residue-level features and circularity of cyclic peptides by deep learning. <i>Bioinformatics</i> , 2022, 38, 1110-1117.	1.8	5
12	A Massively Parallel Sequence Similarity Search for Metagenomic Sequencing Data. <i>International Journal of Molecular Sciences</i> , 2017, 18, 2124.	1.8	4
13	Taxonomic and Gene Category Analyses of Subgingival Plaques from a Group of Japanese Individuals with and without Periodontitis. <i>International Journal of Molecular Sciences</i> , 2021, 22, 5298.	1.8	3
14	Multiple HPC Environments-Aware Container Image Configuration Workflow for Large-Scale All-to-All Protein-Protein Docking Calculations. <i>Lecture Notes in Computer Science</i> , 2020, , 23-39.	1.0	3
15	Optimization of memory use of fragment extension-based protein-ligand docking with an original fast minimum cost flow algorithm. <i>Computational Biology and Chemistry</i> , 2018, 74, 399-406.	1.1	2
16	Learning-to-rank technique based on ignoring meaningless ranking orders between compounds. <i>Journal of Molecular Graphics and Modelling</i> , 2019, 92, 192-200.	1.3	1
17	Improved Large-Scale Homology Search by Two-Step Seed Search Using Multiple Reduced Amino Acid Alphabets. <i>Genes</i> , 2021, 12, 1455.	1.0	1