

# Susumu Goto

## List of Publications by Year in descending order

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Version: 2024-02-01

37  
papers

22,840  
citations

218381

26  
h-index

377514

34  
g-index

38  
all docs

38  
docs citations

38  
times ranked

35467  
citing authors

#	ARTICLE	IF	CITATIONS
1	KEGG for linking genomes to life and the environment. <i>Nucleic Acids Research</i> , 2007, 36, D480-D484.	6.5	5,451
2	KEGG for integration and interpretation of large-scale molecular data sets. <i>Nucleic Acids Research</i> , 2012, 40, D109-D114.	6.5	4,174
3	The KEGG resource for deciphering the genome. <i>Nucleic Acids Research</i> , 2004, 32, 277D-280.	6.5	4,093
4	Data, information, knowledge and principle: back to metabolism in KEGG. <i>Nucleic Acids Research</i> , 2014, 42, D199-D205.	6.5	2,846
5	KEGG for representation and analysis of molecular networks involving diseases and drugs. <i>Nucleic Acids Research</i> , 2010, 38, D355-D360.	6.5	1,999
6	The KEGG databases at GenomeNet. <i>Nucleic Acids Research</i> , 2002, 30, 42-46.	6.5	1,202
7	Development of a Chemical Structure Comparison Method for Integrated Analysis of Chemical and Genomic Information in the Metabolic Pathways. <i>Journal of the American Chemical Society</i> , 2003, 125, 11853-11865.	6.6	423
8	ViPTree: the viral proteomic tree server. <i>Bioinformatics</i> , 2017, 33, 2379-2380.	1.8	415
9	Drug-target interaction prediction from chemical, genomic and pharmacological data in an integrated framework. <i>Bioinformatics</i> , 2010, 26, i246-i254.	1.8	396
10	Linking Virus Genomes with Host Taxonomy. <i>Viruses</i> , 2016, 8, 66.	1.5	295
11	LIGAND: database of chemical compounds and reactions in biological pathways. <i>Nucleic Acids Research</i> , 2002, 30, 402-404.	6.5	255
12	SIMCOMP/SUBCOMP: chemical structure search servers for network analyses. <i>Nucleic Acids Research</i> , 2010, 38, W652-W656.	6.5	125
13	Environmental Viral Genomes Shed New Light on Virus-Host Interactions in the Ocean. <i>MSphere</i> , 2017, 2, .	1.3	114
14	Drug target prediction using adverse event report systems: a pharmacogenomic approach. <i>Bioinformatics</i> , 2012, 28, i611-i618.	1.8	107
15	The KEGG Databases and Tools Facilitating Omics Analysis: Latest Developments Involving Human Diseases and Pharmaceuticals. <i>Methods in Molecular Biology</i> , 2012, 802, 19-39.	0.4	104
16	Type III Polyketide Synthases: Functional Classification and Phylogenomics. <i>ChemBioChem</i> , 2017, 18, 50-65.	1.3	104
17	DINIES: drug-target interaction network inference engine based on supervised analysis. <i>Nucleic Acids Research</i> , 2014, 42, W39-W45.	6.5	97
18	Taxon Richness of Megaviridae Exceeds those of Bacteria and Archaea in the Ocean. <i>Microbes and Environments</i> , 2018, 33, 162-171.	0.7	83

#	ARTICLE	IF	CITATIONS
19	MAPLE 2.3.0: an improved system for evaluating the functionomes of genomes and metagenomes. <i>Bioscience, Biotechnology and Biochemistry</i> , 2018, 82, 1515-1517.	0.6	80
20	Modular Architecture of Metabolic Pathways Revealed by Conserved Sequences of Reactions. <i>Journal of Chemical Information and Modeling</i> , 2013, 53, 613-622.	2.5	73
21	An automated system for evaluation of the potential functionome: MAPLE version 2.1.0. <i>DNA Research</i> , 2016, 23, 467-475.	1.5	66
22	Evaluation method for the potential functionome harbored in the genome and metagenome. <i>BMC Genomics</i> , 2012, 13, 699.	1.2	65
23	Extraction of phylogenetic network modules from the metabolic network. <i>BMC Bioinformatics</i> , 2006, 7, 130.	1.2	44
24	Effects of post-electrophoretic analysis on variance in gel-based proteomics. <i>Expert Review of Proteomics</i> , 2006, 3, 129-142.	1.3	36
25	EGENES: Transcriptome-Based Plant Database of Genes with Metabolic Pathway Information and Expressed Sequence Tag Indices in KEGG. <i>Plant Physiology</i> , 2007, 144, 857-866.	2.3	35
26	Prediction of missing enzyme genes in a bacterial metabolic network. <i>FEBS Journal</i> , 2007, 274, 2262-2273.	2.2	30
27	KEGG Bioinformatics Resource for Plant Genomics Research. , 2007, 406, 437-458.		27
28	Regulation of metabolic networks by small molecule metabolites. <i>BMC Bioinformatics</i> , 2007, 8, 88.	1.2	23
29	Bioinformatics Strategies for the Analysis of Lipids. , 2009, 580, 339-368.		23
30	varDB: a pathogen-specific sequence database of protein families involved in antigenic variation. <i>Bioinformatics</i> , 2008, 24, 2564-2565.	1.8	18
31	Extraction of phylogenetic network modules from prokaryote metabolic pathways. <i>Genome Informatics</i> , 2004, 15, 249-58.	0.4	12
32	Pharmacoepidemiological characterization of drug-induced adverse reaction clusters towards understanding of their mechanisms. <i>Computational Biology and Chemistry</i> , 2014, 50, 50-59.	1.1	10
33	Discriminating the reaction types of plant type III polyketide synthases. <i>Bioinformatics</i> , 2017, 33, 1937-1943.	1.8	7
34	varDB: A database of antigenic variant sequencesâ€”Current status and future prospects. <i>Acta Tropica</i> , 2010, 114, 144-151.	0.9	5
35	Virus proteins similar to human proteins as possible disturbance on human pathways. <i>Systems and Synthetic Biology</i> , 2014, 8, 283-295.	1.0	2
36	MUCHA: multiple chemical alignment algorithm to identify building block substructures of orphan secondary metabolites. <i>BMC Bioinformatics</i> , 2011, 12, S1.	1.2	1

#	ARTICLE	IF	CITATIONS
37	Revealing phenotype-associated functional differences by genome-wide scan of ancient haplotype blocks. PLoS ONE, 2017, 12, e0176530.	1.1	0