Gajendra P S Raghava

List of Publications by Year in Descending Order

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

14,589 63 296 111 h-index g-index citations papers 18,164 6.99 336 5.5 L-index avg, IF ext. papers ext. citations

#	Paper	IF	Citations
296	IL13Pred: A method for predicting immunoregulatory cytokine IL-13 inducing peptides <i>Computers in Biology and Medicine</i> , 2022 , 143, 105297	7	O
295	In Silico Design of Chemically Modified Cell-Penetrating Peptides. <i>Methods in Molecular Biology</i> , 2022 , 2383, 63-71	1.4	1
294	In silico tools and databases for designing cancer immunotherapy <i>Advances in Protein Chemistry and Structural Biology</i> , 2022 , 129, 1-50	5.3	
293	Tool for Predicting, Scanning, and Designing Defensins. <i>Frontiers in Immunology</i> , 2021 , 12, 780610	8.4	1
292	Prognostic biomarkers for predicting papillary thyroid carcinoma patients at high risk using nine genes of apoptotic pathway. <i>PLoS ONE</i> , 2021 , 16, e0259534	3.7	1
291	Universal and cross-cancer prognostic biomarkers for predicting survival risk of cancer patients from expression profile of apoptotic pathway genes. <i>Proteomics</i> , 2021 , e2000311	4.8	
2 90	AlgPred 2.0: an improved method for predicting allergenic proteins and mapping of IgE epitopes. <i>Briefings in Bioinformatics</i> , 2021 , 22,	13.4	36
289	In-silico identification of subunit vaccine candidates against lung cancer-associated oncogenic viruses. <i>Computers in Biology and Medicine</i> , 2021 , 130, 104215	7	3
288	Crowdsourcing digital health measures to predict Parkinsonß disease severity: the Parkinsonß Disease Digital Biomarker DREAM Challenge. <i>Npj Digital Medicine</i> , 2021 , 4, 53	15.7	8
287	FermFooDb: A database of bioactive peptides derived from fermented foods. <i>Heliyon</i> , 2021 , 7, e06668	3.6	10
286	Computational resources for identification of cancer biomarkers from omics data. <i>Briefings in Functional Genomics</i> , 2021 , 20, 213-222	4.9	1
285	Prognostic Biomarker-Based Identification of Drugs for Managing the Treatment of Endometrial Cancer. <i>Molecular Diagnosis and Therapy</i> , 2021 , 25, 629-646	4.5	3
284	Crowdsourcing assessment of maternal blood multi-omics for predicting gestational age and preterm birth. <i>Cell Reports Medicine</i> , 2021 , 2, 100323	18	5
283	SAPdb: A database of short peptides and the corresponding nanostructures formed by self-assembly. <i>Computers in Biology and Medicine</i> , 2021 , 133, 104391	7	5
282	DenvInD: dengue virus inhibitors database for clinical and molecular research. <i>Briefings in Bioinformatics</i> , 2021 , 22,	13.4	2
281	AntiCP 2.0: an updated model for predicting anticancer peptides. <i>Briefings in Bioinformatics</i> , 2021 , 22,	13.4	45
280	Computer-aided prediction and design of IL-6 inducing peptides: IL-6 plays a crucial role in COVID-19. <i>Briefings in Bioinformatics</i> , 2021 , 22, 936-945	13.4	27

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279	B3Pdb: an archive of blood-brain barrier-penetrating peptides. <i>Brain Structure and Function</i> , 2021 , 226, 2489-2495	4	7
278	B3Pred: A Random-Forest-Based Method for Predicting and Designing Blood-Brain Barrier Penetrating Peptides. <i>Pharmaceutics</i> , 2021 , 13,	6.4	3
277	ChAlPred: A web server for prediction of allergenicity of chemical compounds. <i>Computers in Biology and Medicine</i> , 2021 , 136, 104746	7	3
276	Computational resources in the management of antibiotic resistance: Speeding up drug discovery. <i>Drug Discovery Today</i> , 2021 , 26, 2138-2151	8.8	4
275	Computer-aided prediction of inhibitors against STAT3 for managing COVID-19 associated cytokine storm. <i>Computers in Biology and Medicine</i> , 2021 , 137, 104780	7	3
274	ProCanBio: A Database of Manually Curated Biomarkers for Prostate Cancer <i>Journal of Computational Biology</i> , 2021 , 28, 1248-1257	1.7	
273	A Web-Based Platform on Coronavirus Disease-19 to Maintain Predicted Diagnostic, Drug, and Vaccine Candidates. <i>Monoclonal Antibodies in Immunodiagnosis and Immunotherapy</i> , 2020 , 39, 204-216	1.9	17
272	CancerEnD: A database of cancer associated enhancers. <i>Genomics</i> , 2020 , 112, 3696-3702	4.3	7
271	OvirusTdb: A database of oncolytic viruses for the advancement of therapeutics in cancer. <i>Virology</i> , 2020 , 548, 109-116	3.6	10
270	A Method for Predicting Hemolytic Potency of Chemically Modified Peptides From Its Structure. <i>Frontiers in Pharmacology</i> , 2020 , 11, 54	5.6	11
269	CancerLivER: a database of liver cancer gene expression resources and biomarkers. <i>Database: the Journal of Biological Databases and Curation</i> , 2020 , 2020,	5	7
268	A Hybrid Model for Predicting Pattern Recognition Receptors Using Evolutionary Information. <i>Frontiers in Immunology</i> , 2020 , 11, 71	8.4	7
267	Computing Skin Cutaneous Melanoma Outcome From the HLA-Alleles and Clinical Characteristics. <i>Frontiers in Genetics</i> , 2020 , 11, 221	4.5	14
266	Expression based biomarkers and models to classify early and late-stage samples of Papillary Thyroid Carcinoma. <i>PLoS ONE</i> , 2020 , 15, e0231629	3.7	8
265	Computer-aided designing of oncolytic viruses for overcoming translational challenges of cancer immunotherapy. <i>Drug Discovery Today</i> , 2020 , 25, 1198-1205	8.8	8
264	NAGbinder: An approach for identifying N-acetylglucosamine interacting residues of a protein from its primary sequence. <i>Protein Science</i> , 2020 , 29, 201-210	6.3	10
263	Potential Challenges for Coronavirus (SARS-CoV-2) Vaccines Under Trial. <i>Frontiers in Immunology</i> , 2020 , 11, 561851	8.4	2
262	Identification of prognostic biomarkers for major subtypes of non-small-cell lung cancer using genomic and clinical data. <i>Journal of Cancer Research and Clinical Oncology</i> , 2020 , 146, 2743-2752	4.9	7

261	Risk prediction in cutaneous melanoma patients from their clinico-pathological features: superiority of clinical data over gene expression data. <i>Heliyon</i> , 2020 , 6, e04811	3.6	6
260	Community Assessment of the Predictability of Cancer Protein and Phosphoprotein Levels from Genomics and Transcriptomics. <i>Cell Systems</i> , 2020 , 11, 186-195.e9	10.6	11
259	Analysis and prediction of cholangiocarcinoma from transcriptomic profile of patients. <i>Journal of Hepatology</i> , 2020 , 73, S16-S17	13.4	2
258	Docking-based approach for identification of mutations that disrupt binding between Bcl-2 and Bax proteins: Inducing apoptosis in cancer cells. <i>Molecular Genetics & Camp; Genomic Medicine</i> , 2019 , 7, e910	2.3	16
257	Classification of early and late stage liver hepatocellular carcinoma patients from their genomics and epigenomics profiles. <i>PLoS ONE</i> , 2019 , 14, e0221476	3.7	16
256	Prediction of risk scores for colorectal cancer patients from the concentration of proteins involved in mitochondrial apoptotic pathway. <i>PLoS ONE</i> , 2019 , 14, e0217527	3.7	11
255	HumCFS: a database of fragile sites in human chromosomes. <i>BMC Genomics</i> , 2019 , 19, 985	4.5	26
254	APBioNetß annual International Conference on Bioinformatics (InCoB) returns to India in 2018. BMC Genomics, 2019 , 19, 266	4.5	0
253	NeuroPlpred: a tool to predict, design and scan insect neuropeptides. Scientific Reports, 2019, 9, 5129	4.9	21
252	ccPDB 2.0: an updated version of datasets created and compiled from Protein Data Bank. <i>Database:</i> the Journal of Biological Databases and Curation, 2019 , 2019,	5	5
251	Benchmarking of different molecular docking methods for protein-peptide docking. <i>BMC Bioinformatics</i> , 2019 , 19, 426	3.6	52
250	SAMbinder: A Web Server for Predicting S-Adenosyl-L-Methionine Binding Residues of a Protein From Its Amino Acid Sequence. <i>Frontiers in Pharmacology</i> , 2019 , 10, 1690	5.6	1
249	PRRDB 2.0: a comprehensive database of pattern-recognition receptors and their ligands. <i>Database: the Journal of Biological Databases and Curation</i> , 2019 , 2019,	5	16
248	In Silico Analysis of Gene Expression Change Associated with Copy Number of Enhancers in Pancreatic Adenocarcinoma. <i>International Journal of Molecular Sciences</i> , 2019 , 20,	6.3	5
247	Prediction and Analysis of Skin Cancer Progression using Genomics Profiles of Patients. <i>Scientific Reports</i> , 2019 , 9, 15790	4.9	32
246	RareLSD: a manually curated database of lysosomal enzymes associated with rare diseases. <i>Database: the Journal of Biological Databases and Curation</i> , 2019 , 2019,	5	2
245	Identification of Platform-Independent Diagnostic Biomarker Panel for Hepatocellular Carcinoma Using Large-Scale Transcriptomics Data. <i>Frontiers in Genetics</i> , 2019 , 10, 1306	4.5	18
244	Large expert-curated database for benchmarking document similarity detection in biomedical literature search. <i>Database: the Journal of Biological Databases and Curation</i> , 2019 , 2019,	5	4

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243	Human Opinion Inspired Feature Selection Strategy for Predicting the Pleasantness of a Molecule. <i>Advances in Intelligent Systems and Computing</i> , 2018 , 197-205	0.4	
242	Computer-aided prediction of antigen presenting cell modulators for designing peptide-based vaccine adjuvants. <i>Journal of Translational Medicine</i> , 2018 , 16, 181	8.5	40
241	Approach for Prediction of Antifungal Peptides. Frontiers in Microbiology, 2018, 9, 323	5.7	60
240	Prediction of Cell-Penetrating Potential of Modified Peptides Containing Natural and Chemically Modified Residues. <i>Frontiers in Microbiology</i> , 2018 , 9, 725	5.7	33
239	Prediction of Antitubercular Peptides From Sequence Information Using Ensemble Classifier and Hybrid Features. <i>Frontiers in Pharmacology</i> , 2018 , 9, 954	5.6	39
238	In silico approaches for predicting the half-life of natural and modified peptides in blood. <i>PLoS ONE</i> , 2018 , 13, e0196829	3.7	35
237	In Silico Tools and Databases for Designing Peptide-Based Vaccine and Drugs. <i>Advances in Protein Chemistry and Structural Biology</i> , 2018 , 112, 221-263	5.3	32
236	TopicalPdb: A database of topically delivered peptides. <i>PLoS ONE</i> , 2018 , 13, e0190134	3.7	6
235	Overview of Free Software Developed for Designing Drugs Based on Protein-Small Molecules Interaction. <i>Current Topics in Medicinal Chemistry</i> , 2018 , 18, 1146-1167	3	6
234	Prediction of Antimicrobial Potential of a Chemically Modified Peptide From Its Tertiary Structure. <i>Frontiers in Microbiology</i> , 2018 , 9, 2551	5.7	30
233	AntiTbPdb: a knowledgebase of anti-tubercular peptides. <i>Database: the Journal of Biological Databases and Curation</i> , 2018 , 2018,	5	26
232	A Web Resource for Designing Subunit Vaccine Against Major Pathogenic Species of Bacteria. <i>Frontiers in Immunology</i> , 2018 , 9, 2280	8.4	22
231	Computer-aided designing of immunosuppressive peptides based on IL-10 inducing potential. <i>Scientific Reports</i> , 2017 , 7, 42851	4.9	96
230	Anticancer properties of a defensin like class IId bacteriocin Laterosporulin10. <i>Scientific Reports</i> , 2017 , 7, 46541	4.9	54
229	Assessing therapeutic potential of molecules: molecular property diagnostic suite for tuberculosis ((mathbf{MPDS}^{mathbf{TB}})). <i>Journal of Chemical Sciences</i> , 2017 , 129, 515-531	1.8	12
228	Gene expression-based biomarkers for discriminating early and late stage of clear cell renal cancer. <i>Scientific Reports</i> , 2017 , 7, 44997	4.9	54
227	Novel in silico tools for designing peptide-based subunit vaccines and immunotherapeutics. <i>Briefings in Bioinformatics</i> , 2017 , 18, 467-478	13.4	49
226	CancerPDF: A repository of cancer-associated peptidome found in human biofluids. <i>Scientific Reports</i> , 2017 , 7, 1511	4.9	20

225	In silico analysis to identify vaccine candidates common to multiple serotypes of Shigella and evaluation of their immunogenicity. <i>PLoS ONE</i> , 2017 , 12, e0180505	3.7	12
224	THPdb: Database of FDA-approved peptide and protein therapeutics. <i>PLoS ONE</i> , 2017 , 12, e0181748	3.7	243
223	Computational Prediction of the Immunomodulatory Potential of RNA Sequences. <i>Methods in Molecular Biology</i> , 2017 , 1632, 75-90	1.4	9
222	BIS-CATTLE: A Web Server for Breed Identification using Microsatellite DNA Markers. <i>Current Research in Bioinformatics</i> , 2016 , 5, 10-17		1
221	Topical Delivery of Protein and Peptide Using Novel Cell Penetrating Peptide IMT-P8. <i>Scientific Reports</i> , 2016 , 6, 26278	4.9	39
220	ZikaVR: An Integrated Zika Virus Resource for Genomics, Proteomics, Phylogenetic and Therapeutic Analysis. <i>Scientific Reports</i> , 2016 , 6, 32713	4.9	40
219	Prediction of Immunomodulatory potential of an RNA sequence for designing non-toxic siRNAs and RNA-based vaccine adjuvants. <i>Scientific Reports</i> , 2016 , 6, 20678	4.9	12
218	ApoCanD: Database of human apoptotic proteins in the context of cancer. <i>Scientific Reports</i> , 2016 , 6, 20797	4.9	7
217	A Web Server and Mobile App for Computing Hemolytic Potency of Peptides. <i>Scientific Reports</i> , 2016 , 6, 22843	4.9	78
216	A web-based resource for designing therapeutics against Ebola Virus. <i>Scientific Reports</i> , 2016 , 6, 24782	4.9	12
215	A web server for analysis, comparison and prediction of protein ligand binding sites. <i>Biology Direct</i> , 2016 , 11, 14	7.2	11
214	BLAST-based structural annotation of protein residues using Protein Data Bank. <i>Biology Direct</i> , 2016 , 11, 4	7.2	7
213	CPPsite 2.0: a repository of experimentally validated cell-penetrating peptides. <i>Nucleic Acids Research</i> , 2016 , 44, D1098-103	20.1	167
212	SATPdb: a database of structurally annotated therapeutic peptides. <i>Nucleic Acids Research</i> , 2016 , 44, D1119-26	20.1	89
211	Cell-penetrating peptide and antibiotic combination therapy: a potential alternative to combat drug resistance in methicillin-resistant Staphylococcus aureus. <i>Applied Microbiology and Biotechnology</i> , 2016 , 100, 4073-83	5.7	13
210	Managing Drug Resistance in Cancer: Role of Cancer Informatics. <i>Methods in Molecular Biology</i> , 2016 , 1395, 299-312	1.4	9
209	A Web-Based Platform for Designing Vaccines against Existing and Emerging Strains of Mycobacterium tuberculosis. <i>PLoS ONE</i> , 2016 , 11, e0153771	3.7	15
208	A Platform for Designing Genome-Based Personalized Immunotherapy or Vaccine against Cancer. <i>PLoS ONE</i> , 2016 , 11, e0166372	3.7	11

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207	Prioritization of anticancer drugs against a cancer using genomic features of cancer cells: A step towards personalized medicine. <i>Scientific Reports</i> , 2016 , 6, 23857	4.9	38	
206	PEPlife: A Repository of the Half-life of Peptides. <i>Scientific Reports</i> , 2016 , 6, 36617	4.9	76	
205	dbEM: A database of epigenetic modifiers curated from cancerous and normal genomes. <i>Scientific Reports</i> , 2016 , 6, 19340	4.9	34	
204	ProCarDB: a database of bacterial carotenoids. <i>BMC Microbiology</i> , 2016 , 16, 96	4.5	41	
203	Prediction of anticancer molecules using hybrid model developed on molecules screened against NCI-60 cancer cell lines. <i>BMC Cancer</i> , 2016 , 16, 77	4.8	30	
202	ETurn: a novel Eturn mimic in globular proteins stabilized by main-chain to side-chain CHIIIO interaction. <i>Proteins: Structure, Function and Bioinformatics</i> , 2015 , 83, 203-14	4.2	6	
201	In silico platform for predicting and initiating Eturns in a protein at desired locations. <i>Proteins: Structure, Function and Bioinformatics</i> , 2015 , 83, 910-21	4.2	9	
200	QSAR based model for discriminating EGFR inhibitors and non-inhibitors using Random forest. <i>Biology Direct</i> , 2015 , 10, 10	7.2	45	
199	Designing B-Cell Epitopes for Immunotherapy and Subunit Vaccines. <i>Methods in Molecular Biology</i> , 2015 , 1348, 327-40	1.4	1	
198	Identification and characterization of novel protein-derived arginine-rich cell-penetrating peptides. <i>European Journal of Pharmaceutics and Biopharmaceutics</i> , 2015 , 89, 93-106	5.7	27	
197	An in silico platform for predicting, screening and designing of antihypertensive peptides. <i>Scientific Reports</i> , 2015 , 5, 12512	4.9	80	
196	VaccineDA: Prediction, design and genome-wide screening of oligodeoxynucleotide-based vaccine adjuvants. <i>Scientific Reports</i> , 2015 , 5, 12478	4.9	20	
195	AHTPDB: a comprehensive platform for analysis and presentation of antihypertensive peptides. <i>Nucleic Acids Research</i> , 2015 , 43, D956-62	20.1	96	
194	PEPstrMOD: structure prediction of peptides containing natural, non-natural and modified residues. <i>Biology Direct</i> , 2015 , 10, 73	7.2	100	
193	CancerPPD: a database of anticancer peptides and proteins. <i>Nucleic Acids Research</i> , 2015 , 43, D837-43	20.1	169	
192	Assessment of SYBR green I dye-based fluorescence assay for screening antimalarial activity of cationic peptides and DNA intercalating agents. <i>Antimicrobial Agents and Chemotherapy</i> , 2015 , 59, 2886	5- 5 -9	3	
191	Identification of protein-interacting nucleotides in a RNA sequence using composition profile of tri-nucleotides. <i>Genomics</i> , 2015 , 105, 197-203	4.3	17	
190	AntiAngioPred: A Server for Prediction of Anti-Angiogenic Peptides. <i>PLoS ONE</i> , 2015 , 10, e0136990	3.7	33	

189	In Silico Designing and Screening of Antagonists against Cancer Drug Target XIAP. <i>Current Cancer Drug Targets</i> , 2015 , 15, 836-46	2.8	1
188	Peptide toxicity prediction. <i>Methods in Molecular Biology</i> , 2015 , 1268, 143-57	1.4	39
187	Computer-Aided Virtual Screening and Designing of Cell-Penetrating Peptides. <i>Methods in Molecular Biology</i> , 2015 , 1324, 59-69	1.4	35
186	Designing of promiscuous inhibitors against pancreatic cancer cell lines. Scientific Reports, 2014, 4, 466	84.9	12
185	Prediction and classification of ncRNAs using structural information. <i>BMC Genomics</i> , 2014 , 15, 127	4.5	63
184	Designing of peptides with desired half-life in intestine-like environment. <i>BMC Bioinformatics</i> , 2014 , 15, 282	3.6	43
183	PCMdb: pancreatic cancer methylation database. <i>Scientific Reports</i> , 2014 , 4, 4197	4.9	19
182	Herceptin resistance database for understanding mechanism of resistance in breast cancer patients. <i>Scientific Reports</i> , 2014 , 4, 4483	4.9	33
181	QSAR-based models for designing quinazoline/imidazothiazoles/pyrazolopyrimidines based inhibitors against wild and mutant EGFR. <i>PLoS ONE</i> , 2014 , 9, e101079	3.7	27
180	Prediction of uridine modifications in tRNA sequences. <i>BMC Bioinformatics</i> , 2014 , 15, 326	3.6	21
179	ParaPep: a web resource for experimentally validated antiparasitic peptide sequences and their structures. <i>Database: the Journal of Biological Databases and Curation</i> , 2014 , 2014,	5	42
178	Hemolytik: a database of experimentally determined hemolytic and non-hemolytic peptides. <i>Nucleic Acids Research</i> , 2014 , 42, D444-9	20.1	85
177	ToxiPred: A Server for Prediction of Aqueous Toxicity of Small Chemical Molecules in T. Pyriformis 2014 , 1, 21-27		7
176	Evaluation of protein dihedral angle prediction methods. <i>PLoS ONE</i> , 2014 , 9, e105667	3.7	14
175	Tumor homing peptides as molecular probes for cancer therapeutics, diagnostics and theranostics. <i>Current Medicinal Chemistry</i> , 2014 , 21, 2367-91	4.3	60
174	EGFRIndb: epidermal growth factor receptor inhibitor database. <i>Anti-Cancer Agents in Medicinal Chemistry</i> , 2014 , 14, 928-35	2.2	13
173	In silico approaches for designing highly effective cell penetrating peptides. <i>Journal of Translational Medicine</i> , 2013 , 11, 74	8.5	170
172	Designing of inhibitors against drug tolerant Mycobacterium tuberculosis (H37Rv). <i>Chemistry Central Journal</i> , 2013 , 7, 49		17

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171	Prediction of vitamin interacting residues in a vitamin binding protein using evolutionary information. <i>BMC Bioinformatics</i> , 2013 , 14, 44	3.6	18
170	PHDcleav: a SVM based method for predicting human Dicer cleavage sites using sequence and secondary structure of miRNA precursors. <i>BMC Bioinformatics</i> , 2013 , 14 Suppl 14, S9	3.6	31
169	Development of a model webserver for breed identification using microsatellite DNA marker. <i>BMC Genetics</i> , 2013 , 14, 118	2.6	5
168	Designing of interferon-gamma inducing MHC class-II binders. <i>Biology Direct</i> , 2013 , 8, 30	7.2	265
167	DrugMint: a webserver for predicting and designing of drug-like molecules. <i>Biology Direct</i> , 2013 , 8, 28	7.2	34
166	Computational approach for designing tumor homing peptides. Scientific Reports, 2013, 3, 1607	4.9	48
165	NPACT: Naturally Occurring Plant-based Anti-cancer Compound-Activity-Target database. <i>Nucleic Acids Research</i> , 2013 , 41, D1124-9	20.1	170
164	In silico models for B-cell epitope recognition and signaling. <i>Methods in Molecular Biology</i> , 2013 , 993, 129-38	1.4	7
163	Genome Sequence of the "Indian Bison Type" Biotype of Mycobacterium avium subsp. paratuberculosis Strain S5. <i>Genome Announcements</i> , 2013 , 1,		9
162	Draft Genome Sequence of the Type Species of the Genus Citrobacter, Citrobacter freundii MTCC 1658. <i>Genome Announcements</i> , 2013 , 1,		10
161	Draft Genome Sequence of the 2-Chloro-4-Nitrophenol-Degrading Bacterium Arthrobacter sp. Strain SJCon. <i>Genome Announcements</i> , 2013 , 1, e0005813		10
160	Draft Genome Sequence of Rhodococcus qingshengii Strain BKS 20-40. <i>Genome Announcements</i> , 2013 , 1, e0012813		2
159	Draft Genome Sequence of Rhodococcus triatomae Strain BKS 15-14. <i>Genome Announcements</i> , 2013 , 1, e0012913		3
158	Draft Genome Sequence of Acinetobacter baumannii Strain MSP4-16. <i>Genome Announcements</i> , 2013 , 1, e0013713		1
157	Draft Genome Sequence of Amycolatopsis decaplanina Strain DSM 44594T. <i>Genome Announcements</i> , 2013 , 1, e0013813		3
156	Draft Genome Sequence of Rhodococcus ruber Strain BKS 20-38. <i>Genome Announcements</i> , 2013 , 1, e00	13913	0
155	Draft Genome Sequence of Streptomyces gancidicus Strain BKS 13-15. <i>Genome Announcements</i> , 2013 , 1, e0015013		3
154	CancerDR: cancer drug resistance database. <i>Scientific Reports</i> , 2013 , 3, 1445	4.9	79

153	In silico models for designing and discovering novel anticancer peptides. Scientific Reports, 2013, 3, 29	844.9	149
152	lncRNome: a comprehensive knowledgebase of human long noncoding RNAs. <i>Database: the Journal of Biological Databases and Curation</i> , 2013 , 2013, bat034	5	110
151	Identification of B-cell epitopes in an antigen for inducing specific class of antibodies. <i>Biology Direct</i> , 2013 , 8, 27	7.2	59
150	In silico approach for predicting toxicity of peptides and proteins. <i>PLoS ONE</i> , 2013 , 8, e73957	3.7	600
149	Improved method for linear B-cell epitope prediction using antigen® primary sequence. <i>PLoS ONE</i> , 2013 , 8, e62216	3.7	171
148	Genome annotation of Burkholderia sp. SJ98 with special focus on chemotaxis genes. <i>PLoS ONE</i> , 2013 , 8, e70624	3.7	7
147	Genes involved in degradation of para-nitrophenol are differentially arranged in form of non-contiguous gene clusters in Burkholderia sp. strain SJ98. <i>PLoS ONE</i> , 2013 , 8, e84766	3.7	19
146	Prediction of IL4 inducing peptides. Clinical and Developmental Immunology, 2013, 2013, 263952		111
145	Open source software and web services for designing therapeutic molecules. <i>Current Topics in Medicinal Chemistry</i> , 2013 , 13, 1172-91	3	17
144	Hybrid approach for predicting coreceptor used by HIV-1 from its V3 loop amino acid sequence. <i>PLoS ONE</i> , 2013 , 8, e61437	3.7	16
143	In silico platform for prediction of N-, O- and C-glycosites in eukaryotic protein sequences. <i>PLoS ONE</i> , 2013 , 8, e67008	3.7	142
142	On the Development of Vaccine Antigen Databases: Progress, Opportunity, and Challenge 2013 , 117-1	130	
141	ProGlycProt: a repository of experimentally characterized prokaryotic glycoproteins. <i>Nucleic Acids Research</i> , 2012 , 40, D388-93	20.1	22
140	Branching of the p-nitrophenol (PNP) degradation pathway in burkholderia sp. Strain SJ98: Evidences from genetic characterization of PNP gene cluster. <i>AMB Express</i> , 2012 , 2, 30	4.1	23
139	PolysacDB: a database of microbial polysaccharide antigens and their antibodies. <i>PLoS ONE</i> , 2012 , 7, e34613	3.7	8
138	GlycoPP: a webserver for prediction of N- and O-glycosites in prokaryotic protein sequences. <i>PLoS ONE</i> , 2012 , 7, e40155	3.7	44
137	Predicting sub-cellular localization of tRNA synthetases from their primary structures. <i>Amino Acids</i> , 2012 , 42, 1703-13	3.5	7
136	Draft genome sequence of salt-tolerant yeast Debaryomyces hansenii var. hansenii MTCC 234. <i>Eukaryotic Cell</i> , 2012 , 11, 961-2		15

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135	CPPsite: a curated database of cell penetrating peptides. <i>Database: the Journal of Biological Databases and Curation</i> , 2012 , 2012, bas015	5	122
134	Genome sequence of the halotolerant bacterium Imtechella halotolerans K1T. <i>Journal of Bacteriology</i> , 2012 , 194, 3731	3.5	3
133	Genome sequence of the marine bacterium Marinilabilia salmonicolor JCM 21150T. <i>Journal of Bacteriology</i> , 2012 , 194, 3746	3.5	5
132	ccPDB: compilation and creation of data sets from Protein Data Bank. <i>Nucleic Acids Research</i> , 2012 , 40, D486-9	20.1	21
131	Genome sequence of the oleaginous red yeast Rhodosporidium toruloides MTCC 457. <i>Eukaryotic Cell</i> , 2012 , 11, 1083-4		51
130	Draft genome sequence of the nitrophenol-degrading actinomycete Rhodococcus imtechensis RKJ300. <i>Journal of Bacteriology</i> , 2012 , 194, 3543	3.5	11
129	Genome sequence of the nitroaromatic compound-degrading Bacterium Burkholderia sp. strain SJ98. <i>Journal of Bacteriology</i> , 2012 , 194, 3286	3.5	14
128	TumorHoPe: a database of tumor homing peptides. <i>PLoS ONE</i> , 2012 , 7, e35187	3.7	80
127	Open source drug discoverya new paradigm of collaborative research in tuberculosis drug development. <i>Tuberculosis</i> , 2011 , 91, 479-86	2.6	32
126	A web server for predicting inhibitors against bacterial target GlmU protein. <i>BMC Pharmacology</i> , 2011 , 11, 5		23
125	Analysis and prediction of cancerlectins using evolutionary and domain information. <i>BMC Research Notes</i> , 2011 , 4, 237	2.3	17
124	SVM based prediction of RNA-binding proteins using binding residues and evolutionary information. <i>Journal of Molecular Recognition</i> , 2011 , 24, 303-13	2.6	88
123	CCDB: a curated database of genes involved in cervix cancer. <i>Nucleic Acids Research</i> , 2011 , 39, D975-9	20.1	59
122	Designing of highly effective complementary and mismatch siRNAs for silencing a gene. <i>PLoS ONE</i> , 2011 , 6, e23443	3.7	36
121	Identification of mannose interacting residues using local composition. <i>PLoS ONE</i> , 2011 , 6, e24039	3.7	22
120	HIVsirDB: a database of HIV inhibiting siRNAs. <i>PLoS ONE</i> , 2011 , 6, e25917	3.7	34
119	Prediction of Specificity and Cross-Reactivity of Kinase Inhibitors. <i>Letters in Drug Design and Discovery</i> , 2011 , 8, 223-228	0.8	14
118	Bridging innate and adaptive antitumor immunity targeting glycans. <i>Journal of Biomedicine and Biotechnology</i> , 2010 , 2010, 354068		14

117	Identification of conformational B-cell Epitopes in an antigen from its primary sequence. Immunome Research, 2010 , 6, 6	186
116	AntigenDB: an immunoinformatics database of pathogen antigens. <i>Nucleic Acids Research</i> , 2010 , 38, D84 3 e5£	3 48
115	Prediction and classification of aminoacyl tRNA synthetases using PROSITE domains. <i>BMC Genomics</i> , 2010 , 11, 507	12
114	A simple approach for predicting protein-protein interactions. <i>Current Protein and Peptide Science</i> , 2.8	32
113	Prediction of mitochondrial proteins of malaria parasite using split amino acid composition and PSSM profile. <i>Amino Acids</i> , 2010 , 39, 101-10	47
112	KiDoQ: using docking based energy scores to develop ligand based model for predicting antibacterials. <i>BMC Bioinformatics</i> , 2010 , 11, 125	24
111	Identification of NAD interacting residues in proteins. <i>BMC Bioinformatics</i> , 2010 , 11, 160 3.6	50
110	Prediction of GTP interacting residues, dipeptides and tripeptides in a protein from its evolutionary information. <i>BMC Bioinformatics</i> , 2010 , 11, 301	35
109	AntiBP2: improved version of antibacterial peptide prediction. <i>BMC Bioinformatics</i> , 2010 , 11 Suppl 1, S19	132
108	Prediction of FAD interacting residues in a protein from its primary sequence using evolutionary information. <i>BMC Bioinformatics</i> , 2010 , 11 Suppl 1, S48	24
107	Virtual Screening of potential drug-like inhibitors against Lysine/DAP pathway of Mycobacterium tuberculosis. <i>BMC Bioinformatics</i> , 2010 , 11 Suppl 1, S53	16
106	BIAdb: a curated database of benzylisoquinoline alkaloids. <i>BMC Pharmacology</i> , 2010 , 10, 4	34
105	Prediction of cytochrome P450 isoform responsible for metabolizing a drug molecule. <i>BMC Pharmacology</i> , 2010 , 10, 8	49
104	Prediction and classification of chemokines and their receptors. <i>Protein Engineering, Design and Selection</i> , 2009 , 22, 441-4	7
103	Prediction of guide strand of microRNAs from its sequence and secondary structure. <i>BMC Bioinformatics</i> , 2009 , 10, 105	37
102	Identification of ATP binding residues of a protein from its primary sequence. <i>BMC Bioinformatics</i> , 2009 , 10, 434	82
101	Hmrbase: a database of hormones and their receptors. <i>BMC Genomics</i> , 2009 , 10, 307 4.5	22
100	MHCBN 4.0: A database of MHC/TAP binding peptides and T-cell epitopes. <i>BMC Research Notes</i> , 2.3	63

99	OxDBase: a database of oxygenases involved in biodegradation. BMC Research Notes, 2009, 2, 67	2.3	78
98	RSLpred: an integrative system for predicting subcellular localization of rice proteins combining compositional and evolutionary information. <i>Proteomics</i> , 2009 , 9, 2324-42	4.8	35
97	Databases and Web-Based Tools for Innate Immunity 2009 , 67-76		
96	Prediction of nuclear proteins using SVM and HMM models. BMC Bioinformatics, 2009, 10, 22	3.6	30
95	Prediction of Polyadenylation Signals in Human DNA Sequences using Nucleotide Frequencies. <i>In Silico Biology</i> , 2009 , 9, 135-148	2	37
94	Prediction of polyadenylation signals in human DNA sequences using nucleotide frequencies. <i>In Silico Biology</i> , 2009 , 9, 135-48	2	22
93	Identification of proteins secreted by malaria parasite into erythrocyte using SVM and PSSM profiles. <i>BMC Bioinformatics</i> , 2008 , 9, 201	3.6	29
92	ESLpred2: improved method for predicting subcellular localization of eukaryotic proteins. <i>BMC Bioinformatics</i> , 2008 , 9, 503	3.6	41
91	PRRDB: a comprehensive database of pattern-recognition receptors and their ligands. <i>BMC Genomics</i> , 2008 , 9, 180	4.5	27
90	Computer-aided biotechnology: from immuno-informatics to reverse vaccinology. <i>Trends in Biotechnology</i> , 2008 , 26, 190-200	15.1	79
89	CytoPred: a server for prediction and classification of cytokines. <i>Protein Engineering, Design and Selection</i> , 2008 , 21, 279-82	1.9	14
88	CyclinPred: a SVM-based method for predicting cyclin protein sequences. <i>PLoS ONE</i> , 2008 , 3, e2605	3.7	22
87	DPROT: prediction of disordered proteins using evolutionary information. Amino Acids, 2008, 35, 599-6	0 <u>5</u> .5	10
86	Prediction of RNA binding sites in a protein using SVM and PSSM profile. <i>Proteins: Structure, Function and Bioinformatics</i> , 2008 , 71, 189-94	4.2	212
85	AC2DGel: Analysis and Comparison of 2D Gels. <i>Journal of Proteomics and Bioinformatics</i> , 2008 , 01, 043-0	0 4 61	2
84	COPid: composition based protein identification. <i>In Silico Biology</i> , 2008 , 8, 121-8	2	31
83	A machine learning based method for the prediction of secretory proteins using amino acid composition, their order and similarity-search. <i>In Silico Biology</i> , 2008 , 8, 129-40	2	28
82	Towards a consensus on datasets and evaluation metrics for developing B-cell epitope prediction tools. <i>Journal of Molecular Recognition</i> , 2007 , 20, 75-82	2.6	171

81	Analysis and prediction of antibacterial peptides. <i>BMC Bioinformatics</i> , 2007 , 8, 263	3.6	181
80	Support Vector Machine-based method for predicting subcellular localization of mycobacterial proteins using evolutionary information and motifs. <i>BMC Bioinformatics</i> , 2007 , 8, 337	3.6	104
79	Identification of DNA-binding proteins using support vector machines and evolutionary profiles. <i>BMC Bioinformatics</i> , 2007 , 8, 463	3.6	174
78	A hybrid approach for predicting promiscuous MHC class I restricted T cell epitopes. <i>Journal of Biosciences</i> , 2007 , 32, 31-42	2.3	98
77	PEPstr: a de novo method for tertiary structure prediction of small bioactive peptides. <i>Protein and Peptide Letters</i> , 2007 , 14, 626-31	1.9	152
76	Support vector machine based prediction of glutathione S-transferase proteins. <i>Protein and Peptide Letters</i> , 2007 , 14, 575-80	1.9	12
75	Oxypred: prediction and classification of oxygen-binding proteins. <i>Genomics, Proteomics and Bioinformatics</i> , 2007 , 5, 250-2	6.5	4
74	TAPPred prediction of TAP-binding peptides in antigens. <i>Methods in Molecular Biology</i> , 2007 , 409, 381-6	5 1.4	31
73	Prediction methods for B-cell epitopes. <i>Methods in Molecular Biology</i> , 2007 , 409, 387-94	1.4	56
72	Searching and mapping of T-cell epitopes, MHC binders, and TAP binders. <i>Methods in Molecular Biology</i> , 2007 , 409, 95-112	1.4	2
71	Searching haptens, carrier proteins, and anti-hapten antibodies. <i>Methods in Molecular Biology</i> , 2007 , 409, 125-39	1.4	1
70	Application of machine learning techniques in predicting MHC binders. <i>Methods in Molecular Biology</i> , 2007 , 409, 201-15	1.4	40
69	Predicting virulence factors of immunological interest. <i>Methods in Molecular Biology</i> , 2007 , 409, 407-15	1.4	1
68	Searching and mapping of B-cell epitopes in Bcipep database. <i>Methods in Molecular Biology</i> , 2007 , 409, 113-24	1.4	11
67	Prediction of neurotoxins based on their function and source. <i>In Silico Biology</i> , 2007 , 7, 369-87	2	16
66	BTXpred: prediction of bacterial toxins. <i>In Silico Biology</i> , 2007 , 7, 405-12	2	32
65	Quantification of the variation in percentage identity for protein sequence alignments. <i>BMC Bioinformatics</i> , 2006 , 7, 415	3.6	32
64	Machine learning techniques in disease forecasting: a case study on rice blast prediction. <i>BMC Bioinformatics</i> , 2006 , 7, 485	3.6	89

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63	AlgPred: prediction of allergenic proteins and mapping of IgE epitopes. <i>Nucleic Acids Research</i> , 2006 , 34, W202-9	20.1	440
62	HaptenDB: a comprehensive database of haptens, carrier proteins and anti-hapten antibodies. <i>Bioinformatics</i> , 2006 , 22, 253-5	7.2	39
61	VICMpred: an SVM-based method for the prediction of functional proteins of Gram-negative bacteria using amino acid patterns and composition. <i>Genomics, Proteomics and Bioinformatics</i> , 2006 , 4, 42-7	6.5	68
60	VGIchan: prediction and classification of voltage-gated ion channels. <i>Genomics, Proteomics and Bioinformatics</i> , 2006 , 4, 253-8	6.5	21
59	Prediction of mitochondrial proteins using support vector machine and hidden Markov model. Journal of Biological Chemistry, 2006 , 281, 5357-63	5.4	80
58	Prediction of continuous B-cell epitopes in an antigen using recurrent neural network. <i>Proteins:</i> Structure, Function and Bioinformatics, 2006 , 65, 40-8	4.2	761
57	Prediction of subcellular localization of proteins using pairwise sequence alignment and support vector machine. <i>Pattern Recognition Letters</i> , 2006 , 27, 996-1001	4.7	23
56	GPCRsclass: a web tool for the classification of amine type of G-protein-coupled receptors. <i>Nucleic Acids Research</i> , 2005 , 33, W143-7	20.1	52
55	BhairPred: prediction of beta-hairpins in a protein from multiple alignment information using ANN and SVM techniques. <i>Nucleic Acids Research</i> , 2005 , 33, W154-9	20.1	53
54	FASTA Servers for Sequence Similarity Search 2005 , 503-525		1
53	Pcleavage: an SVM based method for prediction of constitutive proteasome and immunoproteasome cleavage sites in antigenic sequences. <i>Nucleic Acids Research</i> , 2005 , 33, W202-7	20.1	71
52	Correlation and prediction of gene expression level from amino acid and dipeptide composition of its protein. <i>BMC Bioinformatics</i> , 2005 , 6, 59	3.6	38
51	Bcipep: a database of B-cell epitopes. <i>BMC Genomics</i> , 2005 , 6, 79	4.5	145
50	Real value prediction of solvent accessibility in proteins using multiple sequence alignment and secondary structure. <i>Proteins: Structure, Function and Bioinformatics</i> , 2005 , 61, 318-24	4.2	73
49	PSLpred: prediction of subcellular localization of bacterial proteins. <i>Bioinformatics</i> , 2005 , 21, 2522-4	7.2	161
48	Support vector machine-based method for subcellular localization of human proteins using amino acid compositions, their order, and similarity search. <i>Journal of Biological Chemistry</i> , 2005 , 280, 14427-3	32 ^{5.4}	169
47	ESLpred: SVM-based method for subcellular localization of eukaryotic proteins using dipeptide composition and PSI-BLAST. <i>Nucleic Acids Research</i> , 2004 , 32, W414-9	20.1	233
46	EGPred: prediction of eukaryotic genes using ab initio methods after combining with sequence similarity approaches. <i>Genome Research</i> , 2004 , 14, 1756-66	9.7	11

45	Classification of nuclear receptors based on amino acid composition and dipeptide composition. Journal of Biological Chemistry, 2004 , 279, 23262-6	5.4	145
44	Spectral Repeat Finder (SRF): identification of repetitive sequences using Fourier transformation. <i>Bioinformatics</i> , 2004 , 20, 1405-12	7.2	116
43	GPCRpred: an SVM-based method for prediction of families and subfamilies of G-protein coupled receptors. <i>Nucleic Acids Research</i> , 2004 , 32, W383-9	20.1	129
42	A neural network method for prediction of beta-turn types in proteins using evolutionary information. <i>Bioinformatics</i> , 2004 , 20, 2751-8	7.2	80
41	Prediction of transmembrane regions of beta-barrel proteins using ANN- and SVM-based methods. <i>Proteins: Structure, Function and Bioinformatics</i> , 2004 , 56, 11-8	4.2	85
40	Prediction of alpha-turns in proteins using PSI-BLAST profiles and secondary structure information. <i>Proteins: Structure, Function and Bioinformatics</i> , 2004 , 55, 83-90	4.2	44
39	SVM based method for predicting HLA-DRB1*0401 binding peptides in an antigen sequence. <i>Bioinformatics</i> , 2004 , 20, 421-3	7.2	128
38	Analysis and prediction of affinity of TAP binding peptides using cascade SVM. <i>Protein Science</i> , 2004 , 13, 596-607	6.3	114
37	BcePred: Prediction of Continuous B-Cell Epitopes in Antigenic Sequences Using Physico-chemical Properties. <i>Lecture Notes in Computer Science</i> , 2004 , 197-204	0.9	147
36	Prediction of CTL epitopes using QM, SVM and ANN techniques. <i>Vaccine</i> , 2004 , 22, 3195-204	4.1	242
35	Role of evolutionary information in prediction of aromatic-backbone NH interactions in proteins. <i>FEBS Letters</i> , 2004 , 564, 47-57	3.8	8
34	MHCBN: a comprehensive database of MHC binding and non-binding peptides. <i>Bioinformatics</i> , 2003 , 19, 665-6	7.2	167
33	Prediction of beta-turns in proteins from multiple alignment using neural network. <i>Protein Science</i> , 2003 , 12, 627-34	6.3	117
32	A neural-network based method for prediction of gamma-turns in proteins from multiple sequence alignment. <i>Protein Science</i> , 2003 , 12, 923-9	6.3	61
31	OXBench: a benchmark for evaluation of protein multiple sequence alignment accuracy. <i>BMC Bioinformatics</i> , 2003 , 4, 47	3.6	139
30	ProPred1: prediction of promiscuous MHC Class-I binding sites. <i>Bioinformatics</i> , 2003 , 19, 1009-14	7.2	348
29	Intrinsic contributions of polar amino acid residues toward thermal stability of an ABC-ATPase of mesophilic origin. <i>Protein Science</i> , 2003 , 12, 2118-20	6.3	5
28	Prediction of promiscuous and high-affinity mutated MHC binders. <i>Hybridoma</i> , 2003 , 22, 229-34		37

27	BTEVAL: a server for evaluation of beta-turn prediction methods. <i>Journal of Bioinformatics and Computational Biology</i> , 2003 , 1, 495-504	1	3
26	GWFASTA: server for FASTA search in eukaryotic and microbial genomes. <i>BioTechniques</i> , 2002 , 33, 548-50, 552, 554-6	2.5	6
25	Locating probable genes using Fourier transform approach. <i>Bioinformatics</i> , 2002 , 18, 196-7	7.2	45
24	An evaluation of beta-turn prediction methods. <i>Bioinformatics</i> , 2002 , 18, 1508-14	7.2	41
23	BetaTPred: prediction of beta-TURNS in a protein using statistical algorithms. <i>Bioinformatics</i> , 2002 , 18, 498-9	7.2	21
22	Detection of Orientation of MHC Class II Binding Peptides Using Bioinformatics Tools. <i>Biotech Software & Internet Report</i> , 2002 , 3, 146-150		1
21	GWFASTA: Server for FASTA Search in Eukaryotic and Microbial Genomes. <i>BioTechniques</i> , 2002 , 33, 548	8-5 5 5	
20	PDSB: Public Domain Software in Biology. <i>Biotech Software & Internet Report</i> , 2001 , 2, 154-156		2
19	ProPred: prediction of HLA-DR binding sites. <i>Bioinformatics</i> , 2001 , 17, 1236-7	7.2	667
18	A Graphical Web Server for the Analysis of Protein Sequences and Alignment. <i>Biotech Software & Internet Report</i> , 2001 , 2, 254-257		5
17	PDWSB: Public Domain Web Servers in Biology. <i>Biotech Software & Internet Report</i> , 2001 , 2, 152-153		1
16	Fingerprinting method for phylogenetic classification and identification of microorganisms based on variation in 16S rRNA gene sequences. <i>BioTechniques</i> , 2000 , 29, 108-12, 114-6	2.5	14
15	Method for determining the affinity of monoclonal antibody using non-competitive ELISA: a computer program. <i>Journal of Immunoassay</i> , 1994 , 15, 115-28		23
14	Measurement and computation of murine interleukin-4 and interferon-gamma by exploiting the unique abilities of these lymphokines to induce the secretion of IgG1 and IgG2a. <i>Journal of Immunoassay</i> , 1993 , 14, 83-97		6
13	Computational resources in healthcare. Wiley Interdisciplinary Reviews: Data Mining and Knowledge Discovery,e1437	6.9	
12	A web-based platform on COVID-19 to maintain Predicted Diagnostic, Drug and Vaccine candidates		3
11	Prediction and analysis of skin cancer progression using genomics profiles of patients		1
10	AntiCP 2.0: An updated model for predicting anticancer peptides		2

9	Evaluation of protein-ligand docking methods on peptide-ligand complexes for docking small ligands to peptides	5
8	HumCFS: A database of fragile sites in human chromosomes	1
7	A web bench for analysis and prediction of oncological status from proteomics data of urine samples	1
6	Expression based biomarkers and models to classify early and late stage samples of Papillary Thyroid Carcinoma	2
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3	SAMbinder: A web server for predicting SAM binding residues of a protein from its amino acid sequence	2
2	GPSRdocker: A Docker-based Resource for Genomics, Proteomics and Systems biology	4
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