

Chris Bailey-Kellogg

List of Publications by Year in descending order

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Version: 2024-02-01

109
papers

3,099
citations

159585

30
h-index

197818

49
g-index

116
all docs

116
docs citations

116
times ranked

3694
citing authors

#	ARTICLE	IF	CITATIONS
1	Polyfunctional HIV-Specific Antibody Responses Are Associated with Spontaneous HIV Control. <i>PLoS Pathogens</i> , 2016, 12, e1005315.	4.7	220
2	High-throughput, multiplexed IgG subclassing of antigen-specific antibodies from clinical samples. <i>Journal of Immunological Methods</i> , 2012, 386, 117-123.	1.4	197
3	Adjuvant-dependent innate and adaptive immune signatures of risk of SIVmac251 acquisition. <i>Nature Medicine</i> , 2016, 22, 762-770.	30.7	197
4	Multiplexed Fc array for evaluation of antigen-specific antibody effector profiles. <i>Journal of Immunological Methods</i> , 2017, 443, 33-44.	1.4	158
5	Pentavalent HIV-1 vaccine protects against simian-human immunodeficiency virus challenge. <i>Nature Communications</i> , 2017, 8, 15711.	12.8	137
6	Route of immunization defines multiple mechanisms of vaccine-mediated protection against SIV. <i>Nature Medicine</i> , 2018, 24, 1590-1598.	30.7	129
7	The two-faced T cell epitope. <i>Human Vaccines and Immunotherapeutics</i> , 2013, 9, 1577-1586.	3.3	88
8	The NOESY Jigsaw: Automated Protein Secondary Structure and Main-Chain Assignment from Sparse, Unassigned NMR Data. <i>Journal of Computational Biology</i> , 2000, 7, 537-558.	1.6	85
9	Learning context-aware structural representations to predict antigen and antibody binding interfaces. <i>Bioinformatics</i> , 2020, 36, 3996-4003.	4.1	79
10	Protein interaction interface region prediction by geometric deep learning. <i>Bioinformatics</i> , 2021, 37, 2580-2588.	4.1	67
11	Graphical Models of Residue Coupling in Protein Families. <i>IEEE/ACM Transactions on Computational Biology and Bioinformatics</i> , 2008, 5, 183-197.	3.0	59
12	Antibody humanization by structure-based computational protein design. <i>MAbs</i> , 2015, 7, 1045-1057.	5.2	58
13	Structure-Guided Deimmunization of Therapeutic Proteins. <i>Journal of Computational Biology</i> , 2013, 20, 152-165.	1.6	54
14	Machine Learning Methods Enable Predictive Modeling of Antibody Feature:Function Relationships in RV144 Vaccines. <i>PLoS Computational Biology</i> , 2015, 11, e1004185.	3.2	50
15	Accounting for conformational entropy in predicting binding free energies of protein-protein interactions. <i>Proteins: Structure, Function and Bioinformatics</i> , 2011, 79, 444-462.	2.6	49
16	Optimization algorithms for functional deimmunization of therapeutic proteins. <i>BMC Bioinformatics</i> , 2010, 11, 180.	2.6	48
17	Depletion of T Cell Epitopes in Lysostaphin Mitigates Anti-Drug Antibody Response and Enhances Antibacterial Efficacy In Vivo. <i>Chemistry and Biology</i> , 2015, 22, 629-639.	6.0	48
18	CHOPPI: A web tool for the analysis of immunogenicity risk from host cell proteins in CHO-based protein production. <i>Biotechnology and Bioengineering</i> , 2014, 111, 2170-2182.	3.3	47

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19	Divergent Antibody Subclass and Specificity Profiles but Not Protective HLA-B Alleles Are Associated with Variable Antibody Effector Function among HIV-1 Controllers. <i>Journal of Virology</i> , 2014, 88, 2799-2809.	3.4	46
20	Structure-based redesign of lysostaphin yields potent antistaphylococcal enzymes that evade immune cell surveillance. <i>Molecular Therapy - Methods and Clinical Development</i> , 2015, 2, 15021.	4.1	45
21	Diversity of Antiviral IgG Effector Activities Observed in HIV-Infected and Vaccinated Subjects. <i>Journal of Immunology</i> , 2016, 197, 4603-4612.	0.8	44
22	Design and engineering of deimmunized biotherapeutics. <i>Current Opinion in Structural Biology</i> , 2016, 39, 79-88.	5.7	43
23	HCV epitope, homologous to multiple human protein sequences, induces a regulatory T cell response in infected patients. <i>Journal of Hepatology</i> , 2015, 62, 48-55.	3.7	39
24	Structure determination of symmetric homo-oligomers by a complete search of symmetry configuration space, using NMR restraints and van der Waals packing. <i>Proteins: Structure, Function and Bioinformatics</i> , 2006, 65, 203-219.	2.6	38
25	Computationally-driven identification of antibody epitopes. <i>ELife</i> , 2017, 6, .	6.0	37
26	High-resolution definition of humoral immune response correlates of effective immunity against HIV. <i>Molecular Systems Biology</i> , 2018, 14, e7881.	7.2	37
27	Design and analysis of immune-evading enzymes for ADEPT therapy. <i>Protein Engineering, Design and Selection</i> , 2012, 25, 613-624.	2.1	36
28	A divide-and-conquer approach to determine the Pareto frontier for optimization of protein engineering experiments. <i>Proteins: Structure, Function and Bioinformatics</i> , 2012, 80, 790-806.	2.6	34
29	Integrated assessment of predicted MHC binding and cross-conservation with self reveals patterns of viral camouflage. <i>BMC Bioinformatics</i> , 2014, 15, S1.	2.6	34
30	Gaussian Processes for Active Data Mining of Spatial Aggregates. , 2005, , .		33
31	Analysis of self-associating proteins by singular value decomposition of solution scattering data. <i>Biophysical Journal</i> , 2008, 94, 4906-4923.	0.5	32
32	Computationally optimized deimmunization libraries yield highly mutated enzymes with low immunogenicity and enhanced activity. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2017, 114, E5085-E5093.	7.1	32
33	Gene and Protein Sequence Optimization for High-Level Production of Fully Active and Aglycosylated Lysostaphin in <i>Pichia pastoris</i> . <i>Applied and Environmental Microbiology</i> , 2014, 80, 2746-2753.	3.1	30
34	Protein deimmunization via structure-based design enables efficient epitope deletion at high mutational loads. <i>Biotechnology and Bioengineering</i> , 2015, 112, 1306-1318.	3.3	29
35	Optimization of Combinatorial Mutagenesis. <i>Journal of Computational Biology</i> , 2011, 18, 1743-1756.	1.6	28
36	Graphical models of protein-protein interaction specificity from correlated mutations and interaction data. <i>Proteins: Structure, Function and Bioinformatics</i> , 2009, 76, 911-929.	2.6	24

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37	Hypergraph Model of Multi-Residue Interactions in Proteins: Sequentially-Constrained Partitioning Algorithms for Optimization of Site-Directed Protein Recombination. <i>Journal of Computational Biology</i> , 2007, 14, 777-790.	1.6	23
38	Mapping the Pareto Optimal Design Space for a Functionally Deimmunized Biotherapeutic Candidate. <i>PLoS Computational Biology</i> , 2015, 11, e1003988.	3.2	23
39	A Random Graph Approach to NMR Sequential Assignment. <i>Journal of Computational Biology</i> , 2005, 12, 569-583.	1.6	22
40	EpiSweep: Computationally Driven Reengineering of Therapeutic Proteins to Reduce Immunogenicity While Maintaining Function. <i>Methods in Molecular Biology</i> , 2017, 1529, 375-398.	0.9	22
41	Globally deimmunized lysostaphin evades human immune surveillance and enables highly efficacious repeat dosing. <i>Science Advances</i> , 2020, 6, .	10.3	22
42	OPTIMIZATION OF THERAPEUTIC PROTEINS TO DELETE T-CELL EPITOPES WHILE MAINTAINING BENEFICIAL RESIDUE INTERACTIONS. <i>Journal of Bioinformatics and Computational Biology</i> , 2011, 09, 207-229.	0.8	21
43	Computationally driven deletion of broadly distributed T cell epitopes in a biotherapeutic candidate. <i>Cellular and Molecular Life Sciences</i> , 2014, 71, 4869-4880.	5.4	21
44	Site-directed combinatorial construction of chimaeric genes: General method for optimizing assembly of gene fragments. <i>Proteins: Structure, Function and Bioinformatics</i> , 2006, 64, 629-642.	2.6	19
45	Contact replacement for NMR resonance assignment. <i>Bioinformatics</i> , 2008, 24, i205-i213.	4.1	18
46	Structure-based redesign of proteins for minimal T cell epitope content. <i>Journal of Computational Chemistry</i> , 2013, 34, 879-891.	3.3	18
47	Balancing sensitivity and specificity in distinguishing TCR groups by CDR sequence similarity. <i>BMC Bioinformatics</i> , 2019, 20, 241.	2.6	18
48	A complete algorithm to resolve ambiguity for intersubunit NOE assignment in structure determination of symmetric homo-oligomers. <i>Protein Science</i> , 2006, 16, 69-81.	7.6	17
49	Antibody Fab-Fc properties outperform titer in predictive models of <i>SIV</i> vaccine-induced protection. <i>Molecular Systems Biology</i> , 2019, 15, e8747.	7.2	17
50	An efficient randomized algorithm for contact-based NMR backbone resonance assignment. <i>Bioinformatics</i> , 2006, 22, 172-180.	4.1	16
51	A High Throughput MHC II Binding Assay for Quantitative Analysis of Peptide Epitopes. <i>Journal of Visualized Experiments</i> , 2014, , .	0.3	16
52	Development and validation of an epitope prediction tool for swine (PigMatrix) based on the pocket profile method. <i>BMC Bioinformatics</i> , 2015, 16, 290.	2.6	16
53	Fine epitope signature of antibody neutralization breadth at the HIV-1 envelope CD4-binding site. <i>JCI Insight</i> , 2018, 3, .	5.0	16
54	Comprehensive analysis of lectin-glycan interactions reveals determinants of lectin specificity. <i>PLoS Computational Biology</i> , 2021, 17, e1009470.	3.2	16

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55	Ensuring scientific reproducibility in bio-macromolecular modeling via extensive, automated benchmarks. <i>Nature Communications</i> , 2021, 12, 6947.	12.8	16
56	Influence-based model decomposition for reasoning about spatially distributed physical systems. <i>Artificial Intelligence</i> , 2001, 130, 125-166.	5.8	15
57	Probabilistic cross-link analysis and experiment planning for high-throughput elucidation of protein structure. <i>Protein Science</i> , 2009, 13, 3298-3313.	7.6	15
58	Protein Design by Sampling an Undirected Graphical Model of Residue Constraints. <i>IEEE/ACM Transactions on Computational Biology and Bioinformatics</i> , 2009, 6, 506-516.	3.0	15
59	Structure-based design of combinatorial mutagenesis libraries. <i>Protein Science</i> , 2015, 24, 895-908.	7.6	15
60	DisrupPI: structure-based computational redesign algorithm for protein binding disruption. <i>Bioinformatics</i> , 2018, 34, i245-i253.	4.1	15
61	Algorithms for Joint Optimization of Stability and Diversity in Planning Combinatorial Libraries of Chimeric Proteins. <i>Journal of Computational Biology</i> , 2009, 16, 1151-1168.	1.6	14
62	Computationally driven antibody engineering enables simultaneous humanization and thermostabilization. <i>Protein Engineering, Design and Selection</i> , 2016, 29, 419-426.	2.1	14
63	Smarter vaccine design will circumvent regulatory T cell-mediated evasion in chronic HIV and HCV infection. <i>Frontiers in Microbiology</i> , 2014, 5, 502.	3.5	13
64	Reducing Mass Degeneracy in SAR by MS by Stable Isotopic Labeling. <i>Journal of Computational Biology</i> , 2001, 8, 19-36.	1.6	12
65	MHCEpitopeEnergy, a Flexible Rosetta-Based Biotherapeutic Deimmunization Platform. <i>Journal of Chemical Information and Modeling</i> , 2021, 61, 2368-2382.	5.4	12
66	Building blocks and blueprints for bacterial autolysins. <i>PLoS Computational Biology</i> , 2021, 17, e1008889.	3.2	11
67	ALGORITHMS FOR SELECTING BREAKPOINT LOCATIONS TO OPTIMIZE DIVERSITY IN PROTEIN ENGINEERING BY SITE-DIRECTED PROTEIN RECOMBINATION. , 2007, , .		11
68	Robotic Hierarchical Mixing for the Production of Combinatorial Libraries of Proteins and Small Molecules. <i>ACS Combinatorial Science</i> , 2008, 10, 63-68.	3.3	10
69	Analysis of sequence-reactivity space for protein-protein interactions. <i>Proteins: Structure, Function and Bioinformatics</i> , 2004, 58, 661-671.	2.6	9
70	A graphical method for analyzing distance restraints using residual dipolar couplings for structure determination of symmetric protein homo-oligomers. <i>Protein Science</i> , 2011, 20, 970-985.	7.6	9
71	Towards conformational fidelity of a quaternary HIV-1 epitope: computational design and directed evolution of a minimal V1V2 antigen. <i>Protein Engineering, Design and Selection</i> , 2018, 31, 121-133.	2.1	8
72	A subgroup algorithm to identify cross-rotation peaks consistent with non-crystallographic symmetry. <i>Acta Crystallographica Section D: Biological Crystallography</i> , 2004, 60, 1057-1067.	2.5	7

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73	Protein Fragment Swapping: A Method for Asymmetric, Selective Site-Directed Recombination. <i>Journal of Computational Biology</i> , 2010, 17, 459-475.	1.6	7
74	Diverse antiviral IgG effector activities are predicted by unique biophysical antibody features. <i>Retrovirology</i> , 2021, 18, 35.	2.0	7
75	A Geometric Arrangement Algorithm for Structure Determination of Symmetric Protein Homo-Oligomers from NOEs and RDCs. <i>Journal of Computational Biology</i> , 2011, 18, 1507-1523.	1.6	6
76	Algorithms for optimizing cross-overs in DNA shuffling. <i>BMC Bioinformatics</i> , 2012, 13, S3.	2.6	6
77	Hit-and-run, hit-and-stay, and commensal bacteria present different peptide content when viewed from the perspective of the T cell. <i>Vaccine</i> , 2015, 33, 6922-6929.	3.8	6
78	Pareto Optimization of Combinatorial Mutagenesis Libraries. <i>IEEE/ACM Transactions on Computational Biology and Bioinformatics</i> , 2019, 16, 1143-1153.	3.0	6
79	Algorithms for selecting breakpoint locations to optimize diversity in protein engineering by site-directed protein recombination. <i>Computational Systems Bioinformatics / Life Sciences Society Computational Systems Bioinformatics Conference</i> , 2007, 6, 31-40.	0.4	6
80	A Hierarchical Grow-and-Match Algorithm for Backbone Resonance Assignments Given 3D Structure. , 2007, , .		5
81	NMR Structural Inference of Symmetric Homo-Oligomers. <i>Journal of Computational Biology</i> , 2011, 18, 1757-1775.	1.6	5
82	Learning Sequence Determinants of Protein:Protein Interaction Specificity with Sparse Graphical Models. <i>Journal of Computational Biology</i> , 2015, 22, 474-486.	1.6	5
83	Characterizing Epitope Binding Regions of Entire Antibody Panels by Combining Experimental and Computational Analysis of Antibody: Antigen Binding Competition. <i>Molecules</i> , 2020, 25, 3659.	3.8	5
84	Induction of cross-reactive HIV-1 specific antibody responses by engineered V1V2 immunogens with reduced conformational plasticity. <i>Vaccine</i> , 2020, 38, 3436-3446.	3.8	5
85	Hypergraph Model of Multi-residue Interactions in Proteins: Sequentiallyâ€œConstrained Partitioning Algorithms for Optimization of Site-Directed Protein Recombination. <i>Lecture Notes in Computer Science</i> , 2006, , 15-29.	1.3	5
86	Stoichiometries and affinities of interacting proteins from concentration series of solution scattering data: decomposition by least squares and quadratic optimization. <i>Journal of Applied Crystallography</i> , 2014, 47, 899-914.	4.5	4
87	Computer-guided binding mode identification and affinity improvement of an LRR protein binder without structure determination. <i>PLoS Computational Biology</i> , 2020, 16, e1008150.	3.2	4
88	A Chimeric Antigen Receptor That Binds to a Conserved Site on MICA. <i>ImmunoHorizons</i> , 2020, 4, 597-607.	1.8	4
89	Learning Sequence Determinants of Protein: Protein Interaction Specificity with Sparse Graphical Models. <i>Lecture Notes in Computer Science</i> , 2014, 8394, 129-143.	1.3	4
90	Combinatorial Resurfacing of Dengue Envelope Protein Domain III Antigens Selectively Ablates Epitopes Associated with Serotype-Specific or Infection-Enhancing Antibody Responses. <i>ACS Combinatorial Science</i> , 2020, 22, 446-456.	3.8	3

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91	A multi-resolution graph convolution network for contiguous epitope prediction. , 2021, , .		3
92	Structure-Guided Deimmunization of Therapeutic Proteins. Lecture Notes in Computer Science, 2012, , 184-198.	1.3	3
93	OPTIMIZING BAYES ERROR FOR PROTEIN STRUCTURE MODEL SELECTION BY STABILITY MUTAGENESIS. , 2008, , .		3
94	Intelligent simulation tools for mining large scientific data sets. New Generation Computing, 1999, 17, 333-347.	3.3	2
95	Functional evolution within a protein superfamily. Proteins: Structure, Function and Bioinformatics, 2006, 63, 697-708.	2.6	2
96	Bioinformatics-driven discovery of novel <i>Clostridioides difficile</i> lysins and experimental comparison with highly active benchmarks. Biotechnology and Bioengineering, 2021, 118, 2482-2492.	3.3	2
97	A Geometric Arrangement Algorithm for Structure Determination of Symmetric Protein Homo-oligomers from NOEs and RDCs. Lecture Notes in Computer Science, 2011, , 222-237.	1.3	2
98	Algorithms for Joint Optimization of Stability and Diversity in Planning Combinatorial Libraries of Chimeric Proteins. , 2008, , 300-314.		2
99	Computational epitope binning reveals functional equivalence of sequence-divergent paratopes. Computational and Structural Biotechnology Journal, 2022, 20, 2169-2180.	4.1	2
100	Planning combinatorial disulfide cross-links for protein fold determination. BMC Bioinformatics, 2011, 12, S5.	2.6	1
101	Structure Discovery from Massive Spatial Data Sets Using Intelligent Simulation Tools. Lecture Notes in Computer Science, 2007, , 158-174.	1.3	1
102	Simultaneous determination of subunit and complex structures of symmetric homo-oligomers from ambiguous NMR data. , 2013, , .		0
103	OCoM-SOCoM. , 2016, , .		0
104	Extended Abstract: Structure Determination of Symmetric Protein Complexes by a Complete Search of Symmetry Configuration Space Using NMR Distance Restraints. Springer Tracts in Advanced Robotics, 2008, , 335-340.	0.4	0
105	Protein Fragment Swapping: A Method for Asymmetric, Selective Site-Directed Recombination. Lecture Notes in Computer Science, 2009, , 321-338.	1.3	0
106	Title is missing!. , 2020, 16, e1008150.		0
107	Title is missing!. , 2020, 16, e1008150.		0
108	Title is missing!. , 2020, 16, e1008150.		0

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109	Title is missing!. , 2020, 16, e1008150.		0