

# Thomas E Ferrin

## List of Publications by Year in descending order

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64  
papers

54,571  
citations

117453

34  
h-index

143772

57  
g-index

70  
all docs

70  
docs citations

70  
times ranked

67910  
citing authors

#	ARTICLE	IF	CITATIONS
1	UCSF Chimera?A visualization system for exploratory research and analysis. <i>Journal of Computational Chemistry</i> , 2004, 25, 1605-1612.	1.5	37,095
2	<scp>UCSF ChimeraX</scp>: Structure visualization for researchers, educators, and developers. <i>Protein Science</i> , 2021, 30, 70-82.	3.1	4,478
3	UCSF ChimeraX: Meeting modern challenges in visualization and analysis. <i>Protein Science</i> , 2018, 27, 14-25.	3.1	3,377
4	A geometric approach to macromolecule-ligand interactions. <i>Journal of Molecular Biology</i> , 1982, 161, 269-288.	2.0	2,043
5	The MIDAS display system. <i>Journal of Molecular Graphics</i> , 1988, 6, 13-27.	1.7	914
6	Visualizing density maps with UCSF Chimera. <i>Journal of Structural Biology</i> , 2007, 157, 281-287.	1.3	802
7	clusterMaker: a multi-algorithm clustering plugin for Cytoscape. <i>BMC Bioinformatics</i> , 2011, 12, 436.	1.2	541
8	Tools for integrated sequence-structure analysis with UCSF Chimera. <i>BMC Bioinformatics</i> , 2006, 7, 339.	1.2	524
9	ModBase, a database of annotated comparative protein structure models, and associated resources. <i>Nucleic Acids Research</i> , 2011, 39, D465-D474.	6.5	506
10	UCSF Chimera, MODELLER, and IMP: An integrated modeling system. <i>Journal of Structural Biology</i> , 2012, 179, 269-278.	1.3	506
11	Designed divergent evolution of enzyme function. <i>Nature</i> , 2006, 440, 1078-1082.	13.7	414
12	Using Sequence Similarity Networks for Visualization of Relationships Across Diverse Protein Superfamilies. <i>PLoS ONE</i> , 2009, 4, e4345.	1.1	385
13	Evolutionary conservation predicts function of variants of the human organic cation transporter, OCT1. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2003, 100, 5902-5907.	3.3	265
14	MODBASE, a database of annotated comparative protein structure models, and associated resources. <i>Nucleic Acids Research</i> , 2004, 32, 217D-222.	6.5	256
15	Natural variation in human membrane transporter genes reveals evolutionary and functional constraints. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2003, 100, 5896-5901.	3.3	224
16	The Structureâ€“Function Linkage Database. <i>Nucleic Acids Research</i> , 2014, 42, D521-D530.	6.5	210
17	Software Extensions to UCSF Chimera for Interactive Visualization of Large Molecular Assemblies. <i>Structure</i> , 2005, 13, 473-482.	1.6	204
18	Outcome of the First wwPDB Hybrid/Integrative Methods Task Force Workshop. <i>Structure</i> , 2015, 23, 1156-1167.	1.6	159

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19	Leveraging Enzyme Structure~Function Relationships for Functional Inference and Experimental Design:~ The Structure~Function Linkage Database. <i>Biochemistry</i> , 2006, 45, 2545-2555.	1.2	157
20	Conic: A fast renderer for spacefilling molecules with shadows. <i>Journal of Molecular Graphics</i> , 1991, 9, 230-236.	1.7	134
21	Enhancing UCSF Chimera through web services. <i>Nucleic Acids Research</i> , 2014, 42, W478-W484.	6.5	116
22	Actin-based protrusions of migrating neutrophils are intrinsically lamellar and facilitate direction changes. <i>ELife</i> , 2017, 6, .	2.8	107
23	Comparison of human solute carriers. <i>Protein Science</i> , 2010, 19, 412-428.	3.1	99
24	Van der Waals Surfaces in Molecular Modeling: Implementation with Real-Time Computer Graphics. <i>Science</i> , 1983, 222, 1325-1327.	6.0	96
25	Automated site-directed drug design using molecular lattices. <i>Journal of Molecular Graphics</i> , 1992, 10, 66-78.	1.7	95
26	Computer graphics in drug design: molecular modeling of thyroid hormone-prealbumin interactions. <i>Journal of Medicinal Chemistry</i> , 1982, 25, 785-790.	2.9	93
27	Engineering Cotton (+)-Î-Cadinene Synthase to an Altered Function: Germacrene D-4-ol Synthase. <i>Chemistry and Biology</i> , 2006, 13, 91-98.	6.2	83
28	structureViz: linking Cytoscape and UCSF Chimera. <i>Bioinformatics</i> , 2007, 23, 2345-2347.	1.8	71
29	Molecular Visualization on the Holodeck. <i>Journal of Molecular Biology</i> , 2018, 430, 3982-3996.	2.0	70
30	Nucleic acid visualization with UCSF Chimera. <i>Nucleic Acids Research</i> , 2006, 34, e29-e29.	6.5	69
31	RRDistMaps: a UCSF Chimera tool for viewing and comparing protein distance maps. <i>Bioinformatics</i> , 2015, 31, 1484-1486.	1.8	52
32	enhancedGraphics: a Cytoscape app for enhanced node graphics. <i>F1000Research</i> , 2014, 3, 147.	0.8	45
33	Visualization software for molecular assemblies. <i>Current Opinion in Structural Biology</i> , 2007, 17, 587-595.	2.6	40
34	Improving the quality of protein similarity network clustering algorithms using the network edge weight distribution. <i>Bioinformatics</i> , 2011, 27, 326-333.	1.8	36
35	The MIDAS database system. <i>Journal of Molecular Graphics</i> , 1988, 6, 2-12.	1.7	27
36	Interactive computer graphics with the UNIX time-sharing system. <i>Computer Graphics</i> , 1980, 13, 320-331.	0.1	23

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37	Clinical Persistence of Chlamydia trachomatis Sexually Transmitted Strains Involves Novel Mutations in the Functional $\epsilon$ -Tetramer of the Tryptophan Synthase Operon. MBio, 2019, 10, .	1.8	20
38	An Atlas of Peroxiredoxins Created Using an Active Site Profile-Based Approach to Functionally Relevant Clustering of Proteins. PLoS Computational Biology, 2017, 13, e1005284.	1.5	19
39	Multidomain Assembler (MDA) Generates Models of Large Multidomain Proteins. Biophysical Journal, 2015, 108, 2097-2102.	0.2	18
40	A real-time malleable molecular surface. Journal of Molecular Graphics, 1990, 8, 16-24.	1.7	16
41	setsApp: Set operations for Cytoscape Nodes and Edges. F1000Research, 2014, 3, 149.	0.8	14
42	Representing structure-function relationships in mechanistically diverse enzyme superfamilies. Pacific Symposium on Biocomputing Pacific Symposium on Biocomputing, 2005, , 358-69.	0.7	14
43	An approach to functionally relevant clustering of the protein universe: Active site profile-based clustering of protein structures and sequences. Protein Science, 2017, 26, 677-699.	3.1	13
44	CyAnimator: Simple Animations of Cytoscape Networks. F1000Research, 2015, 4, 482.	0.8	11
45	Computer-Assisted Drug Receptor Mapping Analysis. ACS Symposium Series, 1986, , 147-158.	0.5	10
46	CyAnimator: Simple Animations of Cytoscape Networks. F1000Research, 2015, 4, 482.	0.8	10
47	Computational Tools for the Interactive Exploration of Proteomic and Structural Data. Molecular and Cellular Proteomics, 2010, 9, 1703-1715.	2.5	9
48	Integrated Tools for Structural and Sequence Alignment and Analysis. , 1999, , 230-41.		9
49	Annotating PDB files with scene information. Journal of Molecular Graphics, 1995, 13, 153-158.	1.7	8
50	setsApp for Cytoscape: Set operations for Cytoscape Nodes and Edges. F1000Research, 2014, 3, 149.	0.8	8
51	The Structure Superposition Database. Nucleic Acids Research, 2003, 31, 505-510.	6.5	7
52	DASP3: identification of protein sequences belonging to functionally relevant groups. BMC Bioinformatics, 2016, 17, 458.	1.2	6
53	Biocuration in the structure-function linkage database: the anatomy of a superfamily. Database: the Journal of Biological Databases and Curation, 2017, 2017, .	1.4	6
54	Macromolecular Assembly Structures by Comparative Modeling and Electron Microscopy. Methods in Molecular Biology, 2011, 857, 331-350.	0.4	4

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55	cddApp: a Cytoscape app for accessing the NCBI conserved domain database. <i>Bioinformatics</i> , 2015, 31, 134-136.	1.8	4
56	Biocuration in the structureâ€function linkage database: the anatomy of a superfamily. <i>Database: the Journal of Biological Databases and Curation</i> , 2017, 2017, .	1.4	2
57	SNP ANALYSIS AND PRESENTATION IN THE PHARMACOGENETICS OF MEMBRANE TRANSPORTERS PROJECT. , 2002, , .		2
58	Molecular image representations and their uses. <i>Journal of Molecular Graphics</i> , 1992, 10, 47.	1.7	0
59	Reply to Rockey et al., â€œGenomics and Chlamydial Persistence <i>In Vivo</i> â€ MBio, 2019, 10, .	1.8	0
60	High-Performance Computing for Computational Biology. , 2000, , .		0
61	The Structureâ€Function Linkage Database. <i>FASEB Journal</i> , 2006, 20, A899.	0.2	0
62	MOLECULES TO MAPS: TOOLS FOR VISUALIZATION AND INTERACTION IN COMPUTATIONAL BIOLOGY. , 1998, , .		0
63	MOLECULES TO MAPS: TOOLS FOR VISUALIZATION AND INTERACTION IN COMPUTATIONAL BIOLOGY. , 1999, , .		0
64	Enhancing data sharing in collaborative research projects with DASH. <i>Pacific Symposium on Biocomputing Pacific Symposium on Biocomputing</i> , 2005, , 260-71.	0.7	0