## Calvin Yu-Chian Chen

## 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.

118	3,204	31	53
papers	citations	h-index	g-index
131 ext. papers	3,549 ext. citations	3.6 avg, IF	6.15 L-index

#	Paper	IF	Citations
118	Network pharmacology-based approach of novel traditional Chinese medicine formula for treatment of acute skin inflammation in silico. <i>Computational Biology and Chemistry</i> , <b>2017</b> , 71, 70-81	3.6	13
117	Residue-based design of small molecule inhibitor for H1N1, H5N1 and H7N1 mutants. <i>Journal of Molecular Modeling</i> , <b>2016</b> , 22, 4	2	
116	Molecular insight and resolution for tumors harboring the H-ras(G12V) mutation. <i>RSC Advances</i> , <b>2015</b> , 5, 20623-20633	3.7	4
115	Susceptible gene of stasis-stagnation constitution from genome-wide association study related to cardiovascular disturbance and possible regulated traditional Chinese medicine. <i>BMC Complementary and Alternative Medicine</i> , <b>2015</b> , 15, 229	4.7	5
114	Beware of docking!. <i>Trends in Pharmacological Sciences</i> , <b>2015</b> , 36, 78-95	13.2	329
113	Insight into molecular dynamics simulation of BRAF(V600E) and potent novel inhibitors for malignant melanoma. <i>International Journal of Nanomedicine</i> , <b>2015</b> , 10, 3131-46	7.3	10
112	When modern technology meets ancient traditional chinese medicine. <i>Evidence-based Complementary and Alternative Medicine</i> , <b>2015</b> , 2015, 156581	2.3	3
111	Identification of tyrosinase inhibitors from traditional Chinese medicines for the management of hyperpigmentation. <i>SpringerPlus</i> , <b>2015</b> , 4, 184		13
110	Insight into two antioxidants binding to the catalase NADPH binding site from traditional Chinese medicines. <i>RSC Advances</i> , <b>2015</b> , 5, 6625-6635	3.7	4
109	Discovery of novel insomnia leads from screening traditional Chinese medicine database. <i>Journal of Biomolecular Structure and Dynamics</i> , <b>2014</b> , 32, 776-91	3.6	19
108	Molecular level activation insights from a NR2A/NR2B agonist. <i>Journal of Biomolecular Structure and Dynamics</i> , <b>2014</b> , 32, 683-93	3.6	8
107	May disordered protein cause serious drug side effect?. <i>Drug Discovery Today</i> , <b>2014</b> , 19, 367-72	8.8	35
106	Chalcone derivatives inhibit human platelet aggregation and inhibit growth in human bladder cancer cells. <i>Biological and Pharmaceutical Bulletin</i> , <b>2014</b> , 37, 1191-8	2.3	12
105	Developing hypothetical inhibition mechanism of novel urea transporter B inhibitor. <i>Scientific Reports</i> , <b>2014</b> , 4, 5775	4.9	14
104	In silico design for adenosine monophosphate-activated protein kinase agonist from traditional chinese medicine for treatment of metabolic syndromes. <i>Evidence-based Complementary and Alternative Medicine</i> , <b>2014</b> , 2014, 928589	2.3	6
103	In Silico Insight into Potent of Anthocyanin Regulation of FKBP52 to Prevent Alzheimerß Disease. <i>Evidence-based Complementary and Alternative Medicine</i> , <b>2014</b> , 2014, 450592	2.3	5
102	Finding inhibitors of mutant superoxide dismutase-1 for amyotrophic lateral sclerosis therapy from traditional chinese medicine. <i>Evidence-based Complementary and Alternative Medicine</i> , <b>2014</b> , 2014, 156	276 <sup>3</sup>	10

101	Computational design of apolipoprotein E4 inhibitors for Alzheimerß disease therapy from traditional Chinese medicine. <i>BioMed Research International</i> , <b>2014</b> , 2014, 452625	3	8
100	Treatment of acute lymphoblastic leukemia from traditional chinese medicine. <i>Evidence-based Complementary and Alternative Medicine</i> , <b>2014</b> , 2014, 601064	2.3	4
99	Potential smoothened inhibitor from traditional Chinese medicine against the disease of diabetes, obesity, and cancer. <i>BioMed Research International</i> , <b>2014</b> , 2014, 873010	3	4
98	A possible solution for hair loss by inhibiting corticotropin-releasing factor (CRF) receptor from traditional Chinese medicine. <i>Journal of Biomolecular Structure and Dynamics</i> , <b>2014</b> , 32, 1613-23	3.6	3
97	In Silico Identification of Potent PPAR-[Agonists from Traditional Chinese Medicine: A Bioactivity Prediction, Virtual Screening, and Molecular Dynamics Study. <i>Evidence-based Complementary and Alternative Medicine</i> , <b>2014</b> , 2014, 192452	2.3	10
96	Design of glucagon-like Peptide-1 receptor agonist for diabetes mellitus from traditional chinese medicine. <i>Evidence-based Complementary and Alternative Medicine</i> , <b>2014</b> , 2014, 385120	2.3	
95	In silico design of BACE1 inhibitor for Alzheimerß disease by traditional Chinese medicine. <i>BioMed Research International</i> , <b>2014</b> , 2014, 741703	3	7
94	Investigation of anti-infection mechanism of lactoferricin and splunc-1. <i>Evidence-based Complementary and Alternative Medicine</i> , <b>2014</b> , 2014, 907028	2.3	4
93	In silico investigation of potential pyruvate kinase M2 regulators from traditional Chinese medicine against cancers. <i>BioMed Research International</i> , <b>2014</b> , 2014, 189495	3	1
92	Investigation of the novel lead of melanocortin 1 receptor for pigmentary disorders. <i>Evidence-based Complementary and Alternative Medicine</i> , <b>2014</b> , 2014, 254678	2.3	10
91	Potential retinoid x receptor agonists for treating Alzheimerß disease from traditional chinese medicine. <i>Evidence-based Complementary and Alternative Medicine</i> , <b>2014</b> , 2014, 278493	2.3	4
90	Ligand-based and structure-based investigation for Alzheimerß disease from traditional chinese medicine. <i>Evidence-based Complementary and Alternative Medicine</i> , <b>2014</b> , 2014, 364819	2.3	23
89	Lead screening for CXCR4 of the human HIV infection receptor inhibited by traditional Chinese medicine. <i>BioMed Research International</i> , <b>2014</b> , 2014, 809816	3	2
88	Lead Screening for HIV of C-C Chemokine Receptor Type 5 Receptor Inhibited by Traditional Chinese Medicine. <i>Evidence-based Complementary and Alternative Medicine</i> , <b>2014</b> , 2014, 313094	2.3	2
87	Lead screening for HIV-1 integrase (IN) inhibited by traditional Chinese medicine. <i>BioMed Research International</i> , <b>2014</b> , 2014, 479367	3	4
86	Treatment of rheumatoid arthritis with traditional chinese medicine. <i>BioMed Research International</i> , <b>2014</b> , 2014, 528018	3	18
85	The inhibition of folylpolyglutamate synthetase (folC) in the prevention of drug resistance in Mycobacterium tuberculosis by traditional Chinese medicine. <i>BioMed Research International</i> , <b>2014</b> , 2014, 635152	3	2
84	Lead discovery for Alzheimerß disease related target protein RbAp48 from traditional Chinese medicine. <i>BioMed Research International</i> , <b>2014</b> , 2014, 764946	3	8

83	In silico investigation of traditional Chinese medicine compounds to inhibit human histone deacetylase 2 for patients with Alzheimerß disease. <i>BioMed Research International</i> , <b>2014</b> , 2014, 769867	3	9
82	Drug design of cyclin-dependent kinase 2 inhibitor for melanoma from traditional Chinese medicine. <i>BioMed Research International</i> , <b>2014</b> , 2014, 798742	3	8
81	Pharmacological chaperone design for reducing risk factor of Parkinson® disease from traditional chinese medicine. <i>Evidence-based Complementary and Alternative Medicine</i> , <b>2014</b> , 2014, 830490	2.3	18
80	Investigation of estrogen receptor (ESR1) for breast cancer from traditional Chinese medicine. <i>BioMed Research International</i> , <b>2014</b> , 2014, 321486	3	5
79	Lead Screening for Chronic Obstructive Pulmonary Disease of IKK2 Inhibited by Traditional Chinese Medicine. <i>Evidence-based Complementary and Alternative Medicine</i> , <b>2014</b> , 2014, 465025	2.3	3
78	In silico investigation of potential mTOR inhibitors from traditional Chinese medicine for treatment of Leigh syndrome. <i>BioMed Research International</i> , <b>2014</b> , 2014, 139492	3	4
77	In Silico Investigation of Cytochrome P450 2C9 in relation to Aging Using Traditional Chinese Medicine. <i>Evidence-based Complementary and Alternative Medicine</i> , <b>2014</b> , 2014, 404505	2.3	1
76	In Silico Investigation of Potential PARP-1 Inhibitors from Traditional Chinese Medicine. <i>Evidence-based Complementary and Alternative Medicine</i> , <b>2014</b> , 2014, 917605	2.3	3
75	Potential Protein Phosphatase 2A Agents from Traditional Chinese Medicine against Cancer. Evidence-based Complementary and Alternative Medicine, <b>2014</b> , 2014, 436863	2.3	6
74	An investigation of small GTPases in relation to liver tumorigenesis using traditional Chinese medicine. <i>BioMed Research International</i> , <b>2014</b> , 2014, 428210	3	1
73	Evaluation of correlation of cell cycle proteins and Ki-67 interaction in paranasal sinus inverted papilloma prognosis and squamous cell carcinoma transformation. <i>BioMed Research International</i> , <b>2014</b> , 2014, 634945	3	13
72	Investigation of potent lead for acquired immunodeficiency syndrome from traditional Chinese medicine. <i>BioMed Research International</i> , <b>2014</b> , 2014, 205890	3	3
71	Insight into HIV of IFN-induced myxovirus resistance 2 (MX2) expressed by traditional Chinese medicine. <i>BioMed Research International</i> , <b>2014</b> , 2014, 871576	3	
70	Potential mitochondrial isocitrate dehydrogenase R140Q mutant inhibitor from traditional Chinese medicine against cancers. <i>BioMed Research International</i> , <b>2014</b> , 2014, 364625	3	5
69	In silico investigation of potential TRAF6 inhibitor from traditional Chinese medicine against cancers. <i>BioMed Research International</i> , <b>2014</b> , 2014, 429486	3	4
68	Possible inhibitor from traditional Chinese medicine for the Iform of calcium-dependent protein kinase type II in the treatment of major depressive disorder. <i>BioMed Research International</i> , <b>2014</b> , 2014, 761849	3	1
67	Treatment of cardiovascular disease by traditional Chinese medicine against pregnane X receptor. BioMed Research International, <b>2014</b> , 2014, 950191	3	15
66	Osteoponin promoter controlled by DNA methylation: aberrant methylation in cloned porcine genome. <i>BioMed Research International</i> , <b>2014</b> , 2014, 327538	3	8

## (2011-2014)

65	Traditional Chinese medicine application in HIV: an in silico study. <i>Journal of Biomolecular Structure and Dynamics</i> , <b>2014</b> , 32, 1-12	3.6	29
64	How to design a drug for the disordered proteins?. <i>Drug Discovery Today</i> , <b>2013</b> , 18, 910-5	8.8	64
63	Investigation of silent information regulator 1 (Sirt1) agonists from Traditional Chinese Medicine. <i>Journal of Biomolecular Structure and Dynamics</i> , <b>2013</b> , 31, 1207-18	3.6	18
62	Memory enhancement by traditional Chinese medicine?. <i>Journal of Biomolecular Structure and Dynamics</i> , <b>2013</b> , 31, 1411-39	3.6	5
61	Han ethnicity-specific type 2 diabetic treatment from traditional Chinese medicine?. <i>Journal of Biomolecular Structure and Dynamics</i> , <b>2013</b> , 31, 1219-35	3.6	16
60	Drug design for neuropathic pain regulation from traditional Chinese medicine. <i>Scientific Reports</i> , <b>2013</b> , 3, 844	4.9	55
59	A possible strategy against head and neck cancer: in silico investigation of three-in-one inhibitors. Journal of Biomolecular Structure and Dynamics, <b>2013</b> , 31, 1358-69	3.6	7
58	A novel integrated framework and improved methodology of computer-aided drug design. <i>Current Topics in Medicinal Chemistry</i> , <b>2013</b> , 13, 965-88	3	53
57	Three-in-one agonists for PPAR-PPPAR-and PPAR-Ifrom traditional Chinese medicine. <i>Journal of Biomolecular Structure and Dynamics</i> , <b>2012</b> , 30, 662-83	3.6	41
56	Traditional Chinese medicine as dual guardians against hypertension and cancer?. <i>Journal of Biomolecular Structure and Dynamics</i> , <b>2012</b> , 30, 299-317	3.6	11
55	In silico investigation of potential SRC kinase ligands from traditional Chinese medicine. <i>PLoS ONE</i> , <b>2012</b> , 7, e33728	3.7	20
54	In silico identification of potent pancreatic triacylglycerol lipase inhibitors from traditional Chinese medicine. <i>PLoS ONE</i> , <b>2012</b> , 7, e43932	3.7	20
53	Mechanism of BAG1 repair on Parkinson® disease-linked DJ1 mutation. <i>Journal of Biomolecular Structure and Dynamics</i> , <b>2012</b> , 30, 1-12	3.6	12
52	A novel algorithm for identification of activated cryptic 5Psplice sites. <i>Journal of Biomolecular Structure and Dynamics</i> , <b>2012</b> , 29, 1089-99	3.6	3
51	Uroporphyrinogen decarboxylase as a potential target for specific components of traditional Chinese medicine: a virtual screening and molecular dynamics study. <i>PLoS ONE</i> , <b>2012</b> , 7, e50087	3.7	27
50	Structure-based and ligand-based drug design for microsomal prostaglandin E synthase-1 inhibitors. <i>Molecular Simulation</i> , <b>2011</b> , 37, 226-236	2	4
49	Screening from TCM Database@Taiwan and QSAR model for identifying HER2 inhibitors. <i>Molecular Simulation</i> , <b>2011</b> , 37, 884-892	2	7
48	Potent inhibitor design against H1N1 swine influenza: structure-based and molecular dynamics analysis for M2 inhibitors from traditional Chinese medicine database. <i>Journal of Biomolecular Structure and Dynamics</i> , <b>2011</b> , 28, 471-82	3.6	44

47	Key features for designing M2 proton channel anti swine flu inhibitors. <i>Journal of the Taiwan Institute of Chemical Engineers</i> , <b>2011</b> , 42, 701-708	5.3	6
46	TCM Database@Taiwan: the worldß largest traditional Chinese medicine database for drug screening in silico. <i>PLoS ONE</i> , <b>2011</b> , 6, e15939	3.7	504
45	Investigation into potent inflammation inhibitors from traditional Chinese medicine. <i>Chemical Biology and Drug Design</i> , <b>2011</b> , 78, 679-88	2.9	50
44	iScreen: worldß first cloud-computing web server for virtual screening and de novo drug design based on TCM database@Taiwan. <i>Journal of Computer-Aided Molecular Design</i> , <b>2011</b> , 25, 525-31	4.2	108
43	Drug design for hemagglutinin: Screening and molecular dynamics from traditional Chinese medicine database. <i>Journal of the Taiwan Institute of Chemical Engineers</i> , <b>2011</b> , 42, 563-571	5.3	10
42	High performance screening, structural and molecular dynamics analysis to identify H1 inhibitors from TCM Database@Taiwan. <i>Molecular BioSystems</i> , <b>2011</b> , 7, 3366-74		34
41	Stroke prevention by traditional Chinese medicine? A genetic algorithm, support vector machine and molecular dynamics approach. <i>Soft Matter</i> , <b>2011</b> , 7, 4001	3.6	52
40	Lose weight with traditional chinese medicine? Potential suppression of fat mass and obesity-associated protein. <i>Journal of Biomolecular Structure and Dynamics</i> , <b>2011</b> , 29, 471-83	3.6	16
39	Traditional Chinese medicine, a solution for reducing dual stroke risk factors at once?. <i>Molecular BioSystems</i> , <b>2011</b> , 7, 2711-9		39
38	In silico pharmacology suggests ginger extracts may reduce stroke risks. <i>Molecular BioSystems</i> , <b>2011</b> , 7, 2702-10		42
37	Screening from the world® largest TCM database for inhibiting DNA repair protein XRCC4. <i>Molecular Simulation</i> , <b>2011</b> , 37, 503-509	2	1
36	Drug discovery for DNA break repair system by screening from TCM database and molecular dynamics approach. <i>Molecular Simulation</i> , <b>2011</b> , 37, 804-811	2	4
35	iSMART: an integrated cloud computing web server for traditional Chinese medicine for online virtual screening, de novo evolution and drug design. <i>Journal of Biomolecular Structure and Dynamics</i> , <b>2011</b> , 29, 243-50	3.6	52
34	Screening from the worldß largest TCM database against H1N1 virus. <i>Journal of Biomolecular Structure and Dynamics</i> , <b>2011</b> , 28, 773-86	3.6	42
33	Drug design for mPGES-1 from traditional Chinese medicine database: A screening, docking, QSAR, molecular dynamics, and pharmacophore mapping study. <i>Journal of the Taiwan Institute of Chemical Engineers</i> , <b>2011</b> , 42, 580-591	5.3	8
32	Treat Alzheimerß disease by traditional Chinese medicine?. <i>Molecular Simulation</i> , <b>2011</b> , 37, 923-931	2	2
31	Novel hemagglutinin inhibitors for H1N1 influenza virus screening from TCM database. <i>Molecular Simulation</i> , <b>2011</b> , 37, 361-368	2	8
30	Molecular dynamics analysis of potent inhibitors of M2 proton channel against H1N1 swine influenza virus. <i>Molecular Simulation</i> , <b>2011</b> , 37, 250-256	2	7

## (2009-2011)

29	Two birds with one stone? Possible dual-targeting H1N1 inhibitors from traditional Chinese medicine. <i>PLoS Computational Biology</i> , <b>2011</b> , 7, e1002315	5	72	
28	Identification of potent EGFR inhibitors from TCM Database@Taiwan. <i>PLoS Computational Biology</i> , <b>2011</b> , 7, e1002189	5	78	
27	Identifying HER2 inhibitors from natural products database. <i>PLoS ONE</i> , <b>2011</b> , 6, e28793	3.7	32	
26	Structure-based and ligand-based drug design for HER 2 receptor. <i>Journal of Biomolecular Structure and Dynamics</i> , <b>2010</b> , 28, 23-37	3.6	77	
25	Application of bioactivity database of Chinese herbal medicine on the therapeutic prediction, drug development, and safety evaluation. <i>Journal of Ethnopharmacology</i> , <b>2010</b> , 132, 429-37	5	36	
24	A novel strategy for designing the selective PPAR agonist by the "sum of activity" model. <i>Journal of Biomolecular Structure and Dynamics</i> , <b>2010</b> , 28, 187-200	3.6	56	
23	Current developments of computer-aided drug design. <i>Journal of the Taiwan Institute of Chemical Engineers</i> , <b>2010</b> , 41, 623-635	5.3	93	
22	Drug design for Influenza A virus subtype H1N1. <i>Journal of the Taiwan Institute of Chemical Engineers</i> , <b>2010</b> , 41, 8-15	5.3	23	
21	Bioinformatics, chemoinformatics, and pharmainformatics analysis of HER2/HSP90 dual-targeted inhibitors. <i>Journal of the Taiwan Institute of Chemical Engineers</i> , <b>2010</b> , 41, 143-149	5.3	25	
20	Computational screening and QSAR analysis for design of AMP-activated protein kinase agonist. Journal of the Taiwan Institute of Chemical Engineers, <b>2010</b> , 41, 352-359	5.3	14	
19	Chemical engineering research in Taiwan in 2007 2009. <i>Journal of the Taiwan Institute of Chemical Engineers</i> , <b>2010</b> , 41, 244-245	5.3	1	
18	Insights into designing the dual-targeted HER2/HSP90 inhibitors. <i>Journal of Molecular Graphics and Modelling</i> , <b>2010</b> , 29, 21-31	2.8	61	
17	Reducing without Side Effects? A Novel Strategy for Designing the PPAR Agonists 2009,		1	
16	Computational screening and design of traditional Chinese medicine (TCM) to block phosphodiesterase-5. <i>Journal of Molecular Graphics and Modelling</i> , <b>2009</b> , 28, 261-9	2.8	48	
15	Magnolol encapsulated by different acyl chain length of liposomes on inhibiting proliferation of smooth muscle cells. <i>Journal of the Taiwan Institute of Chemical Engineers</i> , <b>2009</b> , 40, 380-386	5.3	4	
14	Discovery of potent inhibitors for phosphodiesterase 5 by virtual screening and pharmacophore analysis. <i>Acta Pharmacologica Sinica</i> , <b>2009</b> , 30, 1186-94	8	40	
13	De novo design of novel selective COX-2 inhibitors: From virtual screening to pharmacophore analysis. <i>Journal of the Taiwan Institute of Chemical Engineers</i> , <b>2009</b> , 40, 55-69	5.3	22	
12	Pharmacoinformatics approach for mPGES-1 in anti-inflammation by 3D-QSAR pharmacophore mapping. <i>Journal of the Taiwan Institute of Chemical Engineers</i> , <b>2009</b> , 40, 155-161	5.3	29	

11	Chemoinformatics and pharmacoinformatics approach for exploring the GABA-A agonist from Chinese herb suanzaoren. <i>Journal of the Taiwan Institute of Chemical Engineers</i> , <b>2009</b> , 40, 36-47	5.3	38
10	Ligand-based dual target drug design for H1N1: swine flua preliminary first study. <i>Journal of Biomolecular Structure and Dynamics</i> , <b>2009</b> , 27, 171-8	3.6	80
9	Weighted equation and rulesa novel concept for evaluating protein-ligand interaction. <i>Journal of Biomolecular Structure and Dynamics</i> , <b>2009</b> , 27, 271-82	3.6	65
8	Exploring 3D-QSAR pharmacophore mapping of azaphenanthrenone derivatives for mPGES-1 inhibition Using HypoGen technique <b>2008</b> ,		2
7	Antiplatelet effect and selective binding to cyclooxygenase by molecular docking analysis of 3-alkylaminopropoxy-9,10-anthraquinone derivatives. <i>Biological and Pharmaceutical Bulletin</i> , <b>2008</b> , 31, 1547-51	2.3	14
6	Inhibiting the vascular smooth muscle cells proliferation by EPC and DPPC liposomes encapsulated magnolol. <i>Journal of the Taiwan Institute of Chemical Engineers</i> , <b>2008</b> , 39, 407-411		15
5	A novel perspective on designing the inhibitor of HER2 receptor. <i>Journal of the Taiwan Institute of Chemical Engineers</i> , <b>2008</b> , 39, 291-299		27
4	Insights into the suanzaoren mechanism <b>E</b> rom constructing the 3D structure of GABA-A receptor to its binding interaction analysis. <i>Journal of the Taiwan Institute of Chemical Engineers</i> , <b>2008</b> , 39, 663-67	71	33
3	Discovery of novel inhibitors for c-Met by virtual screening and pharmacophore analysis. <i>Journal of the Taiwan Institute of Chemical Engineers</i> , <b>2008</b> , 39, 617-624		34
2	Antiplatelet Effect and Selective Binding to Cyclooxygenase (COX) by Molecular Docking Analysis of Flavonoids and Lignans. <i>International Journal of Molecular Sciences</i> , <b>2007</b> , 8, 830-841	6.3	22
1	Novel selective inhibitors of hydroxyxanthone derivatives for human cyclooxygenase-2. <i>Acta Pharmacologica Sinica</i> , <b>2007</b> , 28, 2027-32	8	28