# Judith M Rollinger

List of Publications by Year
in descending order



| 123 | 8,525 | 101543 | 48315 |
|---|---|---|---|
| papers | citations | 36 h-index | 88 g-index |

| 129 | 129 | 129 | 11048 |
|---|---|---|---|
| all docs | docs citations | times ranked | citing authors |

| # | Article | IF | Citations |
|---|---|---|---|
| 1 | Natural products in drug discovery: advances and opportunities. Nature Reviews Drug Discovery, 2021, 20, 200-216. | 46.4 | 1,990 |
| 2 | Discovery and resupply of pharmacologically active plant-derived natural products: A review. Biotechnology Advances, 2015, 33, 1582-1614. | 11.7 | 1,871 |
| 3 | Natural product agonists of peroxisome proliferator-activated receptor gamma (PPARÎ³): a review. Biochemical Pharmacology, 2014, 92, 73-89. | 4.4 | 492 |
| 4 | Energy/Temperature Diagram and Compression Behavior of the Polymorphs of d-Mannitol. Journal of Pharmaceutical Sciences, 2000, 89, 457-468. | 3.3 | 306 |
| 5 | Acetylcholinesterase Inhibitory Activity of Scopolin and Scopoletin Discovered by Virtual Screening of Natural Products. Journal of Medicinal Chemistry, 2004, 47, 6248-6254. | 6.4 | 193 |
| 6 | Influenza neuraminidase: A druggable target for natural products. Natural Product Reports, 2012, 29, 11-36. | 10.3 | 146 |
| 7 | In silicoTarget Fishing for Rationalized Ligand Discovery Exemplified on Constituents ofRuta graveolens. Planta Medica, 2009, 75, 195-204. | 1.3 | 131 |
| 8 | European medicinal polypores â€“ A modern view on traditional uses. Journal of Ethnopharmacology, 2014, 154, 564-583. | 4.1 | 120 |
| 9 | Antiviral Potential and Molecular Insight into Neuraminidase Inhibiting Diarylheptanoids from Alpinia katsumadai. Journal of Medicinal Chemistry, 2010, 53, 778-786. | 6.4 | 114 |
| 10 | Honokiol: A non-adipogenic PPARÎ³ agonist from nature. Biochimica Et Biophysica Acta - General Subjects, 2013, 1830, 4813-4819. | 2.4 | 108 |
| 11 | Virtual screening for the discovery of bioactive natural products. , 2008, 65, 211-249. | | 94 |
| 12 | Strategies for Efficient Lead Structure Discovery from Natural Products. Current Medicinal Chemistry, 2006, 13, 1491-1507. | 2.4 | 89 |
| 13 | Carnosol and Carnosic Acids fromSalvia officinalisInhibit Microsomal Prostaglandin E<sub>2</sub>Synthase-1. Journal of Pharmacology and Experimental Therapeutics, 2012, 342, 169-176. | 2.5 | 84 |
| 14 | Structure-Based Virtual Screening for the Discovery of Natural Inhibitors for Human Rhinovirus Coat Protein. Journal of Medicinal Chemistry, 2008, 51, 842-851. | 6.4 | 83 |
| 15 | Computer-Aided Discovery, Validation, and Mechanistic Characterization of Novel Neolignan Activators of Peroxisome Proliferator-Activated Receptor Î³. Molecular Pharmacology, 2010, 77, 559-566. | 2.3 | 72 |
| 16 | Combining Ethnopharmacology and Virtual Screening for Lead Structure Discovery:â€‰ COX-Inhibitors as Application Example. Journal of Chemical Information and Computer Sciences, 2004, 44, 480-488. | 2.8 | 67 |
| 17 | The coumarin scopoletin potentiates acetylcholine release from synaptosomes, amplifies hippocampal long-term potentiation and ameliorates anticholinergic- and age-impaired memory. Neuroscience, 2011, 197, 280-292. | 2.3 | 65 |
| 18 | Discovery of prenylated flavonoids with dual activity against influenza virus and Streptococcus pneumoniae. Scientific Reports, 2016, 6, 27156. | 3.3 | 63 |

| # | Article | IF | Citations |
|---|---|---|---|
| 19 | Experimentally Validated hERG Pharmacophore Models as Cardiotoxicity Prediction Tools. Journal of Chemical Information and Modeling, 2014, 54, 2887-2901. | 5.4 | 62 |
| 20 | Identification of plumericin as a potent new inhibitor of the <scp>NF</scp>â€<scp>ÎºB</scp> pathway with antiâ€inflammatory activity in vitro and in vivo. British Journal of Pharmacology, 2014, 171, 1676-1686. | 5.4 | 61 |
| 21 | Pharmacophore-based discovery of FXR-agonists. Part II: Identification of bioactive triterpenes from Ganoderma lucidum. Bioorganic and Medicinal Chemistry, 2011, 19, 6779-6791. | 3.0 | 59 |
| 22 | Discovery of Depsides and Depsidones from Lichen as Potent Inhibitors of Microsomal Prostaglandin E2 Synthaseâ€1 Using Pharmacophore Models. ChemMedChem, 2012, 7, 2077-2081. | 3.2 | 58 |
| 23 | Taspine:â€‰Bioactivity-Guided Isolation and Molecular Ligandâ^'Target Insight of a Potent Acetylcholinesterase Inhibitor from Magnolia x soulangiana. Journal of Natural Products, 2006, 69, 1341-1346. | 3.0 | 57 |
| 24 | Lignan Derivatives from Krameria lappacea Roots Inhibit Acute Inflammation in Vivo and Pro-inflammatory Mediators in Vitro. Journal of Natural Products, 2011, 74, 1779-1786. | 3.0 | 56 |
| 25 | Impact of Molecular Flexibility on Double Polymorphism, Solid Solutions and Chiral Discrimination during Crystallization of Diprophylline Enantiomers. Molecular Pharmaceutics, 2013, 10, 3850-3861. | 4.6 | 55 |
| 26 | The human rhinovirus: humanâ€pathological impact, mechanisms of antirhinoviral agents, and strategies for their discovery. Medicinal Research Reviews, 2011, 31, 42-92. | 10.5 | 54 |
| 27 | Pharmacophore Modeling and Virtual Screening for Novel Acidic Inhibitors of Microsomal Prostaglandin E<sub>2</sub> Synthase-1 (mPGES-1). Journal of Medicinal Chemistry, 2011, 54, 3163-3174. | 6.4 | 53 |
| 28 | Discovering COX-Inhibiting Constituents ofMorusRoot Bark: Activity-Guided versus Computer-Aided Methods. Planta Medica, 2005, 71, 399-405. | 1.3 | 52 |
| 29 | Natural products modulating the hERG channel: heartaches and hope. Natural Product Reports, 2017, 34, 957-980. | 10.3 | 51 |
| 30 | 11Î²-Hydroxysteroid dehydrogenase 1 inhibiting constituents from Eriobotrya japonica revealed by bioactivity-guided isolation and computational approaches. Bioorganic and Medicinal Chemistry, 2010, 18, 1507-1515. | 3.0 | 50 |
| 31 | Discovery of a novel IKK-Î² inhibitor by ligand-based virtual screening techniques. Bioorganic and Medicinal Chemistry Letters, 2011, 21, 577-583. | 2.2 | 50 |
| 32 | Accessing target information by virtual parallel screeningâ€"The impact on natural product research. Phytochemistry Letters, 2009, 2, 53-58. | 1.2 | 47 |
| 33 | Morphinans and isoquinolines: Acetylcholinesterase inhibition, pharmacophore modeling, and interaction with opioid receptors. Bioorganic and Medicinal Chemistry, 2010, 18, 5071-5080. | 3.0 | 46 |
| 34 | Pharmacophore-based discovery of FXR agonists. Part I: Model development and experimental validation. Bioorganic and Medicinal Chemistry, 2011, 19, 7168-7180. | 3.0 | 46 |
| 35 | Polymorphism of racemic felodipine and the unusual series of solid solutions in the binary system of its enantiomers. Journal of Pharmaceutical Sciences, 2001, 90, 949-959. | 3.3 | 45 |
| 36 | Binary System of (R)- and (S)-Nitrendipineâ€"Polymorphism and Structure. Journal of Pharmaceutical Sciences, 1997, 86, 674-679. | 3.3 | 40 |

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|---|---|---|---|
| 37 | Computer-Guided Approach to Access the Anti-influenza Activity of Licorice Constituents. Journal of Natural Products, 2014, 77, 563-570. | 3.0 | 38 |
| 38 | Dual Acting Neuraminidase Inhibitors Open New Opportunities to Disrupt the Lethal Synergism between Streptococcus pneumoniae and Influenza Virus. Frontiers in Microbiology, 2016, 7, 357. | 3.5 | 38 |
| 39 | Crystal forms of torasemide: new insights. European Journal of Pharmaceutics and Biopharmaceutics, 2002, 53, 75-86. | 4.3 | 37 |
| 40 | Discovery of Novel CB<sub>2</sub> Receptor Ligands by a Pharmacophore-Based Virtual Screening Workflow. Journal of Medicinal Chemistry, 2009, 52, 369-378. | 6.4 | 37 |
| 41 | Anti-Influenza Triterpene Saponins from the Bark of Burkea africana. Journal of Natural Products, 2018, 81, 515-523. | 3.0 | 37 |
| 42 | The crystal polymorphs of metazachlor. Journal of Thermal Analysis and Calorimetry, 2004, 77, 511-522. | 3.6 | 36 |
| 43 | Pharmacophore Model Refinement for 11Î²â€Hydroxysteroid Dehydrogenase Inhibitors: Search for Modulators of Intracellular Glucocorticoid Concentrations. Molecular Informatics, 2014, 33, 15-25. | 2.5 | 35 |
| 44 | Novel neuraminidase inhibitors: identification, biological evaluation and investigations of the binding mode. Future Medicinal Chemistry, 2011, 3, 437-450. | 2.3 | 34 |
| 45 | Predicting Cyclooxygenase Inhibition by Threeâ€Đimensional Pharmacophoric Profiling. Part I: Model Generation, Validation and Applicability in Ethnopharmacology. Molecular Informatics, 2010, 29, 75-86. | 2.5 | 33 |
| 46 | Physico-chemical Characterization of Hydrated and Anhydrous Crystal Forms of Amlodipine Besylate. Magyar AprÃ³vad KÃ¶zlemÃ©nyek, 2002, 68, 361-372. | 1.4 | 32 |
| 47 | Integrated in Silico Tools for Exploiting the Natural Productsâ€™ Bioactivity. Planta Medica, 2006, 72, 671-678. | 1.3 | 32 |
| 48 | Extracts and constituents of Leontopodium alpinum enhance cholinergic transmission: Brain ACh increasing and memory improving properties. Biochemical Pharmacology, 2008, 76, 236-248. | 4.4 | 32 |
| 49 | Inhibition of 11Î²-hydroxysteroid dehydrogenase type 1 by plant extracts used as traditional antidiabetic medicines. FÃ¬toterapÃ¬Ã¢, 2009, 80, 200-205. | 2.2 | 32 |
| 50 | Antipneumococcal activity of neuraminidase inhibiting artocarpin. International Journal of Medical Microbiology, 2015, 305, 289-297. | 3.6 | 32 |
| 51 | Natural products against acute respiratory infections: Strategies and lessons learned. Journal of Ethnopharmacology, 2020, 248, 112298. | 4.1 | 32 |
| 52 | Accessing biological actions of Ganoderma secondary metabolites by in silico profiling. Phytochemistry, 2015, 114, 114-124. | 2.9 | 31 |
| 53 | Imbricaric Acid and Perlatolic Acid: Multi-Targeting Anti-Inflammatory Depsides from Cetrelia monachorum. PLoS ONE, 2013, 8, e76929. | 2.5 | 30 |
| 54 | Anti-cancer Drug Development: Computational Strategies to Identify and Target Proteins Involved in Cancer Metabolism. Current Pharmaceutical Design, 2013, 19, 532-577. | 1.9 | 30 |

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| 55 | 1H NMR-MS-based heterocovariance as a drug discovery tool for fishing bioactive compounds out of a complex mixture of structural analogues. Scientific Reports, 2019, 9, 11113. | 3.3 | 28 |
| 56 | High-performance thin-layer chromatography/bioautography and liquid chromatography-mass spectrometry hyphenated with chemometrics for the quality assessment of Morus alba samples. Journal of Chromatography A, 2019, 1594, 190-198. | 3.7 | 28 |
| 57 | Neoandrographolide fromAndrographis paniculataas a Potential Natural Chemosensitizer. Planta Medica, 2010, 76, 1698-1700. | 1.3 | 27 |
| 58 | Lignans, Phenylpropanoids and Polyacetylenes from Chaerophyllum aureum L. (Apiaceae). Zeitschrift Fur Naturforschung - Section C Journal of Biosciences, 2003, 58, 553-557. | 1.4 | 25 |
| 59 | Dereplication of depsides from the lichen Pseudevernia furfuracea by centrifugal partition chromatography combined to 13C nuclear magnetic resonance pattern recognition. Analytica Chimica Acta, 2014, 846, 60-67. | 5.4 | 25 |
| 60 | Ginkgolic Acid is a Multi-Target Inhibitor of Key Enzymes in Pro-Inflammatory Lipid Mediator Biosynthesis. Frontiers in Pharmacology, 2019, 10, 797. | 3.5 | 25 |
| 61 | Application of the In Combo Screening Approach For the Discovery of Non-Alkaloid Acetylcholinesterase Inhibitors from Cichorium intybus. Current Drug Discovery Technologies, 2005, 2, 185-193. | 1.2 | 24 |
| 62 | Interface dynamics explain assembly dependency of influenza neuraminidase catalytic activity. Journal of Biomolecular Structure and Dynamics, 2015, 33, 104-120. | 3.5 | 24 |
| 63 | Allspice and Clove As Source of Triterpene Acids Activating the G Protein-Coupled Bile Acid Receptor TGR5. Frontiers in Pharmacology, 2017, 8, 468. | 3.5 | 24 |
| 64 | Lignans from Carthamus tinctorius suppress tryptophan breakdown via indoleamine 2,3-dioxygenase. Phytomedicine, 2013, 20, 1190-1195. | 5.3 | 23 |
| 65 | Human Ether-Ã-go-go Related Gene (hERG) Channel Blocking Aporphine Alkaloids from Lotus Leaves and Their Quantitative Analysis in Dietary Weight Loss Supplements. Journal of Agricultural and Food Chemistry, 2015, 63, 5634-5639. | 5.2 | 23 |
| 66 | Complementary assays helping to overcome challenges for identifying neuraminidase inhibitors. Future Virology, 2015, 10, 77-88. | 1.8 | 23 |
| 67 | 2-(2,4-dihydroxyphenyl)-5-(E)-propenylbenzofuran promotes endothelial nitric oxide synthase activity in human endothelial cells. Biochemical Pharmacology, 2012, 84, 804-812. | 4.4 | 22 |
| 68 | Pistacia lentiscus Oleoresin: Virtual Screening and Identification of Masticadienonic and Isomasticadienonic Acids as Inhibitors of 11β-Hydroxysteroid Dehydrogenase 1. Planta Medica, 2015, 81, 525-532. | 1.3 | 22 |
| 69 | Drugs from nature targeting inflammation (DNTI): a successful Austrian interdisciplinary network project. Monatshefte FÃ¼r Chemie, 2016, 147, 479-491. | 1.8 | 22 |
| 70 | Preparative supercritical fluid chromatography for lipid class fractionationâ€"a novel strategy in high-resolution mass spectrometry based lipidomics. Analytical and Bioanalytical Chemistry, 2020, 412, 2365-2374. | 3.7 | 22 |
| 71 | Natural Products in Structure-Assisted Design of Molecular Cancer Therapeutics. Current Pharmaceutical Design, 2010, 16, 1718-1741. | 1.9 | 20 |
| 72 | hERG Channel Blocking Ipecac Alkaloids Identified by Combined In Silico â€“ In Vitro Screening. Planta Medica, 2016, 82, 1009-1015. | 1.3 | 20 |

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| 73 | Podospermic acid, 1,3,5-tri-O-(7,8-dihydrocaffeoyl)quinic acid from Podospermum laciniatum (Asteraceae). Tetrahedron Letters, 2005, 46, 1291-1294. | 1.4 | 19 |
| 74 | Applications of Integrated Data Mining Methods to Exploring Natural Product Space for Acetylcholinesterase Inhibitors. Combinatorial Chemistry and High Throughput Screening, 2010, 13, 54-66. | 1.1 | 18 |
| 75 | Ratanhiaphenol III from Ratanhiae Radix is a PTP1B Inhibitor. Planta Medica, 2012, 78, 678-681. | 1.3 | 18 |
| 76 | Discovery of Bioactive Natural Products for the Treatment of Acute Respiratory Infections â€“ An Integrated Approach. Planta Medica, 2018, 84, 684-695. | 1.3 | 18 |
| 77 | A robust and miniaturized screening platform to study natural products affecting metabolism and survival in Caenorhabditis elegans. Scientific Reports, 2020, 10, 12323. | 3.3 | 18 |
| 78 | Venturia inaequalis-Inhibiting Dielsâ^'Alder Adducts from Morus Root Bark. Journal of Agricultural and Food Chemistry, 2006, 54, 8432-8436. | 5.2 | 17 |
| 79 | Biological Activity of Flavonoids and Rare Sesquiterpene Lactones Isolated From Centaurea ragusina L.. Frontiers in Pharmacology, 2018, 9, 972. | 3.5 | 17 |
| 80 | Rapid analytical approach for bioprofiling compounds with radical scavenging and antimicrobial activities from seaweeds. Food Chemistry, 2021, 334, 127562. | 8.2 | 17 |
| 81 | An evaluation of the transition temperature range of super-elastic orthodontic NiTi springs using differential scanning calorimetry. European Journal of Orthodontics, 1999, 21, 497-502. | 2.4 | 16 |
| 82 | In Silico Workflow for the Discovery of Natural Products Activating the G Protein-Coupled Bile Acid Receptor 1. Frontiers in Chemistry, 2018, 6, 242. | 3.6 | 16 |
| 83 | A Biochemometric Approach for the Identification of In Vitro Anti-Inflammatory Constituents in Masterwort. Biomolecules, 2020, 10, 679. | 4.0 | 16 |
| 84 | Thermal characterization of torasemide using coupled techniques. Journal of Thermal Analysis and Calorimetry, 2003, 73, 519-526. | 3.6 | 15 |
| 85 | Predicting cyclooxygenase inhibition by three-dimensional pharmacophoric profiling. Part II: Identification of enzyme inhibitors from Prasaplai, a Thai traditional medicine. Phytomedicine, 2011, 18, 119-133. | 5.3 | 15 |
| 86 | Plumeridoid C from the Amazonian traditional medicinal plantHimatanthus sucuuba. Acta Crystallographica Section C: Crystal Structure Communications, 2011, 67, o409-o412. | 0.4 | 14 |
| 87 | Piperine Congeners as Inhibitors of Vascular Smooth Muscle Cell Proliferation. Planta Medica, 2015, 81, 1065-1074. | 1.3 | 14 |
| 88 | Streptomyces spp. From Ethiopia Producing Antimicrobial Compounds: Characterization via Bioassays, Genome Analyses, and Mass Spectrometry. Frontiers in Microbiology, 2018, 9, 1270. | 3.5 | 14 |
| 89 | New Insights into the Acetylcholinesterase Inhibitory Activity ofLycopodium clavatum. Planta Medica, 2005, 71, 1040-1043. | 1.3 | 13 |
| 90 | Pharmacophore-based discovery of a novel cytosolic phospholipase A2Î± inhibitor. Bioorganic and Medicinal Chemistry Letters, 2012, 22, 1202-1207. | 2.2 | 13 |

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| 91 | Fast and Green â€“ CO2 Based Extraction, Isolation, and Quantification of Phenolic Styrax Constituents. Planta Medica, 2017, 83, 1068-1075. | 1.3 | 13 |
| 92 | Discovery and Characterization of Diazenylaryl Sulfonic Acids as Inhibitors of Viral and Bacterial Neuraminidases. Frontiers in Microbiology, 2017, 8, 205. | 3.5 | 13 |
| 93 | Steroid sulfatase inhibiting lanostane triterpenes â€“ Structure activity relationship and in silico insights. Bioorganic Chemistry, 2020, 95, 103495. | 4.1 | 11 |
| 94 | Peucedanum ostruthium Inhibits E-Selectin and VCAM-1 Expression in Endothelial Cells through Interference with NF-ÎºB Signaling. Biomolecules, 2020, 10, 1215. | 4.0 | 10 |
| 95 | Lanostane Triterpenes from Gloeophyllum odoratum and Their Anti-Influenza Effects. Planta Medica, 2019, 85, 195-202. | 1.3 | 9 |
| 96 | Insights into the direct anti-influenza virus mode of action of Rhodiola rosea. Phytomedicine, 2022, 96, 153895. | 5.3 | 9 |
| 97 | Anti-cancer drug development: computational strategies to identify and target proteins involved in cancer metabolism. Current Pharmaceutical Design, 2013, 19, 532-77. | 1.9 | 9 |
| 98 | Effects of the coumarin scopoletin on learning and memory, on release of acetylcholine from brain synaptosomes and on long-term potentiation in hippocampus. BMC Pharmacology, 2008, 8, A36. | 0.4 | 8 |
| 99 | In silico discovery of acylated flavonol monorhamnosides from Eriobotrya japonica as natural, small-molecular weight inhibitors of XIAP BIR3. Bioorganic and Medicinal Chemistry, 2011, 19, 1002-1009. | 3.0 | 8 |
| 100 | Discovery of Sanggenon G as a natural cellâ€permeable smallâ€molecular weight inhibitor of Xâ€linked inhibitor of apoptosis protein (XIAP). FEBS Open Bio, 2014, 4, 659-671. | 2.3 | 8 |
| 101 | Black pepper dietary supplementation increases high-density lipoprotein (HDL) levels in pigs. Current Research in Biotechnology, 2019, 1, 28-33. | 3.7 | 8 |
| 102 | Virtual Screening for the Discovery of Active Principles from Natural Products. , 2018, , 333-364. | | 7 |
| 103 | Pharmacophore-Based Screening for the Successful Identification of Bio-Active Natural Products. Chimia, 2007, 61, 350-354. | 0.6 | 6 |
| 104 | Plant extracts in cell-based anti-inflammatory assaysâ€”Pitfalls and considerations related to removal of activity masking bulk components. Phytochemistry Letters, 2014, 10, xli-xlvii. | 1.2 | 6 |
| 105 | Editorial: Ethnopharmacology in Central and Eastern Europe in the Context of Global Research Developments. Frontiers in Pharmacology, 2019, 10, 341. | 3.5 | 5 |
| 106 | High-performance Countercurrent Chromatography to Access Rhodiola rosea Influenza Virus Inhibiting Constituents. Planta Medica, 2021, 87, 818-826. | 1.3 | 5 |
| 107 | Quantitative Analysis of Prenylated Constituents in Commercial Hops Samples Using Ultrahigh-Performance Supercritical Fluid Chromatography. Planta Medica, 2020, 86, 1140-1147. | 1.3 | 5 |
| 108 | Pharmacokinetics of hERG Channel Blocking Voacangine in Wistar Rats Applying a Validated LC-ESI-MS/MS Method. Planta Medica, 2016, 82, 1030-1038. | 1.3 | 4 |

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| 109 | PPARÎ³ transcription effect on naturally occurring O-prenyl cinnamaldehydes and cinnamyl alcohol derivatives. Future Medicinal Chemistry, 2021, 13, 1175-1183. | 2.3 | 3 |
| 110 | A Strength-Weaknesses-Opportunities-Threats (SWOT) Analysis of Cheminformatics in Natural Product Research. Progress in the Chemistry of Organic Natural Products, 2019, 110, 239-271. | 1.1 | 3 |
| 111 | Editorial [Hot topic: Computational Techniques for Lead Discovery from Nature (Executive Editor:) Tj ETQq1 1 0.784314 rgBT /Overlock | 1.9 | 2 |
| 112 | Workflow for Segmentation of Caenorhabditis elegans from Fluorescence Images for the Quantitation of Lipids. Applied Sciences (Switzerland), 2021, 11, 11420. | 2.5 | 2 |
| 113 | (-)-N,N'-but-2-ene-1,4-diylbimorphinans. Die Pharmazie, 1991, 46, 101-2. | 0.5 | 2 |
| 114 | Azepine-Indole Alkaloids From Psychotria nemorosa Modulate 5-HT2A Receptors and Prevent in vivo Protein Toxicity in Transgenic Caenorhabditis elegans. Frontiers in Neuroscience, 2022, 16, 826289. | 2.8 | 2 |
| 115 | Computational Approaches for the Discovery of Natural Lead Structures. , 2011, , 97-132. | | 1 |
| 116 | In silico Driven Pharmacognosy: Forth, Back and Reverse. Planta Medica, 2015, 81, 427-428. | 1.3 | 1 |
| 117 | Pterocarpus santalinus Selectively Inhibits a Subset of Pro-Inflammatory Genes in Interleukin-1 Stimulated Endothelial Cells. Frontiers in Pharmacology, 2021, 12, 802153. | 3.5 | 1 |
| 118 | Expanding the Biological Properties of Alkannins and Shikonins: Their Impact on Adipogenesis and Life Expectancy in Nematodes. Frontiers in Pharmacology, 0, 13, . | 3.5 | 1 |
| 119 | Combining Ethnopharmacology and Virtual Screening for Lead Structure Discovery: COX-Inhibitors as Application Example.. ChemInform, 2004, 35, no. | 0.0 | 0 |
| 120 | In silicoTarget Fishing for Rationalized Ligand Discovery Exemplified on Constituents ofRuta graveolensL.. Planta Medica, 2009, 75, 293-293. | 1.3 | 0 |
| 121 | Special Issue Dedicated to Professor Dr. Max Wichtl. Planta Medica, 2017, 83, 1108-1109. | 1.3 | 0 |
| 122 | Special Issue Dedicated to Professor Dr. Max Wichtl. Planta Medica, 2017, 83, 960-961. | 1.3 | 0 |
| 123 | Biochemometry-Based Discovery of Phenylpropanoids from Azadirachta indica Fruits as Inhibitors of In Vitro Osteoclast Formation. Molecules, 2022, 27, 3611. | 3.8 | 0 |