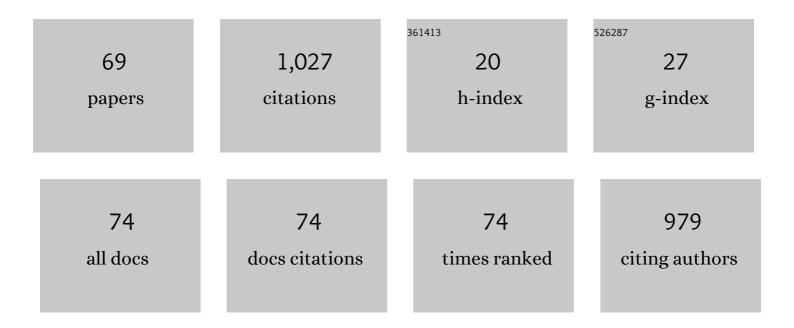
Cristiano Bolchi

List of Publications by Year in descending order

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CRISTIANO ROLCHI

| # | Article | IF | CITATIONS |
|----|---|-------------------|-------------|
| 1 | α9―and α7â€containing receptors mediate the proâ€proliferative effects of nicotine in the A549 adenocarcinoma cell line. British Journal of Pharmacology, 2018, 175, 1957-1972. | 5.4 | 61 |
| 2 | A short entry to enantiopure 2-substituted 1,4-benzodioxanes by efficient resolution methods. Tetrahedron: Asymmetry, 2003, 14, 3779-3785. | 1.8 | 47 |
| 3 | Benzodioxane–benzamides as new bacterial cell division inhibitors. European Journal of Medicinal Chemistry, 2015, 89, 252-265. | 5.5 | 45 |
| 4 | Structure–affinity studies for a novel series of homochiral naphtho and tetrahydronaphtho analogues of α1 antagonist WB-4101. Bioorganic and Medicinal Chemistry, 2004, 12, 4937-4951. | 3.0 | 38 |
| 5 | WB4101-Related Compounds:  New, Subtype-Selective α1-Adrenoreceptor Antagonists (or Inverse) Tj ETQq | 1 1 0.7843 6.4 | 314 rgBT /O |
| 6 | 3-(Benzodioxan-2-ylmethoxy)-2,6-difluorobenzamides bearing hydrophobic substituents at the 7-position of the benzodioxane nucleus potently inhibit methicillin-resistant Sa and Mtb cell division. European Journal of Medicinal Chemistry, 2016, 120, 227-243. | 5.5 | 28 |
| 7 | QSAR study for a novel series of ortho monosubstituted phenoxy analogues of α1-adrenoceptor antagonist WB4101. Bioorganic and Medicinal Chemistry, 2005, 13, 2547-2559. | 3.0 | 26 |
| 8 | Highly efficient resolutions of 1,4-benzodioxane-2-carboxylic acid with para substituted 1-phenylethylamines. Tetrahedron: Asymmetry, 2005, 16, 1639-1643. | 1.8 | 26 |
| 9 | Unichiral 2-(2′-Pyrrolidinyl)-1,4-benzodioxanes: the 2 <i>R</i> ,2′ <i>S</i> Diastereomer of the <i>N</i> -Methyl-7-hydroxy Analogue Is a Potent α4β2- and α6β2-Nicotinic Acetylcholine Receptor Partial Agonist. Journal of Medicinal Chemistry, 2011, 54, 7588-7601. | 6.4 | 26 |
| 10 | Predicting the physicochemical profile of diastereoisomeric histidineâ€containing dipeptides by property space analysis. Chirality, 2012, 24, 566-576. | 2.6 | 26 |
| 11 | 6-Methoxy-7-benzofuranoxy and 6-Methoxy-7-indolyloxy Analogues of 2-[2-(2,6-Dimethoxyphenoxy)ethyl]aminomethyl-1,4-benzodioxane (WB4101):1 Discovery of a Potent and Selective α _{1D} -Adrenoceptor Antagonist. Journal of Medicinal Chemistry, 2013, 56, 6402-6412. | 6.4 | 25 |
| 12 | Ten Years of Fentanyl-like Drugs: a Technical-analytical Review. Analytical Sciences, 2019, 35, 479-491. | 1.6 | 25 |
| 13 | QSAR study forÂaÂnovel series ofÂortho disubstituted phenoxy analogues ofÂα1-adrenoceptor antagonist WB4101. European Journal of Medicinal Chemistry, 2006, 41, 1025-1040. | 5.5 | 24 |
| 14 | One-Pot Racemization Process of 1-Phenyl-1,2,3,4-tetrahydroisoquinoline: A Key Intermediate for the Antimuscarinic Agent Solifenacin. Organic Process Research and Development, 2013, 17, 432-437. | 2.7 | 24 |
| 15 | Chemistry and Pharmacology of a Series of Unichiral Analogues of 2-(2-Pyrrolidinyl)-1,4-benzodioxane, Prolinol Phenyl Ether, and Prolinol 3-Pyridyl Ether Designed as α4β2-Nicotinic Acetylcholine Receptor Agonists. Journal of Medicinal Chemistry, 2015, 58, 6665-6677. | 6.4 | 24 |
| 16 | Resolution of 5-hydroxymethyl-2-oxazolidinone by preferential crystallization and investigations on the nature of the racemates of some 2-oxazolidinone derivatives. Tetrahedron: Asymmetry, 2004, 15, 1659-1665. | 1.8 | 22 |
| 17 | Resolution of 2-substituted 1,4-benzodioxanes by entrainment. Tetrahedron: Asymmetry, 2007, 18, 1038-1041. | 1.8 | 21 |
| 18 | Affinity and activity profiling of unichiral 8-substituted 1,4-benzodioxane analogues of WB4101 reveals a potent and selective α1B-adrenoceptor antagonist. European Journal of Medicinal Chemistry, 2012, 58, 184-191. | 5.5 | 21 |

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|----|--|-----|-----------|
| 19 | Potent Antiglioblastoma Agents by Hybridizing the Onium-Alkyloxy-Stilbene Based Structures of an α7-nAChR, I±9-nAChR Antagonist and of a Pro-Oxidant Mitocan. Journal of Medicinal Chemistry, 2018, 61, 10531-10544. | 6.4 | 21 |
| 20 | Peptidomimetic inhibitors of farnesyltransferase with high in vitro activity and significant cellular potency. Bioorganic and Medicinal Chemistry Letters, 2007, 17, 6192-6196. | 2.2 | 20 |
| 21 | Thiazole- and imidazole-containing peptidomimetic inhibitors of protein farnesyltransferase. Bioorganic and Medicinal Chemistry Letters, 2011, 21, 5408-5412. | 2.2 | 20 |
| 22 | Synthesis and α4β2 nicotinic affinity of unichiral 5-(2-pyrrolidinyl)oxazolidinones and 2-(2-pyrrolidinyl)benzodioxanes. Bioorganic and Medicinal Chemistry Letters, 2006, 16, 5610-5615. | 2.2 | 19 |
| 23 | 5-(2-Pyrrolidinyl)oxazolidinones and 2-(2-pyrrolidinyl)benzodioxanes: Synthesis of all the stereoisomers and α4β2 nicotinic affinity. Bioorganic and Medicinal Chemistry Letters, 2009, 19, 854-859. | 2.2 | 18 |
| 24 | From 2-Aminomethyl-1,4-benzodioxane Enantiomers to Unichiral 2-Cyano- and 2-Carbonyl-Substituted Benzodioxanes via Dichloroamine. Journal of Organic Chemistry, 2014, 79, 6732-6737. | 3.2 | 17 |
| 25 | 1,4-Benzodioxane, an evergreen, versatile scaffold in medicinal chemistry: A review of its recent applications in drug design. European Journal of Medicinal Chemistry, 2020, 200, 112419. | 5.5 | 17 |
| 26 | Influence of (S)-1-phenylethylamine para substitution on the resolution of (±)-1,4-benzodioxane-2-carboxylic acid: a crystallographic, theoretical and morphologic approach. Tetrahedron: Asymmetry, 2005, 16, 2099-2106. | 1.8 | 16 |
| 27 | Modelling of full-length human α4β2 nicotinic receptor by fragmental approach and analysis of its binding modes. Biochemical and Biophysical Research Communications, 2008, 369, 648-653. | 2.1 | 16 |
| 28 | Highly efficient racemisation of a key intermediate of the antibiotic moxifloxacin. Tetrahedron: Asymmetry, 2011, 22, 379-380. | 1.8 | 16 |
| 29 | Farnesyltransferase inhibitors: CAAX mimetics based on different biaryl scaffolds. Bioorganic and Medicinal Chemistry Letters, 2014, 24, 2924-2927. | 2.2 | 15 |
| 30 | From pyrrolidinyl-benzodioxane to pyrrolidinyl-pyridodioxanes, or from unselective antagonism to selective partial agonism at α4β2 nicotinic acetylcholine receptor. European Journal of Medicinal Chemistry, 2017, 125, 1132-1144. | 5.5 | 15 |
| 31 | Influence of Ionization State on the Activation of Temocapril by hCES1: A Molecularâ€Dynamics Study. Chemistry and Biodiversity, 2009, 6, 2092-2100. | 2.1 | 14 |
| 32 | Liver and intestinal protective effects of Castanea sativa Mill. bark extract in high-fat diet rats. PLoS ONE, 2018, 13, e0201540. | 2.5 | 14 |
| 33 | Synthesis and α4β2 nicotinic affinity of 2-pyrrolidinylmethoxyimines and prolinal oxime ethers. Bioorganic and Medicinal Chemistry Letters, 2004, 14, 5827-5830. | 2.2 | 12 |
| 34 | New Ras CAAX mimetics: Design, synthesis, antiproliferative activity, and RAS prenylation inhibition. Bioorganic and Medicinal Chemistry Letters, 2009, 19, 5500-5504. | 2.2 | 12 |
| 35 | Modifications at C(5) of 2-(2-Pyrrolidinyl)-Substituted 1,4-Benzodioxane Elicit Potent α4β2 Nicotinic Acetylcholine Receptor Partial Agonism with High Selectivity over the α3β4 Subtype. Journal of Medicinal Chemistry, 2020, 63, 15668-15692. | 6.4 | 12 |
| 36 | Univocal syntheses of 2- and 3-hydroxymethyl-2,3-dihydro[1,4]dioxino[2,3-b]pyridine enantiomers. Tetrahedron: Asymmetry, 2005, 16, 3380-3384. | 1.8 | 11 |

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|----|--|-----|-----------|
| 37 | Entrainment resolution of carnitinamide chloride. Tetrahedron: Asymmetry, 2008, 19, 1637-1640. | 1.8 | 11 |
| 38 | Antihypertensive phytocomplexes of proven efficacy and well-established use: Mode of action and individual characterization of the active constituents. Phytochemistry, 2020, 170, 112222. | 2.9 | 11 |
| 39 | Behavioural and pharmacological profiles of zebrafish administrated pyrrolidinyl benzodioxanes and prolinol aryl ethers with high affinity for heteromeric nicotinic acetylcholine receptors. Psychopharmacology, 2020, 237, 2317-2326. | 3.1 | 11 |
| 40 | A highly efficient method for the α,β-dehydrogenation of α-amino esters and α-amino-β-diesters. Tetrahedron Letters, 2010, 51, 5540-5542. | 1.4 | 10 |
| 41 | Enantiomer systems of carnitinamide inorganic salts: introductory studies to a successful entrainment resolution. Tetrahedron: Asymmetry, 2007, 18, 906-909. | 1.8 | 9 |
| 42 | Evidence of a dual mechanism of action underlying the anti-proliferative and cytotoxic effects of ammonium-alkyloxy-stilbene-based α7- and α9-nicotinic ligands on glioblastoma cells. Pharmacological Research, 2022, 175, 105959. | 7.1 | 9 |
| 43 | From 2-Triethylammonium Ethyl Ether of 4-Stilbenol (MG624) to Selective Small-Molecule Antagonists of Human α9α10 Nicotinic Receptor by Modifications at the Ammonium Ethyl Residue. Journal of Medicinal Chemistry, 2022, 65, 10079-10097. | 6.4 | 9 |
| 44 | From carnitinamide to 5-aminomethyl-2-oxazolidinones. Tetrahedron: Asymmetry, 2012, 23, 217-220. | 1.8 | 8 |
| 45 | A selective alpha1D-adrenoreceptor antagonist inhibits human prostate cancer cell proliferation and motility "in vitro― Pharmacological Research, 2016, 103, 215-226. | 7.1 | 8 |
| 46 | One-step preparation of enantiopure l- or d-amino acid benzyl esters avoiding the use of banned solvents. Amino Acids, 2017, 49, 965-974. | 2.7 | 8 |
| 47 | Determination of Methyldibromoglutaronitrile (MDBGN) in Skin Care Products by Gaschromatographyâ€Mass Spectrometry Employing an Enhanced Matrix Removal (EMR) Lipid Cleanâ€Up. European Journal of Lipid Science and Technology, 2018, 120, 1700525. | 1.5 | 8 |
| 48 | Efficient Oneâ€Pot Reductive Aminations of Carbonyl Compounds with Aquivionâ€Fe as a Recyclable Catalyst and Sodium Borohydride. European Journal of Organic Chemistry, 2020, 2020, 162-168. | 2.4 | 8 |
| 49 | Characterization of chemotype-dependent terpenoids profile in cannabis by headspace gas-chromatography coupled to time-of-flight mass spectrometry. Journal of Pharmaceutical and Biomedical Analysis, 2021, 203, 114180. | 2.8 | 8 |
| 50 | Highly efficient resolutions with isopropylidene glycerol 3-carboxy-2-naphthoate. Tetrahedron: Asymmetry, 2002, 13, 2277-2282. | 1.8 | 7 |
| 51 | Diastereomeric 2-aminomethyl-1,4-benzodioxane mandelates: phase diagrams and resolution. Tetrahedron: Asymmetry, 2013, 24, 796-800. | 1.8 | 7 |
| 52 | Enantiomerically Pure Dibenzyl Esters of <scp>l</scp> -Aspartic and <scp>l</scp> -Glutamic Acid. Organic Process Research and Development, 2015, 19, 878-883. | 2.7 | 7 |
| 53 | Novel 5-substituted 3-hydroxyphenyl and 3-nitrophenyl ethers of S -prolinol as α4β2-nicotinic acetylcholine receptor ligands. Bioorganic and Medicinal Chemistry Letters, 2016, 26, 5613-5617. | 2.2 | 7 |
| 54 | Determinants for α4β2 vs. α3β4 Subtype Selectivity of Pyrrolidine-Based nAChRs Ligands: A Computational Perspective with Focus on Recent cryo-EM Receptor Structures. Molecules, 2021, 26, 3603. | 3.8 | 7 |

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| # | Article | lF | CITATIONS |
|----|---|-----|-----------|
| 55 | Resolution of 1-phenyl-2-(p-tolyl)ethylamine via diastereomeric salt formation. Tetrahedron: Asymmetry, 2003, 14, 2247-2251. | 1.8 | 6 |
| 56 | Crystallization-based resolution of 1,4-benzodioxane-2-carboxylic acid enantiomers via diastereomeric 1-phenylethylamides. Tetrahedron Letters, 2016, 57, 2009-2011. | 1.4 | 6 |
| 57 | Green Oxidation of Ketones to Lactones with Oxone in Water. Journal of Organic Chemistry, 2021, 86, 15712-15716. | 3.2 | 6 |
| 58 | Preparation of enantiopure methionine, arginine, tryptophan, and proline benzyl esters in green ethers by Fischer–Speier reaction. Amino Acids, 2018, 50, 1261-1268. | 2.7 | 5 |
| 59 | Design, synthesis and binding affinity of acetylcholine carbamoyl analogues. Bioorganic and Medicinal Chemistry Letters, 2013, 23, 6481-6485. | 2.2 | 4 |
| 60 | Preparation and unequivocal identification of the regioisomers of nitrocatechol monobenzyl ether. Synthetic Communications, 2017, 47, 1507-1513. | 2.1 | 4 |
| 61 | Castanea sativa Mill. bark extract cardiovascular effects in a rat model of highâ€fat diet. Phytotherapy Research, 2021, 35, 2145-2156. | 5.8 | 4 |
| 62 | Pyrrolidinyl benzofurans and benzodioxanes: Selective α4β2 nicotinic acetylcholine receptor ligands with different activity profiles at the two receptor stoichiometries. Bioorganic and Medicinal Chemistry Letters, 2022, 65, 128701. | 2.2 | 4 |
| 63 | Development and Early Identification of Cannabis Chemotypes during the Plant Growth: Current Analytical and Chemometric Approaches. Analytical Sciences, 2021, 37, 1665-1673. | 1.6 | 3 |
| 64 | Synthesis of α-Hydroxy Fatty Acids from Fatty Acids by Intermediate α-Chlorination with TCCA under Solvent-Free Conditions: A Way to Valorization of Waste Fat Biomasses. ACS Omega, 2021, 6, 31901-31906. | 3.5 | 3 |
| 65 | Phase Diagrams to Evaluate the Opportunity for Enantiomeric Enrichment of Some Nonracemic Mixtures of Amino Acid Benzyl Esters by Crystallization as p-Toluenesulfonate Salts. Organic Process Research and Development, 2017, 21, 1752-1757. | 2.7 | 2 |
| 66 | Simple route to synthesize (<i>E</i>)-3-propyl-4-oxo-2-butenoic acid esters through the <i>Z</i> isomer. Synthetic Communications, 2018, 48, 85-90. | 2.1 | 1 |
| 67 | 1H NMR spectroscopy in the presence of Mosher acid to rapidly determine the enantiomeric composition of amino acid benzyl esters, chirons susceptible to easy racemization. Amino Acids, 2018, 50, 1759-1767. | 2.7 | 1 |
| 68 | Efficient conversion of d-mannitol into 1,2:5,6-diacetonide with Aquivion-H as a recyclable catalyst. Carbohydrate Research, 2021, 499, 108229. | 2.3 | 1 |
| 69 | Simple Process for the Preparation of Cetyltrimethylammonium Naproxenate (Naprocet). Organic Process Research and Development, 2014, 18, 976-979. | 2.7 | 0 |