Donatella Potenza

List of Publications by Year in descending order

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218592 276775 2,124 78 26 41 h-index citations g-index papers 81 81 81 2119 docs citations times ranked citing authors all docs

#	Article	IF	CITATIONS
1	Synthesis of Sulfonamido-Pseudopeptides: New Chiral Unnatural Oligomers. Angewandte Chemie International Edition in English, 1994, 33, 2067-2069.	4.4	142
2	A Simple Model System for the Study of Carbohydrateâ°'Aromatic Interactions. Journal of the American Chemical Society, 2007, 129, 2890-2900.	6.6	98
3	Conformational Preferences of Peptides Containing Reverse-Turn Mimetic Bicyclic Lactams: Inverse γ-Turns versus Type-ll′ β-Turns – Insights into β-Hairpin Stability. , 1999, 1999, 389-400.		92
4	Stereoselective synthesis of statin analogues Tetrahedron Letters, 1990, 31, 4949-4952.	0.7	78
5	Biological and molecular properties of a new $\hat{l}\pm v\hat{l}^23/\hat{l}\pm v\hat{l}^25$ integrin antagonist. Molecular Cancer Therapeutics, 2005, 4, 1670-1680.	1.9	75
6	Intramolecular Carbohydrate-Aromatic Interactions and Intermolecular van der Waals Interactions Enhance the Molecular Recognition Ability of GM1 Glycomimetics for Cholera Toxin. Chemistry - A European Journal, 2004, 10, 4395-4406.	1.7	69
7	Cyclic RGD Peptidomimetics Containing Bifunctional Diketopiperazine Scaffolds as New Potent Integrin Ligands. Chemistry - A European Journal, 2012, 18, 6195-6207.	1.7	62
8	A new method for the solution and solid phase synthesis of chiral \hat{l}^2 -sulfonopeptides under mild conditions. Tetrahedron Letters, 1996, 37, 8589-8592.	0.7	61
9	Targeting integrins: Insights into structure and activity of cyclic RGD pentapeptide mimics containing azabicycloalkane amino acids. Bioorganic and Medicinal Chemistry, 2006, 14, 169-180.	1.4	61
10	Synthesis of Novel DCâ€SIGN Ligands with an αâ€Fucosylamide Anchor. ChemBioChem, 2008, 9, 1921-1930.	1.3	58
11	Cyclic RGDâ€Peptidomimetics Containing Bifunctional Diketopiperazine Scaffolds as New Potent Integrin Ligands. Chemistry - A European Journal, 2009, 15, 12184-12188.	1.7	58
12	Synthesis and Conformational Analysis of Small Peptides Containing 6-Endo-BT(t)L Scaffolds as Reverse Turn Mimetics. Journal of Organic Chemistry, 2002, 67, 7483-7492.	1.7	51
13	Conformational Studies of Chiral Vinylogous Sulfonamidopeptides. Chemistry - A European Journal, 1996, 2, 644-655.	1.7	50
14	Potent Integrin Antagonists from a Small Library of RGD-Including Cyclic Pseudopeptides. Organic Letters, 2001, 3, 1001-1004.	2.4	49
15	Cyclic RGDâ€Containing Functionalized Azabicycloalkane Peptides as Potent Integrin Antagonists for Tumor Targeting. ChemMedChem, 2009, 4, 615-632.	1.6	44
16	Conformational Analysis of Azabicycloalkane Amino Acid Scaffolds as Reverse-Turn Inducer Dipeptide Mimics. European Journal of Organic Chemistry, 2000, 2000, 2563-2569.	1.2	43
17	STD and trNOESY NMR Study of Receptor–Ligand Interactions in Living Cancer Cells. ChemBioChem, 2011, 12, 695-699.	1.3	39
18	Rational design, synthesis and characterization of potent, non-peptidic Smac mimics/XIAP inhibitors as proapoptotic agents for cancer therapy. Bioorganic and Medicinal Chemistry, 2009, 17, 5834-5856.	1.4	36

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19	Comprehensive analysis of blood group antigen binding to classical and El Tor cholera toxin B-pentamers by NMR. Glycobiology, 2014, 24, 766-778.	1.3	34
20	Stereoselective synthesis of t-butyl 2-amino-2,5-dideoxypentanoate: Formal synthesis of l-daunosamine. Tetrahedron, 1987, 43, 2317-2322.	1.0	32
21	Second-Generation Mimics of Ganglioside GM1 Oligosaccharide: A Three-Dimensional View of Their Interactions with Bacterial Enterotoxins by NMR and Computational Methods. Chemistry - A European Journal, 2002, 8, 4597-4612.	1.7	31
22	Mimics of ganglioside GM1 as cholera toxin ligands: replacement of the GalNAc residueElectronic supplementary information (ESI) available: synthetic details, product characterisations and full NOE contact list. See http://www.rsc.org/suppdata/ob/b2/b210503a/. Organic and Biomolecular Chemistry, 2003, 1, 785-792.	1.5	31
23	Synthesis, Conformational Studies and Mannosidase Stability of a Mimic of 1,2-Mannobioside. European Journal of Organic Chemistry, 2004, 2004, 5119-5225.	1.2	29
24	Dimeric Smac mimetics/IAP inhibitors as in vivo-active pro-apoptotic agents. Part II: Structural and biological characterization. Bioorganic and Medicinal Chemistry, 2012, 20, 6709-6723.	1.4	29
25	A New Isoflavone fromGenistacorsica. Journal of Natural Products, 2000, 63, 504-506.	1.5	28
26	Cyclic <i>iso</i> DGR Peptidomimetics as Lowâ€Nanomolar α _v β ₃ Integrin Ligands. Chemistry - A European Journal, 2013, 19, 3563-3567.	1.7	28
27	A Potent Integrin Antagonist from a Small Library of Cyclic RGD Pentapeptide Mimics Including Benzylâ€ 5 ubstituted Azabicycloalkane Amino Acids. ChemMedChem, 2008, 3, 1589-1603.	1.6	27
28	Stereoselective aldol reactions using ticl4 as stereochemical template. Tetrahedron Letters, 1985, 26, 4129-4132.	0.7	26
29	Structure Revision of the Lantibiotic 97518. Journal of Natural Products, 2009, 72, 605-607.	1.5	25
30	Transferred-NOE NMR experiments on intact human platelets: receptor-bound conformation of RGD-peptide mimics. Organic and Biomolecular Chemistry, 2008, 6, 258-262.	1.5	24
31	Synthese von Sulfonamidâ€Pseudopeptiden: neue chirale synthetische Oligomere. Angewandte Chemie, 1994, 106, 2181-2183.	1.6	23
32	Lupane-triterpenes fromBupleurum flavum. Natural Product Research, 2005, 19, 783-788.	1.0	22
33	Determination of the binding epitope of RGD-peptidomimetics to $\hat{l}\pm v\hat{l}^2$ 3 and $\hat{l}\pm IIb\hat{l}^2$ 3 integrin-rich intact cells by NMR and computational studies. Organic and Biomolecular Chemistry, 2013, 11, 3886.	1.5	22
34	Biosynthesis of citrinin and synthesis of its biogenetic precursors. Journal of the Chemical Society Perkin Transactions 1, 1981, , 2594.	0.9	21
35	Insights into the Binding of Cyclic RGD Peptidomimetics to \hat{l}_{\pm} (sub>5 \hat{l}^{2} (sub>1Integrin by using Live-Cell NMR And Computational Studies. ChemistryOpen, 2017, 6, 128-136.	0.9	21
36	Novel Compounds Targeting the RNA-Binding Protein HuR. Structure-Based Design, Synthesis, and Interaction Studies. ACS Medicinal Chemistry Letters, 2019, 10, 615-620.	1.3	21

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37	Homo- and heterodimeric Smac mimetics/IAP inhibitors as in vivo-active pro-apoptotic agents. Part I: Synthesis. Bioorganic and Medicinal Chemistry, 2012, 20, 6687-6708.	1.4	20
38	(E,E)-10-(1,3-Dihydro-4,6-dihydroxy-7-methyl-3-oxoisobenzofuran-5-yl)4,8-dimethyldeca-4,8-dienoic acid: total synthesis and role in mycophenolic acid biosynthesis. Journal of the Chemical Society Chemical Communications, 1979, , 1021.	2.0	18
39	Stereoconvergent crotylstannane addition to nor-ephedrine-derived 2-methoxy oxazolidines. A clue towards a synclinal transition state geometry. Tetrahedron: Asymmetry, 1990, 1, 429-432.	1.8	18
40	Cyclic RGD Peptides Containing Azabicycloalkane Reverse-Turn Mimics. Helvetica Chimica Acta, 2002, 85, 4353-4368.	1.0	18
41	Synthesis and Characterization of Linkerâ€Armed Fucoseâ€Based Glycomimetics. European Journal of Organic Chemistry, 2013, 2013, 5303-5314.	1.2	18
42	Total synthesis of (±) pseudophrynamine A. Tetrahedron Letters, 1990, 31, 5661-5664.	0.7	17
43	Solid-Phase Synthesis of Peptides Containing Reverse-Turn Mimetic Bicyclic Lactams. , 1999, 1999, 379-388.		17
44	Enantioselective binding of dipeptides using acyclic receptors. Chemical Communications, 2001, , $1358-1359$.	2.2	17
45	Ganglioside GM1 mimics: lipophilic substituents improve affinity for cholera toxin. Bioorganic and Medicinal Chemistry Letters, 2003, 13, 3831-3834.	1.0	17
46	Antitumor Activity of a Novel Homodimeric SMAC Mimetic in Ovarian Carcinoma. Molecular Pharmaceutics, 2014, 11, 283-293.	2.3	17
47	NMR interaction studies of Neu5Ac- \hat{l} ±-(2,6)-Gal- \hat{l} ²-(1-4)-GlcNAc with influenza-virus hemagglutinin expressed in transfected human cells. Glycobiology, 2018, 28, 42-49.	1.3	17
48	Norephedrine derived oxazolidines as chiral acylating agents: An NMR study of the intermediate cations Tetrahedron, 1992, 48, 1343-1352.	1.0	16
49	Rationally Designed Bicyclic Lactams Control Different Turn Motifs and Folding Patterns in Hexapeptide Mimics., 2000, 2000, 695-699.		16
50	Cicloastragenol glycosides from astragalus verrucosus. Phytochemistry, 1998, 49, 2467-2471.	1.4	15
51	Thermodynamically–Weighted Conformational Ensemble of Cyclic RGD Peptidomimetics from NOE Data. Journal of Physical Chemistry B, 2016, 120, 7098-7107.	1.2	15
52	Absolute configuration of A-32'287 [conocandin] and total synthesis of its methyl and tert-butyl esters. Journal of Organic Chemistry, 1987, 52, 5452-5457.	1.7	14
53	A new alpinumisoflavone derivative from Genista pichisermolliana. Phytochemistry Letters, 2011, 4, 342-344.	0.6	14
54	A Combined NMR-Computational Study of the Interaction between Influenza Virus Hemagglutinin and Sialic Derivatives from Human and Avian Receptors on the Surface of Transfected Cells. International Journal of Molecular Sciences, 2018, 19, 1267.	1.8	13

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55	Titanium Ate Enolate Complexes: An NMR Study. Journal of Organic Chemistry, 1994, 59, 3690-3694.	1.7	12
56	Synthesis of a Pseudo Tetrasaccharide Mimic of Ganglioside GM1. European Journal of Organic Chemistry, 1999, 1999, 1311-1317.	1.2	12
57	Solution structure by nuclear magnetic resonance of the two lantibiotics 97518 and NAIâ€107. Journal of Peptide Science, 2012, 18, 129-134.	0.8	12
58	Diffusionâ€Ordered Spectroscopy and Saturation Transfer Difference NMR Spectroscopy Studies of Selective Interactions between ELAV Protein Fragments and an mRNA Target. European Journal of Organic Chemistry, 2014, 2014, 6399-6404.	1.2	12
59	Exploration of ligand binding modes towards the identification of compounds targeting HuR: a combined STD-NMR and Molecular Modelling approach. Scientific Reports, 2018, 8, 13780.	1.6	12
60	Acceleration of hemiacetal cleavage through hydrogen bonding: a new synthetic catalyst with balanced conformational flexibility and preorganization. Journal of Organic Chemistry, 1991, 56, 3201-3203.	1.7	11
61	Chemical composition and volatile constituents of Anthyllis barba-jovis. Natural Product Research, 2007, 21, 418-425.	1.0	11
62	Cyclic DGR-peptidomimetic containing a bicyclic reverse turn inducer as a selective $\hat{l}\pm v\hat{l}^25$ integrin ligand. Amino Acids, 2010, 38, 329-337.	1.2	11
63	Insight to the binding mode of triazole RGD-peptidomimetics to integrin-rich cancer cells by NMR and molecular modeling. Bioorganic and Medicinal Chemistry, 2016, 24, 989-994.	1.4	11
64	A13C and 1H NMR study of diastereomeric \hat{l} ±-methylidene- \hat{l} 2-hydroxy- \hat{l} 3-alkoxy esters. Magnetic Resonance in Chemistry, 1984, 22, 224-227.	0.7	10
65	Solvent-Dependent Conformational Behaviour of Model Tetrapeptides Containing a Bicyclic Proline Mimetic. European Journal of Organic Chemistry, 2004, 2004, 4621-4627.	1.2	10
66	Computer aided design and NMR characterization of an oligopeptide targeting the Ebola virus VP24 protein. New Journal of Chemistry, 2017, 41, 4308-4315.	1.4	10
67	Synthesis, Conformational Studies and Binding Properties of Acyclic Receptors for N-Protected Amino Acids and Dipeptides. European Journal of Organic Chemistry, 2001, 2001, 4625.	1.2	9
68	A NMR and computational study of Smac mimics targeting both the BIR2 and BIR3 domains in XIAP protein. Organic and Biomolecular Chemistry, 2012, 10, 3278.	1.5	8
69	Solution Behavior of Amphiphilic Glycodendrimers with a Rodâ€Like Core. Macromolecular Bioscience, 2016, 16, 896-905.	2.1	8
70	Stereoselective conjugate addition of lithium and titanium enolates to \hat{l}^3 - alkoxy enones. Tetrahedron, 1996, 52, 3497-3508.	1.0	6
71	Exploring E-cadherin-peptidomimetics interaction using NMR and computational studies. PLoS Computational Biology, 2019, 15, e1007041.	1.5	5
72	Cyclic RGD and isoDGR Integrin Ligands Containing cis-2-amino-1-cyclopentanecarboxylic (cis-Î ² -ACPC) Scaffolds. Molecules, 2020, 25, 5966.	1.7	5

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73	Design, Synthesis, Conformational Analysis and Application of AzabicycloÂalkane Amino Acids as Constrained Dipeptide Mimics. Synlett, 2004, 2004, 1449-1471.	1.0	4
74	Selective Acylation of Monosaccharides Using Microbial Cells. Biocatalysis and Biotransformation, 1999, 17, 95-102.	1.1	3
7 5	Chiral 2â€phenylâ€3â€hydroxypropyl esters as PKCâ€alpha modulators: HPLC enantioseparation, NMR absolute configuration assignment, and molecular docking studies. Chirality, 2022, 34, 498-513.	1.3	2
76	Designing Antiviral Substances Targeting the Ebola Virus Viral Protein 24., 2020, , 147-177.		1
77	Hydrogen-Bonding Donor/Acceptor Scales in -Sulfonamidopeptides. Chemistry - A European Journal, 1998, 4, 1924-1931.	1.7	1
78	Towards the Modulation of RNA-Binding Proteins: New Compounds Targeting Protein HuR. Proceedings (mdpi), 2019, 22, .	0.2	0