Laura Belvisi

List of Publications by Year
in descending order

104	2,614	182225	286692
papers	citations	30 h-index	43 g-index

113	113	113	2837
all docs	docs citations	times ranked	citing authors

#	Article	IF	Citations
1	Advanced Pyrrolidineâ€Carbamate Selfâ€Immolative Spacer with Tertiary Amine Handle Induces Superfast Cyclative Drug Release. ChemMedChem, 2022, 17, .	1.6	5
2	Functionalized 2â€Hydroxybenzaldehydeâ€PEG Modules as Portable Tags for the Engagement of Protein Lysine Ïµâ€Amino Groups. European Journal of Organic Chemistry, 2021, 2021, 1763-1767.	1.2	1
3	Prediction and Validation of a Druggable Site on Virulence Factor of Drug Resistant Burkholderia cenocepacia**. Chemistry - A European Journal, 2021, 27, 10341-10348.	1.7	6
4	A trifunctional self-immolative spacer enables drug release with two non-sequential enzymatic cleavages. Chemical Communications, 2021, 57, 7778-7781.	2.2	7
5	Fast Cyclization of a Prolineâ€Derived Selfâ€Immolative Spacer Improves the Efficacy of Carbamate Prodrugs. Angewandte Chemie, 2020, 132, 4205-4210.	1.6	8
6	Fast Cyclization of a Prolineâ€Derived Selfâ€Immolative Spacer Improves the Efficacy of Carbamate Prodrugs. Angewandte Chemie - International Edition, 2020, 59, 4176-4181.	7.2	35
7	Cyclic RGD and isoDGR Integrin Ligands Containing cis-2-amino-1-cyclopentanecarboxylic (cis-Î²-ACPC) Scaffolds. Molecules, 2020, 25, 5966.	1.7	5
8	Side chain effect in the modulation of Î±vÎ²3/Î±5Î²1 integrin activity via clickable isoxazoline-RGD-mimetics: development of molecular delivery systems. Scientific Reports, 2020, 10, 7410.	1.6	4
9	Innovative Linker Strategies for Tumorâ€Targeted Drug Conjugates. Chemistry - A European Journal, 2019, 25, 14740-14757.	1.7	68
10	Bromineâ€Promoted Glycosidation of Conformationally Superarmed Thioglycosides. Chemistry - A European Journal, 2019, 25, 11831-11836.	1.7	10
11	Exploring E-cadherin-peptidomimetics interaction using NMR and computational studies. PLoS Computational Biology, 2019, 15, e1007041.	1.5	5
12	Î²-Glucuronidase triggers extracellular MMAE release from an integrin-targeted conjugate. Organic and Biomolecular Chemistry, 2019, 17, 4705-4710.	1.5	14
13	Synthesis and Biological Evaluation of RGD and isoDGRâ€“Monomethyl Auristatin Conjugates Targeting Integrin Î±_VÎ²₃. ChemMedChem, 2019, 14, 938-942.	1.6	26
14	Rational Design of Antiangiogenic Helical Oligopeptides Targeting the Vascular Endothelial Growth Factor Receptors. Frontiers in Chemistry, 2019, 7, 170.	1.8	10
15	The Importance of Detail: How Differences in Ligand Structures Determine Distinct Functional Responses in Integrin Î± v Î² 3. Chemistry - A European Journal, 2019, 25, 5959-5970.	1.7	10
16	Stem-like Cancer Cells in a Dynamic 3D Culture System: A Model to Study Metastatic Cell Adhesion and Anti-cancer Drugs. Cells, 2019, 8, 1434.	1.8	27
17	A dimeric bicyclic RGD ligand displays enhanced integrin binding affinity and strong biological effects on U-373 MG glioblastoma cells. Organic and Biomolecular Chemistry, 2019, 17, 8913-8917.	1.5	4
18	Frontispiece: Innovative Linker Strategies for Tumorâ€Targeted Drug Conjugates. Chemistry - A European Journal, 2019, 25, .	1.7	0

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19	Neutrophil Elastase Promotes Linker Cleavage and Paclitaxel Release from an Integrinâ€Targeted Conjugate. Chemistry - A European Journal, 2019, 25, 1696-1700.	1.7	29
20	A critical assessment of force field accuracy against NMR data for cyclic peptides containing Î²-amino acids. Physical Chemistry Chemical Physics, 2018, 20, 15807-15816.	1.3	9
21	Tumor Targeting with an isoDGRâ€“Drug Conjugate. Chemistry - A European Journal, 2017, 23, 7910-7914.	1.7	17
22	Insights into the Binding of Cyclic RGD Peptidomimetics to Î±₅Î²₁Integrin by using Live-Cell NMR And Computational Studies. ChemistryOpen, 2017, 6, 128-136.	0.9	21
23	Targeting Integrin Î±_V Î²₃ with Theranostic RGD-Camptothecin Conjugates Bearing a Disulfide Linker: Biological Evaluation Reveals a Complex Scenario. ChemistrySelect, 2017, 2, 4759-4766.	0.7	14
24	Frontispiece: Multivalency Increases the Binding Strength of RGD Peptidomimeticâ€Paclitaxel Conjugates to Integrin Î±_VÎ²₃. Chemistry - A European Journal, 2017, 23, .	1.7	0
25	Multivalency Increases the Binding Strength of RGD Peptidomimeticâ€Paclitaxel Conjugates to Integrin Î±_VÎ²₃. Chemistry - A European Journal, 2017, 23, 14410-14415.	1.7	27
26	Investigating the Interaction of Cyclic RGD Peptidomimetics with Î±VÎ²6 Integrin by Biochemical and Molecular Docking Studies. Cancers, 2017, 9, 128.	1.7	18
27	High Affinity vs. Native Fibronectin in the Modulation of Î±vÎ²3 Integrin Conformational Dynamics: Insights from Computational Analyses and Implications for Molecular Design. PLoS Computational Biology, 2017, 13, e1005334.	1.5	12
28	Crystal Structure of Human E-Cadherin-EC1EC2 in Complex with a Peptidomimetic Competitive Inhibitor of Cadherin Homophilic Interaction. Journal of Medicinal Chemistry, 2016, 59, 5089-5094.	2.9	17
29	New Î²-Lactam Derivatives Modulate Cell Adhesion and Signaling Mediated by RGD-Binding and Leukocyte Integrins. Journal of Medicinal Chemistry, 2016, 59, 9721-9742.	2.9	43
30	Thermodynamicallyâ€“Weighted Conformational Ensemble of Cyclic RGD Peptidomimetics from NOE Data. Journal of Physical Chemistry B, 2016, 120, 7098-7107.	1.2	15
31	New potent Î±_vÎ²₃ integrin ligands based on azabicycloalkane (Î³,Î±)-dipeptide mimics. Organic and Biomolecular Chemistry, 2016, 14, 3221-3233.	1.5	4
32	Synthesis, Characterization, and Biological Evaluation of a Dualâ€Action Ligand Targeting Î±_vÎ²₃ Integrin and VEGF Receptors. ChemistryOpen, 2015, 4, 633-641.	0.9	25
33	Synthesis of Easyâ€toâ€Functionalize AzaÂbicycloalkane Scaffolds as Dipeptide Turn Mimics en Route to cRGDâ€Based Bioconjugates. European Journal of Organic Chemistry, 2015, 2015, 7557-7570.	1.2	12
34	Metadynamics Simulations Rationalise the Conformational Effects Induced by Nâ€Methylation of RGD Cyclic Hexapeptides. Chemistry - A European Journal, 2015, 21, 14165-14170.	1.7	20
35	?v?3 Integrin-Targeted Peptide/Peptidomimetic-Drug Conjugates: In-Depth Analysis of the Linker Technology. Current Topics in Medicinal Chemistry, 2015, 16, 314-329.	1.0	44
36	Cyclic isoDGR and RGD Peptidomimetics Containing Bifunctional Diketopiperazine Scaffolds are Integrin Antagonists. Chemistry - A European Journal, 2015, 21, 6265-6271.	1.7	33

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37	Designing nanomolar antagonists of DC-SIGN-mediated HIV infection: ligand presentation using molecular rods. Chemical Communications, 2015, 51, 3816-3819.	2.2	74
38	New Insights into the Molecular Mechanism of E-Cadherin-Mediated Cell Adhesion by Free Energy Calculations. Journal of Chemical Theory and Computation, 2015, 11, 1354-1359.	2.3	7
39	Synthesis and Biological Evaluation of RGD Peptidomimeticâ€“Paclitaxel Conjugates Bearing Lysosomally Cleavable Linkers. Chemistry - A European Journal, 2015, 21, 6921-6929.	1.7	48
40	Computational design of novel peptidomimetic inhibitors of cadherin homophilic interactions. Organic and Biomolecular Chemistry, 2015, 13, 2570-2573.	1.5	16
41	Antitumor Activity of a Novel Homodimeric SMAC Mimetic in Ovarian Carcinoma. Molecular Pharmaceutics, 2014, 11, 283-293.	2.3	17
42	Cyclic isoDGR Peptidomimetics as Lowâ€Nanomolar Î±_vÎ²₃ Integrin Ligands. Chemistry - A European Journal, 2013, 19, 3563-3567.	1.7	28
43	Determination of the binding epitope of RGD-peptidomimetics to Î±vÎ²3 and Î±IIbÎ²3 integrin-rich intact cells by NMR and computational studies. Organic and Biomolecular Chemistry, 2013, 11, 3886.	1.5	22
44	Modulation of Î±vÎ²3- and Î±5Î²1-integrin-mediated adhesion by dehydro-Î²-amino acids containing peptidomimetics. European Journal of Medicinal Chemistry, 2013, 66, 258-268.	2.6	19
45	Modulation of Î±vÎ²3- and Î±5Î²1-integrin-mediated adhesion by dehydro-Î²-amino acids containing peptidomimetics. , 2013, 66, 258-258.		1
46	Synthesis and Biological Evaluation (in Vitro and in Vivo) of Cyclic Arginineâ€“Glycineâ€“Aspartate (RGD) Peptidomimeticâ€“Paclitaxel Conjugates Targeting Integrin Î±_VÎ²₃. Journal of Medicinal Chemistry, 2012, 55, 10460-10474.	2.9	68
47	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
48	Design, Synthesis, and Biological Evaluation of Novel cRGDâ€“Paclitaxel Conjugates for Integrin-Assisted Drug Delivery. Bioconjugate Chemistry, 2012, 23, 1610-1622.	1.8	41
49	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
50	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
51	Cyclic RGD Peptidomimetics Containing Bifunctional Diketopiperazine Scaffolds as New Potent Integrin Ligands. Chemistry - A European Journal, 2012, 18, 6195-6207.	1.7	62
52	Inside Cover: Cyclic RGD Peptidomimetics Containing Bifunctional Diketopiperazine Scaffolds as New Potent Integrin Ligands (Chem. Eur. J. 20/2012). Chemistry - A European Journal, 2012, 18, 6106-6106.	1.7	0
53	Synthesis of Gd and ⁶⁸Ga Complexes in Conjugation with a Conformationally Optimized RGD Sequence as Potential MRI and PET Tumorâ€Imaging Probes. ChemMedChem, 2012, 7, 1084-1093.	1.6	53
54	Second generation of fucose-based DC-SIGN ligands : affinity improvement and specificity versus Langerin. Organic and Biomolecular Chemistry, 2011, 9, 5778.	1.5	60

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55	Development of Isoxazolineâ€Containing Peptidomimetics as Dual Î±_vÎ²₃ and Î±₅Î²₁ Integrin Ligands. ChemMedChem, 2011, 6, 2264-2272.	1.6	22
56	A new optical imaging probe targeting Î±_VÎ²₃ integrin in glioblastoma xenografts. Contrast Media and Molecular Imaging, 2011, 6, 449-458.	0.4	39
57	STD and trNOESY NMR Study of Receptorâ€“Ligand Interactions in Living Cancer Cells. ChemBioChem, 2011, 12, 695-699.	1.3	39
58	Foldamers of bifunctional diketopiperazines displaying a Î²-bend ribbon structure. Tetrahedron Letters, 2010, 51, 4278-4280.	0.7	15
59	Antiangiogenic Effect of Dual/Selective Î±₅Î²₁/Î±_vÎ²₃ Integrin Antagonists Designed on Partially Modified Retro-Inverso Cyclotetrapeptide Mimetics. Journal of Medicinal Chemistry, 2010, 53, 106-118.	2.9	29
60	Cyclic RGDâ€Peptidomimetics Containing Bifunctional Diketopiperazine Scaffolds as New Potent Integrin Ligands. Chemistry - A European Journal, 2009, 15, 12184-12188.	1.7	58
61	Cyclic RGDâ€Containing Functionalized Azabicycloalkane Peptides as Potent Integrin Antagonists for Tumor Targeting. ChemMedChem, 2009, 4, 615-632.	1.6	44
62	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
63	A Potent Integrin Antagonist from a Small Library of Cyclic RGD Pentapeptide Mimics Including Benzylâ€Substituted Azabicycloalkane Amino Acids. ChemMedChem, 2008, 3, 1589-1603.	1.6	27
64	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
65	Synthesis and Conformational Studies of Peptidomimetics Containing a New Bifunctional Diketopiperazine Scaffold Acting as a Î²-Hairpin Inducer. Journal of Organic Chemistry, 2008, 73, 652-660.	1.7	47
66	Complete Characterization of Extracts of Onopordum illyricum L. (Asteraceae) by HPLC/PDA/ESIMS and NMR. Natural Product Communications, 2008, 3, 1934578X0800301.	0.2	5
67	Nonpeptide Integrin Antagonists: RGD Mimetics Incorporating Substituted Azabicycloalkanes as Amino Acid Replacements. European Journal of Organic Chemistry, 2007, 2007, 1309-1317.	1.2	10
68	Synthesis and conformational analysis of an Î±-d-mannopyranosyl-(1â†'2)-Î±-d-mannopyranosyl-(1â†'6)-Î±-d-mannopyranose mimic. Carbohydrate Research, 2007, 342, 1859-1868.	1.1	18
69	First round of a focused library of cholera toxin inhibitors. Carbohydrate Research, 2007, 342, 1651-1660.	1.1	18
70	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
71	Assessing the influence of electrostatic schemes on molecular dynamics simulations of secondary structure forming peptides. Journal of Physics Condensed Matter, 2006, 18, S329-S345.	0.7	20
72	Synthesis of Functionalized Azabicycloalkane Amino Acids as Dipeptide Mimics. Synthesis, 2006, 2006, 1133-1140.	1.2	5

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73	Grafting Aminocyclopentane Carboxylic Acids onto the RGD Tripeptide Sequence Generates Low Nanomolar Î±VÎ²3/Î±VÎ²5Integrin Dual Binders. Journal of Medicinal Chemistry, 2005, 48, 7675-7687.	2.9	49
74	Biological and molecular properties of a new Î±vÎ²3/Î±vÎ²5 integrin antagonist. Molecular Cancer Therapeutics, 2005, 4, 1670-1680.	1.9	75
75	Functionalized Azabicycloalkane Amino Acids by Nitrone 1,3-Dipolar Intramolecular Cycloaddition. Journal of Organic Chemistry, 2005, 70, 4124-4132.	1.7	39
76	Design, Synthesis, Conformational Analysis and Application of AzabicycloÂalkane Amino Acids as Constrained Dipeptide Mimics. Synlett, 2004, 2004, 1449-1471.	1.0	4
77	Stereoselective synthesis of CÎ±-tetrasubstituted azabicyclo[X.3.0]alkane amino acids. Tetrahedron Letters, 2004, 45, 6311-6315.	0.7	7
78	Synthesis of substituted conformationally constrained 6,5- and 7,5-fused bicyclic lactams as dipeptide mimics. Tetrahedron, 2003, 59, 6241-6250.	1.0	22
79	Structural investigations of isomeric oxidised forms of hyperforin by HPLC-NMR and HPLC-MSn. Phytochemical Analysis, 2003, 14, 290-297.	1.2	45
80	Cyclic RGD Peptides Containing Azabicycloalkane Reverse-Turn Mimics. Helvetica Chimica Acta, 2002, 85, 4353-4368.	1.0	18
81	Potent Integrin Antagonists from a Small Library of RGD-Including Cyclic Pseudopeptides. Organic Letters, 2001, 3, 1001-1004.	2.4	49
82	Practical stereoselective synthesis of conformationally constrained unnatural proline-based amino acids and peptidomimetics. Tetrahedron, 2001, 57, 6463-6473.	1.0	39
83	Simulation of carbohydrate-protein interactions: computer-aided design of a second generation GM1 mimic. Journal of Computer-Aided Molecular Design, 2001, 15, 117-128.	1.3	9
84	Title is missing!. Organic Letters, 2001, 3, 1001-1004.	2.4	2
85	Potent integrin antagonists from a small library of RGD-including cyclic pseudopeptides. Organic Letters, 2001, 3, 1001-4.	2.4	56
86	Rationally Designed Bicyclic Lactams Control Different Turn Motifs and Folding Patterns in Hexapeptide Mimics. , 2000, 2000, 695-699.		16
87	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
88	Synthesis of Azabicycloalkane Amino Acid Scaffolds as Reverse-Turn Inducer Dipeptide Mimics. European Journal of Organic Chemistry, 2000, 2000, 2571-2581.	1.2	27
89	Conformational Preferences of Peptides Containing Reverse-Turn Mimetic Bicyclic Lactams: Inverse Î³-Turns versus Type-IIâ€² Î²-Turns â€“ Insights into Î²-Hairpin Stability. , 1999, 1999, 389-400.		92
90	Stereoselective synthesis of 6,5-bicyclic reverse-turn peptidomimetics. Tetrahedron, 1998, 54, 5325-5336.	1.0	14

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91	Design and synthesis of nonpeptide angiotensin II receptor antagonists featuring acyclic imidazole-mimicking structural units. Bioorganic and Medicinal Chemistry, 1998, 6, 2013-2027.	1.4	4
92	Design and synthesis of conformationally constrained arginal thrombin inhibitors. Bioorganic and Medicinal Chemistry Letters, 1997, 7, 2205-2210.	1.0	15
93	A 3D QSAR CoMFA study of non-peptide angiotensin II receptor antagonists. Journal of Computer-Aided Molecular Design, 1996, 10, 567-582.	1.3	33
94	Conformationally constrained dipeptides: Synthesis of 7,5- and 6,5-fused bicyclic lactams by stereoselective radical cyclizations. Tetrahedron Letters, 1995, 36, 625-628.	0.7	46
95	Non peptide angiotensin II receptor antagonists: A 3D-QSAR comfa-like approach. AIP Conference Proceedings, 1995, , .	0.3	1
96	N-3-Substituted Pyrimidinones as Potent, Orally Active, AT1 Selective Angiotensin II Receptor Antagonists. Journal of Medicinal Chemistry, 1995, 38, 4806-4820.	2.9	59
97	A 3D QSAR approach to the search for geometrical similarity in a series of nonpeptide angiotensin II receptor antagonists. Journal of Computer-Aided Molecular Design, 1994, 8, 211-220.	1.3	15
98	A conformational study of N-tosyl oxazolidines using molecular mechanics and crystallography. Journal of Molecular Structure, 1994, 318, 189-202.	1.8	17
99	Nonpeptide Angiotensin II Receptor Antagonists. Synthesis, in vitro Activity, and Molecular Modeling Studies of N-[(Heterobiaryl)methyl]imidazoles. Journal of Medicinal Chemistry, 1994, 37, 3928-3938.	2.9	19
100	Stereoselective radical-mediated cyclization of norephedrine derived o-bromobenzamides: Enantioselective synthesis of 4-substituted 1,2,3,4-tetrahydroisoquinolines. Tetrahedron: Asymmetry, 1993, 4, 273-280.	1.8	25
101	On the role of the molecular electrostatic potential in modelling the activity of non-peptide angiotensin II receptor antagonists. Computational and Theoretical Chemistry, 1993, 281, 237-252.	1.5	6
102	The first asymmetric synthesis of enantiopure .alpha.-sulfenyl dithioacetals and .alpha.-sulfenyl aldehydes. Journal of Organic Chemistry, 1993, 58, 3165-3168.	1.7	33
103	Stereoselective radical-mediated cyclization of norephdrine derived Î±-iodoamides: synthesis of enantiopure pyrrolidines and trandition state modelling1. Tetrahedron, 1992, 48, 3945-3960.	1.0	32
104	Structure-activity relationship of Ca2+ channel blockers: A study using conformational analysis and chemometric methods. Journal of Computer-Aided Molecular Design, 1991, 5, 571-584.	1.3	8