

Paolo Benedetti

List of Publications by Year in descending order

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Version: 2024-02-01

18
papers

587
citations

687363

13
h-index

752698

20
g-index

20
all docs

20
docs citations

20
times ranked

1011
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|---|-----|-----------|
| 1 | First virtual screening and experimental validation of inhibitors targeting GES-5 carbapenemase. <i>Journal of Computer-Aided Molecular Design</i> , 2019, 33, 295-305. | 2.9 | 9 |
| 2 | Discovering New Casein Kinase 1d Inhibitors with an Innovative Molecular Dynamics Enabled Virtual Screening Workflow. <i>ACS Medicinal Chemistry Letters</i> , 2019, 10, 487-492. | 2.8 | 10 |
| 3 | Human dopamine transporter: the first implementation of a combined in silico/in vitro approach revealing the substrate and inhibitor specificities. <i>Journal of Biomolecular Structure and Dynamics</i> , 2019, 37, 291-306. | 3.5 | 8 |
| 4 | From Experiments to a Fast Easy-to-Use Computational Methodology to Predict Human Aldehyde Oxidase Selectivity and Metabolic Reactions. <i>Journal of Medicinal Chemistry</i> , 2018, 61, 360-371. | 6.4 | 29 |
| 5 | A Pipeline To Enhance Ligand Virtual Screening: Integrating Molecular Dynamics and Fingerprints for Ligand and Proteins. <i>Journal of Chemical Information and Modeling</i> , 2015, 55, 2256-2274. | 5.4 | 65 |
| 6 | Targeting Cystalyisin, a Virulence Factor of <i>Treponema denticola</i> Supported Periodontitis. <i>ChemMedChem</i> , 2014, 9, 1501-1511. | 3.2 | 26 |
| 7 | Exposition and reactivity optimization to predict sites of metabolism in chemicals. <i>Drug Discovery Today: Technologies</i> , 2013, 10, e155-e165. | 4.0 | 40 |
| 8 | Isozyme-Specific Ligands for O-acetylserine sulfhydrylase, a Novel Antibiotic Target. <i>PLoS ONE</i> , 2013, 8, e77558. | 2.5 | 43 |
| 9 | Inhibitor of Ovarian Cancer Cells Growth by Virtual Screening: A New Thiazole Derivative Targeting Human Thymidylate Synthase. <i>Journal of Medicinal Chemistry</i> , 2012, 55, 10272-10276. | 6.4 | 20 |
| 10 | FLAP: GRID Molecular Interaction Fields in Virtual Screening. Validation using the DUD Data Set. <i>Journal of Chemical Information and Modeling</i> , 2010, 50, 1442-1450. | 5.4 | 94 |
| 11 | Non-invasive identification of organic materials in wall paintings by fiber optic reflectance infrared spectroscopy: a statistical multivariate approach. <i>Analytical and Bioanalytical Chemistry</i> , 2009, 395, 2097-2106. | 3.7 | 70 |
| 12 | Fiber-Optic Fourier Transform Mid-Infrared Reflectance Spectroscopy: A Suitable Technique for in Situ Studies of Mural Paintings. <i>Applied Spectroscopy</i> , 2007, 61, 293-299. | 2.2 | 65 |
| 13 | Design, synthesis and in vitro antitumor activity of new trans 2-[2-(heteroaryl)vinyl]-1,3-dimethylimidazolium iodides. <i>Bioorganic and Medicinal Chemistry</i> , 2004, 12, 1689-1695. | 3.0 | 33 |
| 14 | GRIND/ALMOND investigations on CysLT1 receptor antagonists of the quinoliny(bridged)aryl type. <i>Bioorganic and Medicinal Chemistry</i> , 2004, 12, 3607-3617. | 3.0 | 22 |
| 15 | Structure-based rationalization of antitumor drugs mechanism of action by a MIF approach. <i>European Journal of Medicinal Chemistry</i> , 2004, 39, 281-289. | 5.5 | 18 |
| 16 | The CARSO procedure in process optimization. <i>Journal of Chemometrics</i> , 2003, 17, 9-15. | 1.3 | 2 |
| 17 | GBR Compounds and Mepyramines as Cocaine Abuse Therapeutics: A Chemometric Studies on Selectivity Using Grid Independent Descriptors (GRIND). <i>Journal of Medicinal Chemistry</i> , 2002, 45, 1577-1584. | 6.4 | 17 |
| 18 | Distant collaboration in drug discovery: the LINK3D project. <i>Journal of Computer-Aided Molecular Design</i> , 2002, 16, 809-818. | 2.9 | 4 |