# Subhendu Mukherjee

List of Publications by Year
in descending order



| 11 | 144 | 1040056 9 | 1281871 11 |
|---|---|---|---|
| papers | citations | h-index | g-index |

| 12 | 12 | 12 | 197 |
|---|---|---|---|
| all docs | docs citations | times ranked | citing authors |

| # | Article | IF | Citations |
|---|---|---|---|
| 1 | Discovery of O-(3-carbamimidoylphenyl)-l-serine amides as matriptase inhibitors using a fragment-linking approach. Bioorganic and Medicinal Chemistry Letters, 2015, 25, 616-620. | 2.2 | 11 |
| 2 | Structure-guided discovery of 2-aryl/pyridin-2-yl-1H-indole derivatives as potent and selective hepsin inhibitors. Bioorganic and Medicinal Chemistry Letters, 2015, 25, 5309-5314. | 2.2 | 10 |
| 3 | Structure-guided discovery of 1,3,5 tri-substituted benzenes as potent and selective matriptase inhibitors exhibiting in vivo antitumor efficacy. Bioorganic and Medicinal Chemistry, 2014, 22, 3187-3203. | 3.0 | 17 |
| 4 | Discovery of Pyridyl Bis(oxy)dibenzimidamide Derivatives as Selective Matriptase Inhibitors. ACS Medicinal Chemistry Letters, 2013, 4, 1152-1157. | 2.8 | 26 |
| 5 | Pharmacophore mapping of arylbenzothiophene derivatives for MCF cell inhibition using classical and 3D space modeling approaches. Journal of Molecular Graphics and Modelling, 2008, 26, 884-892. | 2.4 | 11 |
| 6 | Pharmacophore Mapping of Selective Binding Affinity of Estrogen Modulators through Classical and Space Modeling Approaches:â€‰ Exploration of Bridged-Cyclic Compounds with Diarylethylene Linkage. Journal of Chemical Information and Modeling, 2007, 47, 475-487. | 5.4 | 12 |
| 7 | Molecular Modeling Studies of Estrogen Receptor Modulators. Current Computer-Aided Drug Design, 2006, 2, 229-253. | 1.2 | 2 |
| 8 | QSAR Studies with E-State Index: Predicting Pharmacophore Signals for Estrogen Receptor Binding Affinity of Triphenylacrylonitriles. Biological and Pharmaceutical Bulletin, 2005, 28, 154-157. | 1.4 | 13 |
| 9 | QSAR modeling on binding affinity of diverse estrogenic flavonoids: electronic, topological and spatial functions in quantitative approximation. Computational and Theoretical Chemistry, 2005, 715, 85-90. | 1.5 | 13 |
| 10 | QSAR of estrogen receptor modulators: exploring selectivity requirements for ERÎ± versus ERÎ² binding of tetrahydroisoquinoline derivatives using E-state and physicochemical parameters. Bioorganic and Medicinal Chemistry Letters, 2005, 15, 957-961. | 2.2 | 20 |
| 11 | Predicting pharmacophore signals for post-coital antifertility activity of 1-trifluoromethyl-1,2,2-triphenylethylene derivatives: a statistical approximation using E-state index. Bioorganic and Medicinal Chemistry Letters, 2004, 14, 897-900. | 2.2 | 9 |