

A comprehensive map of molecular drug targets

Nature Reviews Drug Discovery

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Citation Report

#	ARTICLE	IF	CITATIONS
1	Classical Targets in Drug Discovery. , 2015, , 87-142.		0
2	Thermal proteome profiling: unbiased assessment of protein state through heat-induced stability changes. <i>Proteome Science</i> , 2016, 15, 13.	0.7	101
3	Pharos: Collating protein information to shed light on the druggable genome. <i>Nucleic Acids Research</i> , 2017, 45, D995-D1002.	6.5	271
4	Genetically encoded photocross-linkers determine the biological binding site of exendin-4 peptide in the N-terminal domain of the intact human glucagon-like peptide-1 receptor (GLP-1R). <i>Journal of Biological Chemistry</i> , 2017, 292, 7131-7144.	1.6	41
5	TIN-X: target importance and novelty explorer. <i>Bioinformatics</i> , 2017, 33, 2601-2603.	1.8	27
6	Protein epitope mimetic macrocycles as biopharmaceuticals. <i>Current Opinion in Chemical Biology</i> , 2017, 38, 45-51.	2.8	52
7	Harnessing public domain data to discover and validate therapeutic targets. <i>Expert Opinion on Drug Discovery</i> , 2017, 12, 687-693.	2.5	6
8	Spatial encryption of G protein-coupled receptor signaling in endosomes; Mechanisms and applications. <i>Biochemical Pharmacology</i> , 2017, 143, 1-9.	2.0	25
9	Drug repositioning beyond the low-hanging fruits. <i>Current Opinion in Systems Biology</i> , 2017, 3, 95-102.	1.3	11
10	Phenotypic chemical biology for predicting safety and efficacy. <i>Drug Discovery Today: Technologies</i> , 2017, 23, 53-60.	4.0	29
11	“Big data” approaches for novel anti-cancer drug discovery. <i>Expert Opinion on Drug Discovery</i> , 2017, 12, 599-609.	2.5	9
12	DrugCentral: online drug compendium. <i>Nucleic Acids Research</i> , 2017, 45, D932-D939.	6.5	215
13	Universal and Quantitative Method To Evaluate Inhibitor Potency for Cysteine Proteins Using a Nonspecific Activity-Based Protein Profiling Probe. <i>Biochemistry</i> , 2017, 56, 2921-2927.	1.2	3
14	Novel Structural Insights into GPCR ¹² -Arrestin Interaction and Signaling. <i>Trends in Cell Biology</i> , 2017, 27, 851-862.	3.6	90
15	A combinatorial screen of the CLOUD uncovers a synergy targeting the androgen receptor. <i>Nature Chemical Biology</i> , 2017, 13, 771-778.	3.9	39
16	A resource of potential drug targets and strategic decision-making for obstructive sleep apnoea pharmacotherapy. <i>Respirology</i> , 2017, 22, 861-873.	1.3	50
17	The use of biologics in the management of cardiovascular diseases. <i>Current Opinion in Pharmacology</i> , 2017, 33, 76-80.	1.7	3
18	The utility of target-based discovery. <i>Expert Opinion on Drug Discovery</i> , 2017, 12, 427-429.	2.5	80

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19	The druggable genome and support for target identification and validation in drug development. <i>Science Translational Medicine</i> , 2017, 9, .	5.8	437
20	Trends in GPCR drug discovery: new agents, targets and indications. <i>Nature Reviews Drug Discovery</i> , 2017, 16, 829-842.	21.5	1,773
21	Integrating human and environmental health in antibiotic risk assessment: A critical analysis of protection goals, species sensitivity and antimicrobial resistance. <i>Environment International</i> , 2017, 109, 155-169.	4.8	163
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23	Pharmacological <i>in vivo</i> classical approaches in the design of first in man clinical drug trials. <i>British Journal of Clinical Pharmacology</i> , 2017, 83, 2807-2812.	1.1	3
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32	Natural diversity facilitates the discovery of conserved chemotherapeutic response mechanisms. <i>Current Opinion in Genetics and Development</i> , 2017, 47, 41-47.	1.5	11
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38	Logic Modeling in Quantitative Systems Pharmacology. <i>CPT: Pharmacometrics and Systems Pharmacology</i> , 2017, 6, 499-511.	1.3	25
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