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10 years into the resurgence of covalent drugs

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45	Recent advances in the development of covalent inhibitors. <i>RSC Medicinal Chemistry</i> , 2021 , 12, 1037-10	1 45 .5	7
44	Early career research in medicinal chemistry. Future Medicinal Chemistry, 2021, 13, 91-93	4.1	0
43	Prioritization of antimicrobial targets by CRISPR-based oligo recombineering.		O
42	Design, Synthesis, and Structural Characterization of Lysine Covalent BH3 Peptides Targeting Mcl-1. <i>Journal of Medicinal Chemistry</i> , 2021 , 64, 4903-4912	8.3	5
41	Electrophilic Natural Products as Drug Discovery Tools. <i>Trends in Pharmacological Sciences</i> , 2021 , 42, 434-447	13.2	5
40	Covalent fragment screening. Annual Reports in Medicinal Chemistry, 2021, 56, 243-265	1.6	0
39	Covalent PROTACs: the best of both worlds?. <i>RSC Medicinal Chemistry</i> , 2021 , 12, 1452-1458	3.5	4
38	Mechanism of activation and the rewired network: New drug design concepts. <i>Medicinal Research Reviews</i> , 2021 ,	14.4	2
37	Proteasome Inhibitors and Their Pharmacokinetics, Pharmacodynamics, and Metabolism. <i>International Journal of Molecular Sciences</i> , 2021 , 22,	6.3	4
36	Discovery of 9,10-dihydrophenanthrene derivatives as SARS-CoV-2 3CL inhibitors for treating COVID-19. <i>European Journal of Medicinal Chemistry</i> , 2021 , 228, 114030	6.8	3
35	Welcome to Volume 14 of Future Medicinal Chemistry, 2021,	4.1	
34	High-Throughput Kinetic Characterization of Irreversible Covalent Inhibitors of KRAS by Intact Protein MS and Targeted MRM <i>Analytical Chemistry</i> , 2022 ,	7.8	1
33	Identification of the first structurally validated covalent ligands of the small GTPase RAB27A <i>RSC Medicinal Chemistry</i> , 2022 , 13, 150-155	3.5	O
32	Genetically encoding latent bioreactive amino acids and the development of covalent protein drugs <i>Current Opinion in Chemical Biology</i> , 2021 , 66, 102106	9.7	3
31	Chemical reactivity prediction: current methods and different application areas <i>Molecular Informatics</i> , 2021 ,	3.8	O
30	How neocarcerand Octacid4 self-assembles with guests into irreversible noncovalent complexes and what accelerates the assembly. <i>Communications Chemistry</i> , 2022 , 5,	6.3	
29	Inverse Drug Discovery identifies weak electrophiles affording protein conjugates <i>Current Opinion in Chemical Biology</i> , 2022 , 67, 102113	9.7	1

28	Electrophilic warheads in covalent drug discovery: an overview <i>Expert Opinion on Drug Discovery</i> , 2022 , 1-10	6.2	7
27	Promising reversible protein inhibitors kept on target <i>Nature</i> , 2022 ,	50.4	1
26	In situ identification of cellular drug targets in mammalian tissue Cell, 2022,	56.2	5
25	Therapeutic Targeting the Allosteric Cysteinome of RAS and Kinase Families. <i>Journal of Molecular Biology</i> , 2022 , 167626	6.5	О
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23	Pathobiology and Therapeutic Relevance of GSK-3 in Chronic Hematological Malignancies. <i>Cells</i> , 2022 , 11, 1812	7.9	1
22	Rational identification of small molecules derived from 9,10-dihydrophenanthrene as potential inhibitors of 3CLpro enzyme for COVID-19 therapy: a computer-aided drug design approach. Structural Chemistry,	1.8	3
21	N-Acylamino Saccharin as an Emerging Cysteine-Directed Covalent Warhead and Its Application in the Identification of Novel FBPase Inhibitors toward Glucose Reduction. <i>Journal of Medicinal Chemistry</i> , 2022 , 65, 9126-9143	8.3	
20	Phenotypic screening of low molecular weight compounds is rich ground for repurposed, on-target drugs. 13,		
19	Development of highly potent non-covalent inhibitors of SARS-CoV-2 3CLpro.		O
18	Tideglusib Inhibits Pif1 Helicase of Bacteroides sp. via an Irreversible and Cys-380-Dependent Mechanism.		
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