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P001	Corsini Martina	DESIGN AND SYNTHESIS OF A LIBRARY OF THALIDOMIDE-BASED PROTEOLYSIS TARGETING CHIMERAS (PROTACS) AS ANTIVIRAL AGENTS	I + II
P002	Albano Aurora	DEVELOPMENT OF INNOVATIVE COMPOUNDS TARGETING SARS-COV-2 NSP13	I
P003	Allahi Shahin	DESIGNING A NEW-GENERATION OF GLS1 ALLOSTERIC INHIBITORS BASED ON IN SILICO FRAGMENT-BASED STRATEGY	I
P004	Almerico Anna Maria	INHIBITION OF FTSJ1 BY 1,2,4-TRIAZOLE DERIVATIVES AS A MECHANISM FOR TRANSLATIONAL READTHROUGH: AN INTEGRATED COMPUTATIONAL AND EXPERIMENTAL STUDY	I
P005	Antonini Giulia	DEVELOPMENT OF PFKFB3 KINASE COVALENT INHIBITORS WITH ANTITUMOR ACTIVITY	I
P006	Antonini Giulia	EFFICIENT SYNTHESIS OF THE KINASE INHIBITOR NVS-PI3-4 USING FLOW CHEMISTRY, MICROWAVE IRRADIATION, AND IMMOBILIZED REAGENTS	I
P007	Arrighi Francesca	FROM MOLECULAR DESIGN TO NEUROPROTECTION: A JOURNEY THROUGH BENZO[B]THIOPHENE-BASED MAO-B INHIBITORS	I
P008	Astolfi Andrea	REPURPOSING MEDICINAL CHEMISTRY CONCEPTS FOR TARGET-BASED DISCOVERY OF ANTI-XYLELLA FASTIDIOSA AGROCHEMICALS	I
P009	Barakat Assem	FIRST-IN-CLASS SPIROOXINDOLE-SULFONAMIDE HYBRIDS AS TRIPLE-TARGET EGFR/CDK2/CA IX/XII INHIBITORS: POTENT ANTI-BREAST CANCER AGENTS WITH MULTIMODAL ACTION	I
P010	Barbarossa Alexia	MELATONIN-RELATED COMPOUNDS WITH ANTIOXIDANT AND NEUROPROTECTIVE PROPERTIES	I
P011	Basagni Filippo	EXPLORING DISULFIDE BRIDGE AS A TOOL TO IMPROVE PIOGLITAZONE'S NEUROPROTECTIVE POTENTIAL: TOWARD THE DEVELOPMENT OF PROLONGED-ACTING MAO-B/PPAR γ MODULATORS	I
P012	Bilotta Marina	INTEGRATED MACHINE LEARNING AND STRUCTURE-BASED DISCOVERY OF TANKYRASE 1 INHIBITORS	I
P013	Bono Alessia	B.R.E.A.S.T.: AN INTERPRETABLE MACHINE LEARNING FRAMEWORK FOR PROTEOMICS-DRIVEN THERAPY OUTCOME PREDICTION IN BREAST CANCER AND TRIPLE-NEGATIVE BREAST CANCER	I
P014	Campagna Erica	CATHEPSIN-B ACTIVABLE PRODRUGS OF HISTONE DEACETYLASE INHIBITORS	I
P015	Cannistra' Matilde S.F.	TRANSCRIPTIONAL REPRESSION OF HIV-1 VIA SMALL-MOLECULE STABILIZATION OF THE LTR-III G-QUADRUPLEX	I
P016	Cara Emanuele	IDENTIFICATION OF NOVEL NITROGEN-BASED DERIVATIVES TARGETING SARS-COV-2 RdRp	I

P017	Carbone Giorgio	GRIPkit: A COMPREHENSIVE COMPUTATIONAL SUITE FOR THE CHARACTERIZATION AND QUANTIFICATION OF NON-COVALENT INTERACTIONS IN PROTEIN-LIGAND COMPLEXES	I
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P020	Carradori Simone	COUMARIN- AND DIPICOLYLAMINE-TERPENOID HYBRIDS AS SELECTIVE CARBONIC ANHYDRASES IX AND XII INHIBITORS: MECHANISTIC IN-SIGHTS AND SELECTIVE ANTI-MYELOMA POTENTIAL	I
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