CITATION REPORT List of articles citing

Tracking cancer drugs in living cells by thermal profiling of the proteome

DOI: 10.1126/science.1255784 Science, 2014, 346, 1255784.

Source: https://exaly.com/paper-pdf/59579327/citation-report.pdf

Version: 2024-04-28

This report has been generated based on the citations recorded by exaly.com for the above article. For the latest version of this publication list, visit the link given above.

The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

#	Paper	IF	Citations
679	Nebenwirkungen von Krebsmedikamenten. 2014 , 9, 54-55		
678	Bringing the heat. 2014 , 10, 985-985		
677	Follow your heat. 2014 , 14, 706-706		
676	Know your target, know your molecule. 2015 , 11, 368-72		74
675	Advances in measuring single-cell pharmacology in vivo. 2015 , 20, 1087-92		26
674	2014: signaling breakthroughs of the year. 2015 , 8, eg1		
673	Inhibitors of protein methyltransferases as chemical tools. 2015 , 7, 1327-38		12
672	Hitting the target. 2015 , 12, 1127-8		
671	Target engagement and drug residence time can be observed in living cells with BRET. 2015 , 6, 10091		153
670	Joining time-resolved thermometry and magnetic-induced heating in a single nanoparticle unveils intriguing thermal properties. 2015 , 9, 3134-42		106
669	Extending kinome coverage by analysis of kinase inhibitor broad profiling data. 2015 , 20, 652-8		36
668	A Global Map of Lipid-Binding Proteins and Their Ligandability in Cells. 2015 , 161, 1668-80		140
667	Functional Identification of Target by Expression Proteomics (FITExP) reveals protein targets and highlights mechanisms of action of small molecule drugs. 2015 , 5, 11176		30
666	A Scalable Approach for Protein False Discovery Rate Estimation in Large Proteomic Data Sets. 2015 , 14, 2394-404		210
665	Global analysis of protein folding thermodynamics for disease state characterization. 2015 , 14, 2287-9	7	23
664	Multidimensional proteomics for cell biology. 2015 , 16, 269-80		288
663	CETSA: a target engagement assay with potential to transform drug discovery. 2015 , 7, 975-8		32

(2016-2015)

662	Proteome-wide drug and metabolite interaction mapping by thermal-stability profiling. 2015 , 12, 1055-7	145
661	Thermal proteome profiling monitors ligand interactions with cellular membrane proteins. 2015 , 12, 1129-31	173
660	Assessing the Efficacy of Mdm2/Mdm4-Inhibiting Stapled Peptides Using Cellular Thermal Shift Assays. 2015 , 5, 12116	32
659	CRKL oncogene is downregulated by p53 through miR-200s. 2015 , 106, 1033-40	18
658	Identification of drug candidates and repurposing opportunities through compound-target interaction networks. 2015 , 10, 1333-45	39
657	Limited Proteolysis Combined with Stable Isotope Labeling Reveals Conformational Changes in Protein (Pseudo)kinases upon Binding Small Molecules. 2015 , 14, 4179-93	4
656	Generation of multiple reporter ions from a single isobaric reagent increases multiplexing capacity for quantitative proteomics. 2015 , 87, 9855-63	34
655	Next-generation sequencing: hype and hope for development of personalized radiation therapy?. 2015 , 10, 183	8
654	Thermal proteome profiling for unbiased identification of direct and indirect drug targets using multiplexed quantitative mass spectrometry. 2015 , 10, 1567-93	293
653	Are label-free investigations the best approach to drug discovery?. 2015 , 7, 1561-4	10
652	A Small Molecule that Induces Intrinsic Pathway Apoptosis with Unparalleled Speed. 2015 , 13, 2027-36	48
651	Functional genomics to uncover drug mechanism of action. 2015 , 11, 942-8	58
650	Quantitative proteomics of kinase inhibitor targets and mechanisms. 2015 , 10, 201-12	23
649	Biomarker Development in Targeting Cancer Epigenetic. 2016 , 123-142	
648	Real-Time Biological Annotation of Synthetic Compounds. 2016 , 138, 8920-7	27
647	ECBS & ICBS 2015 Joint Meeting: Bringing Chemistry to Life. 2016 , 17, 447-52	2
646	A Modular Probe Strategy for Drug Localization, Target Identification and Target Occupancy Measurement on Single Cell Level. 2016 , 11, 2541-50	53
645	Chemical Dimerizers in Three-Hybrid Systems for Small Molecule-Target Protein Profiling. 2016 , 11, 2075-90	3

644	Selectivity on-target of bromodomain chemical probes by structure-guided medicinal chemistry and chemical biology. 2016 , 8, 1655-80	41
643	Trapping mammalian protein complexes in viral particles. 2016 , 7, 11416	30
642	Sample Preparation for Mass Spectrometry-Based Proteomics; from Proteomes to Peptides. 2016 , 919, 43-62	20
641	CETSA screening identifies known and novel thymidylate synthase inhibitors and slow intracellular activation of 5-fluorouracil. 2016 , 7, 11040	96
640	Small-Molecule Target Engagement in Cells. 2016 , 23, 435-41	77
639	Novel approaches to map small molecule-target interactions. 2016 , 24, 3232-45	19
638	Cellular thermal shift and clickable chemical probe assays for the determination of drug-target engagement in live cells. 2016 , 14, 6179-83	19
637	Application of Mass Spectrometry Profiling to Establish Brusatol as an Inhibitor of Global Protein Synthesis. 2016 , 15, 1220-31	61
636	Systematic Identification of Protein-Metabolite Interactions in Complex Metabolite Mixtures by Ligand-Detected Nuclear Magnetic Resonance Spectroscopy. 2016 , 55, 2590-600	13
635	Chemical proteomics approaches for identifying the cellular targets of natural products. 2016 , 33, 681-708	225
635 634	Chemical proteomics approaches for identifying the cellular targets of natural products. 2016 , 33, 681-708 An industry perspective on drug target validation. 2016 , 11, 623-5	225 18
		Ĭ
634	An industry perspective on drug target validation. 2016 , 11, 623-5	18
634	An industry perspective on drug target validation. 2016 , 11, 623-5 Medicinal Bioprospecting of the Amazon Rainforest: A Modern Eldorado?. 2016 , 34, 781-790	18
634 633	An industry perspective on drug target validation. 2016 , 11, 623-5 Medicinal Bioprospecting of the Amazon Rainforest: A Modern Eldorado?. 2016 , 34, 781-790 Thermal profiling reveals phenylalanine hydroxylase as an off-target of panobinostat. 2016 , 12, 908-910	18 21 120
634 633 632	An industry perspective on drug target validation. 2016, 11, 623-5 Medicinal Bioprospecting of the Amazon Rainforest: A Modern Eldorado?. 2016, 34, 781-790 Thermal profiling reveals phenylalanine hydroxylase as an off-target of panobinostat. 2016, 12, 908-910 Lanthanides in Luminescent Thermometry. 2016, 49, 339-427 An Unbiased Chemical Proteomics Method Identifies Fabl as the Primary Target of 6-OH-BDE-47.	18 21 120 196
634633632631630	An industry perspective on drug target validation. 2016, 11, 623-5 Medicinal Bioprospecting of the Amazon Rainforest: A Modern Eldorado?. 2016, 34, 781-790 Thermal profiling reveals phenylalanine hydroxylase as an off-target of panobinostat. 2016, 12, 908-910 Lanthanides in Luminescent Thermometry. 2016, 49, 339-427 An Unbiased Chemical Proteomics Method Identifies Fabl as the Primary Target of 6-OH-BDE-47. 2016, 50, 11329-11336	18 21 120 196 20

626	Early Perspective. 2016 , 21, 1019-1033	19
625	Targeted Mass Spectrometry-Based Approach for Protein-Ligand Binding Analyses in Complex Biological Mixtures Using a Phenacyl Bromide Modification Strategy. 2016 , 88, 10987-10993	11
624	Proteome-wide Profiling of Clinical PARP Inhibitors Reveals Compound-Specific Secondary Targets. 2016 , 23, 1490-1503	58
623	Validation and development of MTH1 inhibitors for treatment of cancer. 2016 , 27, 2275-2283	77
622	The diverse and expanding role of mass spectrometry in structural and molecular biology. 2016 , 35, 2634-26	57 160
621	Large-Scale Analysis of Breast Cancer-Related Conformational Changes in Proteins Using Limited Proteolysis. 2016 , 15, 4666-4674	20
620	Trail-blAIZin new directions for conditional proteomics. 2016 , 13, 917-918	
619	Kinetic Insights into the Binding between the nSH3 Domain of CrkII and Proline-Rich Motifs in cAbl. 2016 , 111, 1843-1853	7
618	cAMP-dependent protein kinase (PKA) complexes probed by complementary differential scanning fluorimetry and ion mobility-mass spectrometry. 2016 , 473, 3159-75	33
617	Studying epigenetic complexes and their inhibitors with the proteomics toolbox. 2016 , 8, 76	12
616	Thermal proteome profiling: unbiased assessment of protein state through heat-induced stability changes. 2016 , 15, 13	62
615	Binding Mechanism of the N-Terminal SH3 Domain of CrkII and Proline-Rich Motifs in cAbl. 2016 , 110, 2630-2641	14
614	Functional interdependence of BRD4 and DOT1L in MLL leukemia. 2016 , 23, 673-81	69
613	The Cellular Thermal Shift Assay: A Novel Biophysical Assay for In Situ Drug Target Engagement and Mechanistic Biomarker Studies. 2016 , 56, 141-61	156
612	The contribution of mass spectrometry-based proteomics to understanding epigenetics. 2016 , 8, 429-45	28
611	Characterizing Protein-Protein Interactions Using Mass Spectrometry: Challenges and Opportunities. 2016 , 34, 825-834	99
610	Label-free technologies for target identification and validation. 2016, 7, 769-777	8
609	A Biologist's Field Guide to Multiplexed Quantitative Proteomics. 2016 , 15, 1489-97	44

608	Chemical Proteomics Reveals Ferrochelatase as a Common Off-target of Kinase Inhibitors. 2016 , 11, 1245-54	58
60 7	Strategies in functional proteomics: Unveiling the pathways to precision oncology. 2016 , 382, 86-94	7
606	A Perspective on Implementing a Quantitative Systems Pharmacology Platform for Drug Discovery and the Advancement of Personalized Medicine. 2016 , 21, 521-34	29
605	Characterization of the Saccharomyces cerevisiae ATP-Interactome using the iTRAQ-SPROX Technique. 2016 , 27, 233-43	26
604	Next-generation proteomics faces new challenges in environmental biotechnology. 2016 , 38, 174-82	36
603	Advances in identification and validation of protein targets of natural products without chemical modification. 2016 , 33, 719-30	64
602	Non-stoichiometric inhibition in integrated lead finding - a literature review. 2016 , 11, 149-62	6
601	The resurgence of phenotypic screening in drug discovery and development. 2016 , 11, 121-5	32
600	Whole/Intact Cell MALDI MS Biotyping in Mammalian Cell Analysis. 2016, 249-262	2
599	Identifying compound efficacy targets in phenotypic drug discovery. 2016 , 21, 82-89	103
599 598	Identifying compound efficacy targets in phenotypic drug discovery. 2016 , 21, 82-89 A timeline of stable isotopes and mass spectrometry in the life sciences. 2017 , 36, 58-85	103 47
		, in the second
598	A timeline of stable isotopes and mass spectrometry in the life sciences. 2017 , 36, 58-85	47
598 597	A timeline of stable isotopes and mass spectrometry in the life sciences. 2017 , 36, 58-85 Ligand and Target Discovery by Fragment-Based Screening in Human Cells. 2017 , 168, 527-541.e29	47 231
598597596	A timeline of stable isotopes and mass spectrometry in the life sciences. 2017, 36, 58-85 Ligand and Target Discovery by Fragment-Based Screening in Human Cells. 2017, 168, 527-541.e29 Target engagement: Shining a light. 2017, 13, 133-134 A PDE6EKRas Inhibitor Chemotype with up to Seven H-Bonds and Picomolar Affinity that Prevents	47 231 5
598597596595	A timeline of stable isotopes and mass spectrometry in the life sciences. 2017, 36, 58-85 Ligand and Target Discovery by Fragment-Based Screening in Human Cells. 2017, 168, 527-541.e29 Target engagement: Shining a light. 2017, 13, 133-134 A PDE6EKRas Inhibitor Chemotype with up to Seven H-Bonds and Picomolar Affinity that Prevents Efficient Inhibitor Release by Arl2. 2017, 56, 2423-2428 From genome to proteome: Looking beyond DNA and RNA in chronic lymphocytic leukemia. 2017,	47 231 5
598597596595594	A timeline of stable isotopes and mass spectrometry in the life sciences. 2017, 36, 58-85 Ligand and Target Discovery by Fragment-Based Screening in Human Cells. 2017, 168, 527-541.e29 Target engagement: Shining a light. 2017, 13, 133-134 A PDE6EKRas Inhibitor Chemotype with up to Seven H-Bonds and Picomolar Affinity that Prevents Efficient Inhibitor Release by Arl2. 2017, 56, 2423-2428 From genome to proteome: Looking beyond DNA and RNA in chronic lymphocytic leukemia. 2017, 155, 73-84	47 231 5 63 5

(2017-2017)

590	Cell-wide analysis of protein thermal unfolding reveals determinants of thermostability. <i>Science</i> , 2017 , 355,	33.3	198
589	Technological advances and proteomic applications in drug discovery and target deconvolution: identification of the pleiotropic effects of statins. 2017 , 22, 848-869		19
588	Non-kinase targets of protein kinase inhibitors. 2017 , 16, 424-440		72
587	Plant Chemical Genetics: From Phenotype-Based Screens to Synthetic Biology. 2017 , 174, 5-20		43
586	System-wide detection of protein-small molecule complexes suggests extensive metabolite regulation in plants. 2017 , 7, 42387		26
585	Structure of the mycobacterial ESX-5 type VII secretion system membrane complex by single-particle analysis. 2017 , 2, 17047		74
584	Covalent Protein Labeling at Glutamic Acids. 2017 , 24, 589-597.e5		50
583	In silico polypharmacology of natural products. 2018 , 19, 1153-1171		64
582	Target Engagement Measures in Preclinical Drug Discovery: Theory, Methods, and Case Studies. 2017 , 41-80		2
581	Discovery of Nicotinamide Adenine Dinucleotide Binding Proteins in the Escherichia coli Proteome Using a Combined Energetic- and Structural-Bioinformatics-Based Approach. 2017 , 16, 470-480		8
580	Ferrochelatase is a therapeutic target for ocular neovascularization. 2017, 9, 786-801		30
579	Measuring protein structural changes on a proteome-wide scale using limited proteolysis-coupled mass spectrometry. 2017 , 12, 2391-2410		83
578	Optimization of Experimental Parameters in Data-Independent Mass Spectrometry Significantly Increases Depth and Reproducibility of Results. 2017 , 16, 2296-2309		186
577	Quantitative proteomics of model organisms. 2017 , 6, 58-66		7
576	A Primer on Concepts and Applications of Proteomics in Neuroscience. 2017 , 96, 558-571		41
575	Biomarkers of Cancer. 2017 , 273-462		1
574	Combined CRISPRi/a-Based Chemical Genetic Screens Reveal that Rigosertib Is a Microtubule-Destabilizing Agent. 2017 , 68, 210-223.e6		127
573	Sensitive and Accurate Quantitation of Phosphopeptides Using TMT Isobaric Labeling Technique. 2017 , 16, 4244-4252		16

572	Structural coverage of the proteome for pharmaceutical applications. 2017 , 22, 1792-1799	34
571	ATP alters protein folding and function of Escherichia coli uridine phosphorylase. 2017 , 634, 11-20	3
570	A network integration approach for drug-target interaction prediction and computational drug repositioning from heterogeneous information. 2017 , 8, 573	295
569	Genome-Scale Architecture of Small Molecule Regulatory Networks and the Fundamental Trade-Off between Regulation and Enzymatic Activity. 2017 , 20, 2666-2677	43
568	Mass spectrometry methods to study protein-metabolite interactions. 2017 , 12, 1271-1280	6
567	Nanoparticle Surface Functionality Dictates Cellular and Systemic Toxicity. 2017 , 29, 6578-6595	73
566	Towards detecting regulatory protein-metabolite interactions. 2017, 39, 16-23	28
565	Cellular Assays. 2017 , 313-333	1
564	The target landscape of clinical kinase drugs. <i>Science</i> , 2017 , 358,	389
-(-		
563	Chemical Biology in Drug Discovery. 2017 , 335-370	2
563	Prediction of intracellular exposure bridges the gap between target- and cell-based drug discovery. 2017 , 114, E6231-E6239	60
	Prediction of intracellular exposure bridges the gap between target- and cell-based drug discovery.	
562	Prediction of intracellular exposure bridges the gap between target- and cell-based drug discovery. 2017 , 114, E6231-E6239	60
562 561	Prediction of intracellular exposure bridges the gap between target- and cell-based drug discovery. 2017, 114, E6231-E6239 Biophysics: for HTS hit validation, chemical lead optimization, and beyond. 2017, 12, 897-907 Measurement of drug-target engagement in live cells by two-photon fluorescence anisotropy	60
562 561 560	Prediction of intracellular exposure bridges the gap between target- and cell-based drug discovery. 2017, 114, E6231-E6239 Biophysics: for HTS hit validation, chemical lead optimization, and beyond. 2017, 12, 897-907 Measurement of drug-target engagement in live cells by two-photon fluorescence anisotropy imaging. 2017, 12, 1472-1497	60 12 14
562 561 560 559	Prediction of intracellular exposure bridges the gap between target- and cell-based drug discovery. 2017, 114, E6231-E6239 Biophysics: for HTS hit validation, chemical lead optimization, and beyond. 2017, 12, 897-907 Measurement of drug-target engagement in live cells by two-photon fluorescence anisotropy imaging. 2017, 12, 1472-1497 Emerging Methods in Chemoproteomics with Relevance to Drug Discovery. 2017, 1513, 11-22	60 12 14
562 561 560 559 558	Prediction of intracellular exposure bridges the gap between target- and cell-based drug discovery. 2017, 114, E6231-E6239 Biophysics: for HTS hit validation, chemical lead optimization, and beyond. 2017, 12, 897-907 Measurement of drug-target engagement in live cells by two-photon fluorescence anisotropy imaging. 2017, 12, 1472-1497 Emerging Methods in Chemoproteomics with Relevance to Drug Discovery. 2017, 1513, 11-22 Label-free target identification using in-gel fluorescence difference thermal stability shift. 2017, 8, 1127-113 Detergents: Friends not foes for high-performance membrane proteomics toward precision	60 12 14 15 33 25

554	Novel Molecular Challenges in Targeting Anaplastic Lymphoma Kinase in ALK-Expressing Human Cancers. 2017 , 9,		4
553	Contemporary Considerations for Discovering a Successful Medicine. 2017 , 560-576		
552	Capturing dynamic protein interactions. <i>Science</i> , 2018 , 359, 1105-1106	33.3	35
551	Targeting CDK2 overcomes melanoma resistance against BRAF and Hsp90 inhibitors. 2018 , 14, e7858		35
550	Advancing translational research and precision medicine with targeted proteomics. 2018, 189, 1-10		44
549	Homogeneous Assay for Target Engagement Utilizing Bioluminescent Thermal Shift. 2018 , 9, 546-551		22
548	New use for CETSA: monitoring innate immune receptor stability via post-translational modification by OGT. 2018 , 50, 231-240		10
547	Thermal proteome profiling of breast cancer cells reveals proteasomal activation by CDK4/6 inhibitor palbociclib. 2018 , 37,		52
546	Global profiling of protein-DNA and protein-nucleosome binding affinities using quantitative mass spectrometry. 2018 , 9, 1653		36
545	Modulation of Protein-Interaction States through the Cell Cycle. 2018 , 173, 1481-1494.e13		80
544	Systematic analysis of protein turnover in primary cells. 2018 , 9, 689		145
543	Methods to validate Hsp90 inhibitor specificity, to identify off-target effects, and to rethink approaches for further clinical development. 2018 , 23, 467-482		58
542	Application of targeted mass spectrometry in bottom-up proteomics for systems biology research. 2018 , 189, 75-90		49
541	Structure-guided design of a potent peptide inhibitor targeting the interaction between CRK and ABL kinase. 2018 , 9, 519-524		1
540	An integrative overview of genomic, transcriptomic and proteomic analyses in organohalide respiration research. 2018 , 94,		23
539	ProteomicsDB. 2018 , 46, D1271-D1281		113
538	Targeted NUDT5 inhibitors block hormone signaling in breast cancer cells. 2018 , 9, 250		28
537	Thermal proximity coaggregation for system-wide profiling of protein complex dynamics in cells. <i>Science</i> , 2018 , 359, 1170-1177	33.3	90

536	Metabolism as a signal generator across trans-omic networks at distinct time scales. 2018, 8, 59-66	18
535	New Perspectives, Opportunities, and Challenges in Exploring the Human Protein Kinome. 2018 , 78, 15-29	81
534	CRISPR Approaches to Small Molecule Target Identification. 2018 , 13, 366-375	41
533	A Map of Protein-Metabolite Interactions Reveals Principles of Chemical Communication. 2018 , 172, 358-372.e23	192
532	Emerging Approaches for the Identification of Protein Targets of Small Molecules - A Practitioners' Perspective. 2018 , 61, 8504-8535	38
531	Pervasive Protein Thermal Stability Variation during the Cell Cycle. 2018 , 173, 1495-1507.e18	109
530	Multiplexed Proteome Dynamics Profiling Reveals Mechanisms Controlling Protein Homeostasis. 2018 , 173, 260-274.e25	125
529	Bioactive Compound Collections: From Design to Target Identification. 2018 , 4, 705-730	29
528	Kinase inhibitors: the road ahead. 2018, 17, 353-377	447
527	Toxoplasma Calcium-Dependent Protein Kinase 1 Inhibitors: Probing Activity and Resistance Using Cellular Thermal Shift Assays. 2018 , 62,	11
526	Cell Density Affects the Detection of Chk1 Target Engagement by the Selective Inhibitor V158411. 2018 , 23, 144-153	2
525	A High-Throughput Dose-Response Cellular Thermal Shift Assay for Rapid Screening of Drug Target Engagement in Living Cells, Exemplified Using SMYD3 and IDO1. 2018 , 23, 34-46	28
524	A Vinyl Sulfone-Based Fluorogenic Probe Capable of Selective Labeling of PHGDH in Live Mammalian Cells. 2018 , 57, 579-583	27
523	A Vinyl Sulfone-Based Fluorogenic Probe Capable of Selective Labeling of PHGDH in Live Mammalian Cells. 2018 , 130, 588-592	8
522	Demonstrating In-Cell Target Engagement Using a Pirin Protein Degradation Probe (CCT367766). 2018 , 61, 918-933	61
521	Small-Molecule Kinase Downregulators. 2018 , 25, 30-35	45
520	A Step-Up LC-MS/MS for Proteomics. 2018 , 377-414	
519	Ber bisherige Denkweisen hinaus [heue Wirkstoffe zur Berwindung der Antibiotika-Krise. 2018 , 130, 14642-14682	15

518 Chemo-Selection Strategy for Limited Proteolysis Experiments on the Proteomic Scale. **2018**, 90, 14039-14047 9

5 ¹ 7	An efficient proteome-wide strategy for discovery and characterization of cellular nucleotide-protein interactions. 2018 , 13, e0208273	22
516	Identification of Protein Targets of Bioactive Small Molecules Using Randomly Photomodified Probes. 2018 , 13, 3333-3342	7
515	Reagents for Isobaric Labeling Peptides in Quantitative Proteomics. 2018 , 90, 12366-12371	21
514	Covalent inhibitors of EGFR family protein kinases induce degradation of human Tribbles 2 (TRIB2) pseudokinase in cancer cells. 2018 , 11,	43
513	Thinking Outside the Box-Novel Antibacterials To Tackle the Resistance Crisis. 2018 , 57, 14440-14475	75
512	Proteome-Wide Structural Biology: An Emerging Field for the Structural Analysis of Proteins on the Proteomic Scale. 2018 , 17, 3614-3627	34
511	Small Molecule Targeting of Specific BAF (mSWI/SNF) Complexes for HIV Latency Reversal. 2018 , 25, 1443-1455.e14	25
510	Detection of Chemical Engagement of Solute Carrier Proteins by a Cellular Thermal Shift Assay. 2018 , 13, 1480-1486	20
509	Impact of Mass Spectrometry-Based Technologies and Strategies on Chemoproteomics as a Tool for Drug Discovery. 2018 , 9, 785-791	14
508	Identification of the Natural Product Rotihibin A as a TOR Kinase Signaling Inhibitor by Unbiased Transcriptional Profiling. 2018 , 24, 12500-12504	2
507	Thermal proteome profiling in bacteria: probing protein state. 2018 , 14, e8242	82
506	Natural Products for Drug Discovery in the 21st Century: Innovations for Novel Drug Discovery. 2018 , 19,	380
505	Innovations in improving lipid production: Algal chemical genetics. 2018 , 71, 101-123	20
504	A high content, high throughput cellular thermal stability assay for measuring drug-target engagement in living cells. 2018 , 13, e0195050	14
503	Chemical Denaturation and Protein Precipitation Approach for Discovery and Quantitation of Protein-Drug Interactions. 2018 , 90, 9249-9255	21
502	Chemoproteomics and Chemical Probes for Target Discovery. 2018 , 36, 1275-1286	57
501	Chemical proteomics, an integrated research engine for exploring drug-target-phenotype interactions. 2018 , 16, 1	5

500	A Novel LC System Embeds Analytes in Pre-formed Gradients for Rapid, Ultra-robust Proteomics. 2018 , 17, 2284-2296	133
499	A widely-applicable high-throughput cellular thermal shift assay (CETSA) using split Nano Luciferase. 2018 , 8, 9472	42
498	Polypharmacology by Design: A Medicinal Chemist's Perspective on Multitargeting Compounds. 2019 , 62, 420-444	188
497	Thermal proteome profiling allows quantitative assessment of interactions between tetrachloroethene reductive dehalogenase and trichloroethene. 2019 , 192, 10-17	17
496	An Antitumor Bis(N-Heterocyclic Carbene)Platinum(II) Complex That Engages Asparagine Synthetase as an Anticancer Target. 2019 , 131, 11030-11034	8
495	High throughput discovery of functional protein modifications by Hotspot Thermal Profiling. 2019 , 16, 894-901	52
494	Application of Bioactive Thermal Proteome Profiling to Decipher the Mechanism of Action of the Lipid Lowering 13-Hydroxy-pheophytin Isolated from a Marine Cyanobacteria. 2019 , 17,	10
493	Target engagement approaches for pharmacological evaluation in animal models. 2019 , 55, 9241-9250	1
492	CETSA beyond Soluble Targets: a Broad Application to Multipass Transmembrane Proteins. 2019 , 14, 1913-1920	31
491	Advances in exploring the therapeutic potential of marine natural products. 2019 , 147, 104373	39
490	Mechanistic MALDI-TOF Cell-Based Assay for the Discovery of Potent and Specific Fatty Acid Synthase Inhibitors. 2019 , 26, 1322-1331.e4	7
489	ACT001, a novel PAI-1 inhibitor, exerts synergistic effects in combination with cisplatin by inhibiting PI3K/AKT pathway in glioma. 2019 , 10, 757	56
488	Target Discovery Using Thermal Proteome Profiling. 2019 , 267-291	3
487	ProteomicsDB: a multi-omics and multi-organism resource for life science research. 2020 , 48, D1153-D1163	67
486	Hydralazine targets cAMP-dependent protein kinase leading to sirtuin1/5 activation and lifespan extension in C. elegans. 2019 , 10, 4905	16
485	Differential PROTAC substrate specificity dictated by orientation of recruited E3 ligase. 2019 , 10, 131	188
484	Anticancer Effect of Deuterium Depleted Water - Redox Disbalance Leads to Oxidative Stress. 2019 , 18, 2373-2387	18
483	Epigenetic drug target deconvolution by mass spectrometry-based technologies. 2019 , 26, 854-857	3

482	Chemical proteomics reveals target selectivity of clinical Jak inhibitors in human primary cells. 2019 , 9, 14159	22
481	Nonparametric Analysis of Thermal Proteome Profiles Reveals Novel Drug-binding Proteins. 2019 , 18, 2506-2515	34
480	Proteome Integral Solubility Alteration: A High-Throughput Proteomics Assay for Target Deconvolution. 2019 , 18, 4027-4037	56
479	Illuminating the dark phosphoproteome. 2019 , 12,	125
478	Effect of Sec61 interaction with Mpd1 on endoplasmic reticulum-associated degradation. 2019 , 14, e0211180	4
477	Plasma proteomic and autoantibody profiles reveal the proteomic characteristics involved in longevity families in Bama, China. 2019 , 16, 22	6
476	Mechanistic Basis of Cabotegravir-Glucuronide Disposition in Humans. 2019 , 370, 269-277	3
475	Emerging strategies for the identification of protein-metabolite interactions. 2019 , 70, 4605-4618	15
474	An Antitumor Bis(N-Heterocyclic Carbene)Platinum(II) Complex That Engages Asparagine Synthetase as an Anticancer Target. 2019 , 58, 10914-10918	31
473	Posttranslational Modifications Drive Protein Stability to Control the Dynamic Beer Brewing Proteome. 2019 , 18, 1721-1731	19
472	DO-MS: Data-Driven Optimization of Mass Spectrometry Methods. 2019 , 18, 2493-2500	31
471	Melting Down Protein Stability: PAPS Synthase 2 in Patients and in a Cellular Environment. 2019 , 6, 31	8
470	Quantifying drug-target engagement in live cells using sulfonyl fluoride chemical probes. 2019 , 622, 201-220	2
469	Chemical Epigenetics: The Impact of Chemical and Chemical Biology Techniques on Bromodomain Target Validation. 2019 , 58, 17930-17952	20
468	Global analysis of methionine oxidation provides a census of folding stabilities for the human proteome. 2019 , 116, 6081-6090	38
467	Engagement with tNOX (ENOX2) to Inhibit SIRT1 and Activate p53-Dependent and -Independent Apoptotic Pathways by Novel 4,11-Diaminoanthra[2,3-]furan-5,10-diones in Hepatocellular Carcinoma Cells. 2019 , 11,	9
466	Proteome-wide solubility and thermal stability profiling reveals distinct regulatory roles for ATP. 2019 , 10, 1155	88
465	Chemische Epigenetik: der Einfluss chemischer und chemo-biologischer Techniken auf die Zielstruktur-Validierung von Bromodomfien. 2019 , 131, 18096-18120	1

464	Monitoring structural modulation of redox-sensitive proteins in cells with MS-CETSA. 2019 , 24, 101168	23
463	Advanced proteomics approaches to unravel protein homeostasis. 2019 , 31, 99-108	11
462	Label-free target identification in drug discovery via phenotypic screening. 2019 , 50, 66-72	20
461	Horizontal Cell Biology: Monitoring Global Changes of Protein Interaction States with the Proteome-Wide Cellular Thermal Shift Assay (CETSA). 2019 , 88, 383-408	39
460	Protein corona formed on silver nanoparticles in blood plasma is highly selective and resistant to physicochemical changes of the solution. 2019 , 6, 1089-1098	36
459	Proteomics in Drug Development: The Dawn of a New Era?. 2019 , 13, e1800087	30
458	Mass Spectrometry and Chemical Biology in Epigenetics Drug Discovery. 2019 , 79-106	
457	Chemical Probes. 2019 , 133-152	
456	Identification of bioactive metabolites using activity metabolomics. 2019 , 20, 353-367	258
455	Label-free target identification reveals oxidative DNA damage as the mechanism of a selective cytotoxic agent. 2019 , 10, 3449-3458	14
454	CETSA-based target engagement of taxanes as biomarkers for efficacy and resistance. 2019 , 9, 19384	10
453	PROMIS: Global Analysis of PROtein-Metabolite Interactions. 2019 , 4, e20101	8
452	A new era for proteomics. 2019 , 11, 1731-1735	4
451	Label-Free Techniques for Target Discovery and Validation. 2019 , 131-152	
450	ProTargetMiner as a proteome signature library of anticancer molecules for functional discovery. 2019 , 10, 5715	16
449	Target deconvolution from phenotype-based drug discovery by using chemical proteomics approaches. 2019 , 1867, 22-27	40
448	Screening Strategies and Methods for Better Off-Target Liability Prediction and Identification of Small-Molecule Pharmaceuticals. 2019 , 24, 1-24	17
447	A one-pot analysis approach to simplify measurements of protein stability and folding kinetics. 2019 , 1867, 184-193	13

(2020-2019)

446	Identifying purine nucleoside phosphorylase as the target of quinine using cellular thermal shift assay. 2019 , 11,	81
445	Plasma proteome profiling of high-altitude polycythemia using TMT-based quantitative proteomics approach. 2019 , 194, 60-69	16
444	Monitoring and deciphering protein degradation pathways inside cells. 2019 , 31, 61-68	31
443	Proteome-wide Analysis of Protein Thermal Stability in the Model Higher Plant. 2019 , 18, 308-319	23
442	Perturbation-Based Proteomic Correlation Profiling as a Target Deconvolution Methodology. 2019 , 26, 137-143.e8	5
441	TMT-Based Quantitative Proteomic Analysis Reveals Proteomic Changes Involved in Longevity. 2019 , 13, e1800024	9
440	Lysine-Targeted Inhibitors and Chemoproteomic Probes. 2019 , 88, 365-381	42
439	A Review on Quantitative Multiplexed Proteomics. 2019 , 20, 1210-1224	110
438	Quantitative, Real-Time Measurements of Intracellular Target Engagement Using Energy Transfer. 2019 , 1888, 45-71	19
437	Target Engagement of Small Molecules: Thermal Profiling Approaches on Different Levels. 2019 , 1888, 73-98	5
436	Fluorescence anisotropy imaging in drug discovery. 2019 , 151-152, 262-288	29
435	High-Throughput Cellular Thermal Shift Assays in Research and Drug Discovery. 2020 , 25, 137-147	22
434	The future of cystic fibrosis care: a global perspective. 2020 , 8, 65-124	259
433	Tools of the Ethylene Trade: A Chemical Kit to Influence Ethylene Responses in Plants and Its Use in Agriculture. 2020 , 4, 1900267	5
432	Systematic mapping of protein-metabolite interactions with mass spectrometry-based techniques. 2020 , 64, 24-31	9
431	Modification-free approaches to screen drug targets at proteome level. 2020 , 124, 115574	8
430	Der zytotoxische Naturstoff Vioprolid A interagiert mit dem fil die Ribosomen-Biogenese essentiellen nukleolien Protein 14. 2020 , 132, 1611-1617	2
429	Mass spectrometry-based Cellular Thermal Shift Assay (CETSA[]) for target deconvolution in phenotypic drug discovery. 2020 , 28, 115174	17

428	Perspective on CETSA Literature: Toward More Quantitative Data Interpretation. 2020, 25, 118-126	16
427	The Cytotoxic Natural Product Vioprolide A Targets Nucleolar Protein 14, Which Is Essential for Ribosome Biogenesis. 2020 , 59, 1595-1600	24
426	Multiscale modelling of drug mechanism and safety. 2020 , 25, 519-534	10
425	Drug Target Engagement Using Coupled Cellular Thermal Shift Assay-Acoustic Reverse-Phase Protein Array. 2020 , 25, 207-214	3
424	Label-Free Proteome Profiling as a Quantitative Target Identification Technique for Bioactive Small Molecules. 2020 , 59, 213-215	О
423	Mass spectrometry analysis of the structural proteome. 2020 , 60, 57-65	13
422	Chemoproteomic profiling of protein-metabolite interactions. 2020 , 54, 28-36	7
421	The functional landscape of the human phosphoproteome. 2020 , 38, 365-373	106
420	CETSA in integrated proteomics studies of cellular processes. 2020 , 54, 54-62	16
419	Solvent-Induced Protein Precipitation for Drug Target Discovery on the Proteomic Scale. 2020 , 92, 1363-1371	21
418	Image-Based Morphological Profiling Identifies a Lysosomotropic, Iron-Sequestering Autophagy Inhibitor. 2020 , 59, 5721-5729	27
417	Identification of a Small Compound Targeting PKM2-Regulated Signaling Using 2D Gel Electrophoresis-Based Proteome-wide CETSA. 2020 , 27, 186-196.e4	21
416	The Druggability of Solute Carriers. 2020 , 63, 3834-3867	25
415	An anticancer gold(III)-activated porphyrin scaffold that covalently modifies protein cysteine thiols. 2020 , 117, 1321-1329	26
414	Hit Triage and Validation in Phenotypic Screening: Considerations and Strategies. 2020 , 27, 1332-1346	17
413	Detectives and helpers: Natural products as resources for chemical probes and compound libraries. 2020 , 216, 107688	4
412	Profiling of post-translational modifications by chemical and computational proteomics. 2020 , 56, 13506-1351	 1 %
411	Thermal Proteome Profiling and Meltome Analysis of a Thermophilic Bacterial Strain, ARTRW1: Toward Industrial Applications. 2020 , 24, 756-765	2

(2020-2020)

410	A mass spectrometry-based proteome map of drug action in lung cancer cell lines. 2020, 16, 1111-1119	13
409	The glyphosate formulation Roundup LB plus influences the global metabolome of pig gut microbiota in vitro. 2020 , 745, 140932	8
408	Pooled protein tagging, cellular imaging, and in situ sequencing for monitoring drug action in real time. 2020 , 30, 1846-1855	3
407	A computational method for detection of ligand-binding proteins from dose range thermal proteome profiles. 2020 , 11, 5783	8
406	Mutant thermal proteome profiling for characterization of missense protein variants and their associated phenotypes within the proteome. 2020 , 295, 16219-16238	9
405	Direct label-free methods for identification of target proteins in agrochemicals. 2020 , 164, 1475-1483	1
404	Selective Modulation of Dynamic Protein Complexes. 2020 , 27, 986-997	10
403	COVID-19: Nanomedicine Uncovers Blood-Clot Mystery. 2020 , 19, 4364-4373	9
402	Genetic screens reveal a central role for heme metabolism in artemisinin susceptibility. 2020, 11, 4813	10
401	A machine learning-based chemoproteomic approach to identify drug targets and binding sites in complex proteomes. 2020 , 11, 4200	34
400	Analysis of selective target engagement by small-molecule sphingosine kinase inhibitors using the Cellular Thermal Shift Assay (CETSA). 2020 , 21, 841-852	3
399	Microparticle-Assisted Precipitation Screening Method for Robust Drug Target Identification. 2020 , 92, 13912-13921	5
398	On-Chip Acousto Thermal Shift Assay for Rapid and Sensitive Assessment of Protein Thermodynamic Stability. 2020 , 16, e2003506	4
397	Novel Broad-Spectrum Antiviral Inhibitors Targeting Host Factors Essential for Replication of Pathogenic RNA Viruses. 2020 , 12,	9
396	From Phenotypic Hit to Chemical Probe: Chemical Biology Approaches to Elucidate Small Molecule Action in Complex Biological Systems. 2020 , 25,	8
395	Culture-Independent Omics-Techniques for Microbiome-Based Molecular Therapeutics Against Infectious Diseases. 2020 , 95-114	
394	Thermostability profiling of MHC-bound peptides: a new dimension in immunopeptidomics and aid for immunotherapy design. 2020 , 11, 6305	2
393	The functional proteome landscape of Escherichia coli. 2020 , 588, 473-478	14

392	The GOLIATH Project: Towards an Internationally Harmonised Approach for Testing Metabolism Disrupting Compounds. 2020 , 21,	13
391	New approaches to antibacterial drug discovery. 2020 , 223-248	
390	A Dual-Mechanism Antibiotic Kills Gram-Negative Bacteria and Avoids Drug Resistance. 2020 , 181, 1518-1532	2.e8 8
389	Target identification of natural medicine with chemical proteomics approach: probe synthesis, target fishing and protein identification. 2020 , 5, 72	29
388	Spatiotemporal proteomics uncovers cathepsin-dependent macrophage cell death during Salmonella infection. 2020 , 5, 1119-1133	17
387	TMTpro reagents: a set of isobaric labeling mass tags enables simultaneous proteome-wide measurements across 16 samples. 2020 , 17, 399-404	123
386	Cell-Based Ligand Discovery for the ENL YEATS Domain. 2020 , 15, 895-903	14
385	Interaction profiling methods to map protein and pathway targets of bioactive ligands. 2020, 54, 76-84	2
384	Validation of the Applicability of In-Cell Fast Photochemical Oxidation of Proteins across Multiple Eukaryotic Cell Lines. 2020 , 31, 1372-1379	3
383	Comprehensive chemical proteomics for target deconvolution of the redox active drug auranofin. 2020 , 32, 101491	27
382	Characterizing Drug-Target Interactions: Shifting towards the Clinic. 2020 , 41, 295-297	1
381	Quantifying Target Occupancy of Small Molecules Within Living Cells. 2020 , 89, 557-581	20
380	Multiparametric Assays for Accelerating Early Drug Discovery. 2020 , 41, 318-335	8
379	Selection of Heating Temperatures Improves the Sensitivity of the Proteome Integral Solubility Alteration Assay. 2020 , 19, 2159-2166	8
378	Phenotypic Screening of Chemical Libraries Enriched by Molecular Docking to Multiple Targets Selected from Glioblastoma Genomic Data. 2020 , 15, 1424-1444	3
377	Proteomic and interactomic insights into the molecular basis of cell functional diversity. 2020 , 21, 327-340	68
376	Importance of Quantifying Drug-Target Engagement in Cells. 2020 , 11, 403-406	11
375	Monitoring protein communities and their responses to therapeutics. 2020 , 19, 414-426	19

(2020-2020)

374	Identifying the Target of an Antiparasitic Compound in Using Thermal Proteome Profiling. 2020, 15, 1801-180	7 15
373	A Simplified Thermal Proteome Profiling Approach to Screen Protein Targets of a Ligand. 2020 , 20, e1900372	4
372	Data, Reagents, Assays and Merits of Proteomics for SARS-CoV-2 Research and Testing. 2020 , 19, 1503-1522	52
371	Lysine-specific demethylase 1A restricts ex vivo propagation of human HSCs and is a target of UM171. 2020 , 136, 2151-2161	12
370	Comparative Analysis of Mass-Spectrometry-Based Proteomic Methods for Protein Target Discovery Using a One-Pot Approach. 2020 , 31, 217-226	11
369	Chemical Biology Framework to Illuminate Proteostasis. 2020 , 89, 529-555	12
368	Image-Based Morphological Profiling Identifies a Lysosomotropic, Iron-Sequestering Autophagy Inhibitor. 2020 , 132, 5770-5778	10
367	An isothermal shift assay for proteome scale drug-target identification. 2020 , 3, 75	18
366	Progress and pitfalls of using isobaric mass tags for proteome profiling. 2020 , 17, 149-161	13
365	Click Chemistry in Proteomic Investigations. 2020 , 180, 605-632	92
365 364	Click Chemistry in Proteomic Investigations. 2020 , 180, 605-632 High-throughput quantitative top-down proteomics. 2020 , 16, 91-99	92
364	High-throughput quantitative top-down proteomics. 2020 , 16, 91-99	34
364	High-throughput quantitative top-down proteomics. 2020 , 16, 91-99 Chaperone mediated detection of small molecule target binding in cells. 2020 , 11, 465 Pifithrin-Balters p53 post-translational modifications pattern and differentially inhibits p53 target	34
364 363 362	High-throughput quantitative top-down proteomics. 2020, 16, 91-99 Chaperone mediated detection of small molecule target binding in cells. 2020, 11, 465 Pifithrin-Blters p53 post-translational modifications pattern and differentially inhibits p53 target genes. 2020, 10, 1049 A misprocessed form of Apolipoprotein A-I is specifically associated with recurrent Focal Segmental	34 1 14
364 363 362 361	High-throughput quantitative top-down proteomics. 2020, 16, 91-99 Chaperone mediated detection of small molecule target binding in cells. 2020, 11, 465 Pifithrin-talters p53 post-translational modifications pattern and differentially inhibits p53 target genes. 2020, 10, 1049 A misprocessed form of Apolipoprotein A-I is specifically associated with recurrent Focal Segmental Glomerulosclerosis. 2020, 10, 1159	341147
364 363 362 361 360	High-throughput quantitative top-down proteomics. 2020, 16, 91-99 Chaperone mediated detection of small molecule target binding in cells. 2020, 11, 465 Pifithrin-falters p53 post-translational modifications pattern and differentially inhibits p53 target genes. 2020, 10, 1049 A misprocessed form of Apolipoprotein A-I is specifically associated with recurrent Focal Segmental Glomerulosclerosis. 2020, 10, 1159 Identifying drug targets in tissues and whole blood with thermal-shift profiling. 2020, 38, 303-308 Thermal Proteome Profiling Identifies Oxidative-Dependent Inhibition of the Transcription of	34 1 14 7 46

356	Ethacrynic acid inhibits STAT3 activity through the modulation of SHP2 and PTP1B tyrosine phosphatases in DU145 prostate carcinoma cells. 2020 , 175, 113920	5
355	Towards a functional understanding of the plant metabolome. 2020 , 55, 47-51	20
354	Benchmarking the Orbitrap Tribrid Eclipse for Next Generation Multiplexed Proteomics. 2020 , 92, 6478-6485	24
353	Meltome atlas-thermal proteome stability across the tree of life. 2020 , 17, 495-503	53
352	Cellular thermal shift assay for the identification of drug-target interactions in the Plasmodium falciparum proteome. 2020 , 15, 1881-1921	27
351	Proteome-scale studies of protein stability. 2020 , 71-90	O
350	Loss of N-Glycanase 1 Alters Transcriptional and Translational Regulation in K562 Cell Lines. 2020 , 10, 1585-1597	5
349	Turning liabilities into opportunities: Off-target based drug repurposing in cancer. 2021 , 68, 209-229	11
348	Design of next-generation covalent inhibitors: Targeting residues beyond cysteine. 2021 , 56, 95-134	2
347	Rtpca: an R package for differential thermal proximity coaggregation analysis. 2021 , 37, 431-433	3
346	CETSA MS Profiling for a Comparative Assessment of FDA-Approved Antivirals Repurposed for COVID-19 Therapy Identifies TRIP13 as a Remdesivir Off-Target. 2021 , 26, 336-344	6
345	Isocotoin suppresses hepatitis E virus replication through inhibition of heat shock protein 90. 2021 , 185, 104997	4
344	Protein Folding Stability Changes Across the Proteome Reveal Targets of Cu Toxicity in. 2021 , 16, 214-224	7
343	Target Validation Using PROTACs: Applying the Four Pillars Framework. 2021 , 26, 474-483	7
342	ProThermDB: thermodynamic database for proteins and mutants revisited after 15 years. 2021 , 49, D420-D42	423
341	Dynamic 3D proteomes reveal protein functional alterations at high resolution in situ. 2021 , 184, 545-559.e22	28
340	Thermal Proteome Profiling in Zebrafish Reveals Effects of Napabucasin on Retinoic Acid Metabolism. 2021 , 20, 100033	3
339	Applicability of Chromatographic Co-Elution for Antibiotic Target Identification. 2021 , 21, e2000038	1

338 Target Validation Prosecuting the Target. 2021,

337	Affinity Enrichment Chemoproteomics for Target Deconvolution and Selectivity Profiling. 2021 , 2228, 237-252	1
336	System-Wide Profiling of Protein Complexes Via Size Exclusion Chromatography-Mass Spectrometry (SEC-MS). 2021 , 2259, 269-294	1
335	The right tools for the job: the central role for next generation chemical probes and chemistry-based target deconvolution methods in phenotypic drug discovery. 2021 , 12, 646-665	2
334	Cell surface thermal proteome profiling tracks perturbations and drug targets on the plasma membrane. 2021 , 18, 84-91	12
333	Proteomic Approaches to Study SARS-CoV-2 Biology and COVID-19 Pathology. 2021 , 20, 1133-1152	18
332	Selective cross-linking of coinciding protein assemblies by in-gel cross-linking mass spectrometry. 2021 , 40, e106174	8
331	Mass spectrometry-based protein-protein interaction networks for the study of human diseases. 2021 , 17, e8792	30
330	A Tale of Two Tails: Efficient Profiling of Protein Degraders by Specific Functional and Target Engagement Readouts. 2021 , 26, 534-546	7
329	Mass spectrometry-based protein-protein interaction techniques and their applications in studies of DNA damage repair. 2021 , 22, 1-20	2
328	Dynamic supramolecular self-assembly of platinum(ii) complexes perturbs an autophagy-lysosomal system and triggers cancer cell death 2021 , 12, 15229-15238	4
327	The Relationship between the Misfolding Avoidance Hypothesis and Protein Evolutionary Rates in the Light of Empirical Evidence. 2021 , 13,	4
326	System Biology-Guided Chemical Proteomics to Discover Protein Targets of Monoethylhexyl Phthalate in Regulating Cell Cycle. 2021 , 55, 1842-1851	6
325	"Structuromics": another step toward a holistic view of the cell. 2021 , 184, 301-303	2
324	Reproducibility in the unfolding process of protein induced by an external electric field. 2020 , 12, 2030-2038	1
323	Photoaffinity labelling strategies for mapping the small molecule-protein interactome. 2021 , 19, 7792-7809	8
322	Probing the methotrexate-protein interactions by proteomics and thermostability assay for drug resistance study. 2021 , 13, 411-418	
321	Chemoproteomic methods for covalent drug discovery. 2021 , 50, 8361-8381	5

320	Recent advances in bioanalytical methods to measure proteome stability in cells. 2021, 146, 2097-2109	5
319	Prediction and collection of protein-metabolite interactions. 2021 , 22,	19
318	Global mapping of protein-metabolite interactions in Saccharomyces cerevisiae reveals that Ser-Leu dipeptide regulates phosphoglycerate kinase activity. 2021 , 4, 181	12
317	Improved Proteomics-Based Drug Mechanism-of-Action Studies Using 16-Plex Isobaric Mass Tags. 2021 , 20, 1792-1801	5
316	System-wide identification and prioritization of enzyme substrates by thermal analysis. 2021 , 12, 1296	16
315	Hidden information on protein function in censuses of proteome foldedness.	O
314	Recent advances in proteome-wide label-free target deconvolution for bioactive small molecules. 2021 , 41, 2893-2926	1
313	Reply to Ma and Wang: Reliability of various in vitro activity assays on SARS-CoV-2 main protease inhibitors. 2021 , 118,	6
312	SARS-CoV-2 infection remodels the host protein thermal stability landscape. 2021 , 17, e10188	5
311	TP-MAP - an Integrated Software Package for the Analysis of 1D and 2D Thermal Profiling Data.	O
310	Identification of Celecoxib-Targeted Proteins Using Label-Free Thermal Proteome Profiling on Rat Hippocampus. 2021 , 99, 308-318	2
309	Thermal proteome profiling identifies the membrane-bound purinergic receptor P2X4 as a target of the autophagy inhibitor indophagolin. 2021 ,	9
308	A MALDI-TOF assay identifies nilotinib as an inhibitor of inflammation in acute myeloid leukaemia.	0
307	Chemical Phosphoproteomics Sheds New Light on the Targets and Modes of Action of AKT Inhibitors. 2021 , 16, 631-641	5
306	Global Profiling of Lysine Accessibility to Evaluate Protein Structure Changes in Alzheimer's Disease. 2021 , 32, 936-945	4
305	Chemoproteomic-enabled phenotypic screening. 2021 , 28, 371-393	7
304	Probes for Photoaffinity Labelling of Kinases. 2021 , 22, 2206-2218	2
303	Inflect: Optimizing Computational Workflows for Thermal Proteome Profiling Data Analysis. 2021 , 20, 1874-1888	2

302	Recent advances in identifying protein targets in drug discovery. 2021 , 28, 394-423	17
301	Progress of tubulin polymerization activity detection methods. 2021 , 37, 127698	9
300	Boosting Detection of Low-Abundance Proteins in Thermal Proteome Profiling Experiments by Addition of an Isobaric Trigger Channel to TMT Multiplexes. 2021 , 93, 7000-7010	1
299	Chemoproteomic Profiling of Covalent XPO1 Inhibitors to Assess Target Engagement and Selectivity. 2021 , 22, 2116-2123	1
298	Chd8 regulates X chromosome inactivation in mouse through fine-tuning control of Xist expression. 2021 , 4, 485	О
297	Discovery-Versus Hypothesis-Driven Detection of Protein-Protein Interactions and Complexes. 2021 , 22,	2
296	Subcellular proteomics. 2021 , 1,	13
295	A New Drug Discovery Approach Based on Thermal Proteome Profiling to Develop More Effective Drugs. 2021 , 8,	
294	Quantitative Mass Spectrometry-Based Proteomics for Biomarker Development in Ovarian Cancer. 2021 , 26,	4
293	Chemoproteomic-enabled characterization of small GTPase Rab1a as a target of an N-arylbenzdiimidazole ligand rescue of Parkinson associated cell toxicity.	1
292	Hyperthermia Selectively Destabilizes Oncogenic Fusion Proteins. 2021, 2, 388-401	13
291	Recent progress in mass spectrometry-based strategies for elucidating protein-protein interactions. 2021 , 78, 5325-5339	12
290	Thermal proteome profiling efficiently identifies ribosome destabilizing oxazolidinones. 2021 , 87, 132118	1
289	Why and How to Dig into Plant Metabolite-Protein Interactions. 2021 , 26, 472-483	12
288	Diverse Hotspot Thermal Profiling Methods Detect Phosphorylation-Dependent Changes in Protein Stability.	1
287	Principled decision-making workflow with hierarchical Bayesian models of high throughput dose-response measurements.	
286	Identification of phosphosites that alter protein thermal stability. 2021 , 18, 760-762	12
285	Thermal Analysis of a Mixture of Ribosomal Proteins by vT-ESI-MS: Toward a Parallel Approach for Characterizing the. 2021 , 93, 8484-8492	3

284	New applications of advanced instrumental techniques for the characterization of food allergenic proteins. 2021 , 1-17	3
283	Principled Decision-Making Workflow with Hierarchical Bayesian Models of High-Throughput Dose-Response Measurements. 2021 , 23,	1
282	Chemoproteomics for Plasmodium Parasite Drug Target Discovery. 2021 , 22, 2591-2599	2
281	Quantitative Chemical Proteomics Reveals Interspecies Variations on Binding Schemes of L-FABP with Perfluorooctanesulfonate. 2021 , 55, 9012-9023	O
280	Fundamentals to function: Quantitative and scalable approaches for measuring protein stability. 2021 , 12, 547-560	2
279	Systematic profiling of protein complex dynamics reveals DNA-PK phosphorylation of IFI16 en route to herpesvirus immunity. 2021 , 7,	4
278	Discovery of a Ireceptor antagonist by combination of unbiased cell painting and thermal proteome profiling. 2021 , 28, 848-854.e5	5
277	Receptor tyrosine kinases and cancer: oncogenic mechanisms and therapeutic approaches. 2021 , 40, 4079-4093	16
276	Tyr-Asp inhibition of glyceraldehyde 3-phosphate dehydrogenase affects plant redox metabolism. 2021 , 40, e106800	8
275	Thermal Proteome Profiling Reveals Glutathione Peroxidase 4 as the Target of the Autophagy Inducer Conophylline. 2021 , 100, 181-192	O
274	Conservation of metabolic regulation by phosphorylation and non-covalent small-molecule interactions. 2021 , 12, 538-546	2
273	A Bayesian semi-parametric model for thermal proteome profiling. 2021 , 4, 810	O
272	Thermal proteome profiling identifies PIP4K2A and ZADH2 as off-targets of Polo-like kinase 1 inhibitor volasertib. 2021 , 35, e21741	0
271	Protein mimetic amyloid inhibitor potently abrogates cancer-associated mutant p53 aggregation and restores tumor suppressor function. 2021 , 12, 3962	5
270	Impact of phosphorylation on thermal stability of proteins. 2021 , 18, 757-759	15
269	P38EMAPK phosphorylates Snapin and reduces Snapin-mediated BACE1 transportation in APP-transgenic mice. 2021 , 35, e21691	O
268	Emerging Therapeutic Potential of SIRT6 Modulators. 2021 , 64, 9732-9758	7
267	Assessing target engagement using proteome-wide solvent shift assays.	O

(2021-2021)

266	Rapid Evaluation of Small Molecule Cellular Target Engagement with a Luminescent Thermal Shift Assay. 2021 , 12, 1288-1294	4
265	The Impact of Assay Design on Medicinal Chemistry: Case Studies. 2021 , 26, 1243-1255	O
264	Proteomics in the pharmaceutical and biotechnology industry: a look to the next decade. 2021 , 18, 503-526	5
263	Mechanical stress induced protein precipitation method for drug target screening. 2021 , 1168, 338612	3
262	Characterization of the Heat-Stable Proteome during Seed Germination in Arabidopsis with Special Focus on LEA Proteins. 2021 , 22,	2
261	Intracellular Protein-Drug Interactions Probed by Direct Mass Spectrometry of Cell Lysates. 2021 , 60, 19637-19642	3
260	The rise of proteome-wide biophysics. 2021 , 17, e10442	1
259	Degradation of CCNK/CDK12 is a druggable vulnerability of colorectal cancer. 2021 , 36, 109394	9
258	Proteomics-based target identification of natural products affecting cancer metabolism. 2021 , 74, 639-650	2
257	IFN-Drives Human Monocyte Differentiation into Highly Proinflammatory Macrophages That Resemble a Phenotype Relevant to Psoriasis. 2021 , 207, 555-568	2
256	Intracellular Protein D rug Interactions Probed by Direct Mass Spectrometry of Cell Lysates. 2021 , 133, 19789-19794	
255	Sensitive Measurement of Drug-Target Engagement by a Cellular Thermal Shift Assay with Multiplex Proximity Extension Readout. 2021 , 93, 10999-11009	4
254	A Comparison of Two Stability Proteomics Methods for Drug Target Identification in OnePot 2D Format. 2021 , 16, 1445-1455	2
253	Alectinib treatment improves photodynamic therapy in cancer cell lines of different origin. 2021 , 21, 971	O
252	Characterizing Endogenous Protein Complexes with Biological Mass Spectrometry. 2021,	5
251	A Pan-Cancer Analysis of Transcriptome and Survival Reveals Prognostic Differentially Expressed LncRNAs and Predicts Novel Drugs for Glioblastoma Multiforme Therapy. 2021 , 12, 723725	
250	Global Mapping of Metalloproteomes. 2021 , 60, 3507-3514	1
249	Y box binding protein 1 inhibition as a targeted therapy for ovarian cancer. 2021 , 28, 1206-1220.e6	6

248	Application of omics- and multi-omics-based techniques for natural product target discovery. 2021 , 141, 111833	4
247	Natural-Product-Based Solutions for Tropical Infectious Diseases. 2021 , e0034820	1
246	Chemical Probes for Understudied Kinases: Challenges and Opportunities. 2021,	1
245	Kidney toxicity of the BRAF-kinase inhibitor vemurafenib is driven by off-target ferrochelatase inhibition. 2021 , 100, 1214-1226	2
244	Callyspongiolide kills cells by inducing mitochondrial dysfunction via cellular iron depletion. 2021 , 4, 1123	0
243	Discovery and Pharmacological Characterization of JNJ-64619178, a Novel Small-Molecule Inhibitor of PRMT5 with Potent Antitumor Activity. 2021 , 20, 2317-2328	5
242	Therapeutic Response Assessment of High-Grade Gliomas During Early-Phase Drug Development in the Era of Molecular and Immunotherapies. 2021 , 27, 395-403	
241	Novel perspectives of environmental proteomics. 2021, 788, 147588	O
240	Hsp40 Affinity to Identify Proteins Destabilized by Cellular Toxicant Exposure.	
239	Drug Target Identification in Tissues by Thermal Proteome Profiling. 2021,	2
238	Utilizing thermal proteome profiling to identify the molecular targets of anti-leishmanial compounds. 2021 , 2, 100704	1
237	Systematic analysis of chemical-protein interactions from zebrafish embryo by proteome-wide thermal shift assay, bridging the gap between molecular interactions and toxicity pathways. 2021 , 249, 104382	2
236	Flavonoid 4,4'-dimethoxychalcone suppresses cell proliferation via dehydrogenase inhibition and oxidative stress aggravation. 2021 , 175, 206-215	1
235	Organometallic Receptors and Conjugates With Biomolecules in Bioorganometallic Chemistry. 2021 ,	
234	Nephrotoxicity of the BRAF-kinase inhibitor Vemurafenib is driven by off-target Ferrochelatase inhibition.	1
233	Label-Free Target Identification and Confirmation Using Thermal Stability Shift Assays. 2021 , 2213, 163-173	1
232	Profiling of Small Molecules by Chemical Proteomics. 2016 , 1394, 211-218	10
231	Quantitative, Wide-Spectrum Kinase Profiling in Live Cells for Assessing the Effect of Cellular ATP on Target Engagement. 2018 , 25, 206-214.e11	109

230	Target identification and validation of natural products with label-free methodology: A critical review from 2005 to 2020. 2020 , 216, 107690	12
229	Temporal dynamics of protein complex formation and dissociation during human cytomegalovirus infection. 2020 , 11, 806	33
228	Chapter 8:Fluorescent Thermal Shift Assays for Identifying Small Molecule Ligands. 208-238	2
227	Chapter 5:Contemporary Techniques for Target Deconvolution and Mode of Action Elucidation. 2020 , 83-103	1
226	Mass spectrometry-based methods for structural biology on a proteome-wide scale. 2020 , 48, 945-954	1
225	The small non-coding vault RNA1-1 acts as a riboregulator of autophagy.	1
224	Impact of phosphorylation on thermal stability of proteins.	7
223	Identification of phosphosites that alter protein thermal stability.	4
222	Computational analysis of ligand dose range thermal proteome profiles.	1
221	CETSAI MS profiling for a comparative assessment of FDA approved antivirals repurposed for COVID-19 therapy identifies Trip13 as a Remdesivir off-target.	1
220	Thermal proteome profiling of breast cancer cells reveals proteasomal activation by CDK4/6 inhibitor palbociclib.	О
219	Non-parametric analysis of thermal proteome profiles reveals novel drug-binding proteins.	2
218	System-wide identification and prioritization of enzyme substrates by thermal analysis (SIESTA).	2
217	Rapid discovery of drug target engagement by isothermal shift assay.	2
216	The outer membrane lipoprotein NlpI nucleates hydrolases within peptidoglycan multi-enzyme complexes inEscherichia coli.	5
215	Analysis of Independent Differences (AID) detects complex thermal proteome profiles independent of shape and identifies candidate panobinostat targets.	1
214	Towards a systematic map of the functional role of protein phosphorylation.	6
213	Computational-experimental approach to drug-target interaction mapping: A case study on kinase inhibitors. 2017 , 13, e1005678	51

212	Outer membrane lipoprotein NlpI scaffolds peptidoglycan hydrolases within multi-enzyme complexes in Escherichia coli. 2020 , 39, e102246	36
211	Systematic mapping of protein-metabolite interactions in central metabolism of Escherichia coli. 2019 , 15, e9008	22
210	Drug mechanism-of-action discovery through the integration of pharmacological and CRISPR screens. 2020 , 16, e9405	22
209	Aggregation and disaggregation features of the human proteome. 2020 , 16, e9500	9
208	Phenotypic proteomic profiling identifies a landscape of targets for circadian clock-modulating compounds. 2019 , 2,	9
207	Analysis of Brain Protein Stability Changes in Mouse Models of Normal Aging and Esynucleinopathy Reveals Age- and Disease-Related Differences. 2021 , 20, 5156-5168	2
206	Evaluation of FOXO1 Target Engagement Using a Single-Cell Microfluidic Platform. 2021 , 93, 14659-14666	0
205	Non-conserved metabolic regulation by LKB1 distinguishes human and mouse lung adenocarcinoma.	
204	High-throughput functional characterization of protein phosphorylation sites in yeast. 2021,	3
203	Mapping paths: new approaches to dissect eukaryotic signaling circuitry. 2016 , 5,	
203	Mapping paths: new approaches to dissect eukaryotic signaling circuitry. 2016 , 5, Genome-Scale Architecture of Small Molecule Regulatory Networks and the Fundamental Trade-Off Between Regulation and Enzymatic Activity.	
	Genome-Scale Architecture of Small Molecule Regulatory Networks and the Fundamental	
202	Genome-Scale Architecture of Small Molecule Regulatory Networks and the Fundamental Trade-Off Between Regulation and Enzymatic Activity. Repurposing covalent EGFR/HER2 inhibitors for on-target degradation of human Tribbles 2 (TRIB2)	2
202	Genome-Scale Architecture of Small Molecule Regulatory Networks and the Fundamental Trade-Off Between Regulation and Enzymatic Activity. Repurposing covalent EGFR/HER2 inhibitors for on-target degradation of human Tribbles 2 (TRIB2) pseudokinase.	2 O
202 201 200	Genome-Scale Architecture of Small Molecule Regulatory Networks and the Fundamental Trade-Off Between Regulation and Enzymatic Activity. Repurposing covalent EGFR/HER2 inhibitors for on-target degradation of human Tribbles 2 (TRIB2) pseudokinase. A novel LC system embeds analytes in pre-formed gradients for rapid, ultra-robust proteomics. Post-translational modifications drive protein stability to control the dynamic beer brewing	
202 201 200	Genome-Scale Architecture of Small Molecule Regulatory Networks and the Fundamental Trade-Off Between Regulation and Enzymatic Activity. Repurposing covalent EGFR/HER2 inhibitors for on-target degradation of human Tribbles 2 (TRIB2) pseudokinase. A novel LC system embeds analytes in pre-formed gradients for rapid, ultra-robust proteomics. Post-translational modifications drive protein stability to control the dynamic beer brewing proteome.	O
202 201 200 199	Genome-Scale Architecture of Small Molecule Regulatory Networks and the Fundamental Trade-Off Between Regulation and Enzymatic Activity. Repurposing covalent EGFR/HER2 inhibitors for on-target degradation of human Tribbles 2 (TRIB2) pseudokinase. A novel LC system embeds analytes in pre-formed gradients for rapid, ultra-robust proteomics. Post-translational modifications drive protein stability to control the dynamic beer brewing proteome. ProTargetMiner: A proteome signature library of anticancer molecules for functional discovery.	O

194	The relationship between misfolding avoidance hypothesis and protein evolutionary rates in the light of empirical evidence.	О
193	LiP-Quant, an automated chemoproteomic approach to identify drug targets in complex proteomes.	
192	Temperature sensitive Mutant Proteome Profiling: a novel tool for the characterization of the global impact of missense mutations on the proteome.	О
191	The GFP thermal shift assay for screening ligand and lipid interactions to solute carrier transporters. 2021 , 16, 5357-5376	1
190	Experimental methods for dissecting the terra-incognita of protein-metabolite interactomes. 2021 , 100403	1
189	Boosting detection of low abundance proteins in thermal proteome profiling experiments by addition of an isobaric trigger channel to TMT multiplexes.	О
188	CHAPTER 10:Assays to Characterize the Cellular Pharmacology of a Chemical Probe. 2020 , 247-275	О
187	Aggregation and Disaggregation Features of the Human Proteome.	1
186	In-depth characterization of Staurosporine induced proteome thermal stability changes.	2
185	A dual-mechanism antibiotic targets Gram-negative bacteria and avoids drug resistance.	1
184	Identifying Protein-Drug Interactions in Cell Lysates Using Histidine Hydrogen Deuterium Exchange. 2021 , 93, 14985-14995	0
183	Selective cross-linking of coinciding protein assemblies by in-gel cross-linking mass-spectrometry.	
182	A tale of two tails - efficient profiling of protein degraders by specific functional and target engagement readouts.	
181	Off-target identification by chemical proteomics for the understanding of drug side effects. 2020 , 17, 695-697	1
180	Inflect: Optimizing Computational Workflows for Thermal Proteome Profiling Data Analysis.	
179	A Bayesian semi-parametric model for thermal proteome profiling.	
178	Showing NAFLD, as a key connector disease between Alzheimer's disease and diabetes via analysis of systems biology. 2020 , 13, S89-S97	
177	Chemoproteomic-enabled characterization of small GTPase Rab1a as a target of an -arylbenzimidazole ligand's rescue of Parkinson's-associated cell toxicity 2022 , 3, 96-111	2

176	Technique development of high-throughput and high-sensitivity sample preparation and separation for proteomics. 2022 , 14, 101-111	1
175	An integrative proteomics method identifies a regulator of translation during stem cell maintenance and differentiation. 2021 , 12, 6558	3
174	Illustrative Tutorials for ProThermDB: Thermodynamic Database for Proteins and Mutants. 2021 , 1, e306	
173	Unexpected cause of vemurafenib-induced nephrotoxicity: ferrochelatase. 2021 , 100, 1158-1160	
172	Comprehensive chemical proteomics analyses reveal that the new TRi-1 and TRi-2 compounds are more specific thioredoxin reductase 1 inhibitors than auranofin. 2021 , 48, 102184	3
171	Thermal Proteome Profiling to Identify Protein-ligand Interactions in the Apicomplexan Parasite. 2021 , 11, e4207	O
170	Inhibition of prolyl oligopeptidase: A promising pathway to prevent the progression of age-related macular degeneration. 2021 , 146, 112501	
169	Experimental study of proteome halophilicity using nanoDSF: a proof of concept. 2021 , 26, 1	
168	Quantitative Proteomics Using Isobaric Labeling: A Practical Guide 2022,	6
167	An update of label-free protein target identification methods for natural active products 2022 , 12, 1829-	18541
167 166	An update of label-free protein target identification methods for natural active products 2022 , 12, 1829-7 Precipitate-Supported Thermal Proteome Profiling Coupled with Deep Learning for Comprehensive Screening of Drug Target Proteins 2022 ,	2
Í	Precipitate-Supported Thermal Proteome Profiling Coupled with Deep Learning for Comprehensive	
166	Precipitate-Supported Thermal Proteome Profiling Coupled with Deep Learning for Comprehensive Screening of Drug Target Proteins 2022 ,	
166 165	Precipitate-Supported Thermal Proteome Profiling Coupled with Deep Learning for Comprehensive Screening of Drug Target Proteins 2022, A real-time cellular thermal shift assay (RT-CETSA) to monitor target engagement.	
166 165 164	Precipitate-Supported Thermal Proteome Profiling Coupled with Deep Learning for Comprehensive Screening of Drug Target Proteins 2022, A real-time cellular thermal shift assay (RT-CETSA) to monitor target engagement. A novel role of 3DEAMP in the regulation of actin cytoskeleton in Arabidopsis. Integrated analysis reveals FOXA1 and Ku70/Ku80 as direct targets of ivermectin in prostate	
166 165 164	Precipitate-Supported Thermal Proteome Profiling Coupled with Deep Learning for Comprehensive Screening of Drug Target Proteins 2022, A real-time cellular thermal shift assay (RT-CETSA) to monitor target engagement. A novel role of 3[5](EAMP in the regulation of actin cytoskeleton in Arabidopsis. Integrated analysis reveals FOXA1 and Ku70/Ku80 as direct targets of ivermectin in prostate cancer. SU086, an inhibitor of HSP90, impairs glycolysis and represents a treatment strategy for advanced	2
166 165 164 163	Precipitate-Supported Thermal Proteome Profiling Coupled with Deep Learning for Comprehensive Screening of Drug Target Proteins 2022, A real-time cellular thermal shift assay (RT-CETSA) to monitor target engagement. A novel role of 3[b]EAMP in the regulation of actin cytoskeleton in Arabidopsis. Integrated analysis reveals FOXA1 and Ku70/Ku80 as direct targets of ivermectin in prostate cancer. SU086, an inhibitor of HSP90, impairs glycolysis and represents a treatment strategy for advanced prostate cancer 2022, 3, 100502 The antimicrobial drug pyrimethamine inhibits STAT3 transcriptional activity by targeting the	2

158	High-Throughput Cellular Thermal Shift Assay Using Acoustic Transfer of Protein Lysates 2022,	2
157	Mass Spectrometry-Based Technologies for Probing the Three-Dimensional World of Plant Proteins 2022 ,	O
156	Hsp40 Affinity to Identify Proteins Destabilized by Cellular Toxicant Exposure. 2021,	0
155	A study of protein-drug interaction based on solvent-induced protein aggregation by fluorescence correlation spectroscopy 2022 ,	
154	Kurarinone alleviated Parkinson's disease via stabilization of epoxyeicosatrienoic acids in animal model 2022 , 119,	6
153	Metabolomics paves the way for improved drug target identification 2022 , 18, e10914	O
152	ProSAP: a GUI software tool for statistical analysis and assessment of thermal stability data 2022,	О
151	Past accomplishments and future challenges of the multi-omics characterization of leaf growth 2022 ,	O
150	Deep Neural Network-Assisted Drug Recommendation Systems for Identifying Potential Drug-Target Interactions 2022 , 7, 12138-12146	2
149	Evaluating Client Protein Recovery by the Hsp40s DNAJB8 and DNAJB1 with AP-MS.	Ο
148	The emerging role of mass spectrometry-based proteomics in drug discovery 2022,	10
147	In-Cell Structural Biology by NMR: The Benefits of the Atomic Scale 2022,	3
146	Secretomics-A Key to a Comprehensive Picture of Unconventional Protein Secretion 2022, 10, 878027	Ο
145	Proteome-wide cellular thermal shift assay reveals In expected cross-talk between brassinosteroid and auxin signaling 2022 , 119, e2118220119	3
144	Regulation of Plant Primary Metabolism [How Results From Novel Technologies Are Extending Our Understanding From Classical Targeted Approaches. 1-20	O
143	Learning deep representations of enzyme thermal adaptation.	
142	Drug Mechanism Enrichment Analysis: A tool to link molecular signatures with sensitivity to drug mechanisms of action.	
141	A Recent Update on Small-molecule Kinase inhibitors for Targeted Cancer Therapy and Their	3

140	Metabolite interactions in the bacterial Calvin cycle and implications for flux regulation.	0
139	Thermal Proteome Profiling Reveals the O-GlcNAc-Dependent Meltome 2022,	3
138	NMS-873 Leads to Dysfunctional Glycometabolism in A p97-Independent Manner in HCT116 Colon Cancer Cells 2022 , 14,	О
137	High-Content Screening Pipeline for Natural Products Targeting Oncogenic Signaling in Melanoma 2022 ,	O
136	Validating Small Molecule Chemical Probes for Biological Discovery 2022,	3
135	Photoaffinity labeling and bioorthogonal ligation: Two critical tools for designing "Fish Hooks" to scout for target proteins 2022 , 62, 116721	3
134	Structural Alternation in Heat Shock Proteins of Activated Macrophages 2021, 10,	1
133	Glycolytic flux-signaling controls mouse embryo mesoderm development.	
132	Proteome-wide cellular thermal shift assay reveals novel crosstalk between brassinosteroid and auxin signaling.	
131	Discovery of AL-GDa62 as a Potential Synthetic Lethal Lead for the Treatment of Gastric Cancer. 2021 ,	3
130	A specific inhibitor of ALDH1A3 regulates retinoic acid biosynthesis in glioma stem cells 2021 , 4, 1420	3
129	Assessing target engagement using proteome-wide solvent shift assays. 2021 , 10,	3
128	An overview of kinase downregulators and recent advances in discovery approaches 2021 , 6, 423	3
127	Mechanistic Insights into a CDK9 Inhibitor Via Orthogonal Proteomics Methods 2021,	1
126	Hidden information on protein function in censuses of proteome foldedness 2022 , 13, 1992	1
125	Activity-based annotation: the emergence of systems biochemistry 2022,	2
124	Matrix Thermal Shift Assay for Fast Construction of Multidimensional Ligand-Target Space 2022,	O
123	Proteome-wide identification of amino acid substitutions deleterious for protein function.	O

Unmodified methodologies in target discovery for small molecule drugs: A rising star. **2022**,

121	Chapter 5. Mass Spectrometry in Biophysics: from High Throughput Screening to Structural Biology. 87-119	
120	Ainsliadimer A Induces ROS-Mediated Apoptosis in Colorectal Cancer Cells via Directly Targeting Peroxiredoxin 1 and 2.	
119	Comprehensive Target Screening and Cellular Profiling of the Cancer-Active Compound b-AP15 Indicate Abrogation of Protein Homeostasis and Organelle Dysfunction as the Primary Mechanism of Action 2022 , 12, 852980	Ο
118	PACTS-Assisted Thermal Proteome Profiling for Use in Identifying Peptide-Interacting Proteins 2022 ,	О
117	Integrated genomics and chemical biology herald an era of sophisticated antibacterial discovery, from defining essential genes to target elucidation 2022 ,	1
116	Ion-Based Proteome-Integrated Solubility Alteration Assays for Systemwide Profiling of Protein-Molecule Interactions 2022 , 94, 7066-7074	O
115	Selective chemical probes can untangle the complexity of the plant cell endomembrane system 2022 , 68, 102223	O
114	Xlink Mapping and AnalySis (XMAS) - Smooth Integrative Modeling in ChimeraX.	1
113	Rgs16 promotes antitumor CD8 + T cell exhaustion. 2022 , 7,	1
112	Small Molecule Arranged Thermal Proximity Coaggregation (smarTPCA) A Novel Approach to Characterize Protein Protein Interactions in Living Cells by Similar Isothermal Dose Responses. 2022 , 23, 5605	0
111	De novo mapping of the apicomplexan Ca2+-responsive proteome.	
110	Analytical methods for obtaining binding parameters of drugprotein interactions: A review. 2022 , 340012	О
109	Deep thermal proteome profiling for detection of proteoforms and drug sensitivity biomarkers.	
108	Recent advances in isobaric labeling and applications in quantitative proteomics. 2100256	2
107	Flavonoid 4,4?-dimethoxychalcone induced ferroptosis in cancer cells by synergistically activating Keap1/Nrf2/HMOX1 pathway and inhibiting FECH. 2022 , 188, 14-23	2
106	System-wide profiling by proteome integral solubility alteration assay of drug residence times for target characterization.	
105	Current Advances in CETSA. 9,	1

104	FL-DTD: an integrated pipeline to predict the drug interacting targets by feedback loop-based network analysis. 2022 , 23,	
103	Integrating knowledge and omics to decipher mechanisms via large-scale models of signaling networks. 2022 , 18,	2
102	KOPI: Kinase inhibitOr Proteome Impact analysis. 2022 , 12,	
101	Recent advances in proteomics and metabolomics in plants. 2022 , 2,	O
100	Optimization of Ultrafast Proteomics Using an LC-Quadrupole-Orbitrap Mass Spectrometer with Data-Independent Acquisition.	O
99	Temporal and thermal profiling of the Toxoplasma proteome implicates parasite Protein Phosphatase 1 in the regulation of Ca2+-responsive pathways. 11,	O
98	Exploring direct and indirect targets of current antileishmanial drugs using a novel thermal proteomics profiling approach. 12,	О
97	Quantitative Proteomics Explore the Potential Targets and Action Mechanisms of Hydroxychloroquine. 2022 , 27, 5175	
96	Evaluation of a Pooling Chemoproteomics Strategy with an FDA-Approved Drug Library.	
95	Mapping the GALNT1 substrate landscape with versatile proteomics tools.	O
94	OPDA, more than just a jasmonate precursor. 2022 , 204, 113432	3
93	Polypharmacology and Natural Products. 2022 , 625-646	Ο
92	The In-Cell Western immunofluorescence assay to monitor PROTAC mediated protein degradation. 2022 ,	0
91	Highly effective identification of drug targets at proteome level by pH-dependent protein precipitation.	O
90	Quantitative Detection of Protein Splice Variants by Selected Reaction Monitoring (SRM) Mass Spectrometry. 2022 , 231-246	O
89	Polypharmacology in Drug Design and Discovery B asis for Rational Design of Multitarget Drugs. 2022 , 397-533	O
88	Chapter 1. Chemical Approaches for Beta-cell Biology. 2022 , 1-52	0
87	Discovery and Structural Characterization of Small Molecule Binders of the Human CTLH E3 Ligase Subunit GID4.	1

86	Clinical application of advanced multi-omics tumor profiling: Shaping precision oncology of the future. 2022 , 40, 920-938	2
85	Mass Spectrometry-Based Chemical Proteomics for Drug Target Discoveries. 2022 , 87, 983-994	O
84	Metabolization and sequestration of plant specialized metabolites in insect herbivores: Current and emerging approaches. 13,	О
83	Divergent polo boxes in KKT2 bind KKT1 to initiate the kinetochore assembly cascade in Trypanosoma brucei.	O
82	Integrated analysis reveals FOXA1 and Ku70/Ku80 as targets of ivermectin in prostate cancer. 2022 , 13,	O
81	A Matrix-Assisted Laser Desorption/Ionization Time-of-Flight Assay Identifies Nilotinib as an Inhibitor of Inflammation in Acute Myeloid Leukemia. 2022 , 65, 12014-12030	O
80	Cellular Target Deconvolution of Small Molecules Using a Selection-Based Genetic Screening Platform.	О
79	Coupling cellular drug-target engagement to downstream pharmacology with CeTEAM.	O
78	Real-Time Cellular Thermal Shift Assay to Monitor Target Engagement. 2022 , 17, 2471-2482	1
77	Polypharmacology-based approach for screening TCM against coinfection of Mycoplasma gallisepticum and Escherichia coli. 9,	O
76	Integrated changes in thermal stability and proteome abundance during altered nutrient states in Escherichia coli and human cells. 2100254	1
75	Databases and Tools to Investigate Protein-Metabolite Interactions. 2023 , 231-249	O
74	Monitoring drugEarget interactions through target engagement-mediated amplification on arrays and inlaitu.	О
73	Orthogonally-tunable and ER-targeting fluorophores detect avian influenza virus early infection. 2022 , 13,	O
72	Limited Proteolysis Mass Spectrometry to Identify Metabolite Protein Interactions. 2023, 69-89	0
71	Cellular Thermal Shift Assay for the Detection of Small MoleculeII arget Interactions in Arabidopsis Cells. 2023 , 21-34	O
70	Chemoproteomic Mapping of Glycolytic Targetome in Cancer Cells.	О
69	Label-free LC-MS based assay to characterize small molecule compound binding to cells. 2022 , 27, 405-412	Ο

68	PROMIS: Co-fractionation Mass Spectrometry for Analysis of Protein Metabolite Interactions. 2023 , 141-153	O
67	Proteome Integral Solubility Alteration (PISA) for High-Throughput Ligand Target Deconvolution with Increased Statistical Significance and Reduced Sample Amount. 2023 , 91-106	O
66	Solvent-Induced Protein Precipitation for Drug Target Discovery. 2023, 35-45	0
65	Parallel Analysis of Protein P rotein and ProteinMetabolite Complexes Using a Single-Step Affinity Purification. 2023 , 107-122	1
64	Divergent polo boxes in KKT2 and KKT3 initiate the kinetochore assembly cascade in Trypanosoma brucei.	0
63	Learning deep representations of enzyme thermal adaptation.	O
62	Switching of Photocatalytic Tyrosine/Histidine Labeling and Application to Photocatalytic Proximity Labeling. 2022 , 23, 11622	0
61	Nonionic surfactants can modify the thermal stability of globular and membrane proteins interfering with the thermal proteome profiling principles to identify protein targets.	O
60	Thermal proteome profiling reveals Haemonchus orphan protein HCO_011565 as a target of the nematocidal small molecule UMW-868. 13,	0
59	Rewiring of the proteinproteinmetabolite interactome during the diauxic shift in yeast. 2022 , 79,	O
58	Mitochondrial dysfunction rapidly modulates the abundance and thermal stability of cellular proteins.	0
57	Proteome integral solubility alteration assay combined with multi-criteria decision-making analysis for developing adverse outcome pathways.	O
56	Merits of Diazirine Photo-Immobilization for Target Profiling of Natural Products and Cofactors.	O
55	System-Wide Profiling by Proteome Integral Solubility Alteration Assay of Drug Residence Times for Target Characterization.	O
54	Multi-target-based polypharmacology prediction (mTPP): An approach using virtual screening and machine learning for multi-target drug discovery. 2022 , 368, 110239	0
53	Label-free target protein characterization for small molecule drugs: recent advances in methods and applications. 2023 , 223, 115107	1
52	Proteomic characterization of post-translational modifications in drug discovery.	0
51	The mycotoxin viriditoxin induces leukemia- and lymphoma-specific apoptosis by targeting mitochondrial metabolism. 2022 , 13,	O

50	Current and emerging target identification methods for novel antimalarials. 2022,	1
49	Preclinical efficacy of azacitidine and venetoclax for infant KMT2A-rearranged acute lymphoblastic leukemia reveals a new therapeutic strategy.	1
48	Thermal proteome profiling: Insights into protein modifications, associations, and functions. 2022 , 71, 102225	О
47	On the Feasibility of Using an Ultra-Fast DirectMS1 Method of Proteome-Wide Analysis for Searching Drug Targets in Chemical Proteomics. 2022 , 87, 1342-1353	1
46	Site-Specific Activity-Based Protein Profiling Using Phosphonate Handles. 2023, 22, 100455	O
45	Photoaffinity Labeling Chemistries Used to Map Biomolecular Interactions.	O
44	Proteome-wide structural changes measured with limited proteolysis-mass spectrometry: an advanced protocol for high-throughput applications.	О
43	Glycolytic flux-signaling controls mouse embryo mesoderm development. 11,	O
42	Drug discovery inspired by bioactive small molecules from nature. 1-12	1
41	Mega-scale experimental analysis of protein folding stability in biology and protein design.	O
40	Analysis of Copper-Induced Protein Precipitation across the E. coli Proteome.	О
39	Discovery of a small molecule ligand of FRS2 that inhibits invasion and tumor growth.	O
38	Profiling Protein Targets of Cellular Toxicant Exposure.	О
37	Chemical Biology Approaches Confirm MCT4 as the Therapeutic Target of a Cellular Optimized Hit.	O
36	Massive solubility changes of neuronal proteins upon simulated traumatic brain injury reveal the role of shockwave in irreversible damage.	О
35	Control of protein stability by post-translational modifications. 2023 , 14,	О
34	Elucidating Protein Ligand Interactions in Cell Lysates Using High-Throughput Hydrogen Deuterium Exchange Mass Spectrometry with Integrated Protein Thermal Depletion.	О
33	Tandem mass tag-based thermal proteome profiling for the discovery of drug-protein interactions in cancer cells. 2023 , 4, 102012	O

32	Methods to identify protein targets of metal-based drugs. 2023 , 73, 102257	1
31	Obtaining Increased Functional Proteomics Insights from Thermal Proteome Profiling through Optimized Melt Shift Calculation and Statistical Analysis.	0
30	Mass spectrometric exploration of phytohormone profiles and signaling networks. 2022,	0
29	Multi-comparative Thermal Proteome Profiling Uncovers New O-GlcNAc Proteins in a System-wide Method.	O
28	EFMC ITrends that Link Medicinal Chemistry and Chemical Biology to Translational Drug Discovery.	0
27	Advancing Targeted Protein Degradation via Multiomics Profiling and Artificial Intelligence. 2023 , 145, 2711-2732	O
26	Rhodojaponin VI indirectly targets Cav2.2 channels via N-ethylmaleimide-sensitive fusion protein to alleviate neuropathic pain. 2023 ,	0
25	PerTurboID: A targeted in situ method to measure changes in a local protein environment reveals the impact of kinase deletion on cytoadhesion in malaria causing parasites.	O
24	Solvent-induced proteome profiling for proteomic quantitation and target discovery of small molecular drugs. 2200281	0
23	Decrypting drug actions and protein modifications by dose- and time-resolved proteomics. 2023 , 380, 93-101	O
22	Mass spectrometry for mitochondrial multi-omics. 2023 , 117063	0
21	Extraction of active, contaminant degrading enzymes from soil. 2023 , 187, 104841	Ο
20	Global profiling of AMG510 modified proteins identified tumor suppressor KEAP1 as an off-target. 2023 , 26, 106080	0
19	Concentration-Dependent Enrichment Identifies Primary Protein Targets of Multitarget Bioactive Molecules. 2023 , 22, 802-811	O
18	Bioinformatics toolbox for exploring target mutation-induced drug resistance. 2023 , 24,	0
17	Nonionic Surfactants can Modify the Thermal Stability of Globular and Membrane Proteins Interfering with the Thermal Proteome Profiling Principles to Identify Protein Targets. 2023 , 95, 4033-4042	1
16	Improvedin situCharacterization of Proteome-wide Protein Complex Dynamics with Thermal Proximity Co-Aggregation.	0
15	A Comparison of Quantitative Mass Spectrometric Methods for Drug Target Identification by Thermal Proteome Profiling.	Ο

CITATION REPORT

14	Comparative Analysis of Protein Folding Stability-Based Profiling Methods for Characterization of Biological Phenotypes. 2023 , 34, 383-393	О
13	Prediction of Molecular Initiating Events for Adverse Outcome Pathways Using High-Throughput Identification of Chemical Targets. 2023 , 11, 189	O
12	Challenges and Perspectives in Target Identification and Mechanism Illustration for Chinese Medicine.	O
11	A novel antifolate suppresses growth of FPGS-deficient cells and overcomes methotrexate resistance.	O
10	The Promises of Proteomics and Metabolomics for Unravelling the Mechanism and Side Effect Landscape of Beta-Adrenoceptor Antagonists in Cardiovascular Therapeutics. 2023 , 27, 87-92	0
9	Ainsliadimer A induces ROS-mediated apoptosis in colorectal cancer cells via directly targeting peroxiredoxin 1 and 2. 2023 , 30, 295-307.e5	O
8	Multidimensional proteomics identifies molecular trajectories of cellular aging and rejuvenation.	0
7	Lactate regulates cell cycle by remodelling the anaphase promoting complex.	O
6	A Shift in Thinking: Cellular Thermal Shift Assay-Enabled Drug Discovery. 2023, 14, 369-375	0
5	Mitochondrial dysfunction rapidly modulates the abundance and thermal stability of cellular proteins. 2023 , 6, e202201805	1
4	The toxic natural product tutin causes epileptic seizures in mice by activating calcineurin. 2023, 8,	O
3	Deep thermal profiling for detection of functional proteoform groups.	O
2	Experimental strategies to improve drug-target identification in mass spectrometry-based thermal stability assays. 2023 , 6,	0
	Scability assays. Edes, o,	