

# Comprehensive List of Bioinformatics Tools and Databases

*A complete collection of tools and databases utilized in the field of Bioinformatics, curated by COMBIGS, SASTRA Deemed to be University.*

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[DDMut](#): predicting effects of mutations on protein stability using deep learning

[ARMADiLLO](#): a web server for analyzing antibody mutation probabilities

[DIANA-miRPath v4.0](#): expanding target-based miRNA functional analysis in cell-type and tissue contexts

[PhD-SNPg](#): updating a webserver and lightweight tool for scoring nucleotide variants

[CAID\\_prediction\\_portal](#): a comprehensive service for predicting intrinsic disorder and binding regions in proteins

[sfkit](#): a web-based toolkit for secure and federated genomic analysis

[GePI](#): large-scale text mining, customized retrieval and flexible filtering of gene/protein interactions

[NBBC](#): a non-B DNA burden explorer in cancer

[STellaris](#): a web server for accurate spatial mapping of single cells based on spatial transcriptomics data

[FLUXestimator](#): a webserver for predicting metabolic flux and variations using transcriptomics data

[Conserved unique peptide patterns \(CUPP\) online platform 2.0](#):

implementation of +1000 JGI fungal genomes

[SEPPA-mAb](#): spatial epitope prediction of protein antigens for mAbs

[CRISPRimmunity](#): an interactive web server for CRISPR-associated Important Molecular events and Modulators Used in geNome editing Tool identifying

[Human AGEs](#): an interactive spatio-temporal visualization and database of human archeogenomics

[NORMSEQ](#): a tool for evaluation, selection and visualization of RNA-Seq normalization methods

[The AnnotSV webserver in 2023](#): updated visualization and ranking

[SMDB](#): a Spatial Multimodal Data Browser

[ProAct](#): quantifying the differential activity of biological processes in tissues, cells, and user-defined contexts

[Genome Context Viewer \(GCV\) version 2](#): enhanced visual exploration of multiple annotated genomes

[PanDrugs2](#): prioritizing cancer therapies using integrated individual multi-omics data

[GS-SMD server for steered molecular dynamics of peptide substrates in the active site of the  \$\gamma\$ -secretase complex](#):

[PanDrugs2](#): prioritizing cancer therapies using integrated individual multi-omics data

[FunARTS, the Fungal bioActive compound Resistant Target Seeker, an exploration engine for target-directed genome mining in fungi](#):

[Abalign](#): a comprehensive multiple sequence alignment platform for B-cell receptor immune repertoires

[ProAct](#): quantifying the differential activity of biological processes in tissues, cells, and user-defined contexts

[Genome Context Viewer \(GCV\) version 2](#): enhanced visual exploration of multiple annotated genomes

[ChemMaps.com v2.0](#): exploring the environmental chemical universe

[SEanalysis 2.0](#): a comprehensive super-enhancer regulatory network analysis tool for human and mouse

[PHASTEST](#): faster than PHASTER, better than PHAST

[WebQUAST](#): online evaluation of genome assemblies

[Mol\\* Volumes and Segmentations](#): visualization and interpretation of cell imaging data alongside macromolecular structure data and biological annotations

[nCoVdocking2](#): a docking server to predict the binding modes between COVID-19 targets and its potential ligands

[MBROLE3](#): improved functional enrichment of chemical compounds for metabolomics data analysis

[GenomeFLTR](#): filtering reads made easy

[Breeze 2.0](#): an interactive web-tool for visual analysis and comparison of drug response data

[miEAA 2023](#): updates, new functional microRNA sets and improved enrichment visualizations

[MULocDeep web service for protein localization prediction and visualization at subcellular and suborganellar levels](#):

[RNAincoder](#): a deep learning-based encoder for RNA and RNA-associated interaction

[GeneRanger and TargetRanger](#): processed gene and protein expression levels across cells and tissues for target discovery

[MicrobiomeAnalyst 2.0](#): comprehensive statistical, functional and integrative analysis of microbiome data

[PEP-FOLD4](#): a pH-dependent force field for peptide structure prediction in aqueous solution

[OnTarget](#): in silico design of MiniPromoters for targeted delivery of expression

[MpoxRadar](#): a worldwide MPXV genomic surveillance dashboard

[GeneRanger and TargetRanger](#): processed gene and protein expression levels across cells and tissues for target discovery

[vissE.cloud](#): a webserver to visualise higher order molecular phenotypes from enrichment analysis

[OpenXGR](#): a web-server update for genomic summary data interpretation

[ACFIS 2.0](#): an improved web-server for fragment-based drug discovery via a dynamic screening strategy

[IRSOM2](#): a web server for predicting bifunctional RNAs

[WebTetrado](#): a webserver to explore quadruplexes in nucleic acid 3D structures

[TransCRISPR-sgRNA design tool for CRISPR/Cas9 experiments targeting specific sequence motifs](#):

[3D-GNOME 3.0](#): a three-dimensional genome modelling engine for analysing changes of promoter-enhancer contacts in the human genome

[CAVE](#): a cloud-based platform for analysis and visualization of metabolic pathways

[\$\alpha\$ Charges](#): partial atomic charges for AlphaFold structures in high quality

[GPS 6.0](#): an updated server for prediction of kinase-specific phosphorylation sites in proteins

[g:Profiler—interoperable web service for functional enrichment analysis and gene identifier mapping\\_\(2023 update\)](#):

[ChroKit](#): a Shiny-based framework for interactive analysis, visualization and integration of genomic data

[Tree Visualization By One Table \(tvBOT\)](#): a web application for visualizing, modifying and annotating phylogenetic trees

[antiSMASH 7.0](#): new and improved predictions for detection, regulation, chemical structures and visualisation

[KVFinder-web](#): a web-based application for detecting and characterizing biomolecular cavities

[Updated MS<sup>2</sup>PIP web server supports cutting-edge proteomics applications](#):

[Proksee](#): in-depth characterization and visualization of bacterial genomes

[DEPICTER2](#): a comprehensive webserver for intrinsic disorder and disorder function prediction

[PAE viewer](#): a webserver for the interactive visualization of the predicted aligned error for multimer structure predictions and crosslinks

[The LightDock Server](#): Artificial Intelligence-powered modeling of macromolecular interactions

[TCRmodel2](#): high-resolution modeling of T cell receptor recognition using deep learning

[PrismNet](#): predicting protein–RNA interaction using in vivo RNA structural information

[dbCAN3](#): automated carbohydrate-active enzyme and substrate annotation

[PANGEA](#): a new gene set enrichment tool for Drosophila and common research organisms

[OrthoVenn3](#): an integrated platform for exploring and visualizing orthologous data across genomes

[PASSer](#): fast and accurate prediction of protein allosteric sites

[Prediction of protein structures, functions and interactions using the IntFOLD7, MultiFOLD and ModFOLDdock servers](#):

[PlasMapper 3.0—a web server for generating, editing, annotating and visualizing publication quality plasmid maps](#):

[e-RNA](#): a collection of web-servers for the prediction and visualisation of RNA secondary structure and their functional features

[RNAcanvas](#): interactive drawing and exploration of nucleic acid structures

[DIANA-microT 2023](#): including predicted targets of virally encoded miRNAs

[DeepAlloDriver](#): a deep learning-based strategy to predict cancer driver mutations

[DeepNeo](#): a webserver for predicting immunogenic neoantigens

[Haplogrep 3 - an interactive haplogroup classification and analysis platform](#):

[MyGeneset.info](#): an interactive and programmatic platform for community-curated and user-created collections of genes

[AlloReverse](#): multiscale understanding among hierarchical allosteric regulations

[FuzPred](#): a web server for the sequence-based prediction of the context-dependent binding modes of proteins

[REVERSE](#): a user-friendly web server for analyzing next-generation sequencing data from in vitro selection/evolution experiments

[The Ocean Gene Atlas v2.0](#): online exploration of the biogeography and phylogeny of plankton genes

[HemI 2.0](#): an online service for heatmap illustration

[iBIS2Analyzer](#): a web server for a phylogeny-driven coevolution analysis of protein families

[CRISPRedict](#): a CRISPR-Cas9 web tool for interpretable efficiency predictions

[CircadiOmics](#): circadian omic web portal

[Web-based platform for analysis of RNA folding from high throughput chemical probing data](#):

[NetSurfP-3.0](#): accurate and fast prediction of protein structural features by protein language models and deep learning

[FUNAGE-Pro](#): comprehensive web server for gene set enrichment analysis of prokaryotes

[AllerCatPro 2.0](#): a web server for predicting protein allergenicity potential

[Identification of genome edited cells using CRISPRnano](#):

[Annotation Query \(AnnoQ\)](#): an integrated and interactive platform for large-scale genetic variant annotation

[Shiny GATOM](#): omics-based identification of regulated metabolic modules in atom transition networks

[FABIAN-variant](#): predicting the effects of DNA variants on transcription factor binding

[PRECOGx](#): exploring GPCR signaling mechanisms with deep protein representations

[MDsrv](#): visual sharing and analysis of molecular dynamics simulations

[OmicsNet 2.0](#): a web-based platform for multi-omics integration and network visual analytics

[GeCoViz](#): genomic context visualisation of prokaryotic genes from a functional and evolutionary perspective

[AutoESD](#): a web tool for automatic editing sequence design for genetic manipulation of microorganisms

[PCGA](#): a comprehensive web server for phenotype-cell-gene association analysis

[BioExcel Building Blocks Workflows \(BioBB-Wfs\), an integrated web-based platform for biomolecular simulations](#):

[EPIXplorer](#): A web server for prediction, analysis and visualization of enhancer-promoter interactions

[GrAfSS](#): a webserver for substructure similarity searching and comparisons in the structures of proteins and RNA

[5NosoAE](#): a web server for nosocomial bacterial antibiogram investigation and epidemiology survey

[3DGenBench](#): a web-server to benchmark computational models for 3D Genomics

[ExpressVis](#): a biologist-oriented interactive web server for exploring multi-omics data

[PADLOC](#): a web server for the identification of antiviral defence systems in microbial genomes

[Aquaculture Molecular Breeding Platform \(AMBP\)](#): a comprehensive web server for genotype imputation and genetic analysis in aquaculture

[RaacFold](#): a webserver for 3D visualization and analysis of protein structure by using reduced amino acid alphabets

[CB-Dock2](#): improved protein–ligand blind docking by integrating cavity detection, docking and homologous template fitting

[DrugVirus.info 2.0](#): an integrative data portal for broad-spectrum antivirals (BSA) and BSA-containing drug combinations (BCCs)

[AlignMe](#): an update of the web server for alignment of membrane protein sequences

[AlphaKnot](#): server to analyze entanglement in structures predicted by AlphaFold methods

[PrankWeb 3](#): accelerated ligand-binding site predictions for experimental and modelled protein structures

[PaintOmics 4](#): new tools for the integrative analysis of multi-omics datasets supported by multiple pathway databases

[CSM-Potential](#): mapping protein interactions and biological ligands in 3D space using geometric deep learning

[FuzDrop on AlphaFold](#): visualizing the sequence-dependent propensity of liquid–liquid phase separation and aggregation of proteins

[Dali server](#): structural unification of protein families

[CFM-ID 4.0 – a web server for accurate MS-based metabolite identification](#):

[PIER](#): web-based facilities tailored for genetic target prioritisation harnessing human disease genetics, functional genomics and protein interactions

[WebCSEA](#): web-based cell-type-specific enrichment analysis of genes

[GEOexplorer](#): a webserver for gene expression analysis and visualisation

[FuzDrop on AlphaFold](#): visualizing the sequence-dependent propensity of liquid–liquid phase separation and aggregation of proteins

[Dali server](#): structural unification of protein families

[CFM-ID 4.0 – a web server for accurate MS-based metabolite identification](#):

[Cancer driver drug interaction explorer](#):

[SynergyFinder 3.0](#): an interactive analysis and consensus interpretation of multi-drug synergies across multiple samples

[CalFitter 2.0](#): Leveraging the power of singular value decomposition to analyse protein thermostability

[IBS 2.0](#): an upgraded illustrator for the visualization of biological sequences

[SWORD2](#): hierarchical analysis of protein 3D structures

[GenePlexus](#): a web-server for gene discovery using network-based machine learning

[Secondary Metabolite Transcriptomic Pipeline \(SeMa-Trap\), an expression-based exploration tool for increased secondary metabolite production in bacteria](#):

[rna-tools.online](#): a Swiss army knife for RNA 3D structure modeling workflow

[TIRSF](#): a web server for screening gene signatures to predict Tumor immunotherapy response

[DLEB](#): a web application for building deep learning models in biological research

[RING 3.0](#): fast generation of probabilistic residue interaction networks from structural ensembles

[sRNAbench and sRNAtoolbox 2022 update](#): accurate miRNA and sncRNA profiling for model and non-model organisms

[The Quest for Orthologs orthology benchmark service in 2022](#):

[RSAT 2022](#): regulatory sequence analysis tools

[BeStSel](#): webserver for secondary structure and fold prediction for protein CD spectroscopy

[PhyloCloud](#): an online platform for making sense of phylogenomic data

[CATANA](#): an online modelling environment for proteins and nucleic acid nanostructures

[SubcellularRVis](#): a web-based tool to simplify and visualise subcellular compartment enrichment

[DEMO2](#): Assemble multi-domain protein structures by coupling analogous template alignments with deep-learning inter-domain restraint prediction

[ICARUS, an interactive web server for single cell RNA-seq analysis](#):

[BioUML—towards a universal research platform](#):

[pubmedKB](#): an interactive web server for exploring biomedical entity relations in the biomedical literature

[DEPCOD](#): a tool to detect and visualize co-evolution of protein domains

[BioTransformer 3.0—a web server for accurately predicting metabolic transformation products](#):

[SigCom LINCS](#): data and metadata search engine for a million gene expression signatures

[VRprofile2](#): detection of antibiotic resistance-associated mobilome in bacterial pathogens

[The mitoXplorer 2.0 update](#): integrating and interpreting mitochondrial expression dynamics within a cellular context

[TADeus2](#): a web server facilitating the clinical diagnosis by pathogenicity assessment of structural variations disarranging 3D chromatin structure

[MAPIYA contact map server for identification and visualization of molecular interactions in proteins and biological complexes](#):

[TeachOpenCADD 2022](#): open source and FAIR Python pipelines to assist in structural bioinformatics and cheminformatics research

[BioSimulators](#): a central registry of simulation engines and services for recommending specific tools

[DDGun](#): an untrained predictor of protein stability changes upon amino acid variants

[SuperPred 3.0](#): drug classification and target prediction—a machine learning approach

[iFeatureOmega](#): an integrative platform for engineering, visualization and analysis of features from molecular sequences, structural and ligand data sets

[GRaSP-web](#): a machine learning strategy to predict binding sites based on residue neighborhood graphs

[Deep phenotyping](#): symptom annotation made simple with SAMS

[Multi-CSAR](#): a web server for scaffolding contigs using multiple reference genomes

[BusyBee Web](#): towards comprehensive and differential composition-based metagenomic binning

[ProteinsPlus](#): a comprehensive collection of web-based molecular modeling tools

[AutozygosityMapper](#): Identification of disease-mutations in consanguineous families

[DeepLoc 2.0](#): multi-label subcellular localization prediction using protein language models

[ERMer](#): a serverless platform for navigating, analyzing, and visualizing Escherichia coli regulatory landscape through graph database

[ExPheWas](#): a platform for cis-Mendelian randomization and gene-based association scans

[GraPES](#): The Granule Protein Enrichment Server for prediction of biological condensate constituents

[KmerKeys](#): a web resource for searching indexed genome assemblies and variants

[The Galaxy platform for accessible, reproducible and collaborative biomedical analyses](#): 2022 update

[DNAzymeBuilder, a web application for in situ generation of RNA/DNA-cleaving deoxyribozymes](#):

[ANANASTRA](#): annotation and enrichment analysis of allele-specific transcription factor binding at SNPs

[patchHwork](#): a user-friendly pH sensitivity analysis web server for protein sequences and structures

[LoopGrafter](#): a web tool for transplanting dynamical loops for protein engineering

[LOMETS3](#): integrating deep learning and profile alignment for advanced protein template recognition and function annotation

[WashU Epigenome Browser update 2022](#):

[3DLigandSite](#): structure-based prediction of protein–ligand binding sites

[Search and sequence analysis tools services from EMBL-EBI in 2022](#):

[RNAspider](#): a webserver to analyze entanglements in RNA 3D structures

[ChIP-Atlas 2021 update](#): a data-mining suite for exploring epigenomic landscapes by fully integrating ChIP-seq, ATAC-seq and Bisulfite-seq data

[DAVID](#): a web server for functional enrichment analysis and functional annotation of gene lists (2021 update)

[CRISPRloci](#): comprehensive and accurate annotation of CRISPR–Cas systems

[TIMEOR](#): a web-based tool to uncover temporal regulatory mechanisms from multi-omics data

[DGLinker](#): flexible knowledge-graph prediction of disease–gene associations

[eSkip-Finder](#): a machine learning-based web application and database to identify the optimal sequences of antisense oligonucleotides for exon skipping

[Voronoia 4-ever](#):

[catRAPID omics v2.0](#): going deeper and wider in the prediction of protein–RNA interactions

[KOBAS-i](#): intelligent prioritization and exploratory visualization of biological functions for gene enrichment analysis

[Mechnetor](#): a web server for exploring protein mechanism and the functional context of genetic variants

[GPCRsignal](#): webserver for analysis of the interface between G-protein–coupled receptors and their effector proteins by dynamics and mutations

[pegIT - a web-based design tool for prime editing](#):

[DrugComb update](#): a more comprehensive drug sensitivity data repository and analysis portal

[b2bTools](#): online predictions for protein biophysical features and their conservation

[ncFANs v2.0](#): an integrative platform for functional annotation of non-coding RNAs

[MTR3D](#): identifying regions within protein tertiary structures under purifying selection

[GEPIA2021](#): integrating multiple deconvolution-based analysis into GEPIA

[snpXplorer](#): a web application to explore human SNP-associations and annotate SNP-sets

[GalaxyHeteromer](#): protein heterodimer structure prediction by template-based and ab initio docking

[snpXplorer](#): a web application to explore human SNP-associations and annotate SNP-sets

[DeepFun](#): a deep learning sequence-based model to decipher non-coding variant effect in a tissue- and cell type-specific manner

[Mergeomics 2.0](#): a web server for multi-omics data integration to elucidate disease networks and predict therapeutics

[LipidSig](#): a web-based tool for lipidomic data analysis

[The COVID-19 Data Portal](#): accelerating SARS-CoV-2 and COVID-19 research through rapid open access data sharing

[Aviator](#): a web service for monitoring the availability of web services

[Estimage](#): a webserver hub for the computation of methylation age

[NetGO 2.0](#): improving large-scale protein function prediction with massive sequence, text, domain, family and network information

[Thunor](#): visualization and analysis of high-throughput dose–response datasets

[ProLint](#): a web-based framework for the automated data analysis and visualization of lipid–protein interactions

[BRIO](#): a web server for RNA sequence and structure motif scan

[PlantDeepSEA, a deep learning-based web service to predict the regulatory effects of genomic variants in plants](#):

[CPA](#): a web-based platform for consensus pathway analysis and interactive visualization

[Graphery](#): interactive tutorials for biological network algorithms

[G2PDeep](#): a web-based deep-learning framework for quantitative phenotype prediction and discovery of genomic markers

[Preselector.uni-jena.de](#): optimize your cloning—a resource for identifying restriction enzymes for preselection reactions

[DomainViz](#): intuitive visualization of consensus domain distributions across groups of proteins

[LigAdvisor](#): a versatile and user-friendly web-platform for drug design

[AnnotSV and knotAnnotSV](#): a web server for human structural variations annotations, ranking and analysis

[MyCLADE](#): a multi-source domain annotation server for sequence functional exploration

[RunBioSimulations](#): an extensible web application that simulates a wide range of computational modeling frameworks, algorithms, and formats

[DeepGOWeb](#): fast and accurate protein function prediction on the (Semantic) Web

[MetaboAnalyst 5.0](#): narrowing the gap between raw spectra and functional insights

[KEA3](#): improved kinase enrichment analysis via data integration

[The gutSMASH web server](#): automated identification of primary metabolic gene clusters from the gut microbiota

[pLannotate](#): engineered plasmid annotation

[ProteinTools](#): a toolkit to analyze protein structures

[OmicsAnalyst](#): a comprehensive web-based platform for visual analytics of multi-omics data

[Amino Acid Interactions \(INTAA\) web server v2.0](#): a single service for computation of energetics and conservation in biomolecular 3D structures

[eVITTA](#): a web-based visualization and inference toolbox for transcriptome analysis

[CNVxplorer](#): a web tool to assist clinical interpretation of CNVs in rare disease patients

[SynLeGG](#): analysis and visualization of multiomics data for discovery of cancer 'Achilles Heels' and gene function relationships

[PERCEPTRON](#): an open-source GPU-accelerated proteoform identification pipeline for top-down proteomics

[DeepRefiner](#): high-accuracy protein structure refinement by deep network calibration

[ProteoVision](#): web server for advanced visualization of ribosomal proteins

[PredictProtein - Predicting Protein Structure and Function for 29 Years](#):

[OpenAnnotate](#): a web server to annotate the chromatin accessibility of genomic regions

[DoChAP](#): the domain change presenter

[EDGAR3.0](#): comparative genomics and phylogenomics on a scalable infrastructure

[The Dockstore](#): enhancing a community platform for sharing reproducible and accessible computational protocols

[CoffeeProt](#): an online tool for correlation and functional enrichment of systems genetics data

[InterEvDock3](#): a combined template-based and free docking server with increased performance through explicit modeling of complex homologs and integration of covariation-based contact maps

[antiSMASH 6.0](#): improving cluster detection and comparison capabilities

[The Dockstore](#): enhancing a community platform for sharing reproducible and accessible computational protocols

[LZerD webserver for pairwise and multiple protein-protein docking](#):

[ProteoSign v2](#): a faster and evolved user-friendly online tool for statistical analyses of differential proteomics

[BENZ WS](#): the Bologna ENZyme Web Server for four-level EC number annotation

[ModFOLD8](#): accurate global and local quality estimates for 3D protein models

[Proteo3Dnet](#): a web server for the integration of structural information with interactomics data

[Mol\\* Viewer](#): modern web app for 3D visualization and analysis of large biomolecular structures

[CeLaVi](#): an interactive cell lineage visualization tool

[LipidSuite](#): interactive web server for lipidomics differential and enrichment analysis

[PLIP 2021](#): expanding the scope of the protein–ligand interaction profiler to DNA and RNA

[Trips-Viz](#): an environment for the analysis of public and user-generated ribosome profiling data

[LitSuggest](#): a web-based system for literature recommendation and curation using machine learning

[PE-Designer and PE-Analyzer](#): web-based design and analysis tools for CRISPR prime editing

[miRTargetLink 2.0—interactive miRNA target gene and target pathway networks](#):

[OxDNA.org](#): a public webserver for coarse-grained simulations of DNA and RNA nanostructures

[Vaxign2](#): the second generation of the first Web-based vaccine design program using reverse vaccinology and machine learning

[ReFOLD3](#): refinement of 3D protein models with gradual restraints based on predicted local quality and residue contacts

[VirtualTaste](#): a web server for the prediction of organoleptic properties of chemical compounds

[IPC 2.0](#): prediction of isoelectric point and pKa dissociation constants

[Recognizing and validating ligands with CheckMyBlob](#):

[mmCSM-PPI](#): predicting the effects of multiple point mutations on protein–protein interactions

[ADMETlab 2.0](#): an integrated online platform for accurate and comprehensive predictions of ADMET properties

[MutationTaster2021](#):

[Interactive Tree Of Life \(iTOL\) v5](#): an online tool for phylogenetic tree display and annotation

[Arena3Dweb](#): interactive 3D visualization of multilayered networks

[miRMaster 2.0](#): multi-species non-coding RNA sequencing analyses at scale

[Expasy, the Swiss Bioinformatics Resource Portal, as designed by its users](#):

[iNetModels 2.0](#): an interactive visualization and database of multi-omics data

[TISIGNER.com](#): web services for improving recombinant protein production

[RNAProbe](#): a web server for normalization and analysis of RNA structure probing data

[SNPnexus](#): a web server for functional annotation of human genome sequence variation (2020 update)

[Tox21BodyMap](#): a webtool to map chemical effects on the human body

[HomolWat](#): a web server tool to incorporate ‘homologous’ water molecules into GPCR structures

[miRNet 2.0](#): network-based visual analytics for miRNA functional analysis and systems biology

[mirnaQC](#): a webserver for comparative quality control of miRNA-seq data

[CReSCENT](#): CanceR Single Cell ExpressioN Toolkit

[TopMatch-web](#): pairwise matching of large assemblies of protein and nucleic acid chains in 3D

[The Galaxy platform for accessible, reproducible and collaborative biomedical analyses](#): 2020 update

[FATCAT 2.0](#): towards a better understanding of the structural diversity of proteins

[piNET](#): a versatile web platform for downstream analysis and visualization of proteomics data

[mCSM-membrane](#): predicting the effects of mutations on transmembrane proteins

[PaCRISPR](#): a server for predicting and visualizing anti-CRISPR proteins

[ShiftCrypt](#): a web server to understand and biophysically align proteins through their NMR chemical shift values

[EpiRegio](#): analysis and retrieval of regulatory elements linked to genes

[ASAP 2020 update](#): an open, scalable and interactive web-based portal for (single-cell) omics analyses

[PseudoChecker](#): an integrated online platform for gene inactivation inference

[CVCDAP](#): an integrated platform for molecular and clinical analysis of cancer virtual cohorts

[NanoSPC](#): a scalable, portable, cloud compatible viral nanopore metagenomic data processing pipeline

[SYNERGxDB](#): an integrative pharmacogenomic portal to identify synergistic drug combinations for precision oncology

[Fluxer](#): a web application to compute, analyze and visualize genome-scale metabolic flux networks

[TIMER2.0 for analysis of tumor-infiltrating immune cells](#):

[3D-GNOME 2.0](#): a three-dimensional genome modeling engine for predicting structural variation-driven alterations of chromatin spatial structure in the human genome

[mmCSM-AB](#): guiding rational antibody engineering through multiple point mutations

[ARTS 2.0](#): feature updates and expansion of the Antibiotic Resistant Target Seeker for comparative genome mining

[webPSN v2.0](#): a webserver to infer fingerprints of structural communication in biomacromolecules

[RiboToolkit](#): an integrated platform for analysis and annotation of ribosome profiling data to decode mRNA translation at codon resolution

[mRNALoc](#): a novel machine-learning based in-silico tool to predict mRNA subcellular localization

[mRNALoc](#): a novel machine-learning based in-silico tool to predict mRNA subcellular localization

[ToxicoDB](#): an integrated database to mine and visualize large-scale toxicogenomic datasets

[PlaToLoCo](#): the first web meta-server for visualization and annotation of low complexity regions in proteins

[InterPred](#): a webtool to predict chemical autofluorescence and luminescence interference

[IRIS3](#): integrated cell-type-specific regulon inference server from single-cell RNA-Seq

[AnnoLnc2](#): the one-stop portal to systematically annotate novel lncRNAs for human and mouse

[Conserved unique peptide patterns \(CUPP\) online platform](#): peptide-based functional annotation of carbohydrate active enzymes

[NetMHCpan-4.1 and NetMHCIIpan-4.0](#): improved predictions of MHC antigen presentation by concurrent motif deconvolution and integration of MS MHC eluted ligand data

[AcrFinder](#): genome mining anti-CRISPR operons in prokaryotes and their viruses

[InteractomeSeq](#): a web server for the identification and profiling of domains and epitopes from phage display and next generation sequencing data

[PaccMann](#): a web service for interpretable anticancer compound sensitivity prediction

[Atomic Charge Calculator II](#): web-based tool for the calculation of partial atomic charges

[MISCAST](#): Missense variant to protein Structure Analysis web Suite

[MutaRNA](#): analysis and visualization of mutation-induced changes in RNA structure

[AlloSigMA 2](#): paving the way to designing allosteric effectors and to exploring allosteric effects of mutations

[BIOMEX](#): an interactive workflow for (single cell) omics data interpretation and visualization

[CoCoCoNet](#): conserved and comparative co-expression across a diverse set of species

[CausalMGM](#): an interactive web-based causal discovery tool

[EnzymeMiner](#): automated mining of soluble enzymes with diverse structures, catalytic properties and stabilities

[Oviz-Bio](#): a web-based platform for interactive cancer genomics data visualization

[ARIAweb](#): a server for automated NMR structure calculation

[TeamTat](#): a collaborative text annotation tool

[AWSEM-Suite](#): a protein structure prediction server based on template-guided, coevolutionary-enhanced optimized folding landscapes

[SIB Literature Services](#): RESTful customizable search engines in biomedical literature, enriched with automatically mapped biomedical concepts

[GeneTrail 3](#): advanced high-throughput enrichment analysis

[The omics discovery REST interface](#):

[COVTree](#): Coevolution in OVERlapped sequences by Tree analysis server

[The Quest for Orthologs benchmark service and consensus calls in 2020](#):

[miEAA 2.0](#): integrating multi-species microRNA enrichment analysis and workflow management systems

[miRSwitch](#): detecting microRNA arm shift and switch events

[LIST-S2](#): taxonomy based sorting of deleterious missense mutations across species

[PDBMD2CD](#): providing predicted protein circular dichroism spectra from multiple molecular dynamics-generated protein structures

[MetaPhOrs 2.0](#): integrative, phylogeny-based inference of orthology and paralogy across the tree of life

[VarFish](#): comprehensive DNA variant analysis for diagnostics and research

[MetagenoNets](#): comprehensive inference and meta-insights for microbial correlation networks

[TFmotifView](#): a webserver for the visualization of transcription factor motifs in genomic regions

[MusiteDeep](#): a deep-learning based webserver for protein post-translational modification site prediction and visualization

[NOREVA](#): enhanced normalization and evaluation of time-course and multi-class metabolomic data

[miRViz](#): a novel webserver application to visualize and interpret microRNA datasets

[Zebra2](#): advanced and easy-to-use web-server for bioinformatic analysis of subfamily-specific and conserved positions in diverse protein superfamilies

[SPEED2](#): inferring upstream pathway activity from differential gene expression

[novoPathFinder](#): a webserver of designing novel-pathway with integrating GEM-model

[OligoMinerApp](#): a web-server application for the design of genome-scale oligonucleotide in situ hybridization probes through the flexible OligoMiner environment

[Galaxy HiCExplorer 3](#): a web server for reproducible Hi-C, capture Hi-C and single-cell Hi-C data analysis, quality control and visualization

[ProteinsPlus](#): interactive analysis of protein–ligand binding interfaces

[rMAPS2](#): an update of the RNA map analysis and plotting server for alternative splicing regulation

[TREND](#): a platform for exploring protein function in prokaryotes based on phylogenetic, domain architecture and gene neighborhood analyses

[Prediction of synonymous corrections by the BE-FF computational tool expands the targeting scope of base editing](#):

[SynergyFinder 2.0](#): visual analytics of multi-drug combination synergies

[LINbase](#): a web server for genome-based identification of prokaryotes as members of crowdsourced taxa

[SuperCYPsPred—a web server for the prediction of cytochrome activity](#):

[MTR-Viewer](#): identifying regions within genes under purifying selection

[DNAvisualization.org](#): a serverless web tool for DNA sequence visualization

[WashU Epigenome Browser update 2019](#):

[LnCompare](#): gene set feature analysis for human long non-coding RNAs

[CNIT](#): a fast and accurate web tool for identifying protein-coding and long non-coding transcripts based on intrinsic sequence composition

[ORVAL](#): a novel platform for the prediction and exploration of disease-causing oligogenic variant combinations

[RNAmod](#): an integrated system for the annotation of mRNA modifications

[PRECOG](#): PREDicting COupling probabilities of G-protein coupled receptors

[PatchSearch](#): a web server for off-target protein identification

[BioUML](#): an integrated environment for systems biology and collaborative analysis of biomedical data

[IAMBEE](#): a web-service for the identification of adaptive pathways from parallel evolved clonal populations

[tRNAviz](#): explore and visualize tRNA sequence features

[PSICA](#): a fast and accurate web service for protein model quality analysis

[The Gene Sculpt Suite](#): a set of tools for genome editing

[Web 3DNA 2.0 for the analysis, visualization, and modeling of 3D nucleic acid structures](#):

[mCSM-PPI2](#): predicting the effects of mutations on protein–protein interactions

[Evolview v3](#): a webserver for visualization, annotation, and management of phylogenetic trees

[PubTator central](#): automated concept annotation for biomedical full text articles

[WebGestalt 2019](#): gene set analysis toolkit with revamped UIs and APIs

[Interactive web-based visualization and sharing of phylogenetic trees using phylogeny.IO](#):

[ChEA3](#): transcription factor enrichment analysis by orthogonal omics integration

[DaReUS-Loop](#): a web server to model multiple loops in homology models

[Protein Interaction Z Score Assessment \(PIZSA\)](#): an empirical scoring scheme for evaluation of protein–protein interactions

[Prophage Hunter](#): an integrative hunting tool for active prophages

[MEXPRESS update 2019](#):

[web-rMKL](#): a web server for dimensionality reduction and sample clustering of multi-view data based on unsupervised multiple kernel learning

[BEERE](#): a web server for biomedical entity expansion, ranking and explorations

[HNADOCK](#): a nucleic acid docking server for modeling RNA/DNA–RNA/DNA 3D complex structures

[Simple ClinVar](#): an interactive web server to explore and retrieve gene and disease variants aggregated in ClinVar database

[WebGestalt 2019](#): gene set analysis toolkit with revamped UIs and APIs

[Interactive web-based visualization and sharing of phylogenetic trees using phylogeny.IO](#):

[ChEA3](#): transcription factor enrichment analysis by orthogonal omics integration

[GEPIA2](#): an enhanced web server for large-scale expression profiling and interactive analysis

[ImmuneRegulation](#): a web-based tool for identifying human immune regulatory elements

[IEDB-AR](#): immune epitope database—analysis resource in 2019

[Geneshot](#): search engine for ranking genes from arbitrary text queries

[VOLPES](#): an interactive web-based tool for visualizing and comparing physicochemical properties of biological sequences

[sRNAbench and sRNAtoolbox 2019](#): intuitive fast small RNA profiling and differential expression

[Fido-SNP](#): the first webserver for scoring the impact of single nucleotide variants in the dog genome

[ResponseNet v.3](#): revealing signaling and regulatory pathways connecting your proteins and genes across human tissues

[ProSNEx](#): a web-based application for exploration and analysis of protein structures using network formalism

[SEPPA 3.0—enhanced spatial epitope prediction enabling glycoprotein antigens](#):

[EpiAlignment](#): alignment with both DNA sequence and epigenomic data

[Caver Web 1.0](#): identification of tunnels and channels in proteins and analysis of ligand transport

[QBiC-Pred](#): quantitative predictions of transcription factor binding changes due to sequence variants

[Doc2Hpo](#): a web application for efficient and accurate HPO concept curation

[HawkDock](#): a web server to predict and analyze the protein–protein complex based on computational docking and MM/GBSA

[Drug ReposER](#): a web server for predicting similar amino acid arrangements to known drug binding interfaces for potential drug repositioning

[Pergola-web](#): a web server for the visualization and analysis of longitudinal behavioral data using repurposed genomics tools and standards

[EPIC-TABSAT](#): analysis tool for targeted bisulfite sequencing experiments and array-based methylation studies

[Yvis](#): antibody high-density alignment visualization and analysis platform with an integrated database

[NAPS update](#): network analysis of molecular dynamics data and protein–nucleic acid complexes

[SwissTargetPrediction](#): updated data and new features for efficient prediction of protein targets of small molecules

[RegulationSpotter](#): annotation and interpretation of extratranscriptic DNA variants

[Yosshi](#): a web-server for disulfide engineering by bioinformatic analysis of diverse protein families

[MutationDistiller](#): user-driven identification of pathogenic DNA variants

[MERMAID](#): dedicated web server to prepare and run coarse-grained membrane protein dynamics

[CHOPCHOP v3](#): expanding the CRISPR web toolbox beyond genome editing

[NetGO](#): improving large-scale protein function prediction with massive network information

[iMKT](#): the integrative McDonald and Kreitman test

[LOMETS2](#): improved meta-threading server for fold-recognition and structure-based function annotation for distant-homology proteins

[DOGMA](#): a web server for proteome and transcriptome quality assessment

[VoroMQA web server for assessing three-dimensional structures of proteins and protein complexes](#):

[INGA 2.0](#): improving protein function prediction for the dark proteome

[The RNA workbench 2.0](#): next generation RNA data analysis

[AppA](#): a web server for analysis, comparison, and visualization of contact residues and interfacial waters of antibody–antigen structures and models

[modEnrichr](#): a suite of gene set enrichment analysis tools for model organisms

[MISIM v2.0](#): a web server for inferring microRNA functional similarity based on microRNA-disease associations

[AlloDriver](#): a method for the identification and analysis of cancer driver targets

[SPADE web service for prediction of allergen IgE epitopes:](#)

[MFEprimer-3.0:](#) quality control for PCR primers

[g:Profiler:](#) a web server for functional enrichment analysis and conversions of gene lists (2019 update)

[DrugComb:](#) an integrative cancer drug combination data portal

[CPGAVAS2, an integrated plastome sequence annotator and analyzer:](#)

[MAFFT-DASH:](#) integrated protein sequence and structural alignment

[OrthoVenn2:](#) a web server for whole-genome comparison and annotation of orthologous clusters across multiple species

[Cistrome-GO:](#) a web server for functional enrichment analysis of transcription factor ChIP-seq peaks

[Aggrescan3D \(A3D\) 2.0:](#) prediction and engineering of protein solubility

[MyDGR:](#) a server for identification and characterization of diversity-generating retroelements

[SeqTailor:](#) a user-friendly webserver for the extraction of DNA or protein sequences from next-generation sequencing data

[MRPrimerW2:](#) an enhanced tool for rapid design of valid high-quality primers with multiple search modes for qPCR experiments

[IntFOLD:](#) an integrated web resource for high performance protein structure and function prediction

[antiSMASH 5.0:](#) updates to the secondary metabolite genome mining pipeline

[NGPhylogeny.fr:](#) new generation phylogenetic services for non-specialists

[SEanalysis:](#) a web tool for super-enhancer associated regulatory analysis

[Updated MS<sup>2</sup>PIP web server delivers fast and accurate MS<sup>2</sup> peak intensity prediction for multiple fragmentation methods, instruments and labeling techniques](#):

[The PSIPRED Protein Analysis Workbench](#): 20 years on

[LitSense](#): making sense of biomedical literature at sentence level

[GalaxyRefine2](#): simultaneous refinement of inaccurate local regions and overall protein structure

[AutoMLST](#): an automated web server for generating multi-locus species trees highlighting natural product potential

[The EMBL-EBI search and sequence analysis tools APIs in 2019](#):

[OrganellarGenomeDRAW \(OGDRAW\) version 1.3.1](#): expanded toolkit for the graphical visualization of organellar genomes

[NetworkAnalyst 3.0](#): a visual analytics platform for comprehensive gene expression profiling and meta-analysis

[Interactive Tree Of Life \(iTOL\) v4](#): recent updates and new developments

[CircadiOmics](#): circadian omic web portal

[GeneMANIA update 2018](#):

[KnotGenome](#): a server to analyze entanglements of chromosomes

[Identification and visualization of protein binding regions with the ArDock server](#):

[MISTIC2](#): comprehensive server to study coevolution in protein families

[HMMER web server](#): 2018 update

[cgDNAweb](#): a web interface to the cgDNA sequence-dependent coarse-grain model of double-stranded DNA

[The Microbial Genomes Atlas \(MiGA\) webserver](#): taxonomic and gene diversity analysis of Archaea and Bacteria at the whole genome level

[EasyFRAP-web](#): a web-based tool for the analysis of fluorescence recovery after photobleaching data

[Galaxy HiCExplorer](#): a web server for reproducible Hi-C data analysis, quality control and visualization

[BeStSel](#): a web server for accurate protein secondary structure prediction and fold recognition from the circular dichroism spectra

[EviNet](#): a web platform for network enrichment analysis with flexible definition of gene sets

[OmicsNet](#): a web-based tool for creation and visual analysis of biological networks in 3D space

[TAM 2.0](#): tool for MicroRNA set analysis

[LOLAweb](#): a containerized web server for interactive genomic locus overlap enrichment analysis

[PSSMSearch](#): a server for modeling, visualization, proteome-wide discovery and annotation of protein motif specificity determinants

[Coloc-stats](#): a unified web interface to perform colocalization analysis of genomic features

[IUPred2A](#): context-dependent prediction of protein disorder as a function of redox state and protein binding

[CASTp 3.0](#): computed atlas of surface topography of proteins

[VarAFT](#): a variant annotation and filtration system for human next generation sequencing data

[COACH-D](#): improved protein–ligand binding sites prediction with refined ligand-binding poses through molecular docking

[The BaMM web server for de-novo motif discovery and regulatory sequence analysis](#):

[HSYMDOCK](#): a docking web server for predicting the structure of protein homo-oligomers with Cn or Dn symmetry

[GDA, a web-based tool for Genomics and Drugs integrated analysis](#):

[gRINN](#): a tool for calculation of residue interaction energies and protein energy network analysis of molecular dynamics simulations

[GIANT 2.0](#): genome-scale integrated analysis of gene networks in tissues

[PaintOmics 3](#): a web resource for the pathway analysis and visualization of multi-omics data

[SMARTIV](#): combined sequence and structure de-novo motif discovery for in-vivo RNA binding data

[eXpression2Kinases \(X2K\) Web](#): linking expression signatures to upstream cell signaling networks

[HotSpot Wizard 3.0](#): web server for automated design of mutations and smart libraries based on sequence input information

[ComplexContact](#): a web server for inter-protein contact prediction using deep learning

[CRISPRCasFinder, an update of CRISRFinder, includes a portable version, enhanced performance and integrates search for Cas proteins](#):

[TCRmodel](#): high resolution modeling of T cell receptors from sequence

[Mutalisk](#): a web-based somatic MUTation AnaLyIS toolKit for genomic, transcriptional and epigenomic signatures

[The Galaxy platform for accessible, reproducible and collaborative biomedical analyses](#): 2018 update

[Kinact](#): a computational approach for predicting activating missense mutations in protein kinases

[CellAtlasSearch](#): a scalable search engine for single cells

[SWISS-MODEL](#): homology modelling of protein structures and complexes

[GPCRM](#): a homology modeling web service with triple membrane-fitted quality assessment of GPCR models

[mTM-align](#): a server for fast protein structure database search and multiple protein structure alignment

[Freiburg RNA tools](#): a central online resource for RNA-focused research and teaching

[BAGEL4](#): a user-friendly web server to thoroughly mine RiPPs and bacteriocins

[FragFit](#): a web-application for interactive modeling of protein segments into cryo-EM density maps

[The Ocean Gene Atlas](#): exploring the biogeography of plankton genes online

[WEGO 2.0](#): a web tool for analyzing and plotting GO annotations, 2018 update

[ezTag](#): tagging biomedical concepts via interactive learning

[GlobAI Distribution of GEnetic Traits \(GADGET\) web server](#): polygenic trait scores worldwide

[dbCAN2](#): a meta server for automated carbohydrate-active enzyme annotation

[GWAS4D](#): multidimensional analysis of context-specific regulatory variant for human complex diseases and traits

[MetaboAnalyst 4.0](#): towards more transparent and integrative metabolomics analysis

[CRISPOR](#): intuitive guide selection for CRISPR/Cas9 genome editing experiments and screens

[CalFitter](#): a web server for analysis of protein thermal denaturation data

[LitVar](#): a semantic search engine for linking genomic variant data in PubMed and PMC

[CABS-flex 2.0](#): a web server for fast simulations of flexibility of protein structures

[AAI-profiler](#): fast proteome-wide exploratory analysis reveals taxonomic identity, misclassification and contamination

[AlloFinder](#): a strategy for allosteric modulator discovery and allosterome analyses

[SNPnexus](#): assessing the functional relevance of genetic variation to facilitate the promise of precision medicine

[CavityPlus](#): a web server for protein cavity detection with pharmacophore modelling, allosteric site identification and covalent ligand binding ability prediction

[RepeatsDB-lite](#): a web server for unit annotation of tandem repeat proteins

[INTERSPIA](#): a web application for exploring the dynamics of protein-protein interactions among multiple species

[HPEPDOCK](#): a web server for blind peptide–protein docking based on a hierarchical algorithm

[PANNZER2](#): a rapid functional annotation web server

[xiSPEC](#): web-based visualization, analysis and sharing of proteomics data

[InterEvDock2](#): an expanded server for protein docking using evolutionary and biological information from homology models and multimeric inputs

[SPAR](#): small RNA-seq portal for analysis of sequencing experiments

[CSAR-web](#): a web server of contig scaffolding using algebraic rearrangements

[oriTfinder](#): a web-based tool for the identification of origin of transfers in DNA sequences of bacterial mobile genetic elements

[Patscanui](#): an intuitive web interface for searching patterns in DNA and protein data

[RSAT 2018](#): regulatory sequence analysis tools 20th anniversary

[CoNekT](#): an open-source framework for comparative genomic and transcriptomic network analyses

[geno2pheno\[ngs-freq\]](#): a genotypic interpretation system for identifying viral drug resistance using next-generation sequencing data

[DynaMut](#): predicting the impact of mutations on protein conformation, flexibility and stability

[ProTox-II](#): a webserver for the prediction of toxicity of chemicals

[Oli2go](#): an automated multiplex oligonucleotide design tool

[RNApdbee 2.0](#): multifunctional tool for RNA structure annotation

[BUSCA](#): an integrative web server to predict subcellular localization of proteins

[RNApdbee 2.0](#): multifunctional tool for RNA structure annotation

[PhytoNet](#): comparative co-expression network analyses across phytoplankton and land plants

[iPath3.0](#): interactive pathways explorer v3

[MetExplore](#): collaborative edition and exploration of metabolic networks

[Oli2go](#): an automated multiplex oligonucleotide design tool

[DynaMut](#): predicting the impact of mutations on protein conformation, flexibility and stability

[ProTox-II](#): a webserver for the prediction of toxicity of chemicals

[pirScan](#): a webserver to predict piRNA targeting sites and to avoid transgene silencing in *C. elegans*

[BRepertoire](#): a user-friendly web server for analysing antibody repertoire data

[GPCR-SSFE 2.0—a fragment-based molecular modeling web tool for Class A G-protein coupled receptors](#):

[The RNA workbench](#): best practices for RNA and high-throughput sequencing bioinformatics in Galaxy

[HGVA](#): the Human Genome Variation Archive

[LimTox](#): a web tool for applied text mining of adverse event and toxicity associations of compounds, drugs and genes

[SDM](#): a server for predicting effects of mutations on protein stability

[NOREVA](#): normalization and evaluation of MS-based metabolomics data

[SCENERY](#): a web application for (causal) network reconstruction from cytometry data

[RNA-MoIP](#): prediction of RNA secondary structure and local 3D motifs from sequence data

[GeMSTONE](#): orchestrated prioritization of human germline mutations in the cloud

[CPC2](#): a fast and accurate coding potential calculator based on sequence intrinsic features

[VCF.Filter](#): interactive prioritization of disease-linked genetic variants from sequencing data

[ProteoSign](#): an end-user online differential proteomics statistical analysis platform

[HDOCK](#): a web server for protein–protein and protein–DNA/RNA docking based on a hybrid strategy

[XSuLT](#): a web server for structural annotation and representation of sequence-structure alignments

[SODA](#): prediction of protein solubility from disorder and aggregation propensity

[DEOGEN2](#): prediction and interactive visualization of single amino acid variant deleteriousness in human proteins

[RiPPMiner](#): a bioinformatics resource for deciphering chemical structures of RiPPs based on prediction of cleavage and cross-links

[complexView](#): a server for the interpretation of protein abundance and connectivity information to identify protein complexes

[GenProBiS](#): web server for mapping of sequence variants to protein binding sites

[ThreaDomEx](#): a unified platform for predicting continuous and discontinuous protein domains by multiple-threading and segment assembly

[AMMOS2](#): a web server for protein–ligand–water complexes refinement via molecular mechanics

[GeSeq – versatile and accurate annotation of organelle genomes](#):

[CSTE](#): a webserver for the Cell State Transition Expression Atlas

[PhD-SNPg](#): a webserver and lightweight tool for scoring single nucleotide variants

[Web3DMol](#): interactive protein structure visualization based on WebGL

[mTCTScan](#): a comprehensive platform for annotation and prioritization of mutations affecting drug sensitivity in cancers

[Pathview Web](#): user friendly pathway visualization and data integration

[P4P](#): a peptidome-based strain-level genome comparison web tool

[C-SPADE](#): a web-tool for interactive analysis and visualization of drug screening experiments through compound-specific bioactivity dendrograms

[IntaRNA 2.0](#): enhanced and customizable prediction of RNA–RNA interactions

[ConTra v3](#): a tool to identify transcription factor binding sites across species, update 2017

[PIGSPro](#): prediction of immunoGlobulin structures v2

[Exploring background mutational processes to decipher cancer genetic heterogeneity](#):

[agriGO v2.0](#): a GO analysis toolkit for the agricultural community, 2017 update

[DSSR-enhanced visualization of nucleic acid structures in Jmol](#):

[WebGestalt 2017](#): a more comprehensive, powerful, flexible and interactive gene set enrichment analysis toolkit

[Olelo](#): a web application for intuitive exploration of biomedical literature

[CAFE](#): aCcelerated Alignment-FrEe sequence analysis

[PharmMapper 2017 update](#): a web server for potential drug target identification with a comprehensive target pharmacophore database

[DynOmics](#): dynamics of structural proteome and beyond

[The EBI search engine](#): EBI search as a service—making biological data accessible for all

[The Antibiotic Resistant Target Seeker \(ARTS\), an exploration engine for antibiotic cluster prioritization and novel drug target discovery](#):

[I-TASSER-MR](#): automated molecular replacement for distant-homology proteins using iterative fragment assembly and progressive sequence truncation

[ProteinsPlus](#): a web portal for structure analysis of macromolecules

[BusyBee Web](#): metagenomic data analysis by bootstrapped supervised binning and annotation

[COFACTOR](#): improved protein function prediction by combining structure, sequence and protein–protein interaction information

[BIS2Analyzer](#): a server for co-evolution analysis of conserved protein families

[IslandViewer 4](#): expanded prediction of genomic islands for larger-scale datasets

[BepiPred-2.0](#): improving sequence-based B-cell epitope prediction using conformational epitopes

[RegulatorTrail](#): a web service for the identification of key transcriptional regulators

[COFACTOR](#): improved protein function prediction by combining structure, sequence and protein–protein interaction information

[Interactive microbial distribution analysis using BioAtlas](#):

[The pepATTRACT web server for blind, large-scale peptide–protein docking](#):

[HH-MOTiF](#): de novo detection of short linear motifs in proteins by Hidden Markov Model comparisons

[SpartaABC](#): a web server to simulate sequences with indel parameters inferred using an approximate Bayesian computation algorithm

[kpLogo](#): positional k-mer analysis reveals hidden specificity in biological sequences

[WoPPER](#): Web server for Position Related data analysis of gene Expression in Prokaryotes

[ModFOLD6](#): an accurate web server for the global and local quality estimation of 3D protein models

[GASS-WEB](#): a web server for identifying enzyme active sites based on genetic algorithms

[SBSPKSV2](#): structure-based sequence analysis of polyketide synthases and non-ribosomal peptide synthetases

[antiSMASH 4.0—improvements in chemistry prediction and gene cluster boundary identification](#):

[PRISM 3](#): expanded prediction of natural product chemical structures from microbial genomes

[SeMPI](#): a genome-based secondary metabolite prediction and identification web server

[plantiSMASH](#): automated identification, annotation and expression analysis of plant biosynthetic gene clusters

[PMut](#): a web-based tool for the annotation of pathological variants on proteins, 2017 update

[TraitRateProp](#): a web server for the detection of trait-dependent evolutionary rate shifts in sequence sites

[A PanorOmic view of personal cancer genomes](#):

[The Bologna Annotation Resource \(BAR 3.0\)](#): improving protein functional annotation

[FireProt](#): web server for automated design of thermostable proteins

[MicrobiomeAnalyst](#): a web-based tool for comprehensive statistical, visual and meta-analysis of microbiome data

[GWAB](#): a web server for the network-based boosting of human genome-wide association data

[LigParGen web server](#): an automatic OPLS-AA parameter generator for organic ligands

[Gene ORGANizer](#): linking genes to the organs they affect

[Programmatic access to bioinformatics tools from EMBL-EBI update](#): 2017

[TRAPP webserver](#): predicting protein binding site flexibility and detecting transient binding pockets

[DNAproDB](#): an interactive tool for structural analysis of DNA–protein complexes

[minepath.org](#): a free interactive pathway analysis web server

[NNAlign](#): a platform to construct and evaluate artificial neural network models of receptor–ligand interactions

[GEPIA](#): a web server for cancer and normal gene expression profiling and interactive analyses

[GibbsCluster](#): unsupervised clustering and alignment of peptide sequences

[ReFOLD](#): a server for the refinement of 3D protein models guided by accurate quality estimates

[Omicseq](#): a web-based search engine for exploring omics datasets

[GalaxyHomomer](#): a web server for protein homo-oligomer structure prediction from a monomer sequence or structure

[SLIMSearch](#): a framework for proteome-wide discovery and annotation of functional modules in intrinsically disordered regions

[The Proteins API](#): accessing key integrated protein and genome information

[mCSM–NA](#): predicting the effects of mutations on protein–nucleic acids interactions

[The RING 2.0 web server for high quality residue interaction networks](#):

[Heatmapper](#): web-enabled heat mapping for all

[QuadBase2](#): web server for multiplexed guanine quadruplex mining and visualization

[3D-GNOME](#): an integrated web service for structural modeling of the 3D genome

[incaRNAfbinv](#): a web server for the fragment-based design of RNA sequences

[MoRFchibi SYSTEM](#): software tools for the identification of MoRFs in protein sequences

[tRNAscan-SE On-line](#): integrating search and context for analysis of transfer RNA genes

[PSSweb](#): protein structural statistics web server

[PASMet](#): a web-based platform for prediction, modelling and analyses of metabolic systems

[Breaking-Cas—interactive design of guide RNAs for CRISPR-Cas experiments for ENSEMBL genomes](#):

[MutaBind estimates and interprets the effects of sequence variants on protein–protein interactions](#):

[BetaSCPWeb](#): side-chain prediction for protein structures using Voronoi diagrams and geometry prioritization

[KeyPathwayMinerWeb](#): online multi-omics network enrichment

[NAPS](#): Network Analysis of Protein Structures

[PathwAX](#): a web server for network crosstalk based pathway annotation

[GREAT](#): a web portal for Genome Regulatory Architecture Tools

[PHASTER](#): a better, faster version of the PHAST phage search tool

[The Galaxy platform for accessible, reproducible and collaborative biomedical analyses](#): 2016 update

[Actionable pathways](#): interactive discovery of therapeutic targets using signaling pathway models

[Evolview v2](#): an online visualization and management tool for customized and annotated phylogenetic trees

[Protein Frustratometer 2](#): a tool to localize energetic frustration in protein molecules, now with electrostatics

[3Drefine](#): an interactive web server for efficient protein structure refinement

[Rtools](#): a web server for various secondary structural analyses on single RNA sequences

[Dali server update](#):

[MRE](#): a web tool to suggest foreign enzymes for the biosynthesis pathway design with competing endogenous reactions in mind

[The MPI bioinformatics Toolkit as an integrative platform for advanced protein sequence and structure analysis](#):

[PHYLOViZ Online](#): web-based tool for visualization, phylogenetic inference, analysis and sharing of minimum spanning trees

[CoinFold](#): a web server for protein contact prediction and contact-assisted protein folding

[PSI/TM-Coffee](#): a web server for fast and accurate multiple sequence alignments of regular and transmembrane proteins using homology extension on reduced databases

[USR-VS](#): a web server for large-scale prospective virtual screening using ultrafast shape recognition techniques

[SensiPath](#): computer-aided design of sensing-enabling metabolic pathways

[EXPLoRA-web](#): linkage analysis of quantitative trait loci using bulk segregant analysis

[miRNet - dissecting miRNA-target interactions and functional associations through network-based visual analysis](#):

[SL2](#): an interactive webtool for modeling of missing segments in proteins

[GIANT API](#): an application programming interface for functional genomics

[Pharmit](#): interactive exploration of chemical space

[Interactive tree of life \(iTOL\) v3](#): an online tool for the display and annotation of phylogenetic and other trees

[SimRNAweb](#): a web server for RNA 3D structure modeling with optional restraints

[W-IQ-TREE](#): a fast online phylogenetic tool for maximum likelihood analysis

[RBscore&NBench](#): a high-level web server for nucleic acid binding residues prediction with a large-scale benchmarking database

[DeepBlue epigenomic data server](#): programmatic data retrieval and analysis of epigenome region sets

[MBROLE 2.0—functional enrichment of chemical compounds](#):

[deepTools2](#): a next generation web server for deep-sequencing data analysis

[RNAAssess—a web server for quality assessment of RNA 3D structures](#):

[R3D-2-MSA](#): the RNA 3D structure-to-multiple sequence alignment server

[CATNAP](#): a tool to compile, analyze and tally neutralizing antibody panels

[SNiPlay3](#): a web-based application for exploration and large scale analyses of genomic variations

[I-COMS](#): Interprotein-CORrelated Mutations Server

[RNAiFold 2.0](#): a web server and software to design custom and Rfam-based RNA molecules

[INGA](#): protein function prediction combining interaction networks, domain assignments and sequence similarity

[sRNAtoolbox](#): an integrated collection of small RNA research tools

[DeAnnCNV](#): a tool for online detection and annotation of copy number variations from whole-exome sequencing data

[PatternQuery](#): web application for fast detection of biomacromolecular structural patterns in the entire Protein Data Bank

[StemChecker](#): a web-based tool to discover and explore stemness signatures in gene sets

[LYRA, a webserver for lymphocyte receptor structural modeling](#):

[Primerize](#): automated primer assembly for transcribing non-coding RNA domains

[Multidimensional gene search with Genehopper](#):

[StarScan](#): a web server for scanning small RNA targets from degradome sequencing data

[MS2PIP prediction server](#): compute and visualize MS2 peak intensity predictions for CID and HCD fragmentation

[APPRIS WebServer and WebServices](#):

[MyProteinNet](#): build up-to-date protein interaction networks for organisms, tissues and user-defined contexts

[SIFTER search](#): a web server for accurate phylogeny-based protein function prediction

[CSI 3.0](#): a web server for identifying secondary and super-secondary structure in proteins using NMR chemical shifts

[PrionW](#): a server to identify proteins containing glutamine/asparagine rich prion-like domains and their amyloid cores

[Web-Beagle](#): a web server for the alignment of RNA secondary structures

[PrionW](#): a server to identify proteins containing glutamine/asparagine rich prion-like domains and their amyloid cores

[ZCURVE 3.0](#): identify prokaryotic genes with higher accuracy as well as automatically and accurately select essential genes

[DIANA-miRPath v3.0](#): deciphering microRNA function with experimental support

[NPDock](#): a web server for protein–nucleic acid docking

[ClustVis](#): a web tool for visualizing clustering of multivariate data using Principal Component Analysis and heatmap

[The TOPCONS web server for consensus prediction of membrane protein topology and signal peptides](#):

[IMP 2.0](#): a multi-species functional genomics portal for integration, visualization and prediction of protein functions and networks

[GalaxyPepDock](#): a protein–peptide docking tool based on interaction similarity and energy optimization

[CATH FunFHMMer web server](#): protein functional annotations using functional family assignments

[RNA-Redesign](#): a web server for fixed-backbone 3D design of RNA

[OrthoVenn](#): a web server for genome wide comparison and annotation of orthologous clusters across multiple species

[SCUDO](#): a tool for signature-based clustering of expression profiles

[Pse-in-One](#): a web server for generating various modes of pseudo components of DNA, RNA, and protein sequences

[PhyloGene server for identification and visualization of co-evolving proteins using normalized phylogenetic profiles](#):

[NaviCell Web Service for network-based data visualization:](#)

[xVis](#): a web server for the schematic visualization and interpretation of crosslink-derived spatial restraints

[PockDrug-Server](#): a new web server for predicting pocket druggability on holo and apo proteins

[ProtPhylo](#): identification of protein–phenotype and protein–protein functional associations via phylogenetic profiling

[The MEME Suite](#):

[SELPHI](#): correlation-based identification of kinase-associated networks from global phospho-proteomics data sets

[antiSMASH 3.0—a comprehensive resource for the genome mining of biosynthetic gene clusters](#):

[PACCMIT/PACCMIT-CDS](#): identifying microRNA targets in 3' UTRs and coding sequences

[CCTOP](#): a Consensus Constrained TOPology prediction web server

[FlyNet](#): a versatile network prioritization server for the Drosophila community

[CABS-dock web server for the flexible docking of peptides to proteins without prior knowledge of the binding site](#):

[HMMER web server](#): 2015 update

[\(PS\)2](#): protein structure prediction server version 3.0

[TFmiR](#): a web server for constructing and analyzing disease-specific transcription factor and miRNA co-regulatory networks

[MapMyFlu](#): visualizing spatio-temporal relationships between related influenza sequences

[ENTM](#): a server for predicting functional networks of tissues in mouse

[NFFinder](#): an online bioinformatics tool for searching similar transcriptomics experiments in the context of drug repositioning

[RNAPattMatch](#): a web server for RNA sequence/structure motif detection based on pattern matching with flexible gaps

[Galahad](#): a web server for drug effect analysis from gene expression

[Pathways with PathWhiz](#):

[Localize.pytom](#): a modern webserver for cryo-electron tomography

[PUG-SOAP and PUG-REST](#): web services for programmatic access to chemical information in PubChem

[NGL Viewer](#): a web application for molecular visualization

[PolySearch2](#): a significantly improved text-mining system for discovering associations between human diseases, genes, drugs, metabolites, toxins and more

[i-cisTarget 2015 update](#): generalized cis-regulatory enrichment analysis in human, mouse and fly

[IslandViewer 3](#): more flexible, interactive genomic island discovery, visualization and analysis

[ChIP-Array 2](#): integrating multiple omics data to construct gene regulatory networks

[Stock-based detection of protein oligomeric states in jsPISA](#):

[Introducing the PRIDE Archive RESTful web services](#):

[BetaCavityWeb](#): a webserver for molecular voids and channels

[RSAT 2015](#): Regulatory Sequence Analysis Tools

[Babelomics 5.0](#): functional interpretation for new generations of genomic data

[The BioMart community portal](#): an innovative alternative to large, centralized data repositories

[MetaboAnalyst 3.0—making metabolomics more meaningful](#):

[miRiadne](#): a web tool for consistent integration of miRNA nomenclature

[RBO Aleph](#): leveraging novel information sources for protein structure prediction

[pyDockSAXS](#): protein–protein complex structure by SAXS and computational docking

[The iceLogo web server and SOAP service for determining protein consensus sequences](#):

[Assessing the impact of mutations found in next generation sequencing data over human signaling pathways](#):

[webSDA](#): a web server to simulate macromolecular diffusional association

[I-TASSER server](#): new development for protein structure and function predictions

[ChEMBL web services](#): streamlining access to drug discovery data and utilities

[FAF-Drugs3](#): a web server for compound property calculation and chemical library design

[AGGRESCAN3D \(A3D\)](#): server for prediction of aggregation properties of protein structures

[ENCoM server](#): exploring protein conformational space and the effect of mutations on protein function and stability

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[Assessing the impact of mutations found in next generation sequencing data over human signaling pathways](#):

[webSDA](#): a web server to simulate macromolecular diffusional association

[I-TASSER server](#): new development for protein structure and function predictions

[ChEMBL web services](#): streamlining access to drug discovery data and utilities

[NGS-eval](#): NGS Error analysis and novel sequence VArant detection tool

[PheNetic](#): network-based interpretation of molecular profiling data

[PLIP](#): fully automated protein–ligand interaction profiler

[MTiOpenScreen](#): a web server for structure-based virtual screening

[SANSParallel](#): interactive homology search against Uniprot

[TCS](#): a web server for multiple sequence alignment evaluation and phylogenetic reconstruction

[The EBI Search engine](#): providing search and retrieval functionality for biological data from EMBL-EBI

[WAXSiS](#): a web server for the calculation of SAXS/WAXS curves based on explicit-solvent molecular dynamics

[The EMBL-EBI bioinformatics web and programmatic tools framework](#):

[IntFOLD](#): an integrated server for modelling protein structures and functions from amino acid sequences

[RiceNet v2](#): an improved network prioritization server for rice genes