






Ранитидин: Ранитидин (Ranitidine) — это лекарственное вещество, которое относится к группе **H₂-гистаминоблокаторов**. Оно **снижает выработку соляной кислоты в желудке**. Отозван из-за обнаружения в некоторых партиях канцерогенной примеси NDMA.

 pubchem.ncbi.nlm.nih.gov/#query=ranitidine 

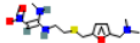
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SEARCH FOR

ranitidine  

Treating this as a text search.

BEST MATCH



ranitidine; 66357-35-5; Ranitidine Base; ZANTAC; Coralen; Gastrial; Quantor; Rantidine; ...

Compound CID: 3001055

MF: [C₁₃H₂₂N₄O₃S](#) MW: 314.41 g/mol

IUPAC Name: (E)-1-N'-[2-[[5-[(dimethylamino)methyl]furan-2-yl]methylsulfanyl]ethyl]-1-N-methyl-2-nitroethene-1,1-diamine

SMILES: [CN/C\(=C\\[N+\]\(=O\)\[O-\]\)/NCCSCC1=CC=C\(O1\)CN\(C\)C](#)

InChIKey: [VMXUWOKSQNHOCA-UKTHLTGXSA-N](#)

InChI: [InChI=1S/C13H22N4O3S/c1-14-13\(9-17\(18\)19\)15-6-7-21-10-12-5-4-11\(20-12\)8-16\(2\)3/h4-5,9,14-15H,6-8,10H2,1-3H3/b13-9+](#)

Create Date: 2005-03-27

[Summary](#) [Similar Structures Search](#) [Related Records](#) [PubMed \(MeSH Keyword\)](#)

Compounds (53)	Substances (876)	Pathways (1)	BioAssays (266)	Literature (7,487)	Patents (877)
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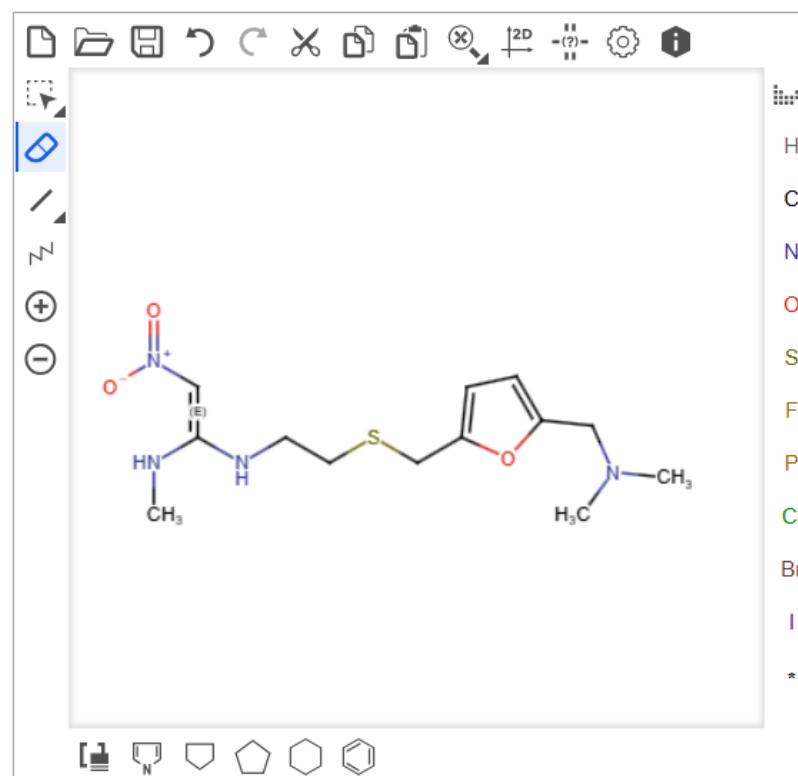
For information: We have changed the look and feel of our tool. However, we have **NOT** changed the underlying technologies and parameters. Consequently, this updated Web tool provides exactly the same results as the previous version.

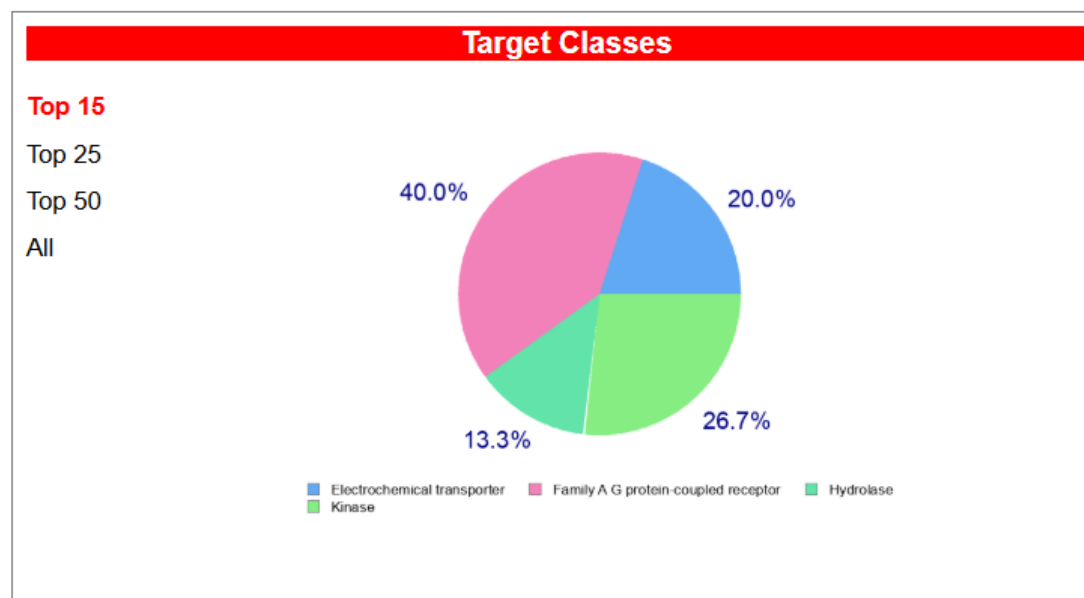
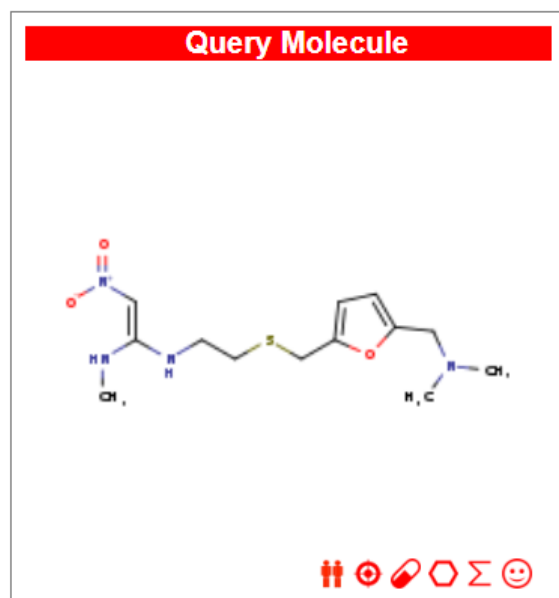
Select a species

- ☒ Homo sapiens
☐ Mus musculus
☐ Rattus norvegicus

Paste a SMILES in this box, or draw a moleculeExamples: ▼

=

(Can take up to one minute)

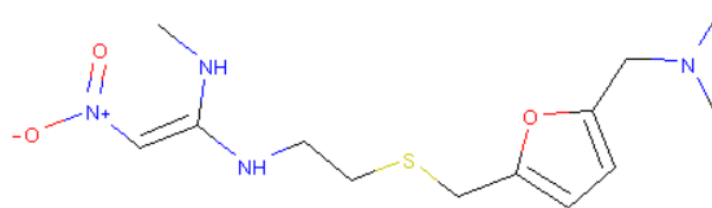


Export results:

Show entries

Search:

Target	Common name	Uniprot ID	ChEMBL ID	Target Class	Probability*	Known actives (3D/2D)
Multidrug and toxin extrusion protein 1	SLC47A1	Q96FL8	CHEMBL1743126	Electrochemical transporter	<div><div></div></div>	6 / 1
Histamine H2 receptor	HRH2	P25021	CHEMBL1941	Family A G protein-coupled receptor	<div><div></div></div>	18 / 1
Acetylcholinesterase	ACHE	P22303	CHEMBL220	Hydrolase	<div><div></div></div>	234 / 41
Muscarinic acetylcholine receptor M2 (by homology)	CHRM2	P08172	CHEMBL211	Family A G protein-coupled receptor	<div><div></div></div>	279 / 30

Input structure **RANITIDINE**

is available in ChEMBL.

Known strong binders

[Copy](#) [CSV](#) [Excel](#) [PDF](#) [Print](#)Search:

Target Name	ChEMBL-ID	UniProt ID	PDB Visualization	TTD ID	Min Activity	Assay type
Histamine H2 receptor	CHEMBL1941	P25021	Not Available	Not Available	63 nm	Kd
Prelamin-A/C	CHEMBL1293235	P02545	6JLB	Not Available	79.4 nm	Potency

All-In-One Tools 2

Software Categories

» Known miRNA Identification 33

» isomiRs Identification 13

» Novel miRNA/Precursor Analysis 45

» Differential Expression Analysis 20

» Target Prediction 50

» Target Functional Analysis 21

» miRNA-SNP Analysis 6

Other Tools 18

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Prediction Features ?

3' Compensatory PairingConservationLocal AU Content

Target-site AbundanceSite AccessibilitySeed Match

Machine LearningG:U Pairs Allowed In SeedFree Energy

Target ?

Region ?

3' UTR5' UTRCDS

AnnotationEnrichment Analysis

Tool Availability

OnlineLocalPrecomputed Data Available For Download

NGS Data Needed ?

YesNo

Metaserver ?

YesNo

Clear Filtering

Filter

Number of tools found: 38

Mode

Browse

Search

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#	Name ?	Description	Online/Local ?	Organism Specific	Algorithm Features	NGS Data Needed ?	Target Annotation ?	References ?
		BioVLAB-MMIA-NGS is Cloud-based miRNA mRNA integrated analysis system			<ul style="list-style-type: none"> seed match conservation free energy 			