RSAT	WSP	ortT	vpe

Service Documentation

Web services for the Regulatory Sequence Analysis Tools (RSAT). Tools developed by Jacques van Helden (jvanheld@scmbb.ulb.ac.be), SOAP/WSDL interface developed by Olivier Sand (oly@scmbb.ulb.ac.be).

Server Address

http://rsat.scmbb.ulb.ac.be/rsat/web_services/RSATWS.cgi

Method retrieve_seq

Description Returns upstream, downstream or coding DNA sequences for list of query genes.

Parameters Input Parameters

> Return type. Accepted values: 'server' (result is stored on a file on the server), 'client' (result is directly transferred to the client), or 'both'. Default is 'both'. Type = String output

organism Organism. Words need to be underscore separated (example: Escherichia_coli_K12).

(required) Type = String

A list of query genes Type = String query

Return sequences for all the genes of the organism if value = 1. Incompatible with query. Type = Integer all

noorf Prevent overlap with upstream open reading frames (ORF) if value = 1.

Inferior limit of the region to retrieve. Default is organism dependant (example: Saccharomyces cerevisiae = -800). from

Type = Integer

Superior limit of the region to retrieve. Default is '-1'. Type = Integer to

feattype Type of genome features to load. Supported: CDS, mRNA, tRNA, rRNA.

Sequence type. Supported: upstream, downstream, ORF (unspliced open reading frame). Type = String $\,$ type

Sequence format. Supported: IG (Intelligenetics), WC (wconsensus), raw, FastA format Type = String

lw Line width (0 for whole sequence on one line).

Type = Integer

 $Field(s) \ to \ be \ used \ in \ the \ sequence \ label. \ Multiple \ fields \ can \ be \ specified, \ separated \ by \ commas. \\ Supported: \ id, \ name, \ organism_name, \ sequence_type, \ current_from, \ current_to, \ ctg, \ orf_strand, \ reg_left, \ reg_right. \ Default: \ name. \\ Type = String$ label

label_sep Separator between the label fields. Default: I (pipe character). Type = String

No comments if value = 1. Only the identifier and the sequence are returned. By default, the comment indicates the ORF and upstream sequence coordinates. nocom

Use the repeat masked version of the genome if value = 1. Attention: repeated regions are annotated for some genomes only. repeat

Type = Integer

Admit imprecise positions if value = 1. Type = Integer imp_pos

Output Parameters

Location of the result file on the server. This can be used as input for a further request. server

Type = String

command The stand alone command executed on the server.

 $\mathbf{Type} = \mathbf{String}$

client The results.

Type = String

Method purge seq

Description Mask repeated fragments of an input sequence.

Parameters Input Parameters

> output Return type. Accepted values: 'server' (result is stored on a file on the server), 'client' (result is directly transferred to the client), or 'both'. Default is 'both'.

Type = String

Sequence to purge. You need to supply either this parameter or the next one (tmp_infile). Type = String $\,$ sequence

tmp_infile Name of the file with input sequence on the server. You need to supply either this parameter or

the previous one (sequence).

Type = String

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Sequence format. Supported: IG (Intelligenetics), WC (w
consensus), raw, Fast
A. Type = String format

Minimal match length. Default is 40. match length

Type = Integer

mismatch Number of mismatches allowed. Default is 3.

Type = Integer

Discard duplications on the direct strand only (1) or on the reverse complement as well (2). Default is 2. str

Type = Integer

Delete repeats instead of masking them if value = 1. Type = Integer delete

mask_short Mask (replace by N characters) sequences shorter than the specified length.

Type = Integer

Output Parameters

Location of the result file on the server. This can be used as input for a further request.

Type = String

command The stand alone command executed on the server.

Type = String

The results. Type = String client

Method oligo_analysis

Analysis of the statistical significance of all the oligomers of a given size in a sequence. Commonly used to detect over-represented oligonucleotides in a set of promoter sequences. Description

Parameters Input Parameters

Return type. Accepted values: 'server' (result is stored on a file on the server), 'client' (result is directly transferred to the client), or 'both'. Default is 'both'.

Type = String output

Input sequence. You need to supply either this parameter or the next one (tmp_infile). sequence

Type = String

tmp_infile Name of the file with input sequence on the server. You need to supply either this parameter or

the previous one (sequence). Type = String

format Input sequence format. Supported: IG (Intelligenetics), WC (wconsensus), raw, fasta. Default is

Type = String

Oligomer length. length

(required) Type = Integer

organism Organism. Words need to be underscore separated (example: Escherichia_coli_K12).

Type = String (required)

Background model: Type of sequences used as background model for estimating expected background

oligonucleotide frequencies. Supported: upstream, upstreamL, upstream-noorf, intergenic,

stats List of statistics to return. Supported:occ, mseq, freq, proba, ratio, zscore, like, pos, rank.

No overlapping of oligos allowed if value = 1. Disable the detection of overlapping matches for self-overlapping patterns (ex TATATA, GATAGA). noov

Type = Integer

Oligonucleotide occurrences found on both stands are summed (2) or not (1). Default is 2. str

Type = Integer

Sort oligomers according to overrepresentation if value = 1. Type = Integer sort

lth Lower threshold on some parameter. Format='param value'.

Type = String

Output Parameters

Location of the result file on the server. This can be used as input for a further request. Type = String

command The stand alone command executed on the server.

Type = String

client The results.

Type = String

Method dyad_analysis

Analysis of the statistical significance of all the spaced dyads of a given size in a sequence. Commonly used to detect over-represented spaced dyads in a set of promoter sequences. Description

Parameters	Input Parame	Input Parameters			
	output	Return type. Accepted values: 'server' (result is stored on a file on the server), 'client' (result i directly transferred to the client), or 'both'. Default is 'both'. Type = String			
	sequence	Input sequence. You need to supply either this parameter or the next one (tmp_infile). Type = String			
	tmp_infile	Name of the file with input sequence on the server. You need to supply either this parameter or the previous one (sequence). Type = String			
l	format	Input sequence format. Supported: IG (Intelligenetics), WC (wconsensus), raw, fasta. Default is fasta. Type = String			
	length (required)	Dyad length. Type = Integer			
	spacing (required)	Spacing between elements of the dyads. Type = String			
	organism (required)	Organism. Words need to be underscore separated (example: Escherichia_coli_K12). Type = String			
	background	Background model: Type of sequences used as background model for estimating expected oligonucleotide frequencies. Supported: upstream, upstreamL, upstream-noorf, intergenic, input. Type = String			
	stats	List of statistics to return. Supported:occ, mseq, freq, proba, ratio, zscore, like, pos, rank. Type = String			
	type	dyad_type (dr ir rep any) In order to fasten execution, the program can be asked to restrict its analysis to symmetric dyads. Four types are accepted: dr - direct repeats: the second element is the same as the first one; ir - inverted repeats: the second element is the reverse complement of the first one; rep - repeats: direct and inverted repeats are evaluated, any - (default) When selecting the option any, the analysis is performed on all non-symmetric dyads as well. Type = String			
	noov	No overlapping of dyads allowed if value = 1. Disable the detection of overlapping matches for self-overlapping patterns (ex TATATA, GATAGA). Type = Integer			
	str	Dyad occurrences found on both stands are summed (2) or not (1). Default is 2. Type = Integer			
	sort	Sort dyads according to overrepresentation if value = 1. Type = Integer			
	under	Detect under-represented instead of over-represented dyads (left tail significance test) if value = 1. Type = Integer			
	two_tails	Detect under-represented and over-represented dyads (two-fail significance test) if value = 1. Type = Integer			
	zeroocc	Report also dyads with zero occurrences (provided they fit the other thresholds) if value = 1. By default, the program reports only patterns present in the sequence. If the left tail or two-tail test is applied, patterns with zero occurrences are automatically taken into account. In some other cases, one would also like to detect patterns absent from the sequence. This is the function of the option -zeroocc. Type = Integer			
	lth	Lower threshold on some parameter. Format='param value'. Type = String			
	uth	Upper threshold on some parameter. Format='param value'. Type = String			
	Outnut Param	Output Parameters			
	server L	•			
		The stand alone command executed on the server. Type = String			
		The results. Type = String			
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Method	dna_pattern				

Searches all occurrences of a pattern within DNA sequences. Description

Parameters Input Parameters

output

Return type. Accepted values: 'server' (result is stored on a file on the server), 'client' (result is directly transferred to the client), or 'both'. Default is 'both'. Type = String

Input sequence. You need to supply either this parameter or the next one (tmp_infile). Type = Stringsequence

tmp_infile Name of the file with input sequence on the server. You need to supply either this parameter or the previous one (sequence).

Type = String

3 of 7 9/13/07 4:04 PM format Input sequence format. Supported: IG (Intelligenetics), WC (wconsensus), raw, fasta. Default is

Type = String

Number of substitutions allowed. Type = Integer

Pattern to match. pattern (required) Type = String

Pattern identifier. id

Origin for the calculation of positions (0 for end of sequence). origin

No overlapping of oligos allowed if value = 1. Disable the detection of overlapping matches for noov

self-overlapping patterns (ex TATATA, GATAGA).

Type = Integer

Oligonucleotide occurrences found on both stands are summed (2) or not (1). Default is 2. str

Type = Integer

Sort oligomers according to overrepresentation if value = 1.

Type = Integer

th Threshold on match count.

Type = Integer

Output Parameters

Location of the result file on the server. This can be used as input for a further request. Type = String

command The stand alone command executed on the server. Type = String

The results client Type = String

Method gene_info

Description Get information about genes

Return type. Accepted values: 'server' (result is stored on a file on the server), 'client' (result is directly transferred to the client), or 'both'. Default is 'both'. Type = String

Organism. Words need to be underscore separated (example: Escherichia coli K12). organism

Type = String (required)

List of gene(s) for which you want info on or list of keywords to search for (can be regular query

expressions).
Type = String (required)

Full match only (no substring matching) if value = 1. Type = Integer full

Do not print the query at the begining of each line if value = 1. Type = Integer $\,$ noquery

Match query against the description, too, not just against gene ID and name if value = 1.

Feature type (CDS, mRNA, tRNA, rRNA, scRNA). Type = String feattype

Output Parameters

Location of the result file on the server. This can be used as input for a further request.

command The stand alone command executed on the server.

Type = String

client The results

Type = String

Method supported_organisms

Description Get a list of supported organisms.

Input Parameters Parameters

output Return type. Accepted values: 'server' (result is stored on a file on the server), 'client' (result is directly transferred to the client), or 'both'. Default is 'both'.

Type = String

 $Output\ format\ (supported:\ html_list,html_table,array,text,keys,names,sizes,full,tree,html_tree).$ Type=String

Root taxon.

Type = String

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Location of the result file on the server. This can be used as input for a further request. server

Type = String

command The stand alone command executed on the server.

The results.
Type = String client

Method convert_seq

Description Converts a sequence between two formats (e.g. fasta -> raw).

Parameters Input Parameters

> Return type. Accepted values: 'server' (result is stored on a file on the server), 'client' (result is directly transferred to the client), or 'both'. Default is 'both'. output

Type = String

Sequence to convert. Type = String sequence

Name of the file with input sequence on the server. You need to supply either this parameter or the tmp_infile

previous one (sequence). Type = String

from Format of input sequence (embl, fasta, filelist, gcg, genbank, ig, maf, multi, ncbi, raw, tab, wc,

(required) wconsensus). Type = String

Format of output sequence (fasta, filelist, ft, ig, multi, raw, tab, wc, wconsensus),

(required) Type = String

Output Parameters

Location of the result file on the server. This can be used as input for a further request. Type = ${\bf String}$

The stand alone command executed on the server.

Type = String

The results.

Type = String

Method compare_classes

Compare two class files(the query file and the reference file). Each class of the query file is compared to each class of the reference Description

file.The number of common elements is reperted, as well as the probability to observe at least this number of common elements by chance alone.

Parameters Input Parameters Return type. Accepted values: 'server' (result is stored on a file on the server), 'client' (result is directly transferred to the client), or 'both'. Default is 'both'. output

Type = String

ref classes A tab-delimited text file containing the description of reference classes

A tab-delimited text file containing the description of query classes Type = String query_classes

return_fields List of fields to return. Supported field: dotprod,entropy,freq,jac_sim,members,occ,proba,rank Type = Integer

Specify a column of the input file containing a score associated to each member Type = Integer score_column

This file will be used as both reference and query. This is equivalent to -q input_file -r input_file. Type = String input classes

upper_threshold field

 $\label{eq:control_equation} Supported fields: E(QR), E_val, F(!Q!R), F(Q!R), F(Q), F(QR), F(R!Q), F(R), H(Q), H(Q), H(R), H(R), H(R), I(Q,R), I(C,P(QR),P(QR),P(R|Q),P_val,Q,QR,QvR,R,U(QR),U(R|Q), H(Q,R), dotprod,jac_sim,rank, sig Type = String$

upper threshold value for a supported field Type = Float upper_threshold_value

Supported fields: same fields as upper_threshold_field. Type = String lower_threshold_field

lower threshold value for a given fields. Type = Float lower_threshold_value

sort on the basis of the specified key.

Type = String

distinct Prevent to compare each class with itself (when the reference and query files contain the same classes)

triangle (only valid if query file and reference file are the same) Do not perform the reciprocal comparisons.

Type = Integer

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Return a pairwise matrix, where each row corresponds to a reference class, each column to a query class, and each cell contains a comparison between the two classes. The next argument indicates which statistics has to be return in the matrix (default=sig)Supported: $E(QR)_E = val,F(Q)R,F(Q)R,F(Q)R,F(R)Q,F(R)Q,H(Q)R,H(R),H(R)Q,H(R),H(R)Q,R),IC,P(QR),P(Q)R,P(R)Q,P_{Q}R,Q_{Q}R,Q_{Q}R,R_{Q}R,U(Q)R_{Q}R,U(R)Q,H(Q,R),dotprod,jac_sim,rank,sig$ matrix

Type = String

Output Parameters

Location of the result file on the server. This can be used as input for a further request. Type = String

The stand alone command executed on the server.

Type = String

client

Type = String

Method matrix scan

 $Scan \ sequences \ with \ one \ or \ several \ position-specific \ scoring \ matrices \ (PSSM) \ to \ identify \ instances \ of \ the \ corresponding \ motifs (putative \ sites). \ This \ program \ supports \ a \ variety \ of \ background \ models \ (Bernoulli,$ Description

Markov chains of any order).

Input Parameters Parameters

Return type. Accepted values: 'server' (result is stored on a file on the server), 'client' (result is directly transferred to the client), or 'both'. Default is 'both'. Type = String output

sequence file Sequence file - all the formats supported in RSAT can be used as input (default:

Type = String

matrix file The matrix file is specified with the option "matrix_format" (see below) Default

format: tab . Type = String

Supported field: tab, cb, consensus, gibbs, meme, assembly. Type = String matrix format

matrix_list

Indicate a file containing a list of matrices to be used for scanning the region. This facilitates the scanning of a sequence with a library of matrices (e.g. all the matrices from RegulonDB or TRANSFAC) Format: the matrix list file is a text file. The first word of each row is suppose to indicate a file name. Any further information on the

same row is ignored. Type = String

Only scan with the top # matrices per matrix file. This option is valid for some file formats containing multiple matrices where top matrices are generally more top_matrices

informative. Type = Integer

background Background model file is a tab-delimited file containing the specification of

oligonucleootide frequencies.

Calculate background model from the input sequence set. This option requires to specify the order of the background model with the option markov . Type = Integer $\,$ background_input

background window

Size of the sliding window for the background model calculation. This option requires to specify the order of the background model with the option markov (suitable for short order model only markov 0 or 1)

Type = Integer

Order of the markov chain for the background model. This option is incompatible markov

with the option background Type = Integer

background_pseudo

Pseudo frequency for the background models. Value must be a real between 0 and 1. If this option is not specified, the pseudo-frequency value depends on the background calculation. For -bginput and -window, the pseudo frequency is automatically calculated with the length (L) of the sequence following this formula : square-root of L divided by L+squareroot of L. For -bgfile, default value is 0.01. If the training sequence length (L) is known, the value can be set by -bg_pseudo option to square-root of L divided by L+squareroot of L. Type = Float

List of fields to return. Supported fields: sites, rank, limits, normw, bg_model, matrix, freq_matrix, weight_matrix, distrib . return_fields

Type = String

upper threshold field

Type = String

upper_threshold_value

Type = Float

lower threshold field

Type = String

lower_threshold_value

Type = Float

both_strand

Type = Integer

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single strand

Type = Integer

Output Parameters

Location of the result file on the server. This can be used as input for a further request.

The stand alone command executed on the server. command

Type = String

client The results. Type = String

Method random_seq

Description Generates random sequences.

Input Parameters Parameters

Return type. Accepted values: 'server' (result is stored on a file on the server), 'client' (result is directly transferred to the client), or 'both'. Default is 'both'.

Type = String output

Length of sequence to generate. sequence length

repetition Number of sequences to generate.

Type = Integer

Format of sequence(s) to generate. Type = String format

A newline character will be inserted in the sequence every # bases, where # is the number provided. Default is 70. A value of 0 will prevent newline insertion. Type = Integer line_width

Type of sequence(s) to generate (protein \mid DNA \mid other). Type = String type

Seed for the random generator. seed

Type = Integer

alphabet alphabet must be followed by residue frequencies expressed precisely this way: -a a:t # c:g

Type = String

expfreq

Expected frequencies of oligomers in sequence(s) to generate. Indicate the file that contains expected oligomer frequencies. When this option is used, the sequences are generated according to a Markov chain.

Type = String

Name of the file with expected frequencies on the server. tmp_expfreq_file

Type = String

bg_model Background model. Automatically load a pre-calibrated exected frequency file from the

RSAT genome distribution. When this option is used, the options organism and oligo_length are also required, to indicate the organism and the oligonucleotide length, olgo_length are also required, to indicate the organism and the oligonucleotide length, respectively. This option is incompatible with the option expfreq. Type of sequences used as background model for estimating expected oligonucleotide frequencies (supported models): equi (equiprobable residue frequencies [default]), upstream (all upstream sequences, allowing overlap with upstream ORFs. Requires to specify a model organism), upstream-noorf (all upstream sequences, preventing overlap with upstream ORFs. Requires to specify a model organism), and intergenic (intergenic frequencies. Whole set of intergenic regions including upstream and downstream sequences. Pocuring to specify a model organism. regions, including upstream and downstream sequences. Requires to specify a model

organism). Type = String

organism Name of the organism when using a background model.

Type = String

Length of oligomer when using a background model. oligo length

Type = Integer

length file Allows to generate random sequences with the same lengths as a set of reference sequences. The length file contains two columns: sequence ID (ignored) and sequence $\,$ length file

length. Type = String

Output Parameters

Location of the result file on the server. This can be used as input for a further request. Type = ${\sf String}$ server

The stand alone command executed on the server. command

Type = String

client The results.

Type = String

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