

---

# TREMSUCS

*Release 1.0*

Gabor Balogh

Jun 20, 2024



**CONTENTS:**

<b>1</b>	<b>Example report:</b>	<b>3</b>
<b>2</b>	<b>Installing from github.com:</b>	<b>5</b>
<b>3</b>	<b>Help Page of the pipeline:</b>	<b>7</b>
3.1	TREMSUCS . . . . .	7
<b>4</b>	<b>Short tutorial:</b>	<b>9</b>
4.1	Usage of the interactive mode: . . . . .	9
	<b>Index</b>	<b>13</b>



“TREMSUCS” a tool to choose, harvest and analyse expression and methylation data of the TCGA-projects for revealing Biomarkers which indicate treatment success.



## **EXAMPLE REPORT:**

An example report can be downloaded [here](#). Be aware that this report has a size of about 300 MB.





## INSTALLING FROM GITHUB.COM:

```
$ git clone https://github.com/dendemayer/TREMSUCS-TCGA.git  
$ cd TREMSUCS-TCGA  
$ pip install .
```

To start the analysis with help of the interactive mode, call the pipeline without any argument:

```
$ TREMSUCS
```

Calling the help or the manual page:

```
$ TREMSUCS --help  
$ man TREMSUCS
```



## HELP PAGE OF THE PIPELINE:

### 3.1 TREMSUCS

“TREMSUCS” a tool to choose, harvest and analyse expression and methylation data of the TCGA-projects for revealing Biomarkers which indicate therapy specific treatment success predictions.

Calling the pipeline without any argument starts the interactive mode to help setting all needed parameters for the analysis.

TREMSUCS [OPTIONS]

#### Options

**-o, --out\_path** <out\_path>  
path to save the result files

**Default**  
'/home/dende/TREMSUCS'

**-p, --project** <project>  
TCGA project(s) to be applied. Any TCGA project can be chosen, like: -p TCGA-CESC -p TCGA-HNSC ...

**-d, --drugs** <drugs>  
drug(s), like: -d drug1 -d drug2 or drugcombination(s), like: -d drug1,drug2

**-c, --cores** <cores>  
number of cores provided to snakemake

**Default**  
1

**-C, --cutoff** <cutoff>  
Cut-off parameter

**Default**  
0

**-t, --threshold** <threshold>  
threshold parameter

**Default**  
0

**-e, --execute <execute>**

choose which pipeline shall be executed

**Default**

'DESeq2', 'metilene'

**-N, --dryrun**

snakemake dryrun

**Default**

False

**-D, --download**

if set, just download raw and meta data for given projects and analysis types, revise them, link them, but do not run any analysis

**Default**

False

**-u, --unlock**

in case the analysis crashes, snakemake locks the output directory, run with -u to unlock, then repeat the analysis

**Default**

False

**-v, --version**

printing out version information: Version 1.0

## SHORT TUTORIAL:

### 4.1 Usage of the interactive mode:

The following example composition of projects, drugs and parameters creates the configuration given in the example report [here](#).

The same configuration can be applied by issuing the following command (the number of cores hereby can be adjusted and would also give the same results):

```
$ TREMSUCS -p TCGA-CESC -p TCGA-HNSC -p TCGA-LUSC -d cisplatin -d carboplatin,paclitaxel
↪ ↵
-d carboplatin -o /scr/TREMSUCS_out -c 40 -t 5 -t 10 -t 20 -C 5 -C 8
```

Calling the pipeline without any argument starts the interactive mode:

```
$ TREMSUCS

OUTPUT_PATH:          /homes/biertruck/gabor/TREMSUCS
SCRIPT_PATH:          /homes/biertruck/gabor/phd/test_git_doc/TREMSUCS/src/shared/
↪ modules
PIPELINES executed:   ['DESeq2', 'metilene']

which projects do you want to include in your analysis:

0:    TCGA-CESC          Cervical Squamous Cell Carcinoma and Endocervical
↪ Adenocarcinoma
1:    TCGA-HNSC          Head and Neck Squamous Cell Carcinoma
2:    TCGA-LUSC          Lung Squamous Cell Carcinoma
3:    TCGA-ESCA          Esophageal Carcinoma
4:    TCGA-BRCA          Breast Invasive Carcinoma
5:    TCGA-GBM           Glioblastoma Multiforme
6:    TCGA-OV            Ovarian Serous Cystadenocarcinoma
7:    TCGA-LUAD          Lung Adenocarcinoma
8:    TCGA-UCEC          Uterine Corpus Endometrial Carinoma
9:    TCGA-KIRC          kidney renal clear cell carcinoma
10:   TCGA-LGG           brain lower grade glioma
11:   TCGA-THCA          thyroid carcinoma
12:   TCGA-PRAD          prostate adenocarcinoma
13:   TCGA-SKCM          skin cutaneous melanoma
14:   TCGA-COAD          colon adenocarcinoma
15:   TCGA-STAD          stomach adenocarcinoma
```

(continues on next page)

(continued from previous page)

```

16:      TCGA-BLCA      bladder urothelial carcinoma
17:      TCGA-LIHC      liver hepatocellular carcinoma
18:      TCGA-KIRP      kidney renal papillary cell carcinoma
19:      TCGA-SARC      sarcoma
20:      TCGA-PAAD      pancreatic adenocarcinoma
21:      TCGA-PCPG      pheochromocytoma and paraganglioma
22:      TCGA-READ      rectum adenocarcinoma
23:      TCGA-TGCT      testicular germcelltumors
24:      TCGA-THYM      thymoma
25:      TCGA-KICH      kidney chromophobe
26:      TCGA-ACC      adrenochordical carcinoma
27:      TCGA-MESO      mesothelioma
28:      TCGA-UVM      uveal melanoma
29:      TCGA-DLBC      lymphoid neoplasm diffuse large b-cell lymphoma
30:      TCGA-UCS      uterine carcinoma
31:      TCGA-CHOL      cholangiocarcinoma
enter your choices one by one, when you are done, simply press "Enter":

```

As suggested, you can now, one by one include the projects you are interested in. A default OUTPUT\_PATH is also already given together with the default analysis types “DESeq” and “metilene”. Those defaults can also be adjusted in next steps with help of the interactive mode.

To recreate the example set, the first three projects have to be selected, afterwards the following prompt is given:

```

you choose:
PROJECTS:      ['TCGA-CESC', 'TCGA-HNSC', 'TCGA-LUSC']

which therapy approach do you want to include in your analysis:

0: cisplatin                TCGA-CESC: 103 TCGA-HNSC: 64 TCGA-LUSC: 1
1: carboplatin,paclitaxel   TCGA-CESC: 5 TCGA-HNSC: 26 TCGA-LUSC: 14
2: 5-fluorouracil,cisplatin TCGA-CESC: 5 TCGA-HNSC: 2 TCGA-LUSC: 0
3: carboplatin              TCGA-CESC: 3 TCGA-HNSC: 6 TCGA-LUSC: 3
4: carboplatin,cisplatin,paclitaxel TCGA-CESC: 3 TCGA-HNSC: 0 TCGA-LUSC: 1
5: cisplatin,gemcitabine    TCGA-CESC: 3 TCGA-HNSC: 0 TCGA-LUSC: 9
6: paclitaxel               TCGA-CESC: 2 TCGA-HNSC: 1 TCGA-LUSC: 0
7: erbitux                  TCGA-CESC: 1 TCGA-HNSC: 9 TCGA-LUSC: 0
8: cisplatin,vectibix       TCGA-CESC: 0 TCGA-HNSC: 5 TCGA-LUSC: 0
9: carboplatin,erbitux,paclitaxel TCGA-CESC: 0 TCGA-HNSC: 4 TCGA-LUSC: 0
10: cisplatin,erbitux       TCGA-CESC: 0 TCGA-HNSC: 3 TCGA-LUSC: 0
11: carboplatin,cisplatin,erbitux,paclitaxel TCGA-CESC: 0 TCGA-HNSC: 3 TCGA-LUSC: 0
12: carboplatin,cisplatin   TCGA-CESC: 0 TCGA-HNSC: 2 TCGA-LUSC: 0
13: docetaxel,erbitux       TCGA-CESC: 0 TCGA-HNSC: 2 TCGA-LUSC: 0
14: cisplatin,docetaxel     TCGA-CESC: 0 TCGA-HNSC: 1 TCGA-LUSC: 10
15: carboplatin,docetaxel   TCGA-CESC: 0 TCGA-HNSC: 1 TCGA-LUSC: 3
16: cisplatin,vinorelbine    TCGA-CESC: 0 TCGA-HNSC: 0 TCGA-LUSC: 21
17: carboplatin,vinorelbine TCGA-CESC: 0 TCGA-HNSC: 0 TCGA-LUSC: 8
18: cisplatin,etoposide     TCGA-CESC: 0 TCGA-HNSC: 0 TCGA-LUSC: 7
19: carboplatin,gemcitabine TCGA-CESC: 0 TCGA-HNSC: 0 TCGA-LUSC: 5
20: cisplatin,pemetrexed    TCGA-CESC: 0 TCGA-HNSC: 0 TCGA-LUSC: 3
21: cisplatin,docetaxel,gemcitabine TCGA-CESC: 0 TCGA-HNSC: 0 TCGA-LUSC: 2
22: carboplatin,gemcitabine,paclitaxel TCGA-CESC: 0 TCGA-HNSC: 0 TCGA-LUSC: 2
23: carboplatin,cisplatin,vinorelbine TCGA-CESC: 0 TCGA-HNSC: 0 TCGA-LUSC: 2

```

(continues on next page)

(continued from previous page)

```

24: carboplatin,docetaxel,gemcitabine      TCGA-CESC: 0 TCGA-HNSC: 0 TCGA-LUSC: 2
25: carboplatin,docetaxel,paclitaxel      TCGA-CESC: 0 TCGA-HNSC: 0 TCGA-LUSC: 2
26: gemcitabine                          TCGA-CESC: 0 TCGA-HNSC: 0 TCGA-LUSC: 2

```

enter your choices one by one, when you are done, simply press "Enter":

Here are therapies listed where the maximum of a row is greater than 1. We apply row 0, 1 and 3 to include cisplatin, the combination of carboplatin and paclitaxel and cases which got solely treated with carboplatin. In the following, every other parameter is requested. With the next prompt, the default OUTPUT\_PATH can be confirmed or replaced:

```

do you want to keep the default OUTPUT_PATH of:
/homes/biertruck/gabor/TREMSUCS
if so, press ENTER, if not, enter your custom output path:

```

**In this example, we confirm the suggested OUTPUT\_PATH and are asked to confirm**  
or set the number of cores which shall be invoked into the analyses:

```

do you want to keep the default number of cores invoked of 1?
if so, press ENTER, if not, enter the number of cores:
40

```

We set the cores to 40 and then can decide which analysis approaches shall be triggered, per default, DESeq2 and metilene based biomarker predictions are produced:

```

which pipeline do you want to include into your analysis
press ENTER if DESeq2 and metilene (default) or
1 for DESeq2 or
2 for metilene

```

We confirm the default of those two analyses and can set the cutoff values, if we want to add those at all:

```

do you want to add one or multiple cutoffs?
it is recommend to choose cutoff values between 5 and 10 years
if not, just press ENTER, if so enter the coutoffs one by one:
5
8

```

Like the example set, we add here a cutoff of 5 and 8. Then the thresholds are requested:

```

do you want to add one or multiple thresholds?
it is recommend to choose threshold values which do not exceed a value of 50
if not, just press ENTER, if so enter the thresholds one by one:
5
10
20

```

We apply thresholds of 5, 10 and 20. All mandatory and optional parameters are set with that and are finally listed before the whole approach is started:

```

OUTPUT_PATH:      /homes/biertruck/gabor/TREMSUCS
PROJECT:          ['TCGA-CESC', 'TCGA-HNSC', 'TCGA-LUSC']
DRUGS:            ['carboplatin', 'carboplatin,paclitaxel', 'cisplatin']
pipelines executed: ['DESeq2', 'metilene']

```

(continues on next page)

(continued from previous page)

```
cores:          40
cutoff:         [0, 5, 8]
threshold:      [0, 5, 10, 20]
press ENTER to start or q to quit:
```

If something went wrong, you can quit now and start over, or of course start the analysis.



## Symbols

-C  
TREMSUCS command line option, 7

-D  
TREMSUCS command line option, 8

-N  
TREMSUCS command line option, 8

--cores  
TREMSUCS command line option, 7

--cutoff  
TREMSUCS command line option, 7

--download  
TREMSUCS command line option, 8

--drugs  
TREMSUCS command line option, 7

--dryrun  
TREMSUCS command line option, 8

--execute  
TREMSUCS command line option, 7

--out\_path  
TREMSUCS command line option, 7

--project  
TREMSUCS command line option, 7

--threshold  
TREMSUCS command line option, 7

--unlock  
TREMSUCS command line option, 8

--version  
TREMSUCS command line option, 8

-c  
TREMSUCS command line option, 7

-d  
TREMSUCS command line option, 7

-e  
TREMSUCS command line option, 7

-o  
TREMSUCS command line option, 7

-p  
TREMSUCS command line option, 7

-t  
TREMSUCS command line option, 7

-u

TREMSUCS command line option, 8

-v

TREMSUCS command line option, 8

## T

TREMSUCS command line option

-C, 7

-D, 8

-N, 8

--cores, 7

--cutoff, 7

--download, 8

--drugs, 7

--dryrun, 8

--execute, 7

--out\_path, 7

--project, 7

--threshold, 7

--unlock, 8

--version, 8

-c, 7

-d, 7

-e, 7

-o, 7

-p, 7

-t, 7

-u, 8

-v, 8