

***NEAT***

## **Adding modules**

Patrick Schorderet

[Patrick.schorderet@molbio.mgh.harvard.edu](mailto:Patrick.schorderet@molbio.mgh.harvard.edu)

Jan 2015

<b>1</b>	<b>INTRODUCTION.....</b>	<b>4</b>
<b>2</b>	<b>ARCHITECTURE.....</b>	<b>5</b>
2.1	GENERAL ARCHITECTURE.....	5
2.2	CODE ARCHITECTURE.....	6

# 1 Introduction

---

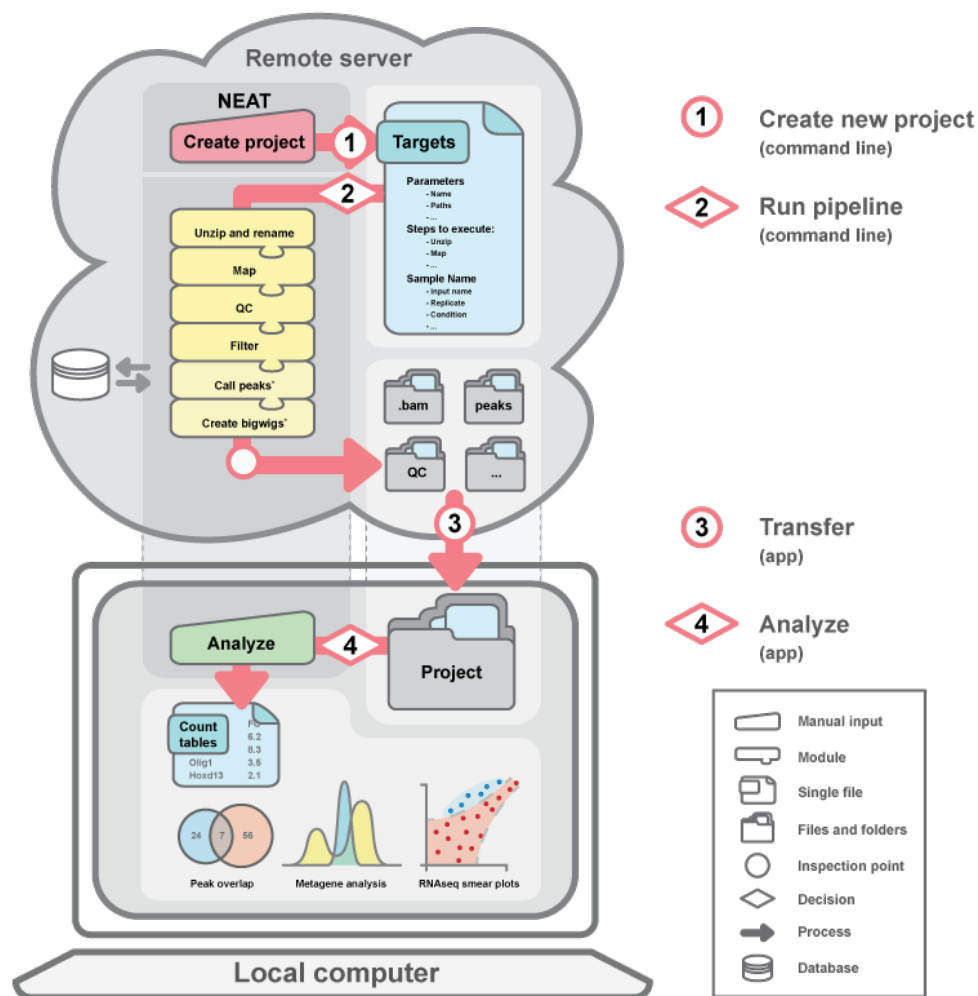
***N***ext generation ***A***nalysis ***T***oolbox (NEAT) is a perl/R package that supports users during the analysis of next generation sequencing (NGS)

NEAT is versatile and easy to modify. In this tutorial, we will show how to add a custom module to NEAT. Adding a new module has been made very easy by automating all the repetitive tasks such as job creation, batch submission and queuing. Adding a new module usually falls down to a single line of code.

## 2 Architecture

### 2.1 General architecture

NEAT contains different modules (yellow boxes) than can be modified / added.



**Fig.1 NEAT architecture.** NGS data can be analyzed using NEAT in less than a day. Users follow a logical 4-step process, including the creation of a new project, running the pipeline on a remote server or in the cloud, transferring the data to a local computer and proceeding to the analysis.

## 2.2 Code architecture

The main code of NEAT (in the example of ChIPseq projects) is found in *./NEAT/ChIPpip/scripts/ChIPpip.pl*. The code is well annotated, highly redundant and should be self-explanatory to advanced users. Main modules are easily identifiable and customizable. A brief summary of how each module is built is depicted below.



Figure 1. Schematic representation of the experimental design. The subjects were divided into two groups: the control group (n = 10) and the experimental group (n = 10). The control group received a placebo (P) and the experimental group received a 10% solution of the active ingredient (A). The subjects were divided into two groups: the control group (n = 10) and the experimental group (n = 10). The control group received a placebo (P) and the experimental group received a 10% solution of the active ingredient (A). The subjects were divided into two groups: the control group (n = 10) and the experimental group (n = 10). The control group received a placebo (P) and the experimental group received a 10% solution of the active ingredient (A).

\* Mapping requests with base

```
#[ Smap == "TRUE" ](
```

```
print "\n*****\n";
print "\n Mapping files\n";
```

```
my $iterJobName = "Iterate.<YOUR_JOB_NAME_HERE>";
my $seqJobName  = "<YOUR_JOB_NAME_HERE>";
my $splitJobName = "Inspect/SeqJobName/split";
```

```

n = Create file to store jobs in
n = f"Storage/{StorageName}"      (mkdir Storage/{StorageName} ;
mkdir Spd2Spd2)
n["-d"] = "Storage/{StorageName}/{StorageName}.n"
open $GSLIB ">" ; $GSLIB or die "Can't open '$GSLIB'";
print $GSLIB "n {n}/bin/{n}"
close $GSLIB
chmod 777 $GSLIB ;
print "I've Store all of the following '{StorageName}' jobs in $GSLIB '{n}'"
my $n { $n }

```

```
foreach my $i (0..$#sampleInputs) {
```

```
# Prepare a personal quabscript
my $QSUDir = "Storage/StorageName/StorageInputs/$I", $StorageName', sh')
'qp SecHead $QSUDir';
```

```
# Create a directory
my $path2currentSampleDir = "$path2aligned/$samplesInputs[$i]"
unless (-d "$path2currentSampleDir") { mkdir $path2currentSampleDir; }
```

```

#-----+-----+-----+-----+-----+-----+-----+-----+-----+-----+
#-----+-----+-----+-----+-----+-----+-----+-----+-----+-----+
my $cmd = "<YOUR_CUSTOM_COMMAND_HERE>"
"echo \"$cmd\" >> $SQL_Bin";
#-----+-----+-----+-----+-----+-----+-----+-----+-----+-----+

```

```
A Keep track of the jobs in @myjobs
my $jobName = "SentryJobName";
push(@myjobs, $jobName);
$cmd = "$jobName.cmd" >> $path2cmd >> $path2cmd && $Q$B$E";
open $Q$B$E ">>", "$Q$B$E" or die "Can't open '$Q$B$E'";
print $Q$B$E "$cmd\n";
close $Q$B$E;
```

```

# Change Targets list file for next iteration
print "\n"
print "\n Changing '$objName' variable to FALSE and proceed"
"/usr/bin/perl -p -i -e 's/$objName/$objName/,DONE/g' $Targets";

```

0 Prepare file containing the jobs to run

[illegible]

```
@ Submit jobs to run
```

```
print "\n")
print "\n Submitting job to cluster: 't' with $QSUB" "\n"
job $QSUB
```

@ Edit script.

```
print "\n";
print "\n Exiting fnyjshName section with no known error 'a'";
print "\n";
```

edit:

1

If job is TRUE, enter module  
(modify the 'Steps\_to\_execute' variables at the  
top of the code)

**Name your job** (flex map)

A file is created with YOUR\_JOB\_NAME. This file will contain the flow of each individual-job related to this job.

For loop that iterates through all samples and inputs. To iterate through samples only, use the @samples table instead of the @samplesInputs table.

For each individual job, create a qsub-like file (in which the job code will be) by copying the qsub header (\$scrhead) into the intermediary \$QSUBint file

**YOUR CUSTOMIZE CODE GOES HERE**

**Keep track of job names to ensure job dependencies (see below)**

**Modify the Targets.txt file to mark the job as '\_DONE'**

**Prepare code for the next iteration (\$IterateSH).** This iteration depends on the completion of all the current jobs (\$myJobsVec)

**Submit job to cluster and exit**