# Building IaaS infrastructures on the AWS Cloud

Saul Pierotti

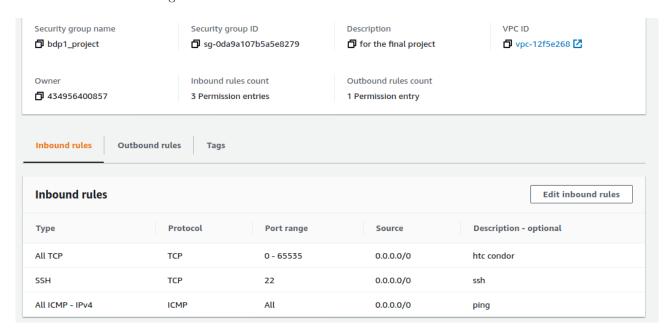
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### 1 Creation of an HTCondor Cluster for the Alignment of NGS Reads

The demonstrative infrastructure described in this project consists of an HTCondor cluster of three nodes. One node is configured as Master Node, while two nodes are configured as Worker Nodes. The infrastructure can be easily expanded by replicating the Worker Node instances. The Master Node was not used also as a Worker Node since the performance benefits and cost savings would be marginal. This was done also to avoid overloading the Master Node, on which the entire cluster depends. A shared storage space directly attached to the Master Node but available to all the Worker Nodes was also implemented using the distributed file system NFS.

### 1.1 Initialization of the Instances on the AWS Cloud

Worker Nodes and the Master Node were both built on similar machines. For the Master Node, the t2.medium instance type was used with a 50 Gb SSD as root storage. For the Worker Nodes, the t2.large instance type was used with a 50 Gb SSD as root storage. The operating system chosen for both machine types is Ubuntu Server 18.04.4 LTS. The Master Node and the Worker Nodes were all instantiated in the same availability zone (us-east-la) so that they would be able to communicate through private IPv4 addresses. The security group for the instances was configured as follows:



All the TCP ports were opened to the other members of the same security group since HTCondor daemons use a dynamically assigned port. The TCP ports were also needed for the setup of a shared NFS volume. ICMP ports were opened for accepting incoming ping requests for testing purposes. TCP port 22 was opened for allowing remote control of the machines via ssh.

#### 1.2 Configuration of the Master Node

The PS1 prompt of the Master Node was changed so to make the node easily identifiable from the command line. HTCondor was then installed with the following commands:

```
ubuntu@bdp1-master-node:~$ echo $PS1
\[\e]0;\u@\h: \w\a\]${debian_chroot:+($debian_chroot)}\[\033[01;32m\]\u@bdp1-master-node\[\033[00m\]:\[\033[01;34m\]\w\[\033[00m\]\$
```

```
sudo su
wget -q0 - https://research.cs.wisc.edu/htcondor/ubuntu/HTCondor-Release.gpg.key | apt-key add - #
    import the gpg key of HTCondor
echo "deb http://research.cs.wisc.edu/htcondor/ubuntu/8.8/bionic bionic contrib" >> /etc/apt/sources.
    list # add the repository
echo "deb-src http://research.cs.wisc.edu/htcondor/ubuntu/8.8/bionic bionic contrib" >> /etc/apt/
    sources.list
apt update
apt install htcondor
systemctl start condor # start and enable the condor service
systemctl enable condor
```

The correct proceeding of the installation and the start of the condor service where checked with the following commands:

```
condor.service - Condor Distributed High-Throughput-Computing
  Loaded: loaded (/lib/systemd/system/condor.service; enabled; vendor preset: enabled)
  Active: active (running) since Tue 2020-06-16 10:31:25 UTC; 1min 16s ago
Main PID: 15225 (condor_master)
  Status: "All daemons are responding"
  CGroup: /system.slice/condor.service
            -15266 condor_procd -A /var/run/condor/procd_pipe -L /var/log/condor/ProcLog -R 1000000 <u>-S 60 -C 111</u>
           lue_{15268} condor_shared_port -f
Jun 16 10:31:25 ip-172-31-8-109 systemd[1]: Started Condor Distributed High-Throughput-Computing.
ubuntu@bdp1-master-node:~$ ps ax | grep condor
                                           _master -f
15266 ?
                     0:00
                                                        dor/procd_pipe -L /var/log/condor/ProcLog -R 1000000 -S 60 -C 111
15268 ?
                     0:00
                                 _shared_port -f
                      0:00 grep --color=auto
15286 pts/0
```

The following lines where appended at the end of the main HTCondor configuration file, located at /etc/condor\_config:

```
# Master Node IP
CONDOR_HOST = <Master_Node_private_IP>
# Master Node config
DAEMON_LIST = COLLECTOR, MASTER, NEGOTIATOR, STARTD
```

Finally, the condor service was restarted with the following command:

```
sudo systemctl restart condor
```

The NFS server was then implemented in the Master Node. A new 100 Gb standard magnetic volume was created from the AWS interface and attached to the Master Node. Fom the server, a primary partiton was initialized on the volume using fdisk and an Ext4 file system was created onto it using mkfs.ext4. The file /etc/fstab of the Master Node was modified so that the machine would mount the volume automatically at boot under the newly created directory /data. The following line was appended to /etc/fstab:

The following commands were then issued, so to install the appropriate packages:

```
sudo apt install nfs-kernel-server
```

The following line was appended to the NFS configuration file /etc/exports:

```
/data 172.31.0.0/16(rw,sync,no_wdelay)
```

Finally, the owner and group of the shared folder were changed to nobody:nogroup and the permissions of the folder were edited so to grant unlimited access to it:

```
sudo chown nobody:nogroup /data
sudo chmod 777 /data
```

The /data folder was so made available to all the Worker Nodes on the address range 172.31.0.0/16. This configuration does not pose a security risk since the Master Node and Worker Nodes machines belong to a Virtual Private Cloud (VPC), and so only machines instantiated on it will be able to access the volume. Moreover, this configuration grants immediate access to the volume to newly instantiated Worker Nodes. A mock file was created on the /data folder so to be able to recognize it when mounted.

```
touch /data/this_is_a_shared_NFS_volume
```

### 1.3 Configuration of the Worker Nodes

A virtual machine identical to the one used for the Master Node was instantiated. The PS1 prompt of this first Worker Node was changed so to make the node easily identifiable from the command line.

```
ubuntu@bpd1-worker-node:~$ echo $PS1
\[\e]0;\u@\h: \w\a\]${debian_chroot:+($debian_chroot)}\[\033[01;32m\]\u@bpd1-worker-node\[\033[00m\]:\[\
033[01;34m\]\w\[\033[00m\]\$
```

HTCondor was installed in this system with the same procedure used for the Master Node. Only the /etc/condor\_config file was configured differently, by appending the following lines to it:

```
# Master Node IP
CONDOR_HOST = <Master_Node_private_IP>

# Worker Node config
DAEMON_LIST = MASTER, STARTD

HOSTALLOW_READ = *
HOSTALLOW_WRITE = *
HOSTALLOW_ADMINISTRATOR = *
```

On the Worker Node access to the shared NFS volume was then set up. The following command was issued to install the required packages and enable the respective services:

```
sudo apt install nfs-common
sudo systemctl start nfs-server
sudo systemctl enable nfs-server
```

A new directory was then created at /data using the mkdir command. The /etc/fstab file was edited by appending the following line, so that the shared volume would be automatically mounted at boot under the directory /data:

```
<Master_Node_private_IP>:/data /data nfs defaults 0 0
```

It was then verified that the shared volume was accessible from the Worker Node.

```
ubuntu@bpd1-worker-node:~$ ll /data
total 24
drwxrwxrwx 3 nobody nogroup 4096 Jun 17 07:16  /
drwxr-xr-x 24 root root 4096 Jun 17 06:46 ../
drwx----- 2 root root 16384 Jun 16 17:38 lost+found/
-rw-rw-r-- 1 ubuntu ubuntu 0 Jun 17 07:16 this_is_a_shared_NFS_volume
```

On the Worker Nodes the application BWA was installed with the command:

```
sudo apt install -y bwa
```

An image of the virtual machine was taken through the AWS interface. In this way, the Worker Nodes would be easily replicable when more computational power is needed by instantiating new virtual machines from the image, without the need for manual configuration.

#### 1.4 Submission of a Test Job to the HTCondor Cluster

For demonstrating the use of the newly created cluster, a new Worker Node was instantiated from the Worker Node image, so that the cluster would contain the Master Node and 2 Worker Nodes.

Subsequently, a new volume was created in the AWS interface from a snapshot containing test data used for the BDP1 course (snap-09ee52d8038fb8094, BDP1\_2020). This snapshot contains NGS reads from 3 different patients. Each patient has a folder with around 500 fasta files, with 1000 reads each. Part of this data will be used for testing the infrastructure. The new volume was mounted on the Master Node under /data, substituting the empty volume used before. The /etc/fstab file was update accordingly and the nfs-server service restarted. It was then tested that the new volume was accessible from the Worker Nodes.

```
ubuntu@bpd1-worker-node:~$ ll /data
total 28
drwxrwxrwx 4 root root 4096 Apr 26 13:38 //
drwxr-xr-x 24 root root 4096 Jun 17 06:46 ../
drwxr-xr-x 5 root root 4096 Apr 19 14:39 BDP1_2020/
drwx----- 2 root root 16384 Apr 26 09:37 lost+found/
```

The test job consisted in aligning ten fasta files (read\_1.fa to read\_10.fa from patient 1) to the human genome build hg19, also stored on the shared volume. The BWA alignment tool was used for the scope. This tool takes advantage of indexing the genome for speeding up the alignment of low-divergent reads. The index for the hg19 build was already present in the volume snapshot, so there was no need to create it from scratch. The test fasta files were copied from the shared volume to the home folder of the Master Node. The test job file alignment\_test.job was created, with the following content:

```
######## The program that will be executed ######
Executable = alignment_test.py
readnum = $(Process)+1
arguments = read_$INT(readnum).fa
= read_$INT(readnum).fa
transfer_input_files = read_$INT(readnum).fa
= read_$INT(readnum).log
Log
# will contain condor log
Output
     = read_$INT(readnum).out
# will contain the standard output
     = read_$INT(readnum).error
Error
# will contain the standard error
should_transfer_files = YES
```

The script alignment\_test.py called as executable in alignment\_test.job was the following:

```
#!/usr/bin/python
import sys, os
from timeit import default_timer as timer
start = timer()
dbpath = "/data/BDP1_2020/hg19/"
dbname = "hg19bwaidx"
queryname = sys.argv[1]
out_name = queryname[:-3]
fafile = queryname
samfile = out_name + ".sam"
gzipfile = out_name + ".sam.gz"
saifile = out_name + ".sai"
md5file = out_name + ".md5"
print "Input: ", queryname
command = "bwa aln -t 1 " + dbpath + dbname + " " + fafile + " > " + saifile
print "launching command: " , command
os.system(command)
\texttt{command = "bwa samse -n 10 " + dbpath + dbname + " " + saifile + " " + fafile + " > " + samfile}
print "launching command: " , command
os.system(command)
# Checksums
print "Creating md5sums"
os.system("md5sum " + samfile + " > " + md5file)
print "gzipping out text file"
command = "gzip " + samfile
print "launching command: " , command
os.system(command)
# Transfer files to shared volume and clean the Output Sandbox
print "Moving files and clearing the Output Sandbox"
os.system("mv "+ gzipfile + " /data/outputs/"+ gzipfile)
os.system("mv "+ md5file + " /data/outputs/"+ md5file)
os.system("rm "+ saifile)
execution_time = timer() - start
print "Total execution time: " + str(execution_time)
print "exiting"
exit(0)
```

Ten instances of this test job were run on the cluster. The following outputs of condor\_q and condor\_status were recorded after submission:

```
ubuntu@bdp1-master-node:~$ condor_q
-- Schedd: ip-172-31-8-109.ec2.internal : <172.31.8.109:9618?... @ 06/18/20 08:03:52
OWNER BATCH NAME
                   SUBMITTED
                               DONE
                                             IDLE TOTAL JOB_IDS
                                      RUN
                   6/18 08:03
ubuntu ID: 24
                                                     10 24.0-9
Total for query: 10 jobs; 0 completed, 0 removed, 10 idle, 0 running, 0 held, 0 suspended
Total for ubuntu: 10 jobs; 0 completed, 0 removed, 10 idle, 0 running, 0 held, 0 suspended
Total for all users: 10 jobs; 0 completed, 0 removed, 10 idle, 0 running, 0 held, 0 suspended
ubuntu@bdp1-master-node:~$ condor_status
Name
                                                 Arch
                                                        State
                                                                   Activity LoadAv Mem
                                                 X86_64 Claimed
slot1@ip-172-31-8-22.ec2.internal
                                     LINUX
                                                                   Busy
                                                                              0.000 3979
slot2@ip-172-31-8-22.ec2.internal LINUX
                                                 X86_64 Claimed
                                                                   Busy
                                                                              0.000 3979
slot1@ip-172-31-14-66.ec2.internal LINUX
                                                 X86_64 Claimed
                                                                   Busy
                                                                              0.000 3979
slot2@ip-172-31-14-66.ec2.internal LINUX
                                                 X86_64 Claimed
                                                                   Busy
                                                                              0.000 3979
                Total Owner Claimed Unclaimed Matched Preempting Backfill Drain
  X86_64/LINUX
                                                      0
                                                                           0
                                                                                   0
         Total
                                                                                   Θ
```

The cluster was able to successfully complete the task and the following output files were produced:

```
ubuntu@bdp1-master-node:/data/outputs$ ll
total 408
drwxrwxrwx 2 nobody nogroup
                             4096 Jun 18 09:12
drwxrwxrwx 5 root
                             4096 Jun 17 15:11
                    root
-rw-r--r-- 1 nobody nogroup
                               45 Jun 18 08:57 read_1.md5
-rw-r--r-- 1 nobody nogroup
                               31 Jun 18 08:56
-rw-r--r-- 1 nobody nogroup
                               46 Jun 18 09:07
                                                read_10.md5
rw-r--r-- 1 nobody nogroup 55074 Jun 18 09:07
rw-r--r-- 1 nobody nogroup
                               45 Jun 18 08:57
                                                read_2.md5
-rw-r--r-- 1 nobody nogroup 41129 Jun 18 08:57
rw-r--r-- 1 nobody nogroup
                               45 Jun 18 08:57
                                                read_3.md5
rw-r--r-- 1 nobody nogroup
                               31 Jun 18 08:56
rw-r--r-- 1 nobody nogroup
                                                read_4.md5
                               45 Jun 18 08:57
-rw-r--r-- 1 nobody nogroup 54914 Jun 18 08:57
rw-r--r-- 1 nobody nogroup
                               45 Jun 18 09:03
                                                read_5.md5
rw-r--r-- 1 nobody nogroup 77299 Jun 18 09:03
rw-r--r-- 1 nobody nogroup
                               45 Jun 18 09:03
                                                read_6.md5
rw-r--r-- 1 nobody nogroup 55883 Jun 18 09:03
rw-r--r-- 1 nobody nogroup
                               45 Jun 18 09:00 read_7.md5
rw-r--r-- 1 nobody nogroup
                               31 Jun 18 08:58
rw-r--r-- 1 nobody nogroup
                               45 Jun 18 09:00 read_8.md5
rw-r--r-- 1 nobody nogroup
                               31 Jun 18 08:58
rw-r--r-- 1 nobody nogroup
                               45 Jun 18 09:07 read_9.md5
 rw-r--r-- 1 nobody nogroup 53760 Jun 18 09:07
```

The time required to complete the task on a single fasta file ranged from 62.77 to 422.54 seconds, with an average required time of 227.7 seconds.

```
ubuntu@bdp1-master-node:~$ cat read_*.out |grep "Total execution ti
me:"
Total execution time: 62.765912056
Total execution time: 417.533436775
Total execution time: 116.913583994
Total execution time: 66.0304908752
Total execution time: 115.42266202
Total execution time: 383.357795
Total execution time: 377.316680908
Total execution time: 158.487968922
Total execution time: 161.080944061
Total execution time: 422.54254508
```

### 1.5 Data Management Model

The general data management model followed would be that of transferring the executables and the input fasta files using the HTCondor Input Sandbox. This is because those files are generally small and their transfer will not overload the Master Node. Moreover, input data are likely to be uploaded dynamically when new sequencing experiments are performed. This system will allow sending easily new inputs to the cluster. On the contrary, the large hg19 genome and index files will be made available to the Worker Nodes through the shared NFS volume. The Output Sandbox will contain the condor log, the standard output, and the standard error. These files are really small and need to be immediately available to the submitter for inspecting the proceeding of the job. The aligned reads will instead be compressed and put in the shared volume, to avoid moving large data on the Output Sandbox.

## 2 Creation of a Container Image for the Alignment of NGS reads

A containerized version of the BWA application was built using docker. A container image for the BWA application is already available on DockerHub (biocontainers/bwa), but for educational purposes, a new image was built from scratch. The Ubuntu docker image was used as a base. The local docker image bwa was built from the following Dockerfile:

```
FROM ubuntu
RUN apt update
RUN apt install -y bwa
RUN apt install -y python
COPY ./alignment_test_docker.py
```

The container was then run using a bind-mount for the /data volume and starting an interactive shell.

```
docker run -v /data:/data -ti bwa /bin/bash
```

The script alignment\_test.py was slightly modified so to omit the final data movement to the shared volume, and renamed alignment\_test\_docker.py. Inside the container, the following command was run so to align a sample fasta file which had been copied from the shared volume:

```
./alignment_test_docker.py read_1.fa
```

The application ran successfully, producing the following output in 213.33 seconds (the container was run on a Ubuntu Server 18.04.4 LTS t2.large AWS virtual machine):

# 3 Evaluation of the Resources Needed for Solving a Real Use-Case

A plausible non-trivial use-case that can be addressed using the HTCondor cluster developped in Section 1 is the allignment of a big number of fasta files to the human reference genome hg19. The scientific challenge proposed

```
oot@871e3f6d9c26:/# cp /data/BDP1_2020/hg19/Patients/patient1/read_1.fa read_1.fa
root@871e3f6d9c26:/# ./alignment_test_docker.py read_1.fa
Input: read_1.fa
launching command: bwa aln -t 1 /data/BDP1_2020/hg19/hg19bwaidx read_1.fa > read_1.sai
[bwa_aln] 17bp reads: max_diff = 2
[bwa_aln] 38bp reads: max_diff = 3
[bwa_aln] 64bp reads: max_diff = 4
[bwa_aln] 93bp reads: max_diff
[bwa_aln] 124bp reads: max_diff = 6
[bwa_aln] 157bp reads: max_diff = 7
[bwa_aln] 190bp reads: max_diff = 8
[bwa_aln] 225bp reads: max_diff = 9
[bwa_aln_core] calculate SA coordinate... 0.21 sec
[bwa_aln_core] write to the disk... 0.00 sec
[bwa_aln_core] 1000 sequences have been processed.
[main] Version: 0.7.17-r1188
[main] CMD: bwa aln -t 1 /data/BDP1_2020/hg19/hg19bwaidx read_1.fa
[main] Real time: 121.035 sec; CPU: 4.582 sec
launching command:  bwa samse -n 10 /data/BDP1_2020/hg19/hg19bwaidx read_1.sai read_1.fa > read_1.sam
[bwa_aln_core] convert to sequence coordinate... 4.28 sec
[bwa_aln_core] refine gapped alignments... 1.13 sec
[bwa_aln_core] print alignments... 0.00 sec
[bwa_aln_core] 1000 sequences have been processed.
[main] Version: 0.7.17-r1188
[main] CMD: bwa samse -n 10 /data/BDP1_2020/hg19/hg19bwaidx read_1.sai read_1.fa
[main] Real time: 92.240 sec; CPU: 5.416 sec
Creating md5sums
gzipping out text file
launching command: gzip read_1.sam
Total execution time: 213.335064888
exiting
```

during the BDP1 course is an example of a task that can be tackled with the aforementioned infrastructure. This task consists in aligning 554000 NGS reads from one individual against the hg19 genome. Given that a single NGS-produced fasta file typically contains 1000 reads, the task consists in aligning 554 fasta files. From the tests performed, I roughly determined that the time needed to align one fasta file is 227.3 seconds. Of course this can vary based on the length and composition of the reads themselves. Using this rough average, we obtain that aligning the 554 fasta files from one individual will require 125924.2 seconds, around 35 hours of CPU time on a single core like the ones provisioned with the AWS t2.large machine.

If I assume that I will use the cluster detailed in this project, with two Worker Nodes and one Master Node, I will have four cores at my disposal (two per Worker Node). This means that the time required on the cluster will be of around 9 hours.

I will now describe a breakdown of the expected costs for aligning the reads from a single patient using the AWS EC2 system with On-Demand pricing on the US East N. Virginia availability zone, on a Linux system. The t2.large machine costs \$ 0.0928 per hour, and the cluster will require two of them for the two Worker Nodes. The t2.medium machine costs \$ 0.0464 per hour, and the cluster will require one of them for the Master Node. With an expected time needed for the task of 9 hours, the costs for the computation for one patient will amount to \$ 2.088. Storage on general purpose SSD EBS devices costs \$ 0.1 per Gb per month. A 100 Gb disk was used for storing the hg19 genome and index, as well as the reads for the patients. Therefore, storage will have a cost of \$ 10 per month. Note that costs for data movements in and out of the AWS system have been omitted.