

Bismark on HPC cluster:

What is Bismark?

Bismark is a program to map bisulfite treated sequencing reads to a genome of interest and perform methylation calls in a single step. The output can be easily imported into a genome viewer, such as Seq Monk, and enables a researcher to analyze the methylation levels of their samples straight away.

Features:

- Bisulfite mapping and methylation calling in one single step
- Supports single-end and paired-end read alignments
- Supports ungapped, gapped or spliced alignments

The official documentation can be found on:

[GitHub- Bismark](#)

[Documentation](#)

Versions Available:

- Bismark = Version: v0.19.0

How to load a version of Bismark?

To load a version of Bismark on the HPC, use the following command:

```
module load bio/bismark
```

Verify by using this command:

```
module list
```

It should load bowtie as dependency.

Note: User are recommended to load Sam tools while working with bismark.

How to use Bismark on the cluster?

To see all the options available, use the following command available,

```
bismak --help
```

For this test run download the required file from following links:

```
# Download fastq file
wget
https://github.com/FelixKrueger/Bismark/raw/master/test_data.fastq

# Download fasta file
wget
https://ftp.ensembl.org/pub/release-
108/fasta/homo_sapiens/dna/Homo_sapiens.GRCh38.dna.chromosome.1.fa.gz

# Unzip the file
gunzip *.fa.gz
```

To run the analysis on this dataset, use the following commands

Gene preparation:

```
# Run bismark gene preparation
# $(pwd) is my current data folder
bismark_genome_preparation $(pwd)
```

Run Bismark:

```
# Run bismark on the test data parallely over 16 cores  
bismark --parallel 16 --genome $(pwd) test_data.fastq
```

Dedublication:

```
deduplicate_bismark --bam test_data_bismark_bt2.bam
```

Running Methyl Extractor

```
bismark_methylation_extractor --parallel 16 --gzip --bedGraph  
test_data_bismark_bt2.bam
```

To see the report, use the following command

```
bismark2report
```

These commands will process necessary data and write files in the current working directory. If user needs to run a batch job, user can wrap commands into a sbatch script and submit it to scheduler.

To learn more, please refer to the documentation,

[Documentation](#)

Where to find help?

If you are stuck on some part or need help at any point, please contact OIT at the following address.

<https://ua-app01.ua.edu/researchComputingPortal/public/oitHelp>

