## ▼ BPL\_TEST2\_Chemostat script with FMPy ver 0.3.15

The key library FMPy v0.3.15 is installed.

After the installation a small application BPL\_TEST2\_Chemostat is loaded and run. You can continue with this example if you like.

```
!lsb_release -a # Actual VM Ubuntu version used by Google
     No LSB modules are available.
     Distributor ID: Ubuntu
     Description: Ubuntu 20.04.5 LTS
                        20.04
     Release:
     Codename:
                        focal
%env PYTHONPATH=
     env: PYTHONPATH=
!wget https://repo.anaconda.com/miniconda/Miniconda3-py38_22.11.1-1-Linux-x86_64.sh
!chmod +x Miniconda3-py38_22.11.1-1-Linux-x86_64.sh
!bash ./Miniconda3-py38_22.11.1-1-Linux-x86_64.sh -b -f -p /usr/local
import sys
sys.path.append('/usr/local/lib/python3.8/site-packages/')
     --2023-03-23 14:05:36-- <a href="https://repo.anaconda.com/miniconda/Miniconda3-py38-22.11.1-1-Linux-x86-64.sh">https://repo.anaconda.com/miniconda/Miniconda3-py38-22.11.1-1-Linux-x86-64.sh</a> Resolving repo.anaconda.com (repo.anaconda.com)... 104.16.130.3, 104.16.131.3, 2606:4700::6810:8203, ...
     Connecting to repo.anaconda.com (repo.anaconda.com) | 104.16.130.3 | :443... connected.
     HTTP request sent, awaiting response... 200 OK
     Length: 64630241 (62M) [application/x-sh]
     Saving to: 'Miniconda3-py38_22.11.1-1-Linux-x86_64.sh'
     Miniconda3-py38_22. 100%[=========>] 61.64M 144MB/s
     2023-03-23 14:05:37 (144 MB/s) - 'Miniconda3-py38_22.11.1-1-Linux-x86_64.sh' saved [64630241/64630241]
     PREFIX=/usr/local
     Unpacking payload ...
     Installing base environment...
     Downloading and Extracting Packages
     Downloading and Extracting Packages
     Preparing transaction: done
     Executing transaction: done
     installation finished.
!conda update -n base -c defaults conda --yes
```

Preparing transaction: done Verifying transaction: done Executing transaction: done

!conda --version !python --version

> conda 23.1.0 Python 3.8.15

!conda install -c conda-forge fmpy --yes # Install the key package

```
Preparing transaction: done
Verifying transaction: done
Executing transaction: done
```

## Now specific installation run a simulation and notebook for that

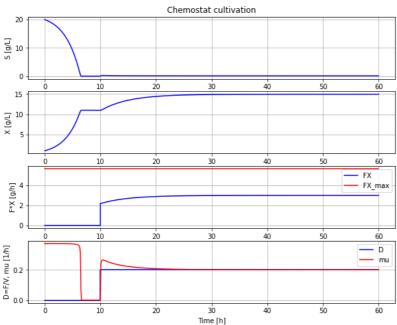
Start with connecting to Github. Then upload the two files:

- FMU BPL\_TEST2\_Chemostat\_linux\_om\_me.fmu
- Setup-file BPL\_TEST2\_Chemostat\_fmpy\_explore.py

```
git clone https://github.com/janpeter19/BPL_TEST2_Chemostat
     Cloning into 'BPL TEST2 Chemostat'...
%cd BPL TEST2 Chemostat
     /content/BPL_TEST2_Chemostat
```

## ▼ BPL\_TEST2\_Chemostat - demo

```
run -i BPL_TEST2_Chemostat_fmpy_explore.py
                             Linux - run FMU pre-compiled OpenModelica 1.21.0
                             Model for bioreactor has been setup. Key commands:
                                                                                                                  - change of parameters and initial values
                                    - par()
                                      - init()
                                                                                                                      - change initial values only
                                    - simu()
                                                                                                                     - simulate and plot
                                    - newplot()
                                                                                                                  - make a new plot
                                                                                                                      - show plot from previous simulation
                                    - show()
                                                                                                                     - display parameters and initial values from the last simulation % \left( 1\right) =\left( 1\right) \left( 1\right) \left(
                                    - disp()
                                    - describe() - describe culture, broth, parameters, variables with values/units
                             Note that both disp() and describe() takes values from the last simulation
                             Brief information about a command by help(), eg help(simu)
                             Key system information is listed with the command system_info()
%matplotlib inline
plt.rcParams['figure.figsize'] = [25/2.54, 20/2.54]
newplot()
par(Y=0.50, qSmax=0.75, Ks=0.1)
                                                                                                                                                                                                                                                 # Culture parameters
init(V_0=1.0, VX_0=1.0, VS_0=20)
                                                                                                                                                                                                                                                 # Bioreactor startup
par(S_in=30, t0=0, F0=0, t1=10, F1=0.2) \# Substrate feeding
```



```
describe('parts')
     ['bioreactor', 'bioreactor.culture', 'D', 'dosagescheme', 'feedtank', 'harvesttank']
describe('MSL')
     MSL: 3.2.3 - used components: RealInput, RealOutput, CombiTimeTable, Types
system_info()
     System information
      -OS: Linux
      -Python: 3.9.16
-Scipy: not installed in the notebook
      -FMPy: 0.3.15
      -FMU by: OpenModelica Compiler OpenModelica 1.21.0~dev-185-g9d983b8
      -FMI: 2.0
      -Type: ME
      -Name: BPL_TEST2.Chemostat
      -Generated: 2023-02-13T10:27:14Z
      -MSL: 3.2.3
      -Description: Bioprocess Library version 2.1.1-beta
      -Interaction: FMU-explore for FMPy version 0.9.7c
```

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