

## ✓ BPL\_TEST2\_Chemostat script with FMPy

The key library FMPy 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 22.04.4 LTS
Release:      22.04
Codename:     jammy

```

```
!python --version
```

```
Python 3.11.11
```

```
!pip install fmpy
```

```

Collecting fmpy
  Downloading FMPy-0.3.22-py3-none-any.whl.metadata (1.9 kB)
Requirement already satisfied: attrs in /usr/local/lib/python3.11/dist-packages (from fmpy) (25.1.0)
Requirement already satisfied: Jinja2 in /usr/local/lib/python3.11/dist-packages (from fmpy) (3.1.5)
Collecting lark (from fmpy)
  Downloading lark-1.2.2-py3-none-any.whl.metadata (1.8 kB)
Requirement already satisfied: lxml in /usr/local/lib/python3.11/dist-packages (from fmpy) (5.3.1)
Requirement already satisfied: msgpack in /usr/local/lib/python3.11/dist-packages (from fmpy) (1.1.0)
Requirement already satisfied: numpy in /usr/local/lib/python3.11/dist-packages (from fmpy) (1.26.4)
Requirement already satisfied: MarkupSafe>=2.0 in /usr/local/lib/python3.11/dist-packages (from Jinja2->fmpy) (2.1.5)
Downloading FMPy-0.3.22-py3-none-any.whl (4.9 MB)
----- 4.9/4.9 MB 42.7 MB/s eta 0:00:00
Downloading lark-1.2.2-py3-none-any.whl (111 kB)
----- 111.0/111.0 kB 8.8 MB/s eta 0:00:00
Installing collected packages: lark, fmpy
Successfully installed fmpy-0.3.22 lark-1.2.2

```

## ✓ 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

```
%bash
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
```

```

Model for bioreactor has been setup. Key commands:
- par()      - change of parameters and initial values
- init()     - change initial values only
- simu()     - simulate and plot
- newplot()  - make a new plot
- show()     - show plot from previous simulation

```


```
- disp()      - display parameters and initial values from the last simulation
- describe()  - describe culture, broth, parameters, variables with values/units
```

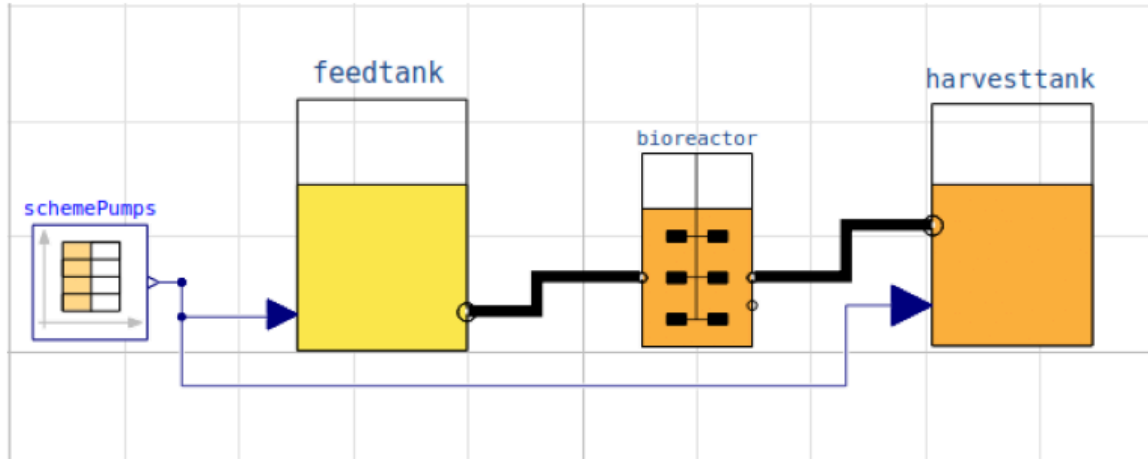
Note that both `disp()` and `describe()` takes values from the last simulation and the command `process_diagram()` brings up the main configuration

Brief information about a command by `help()`, eg `help(simu)`  
Key system information is listed with the command `system_info()`

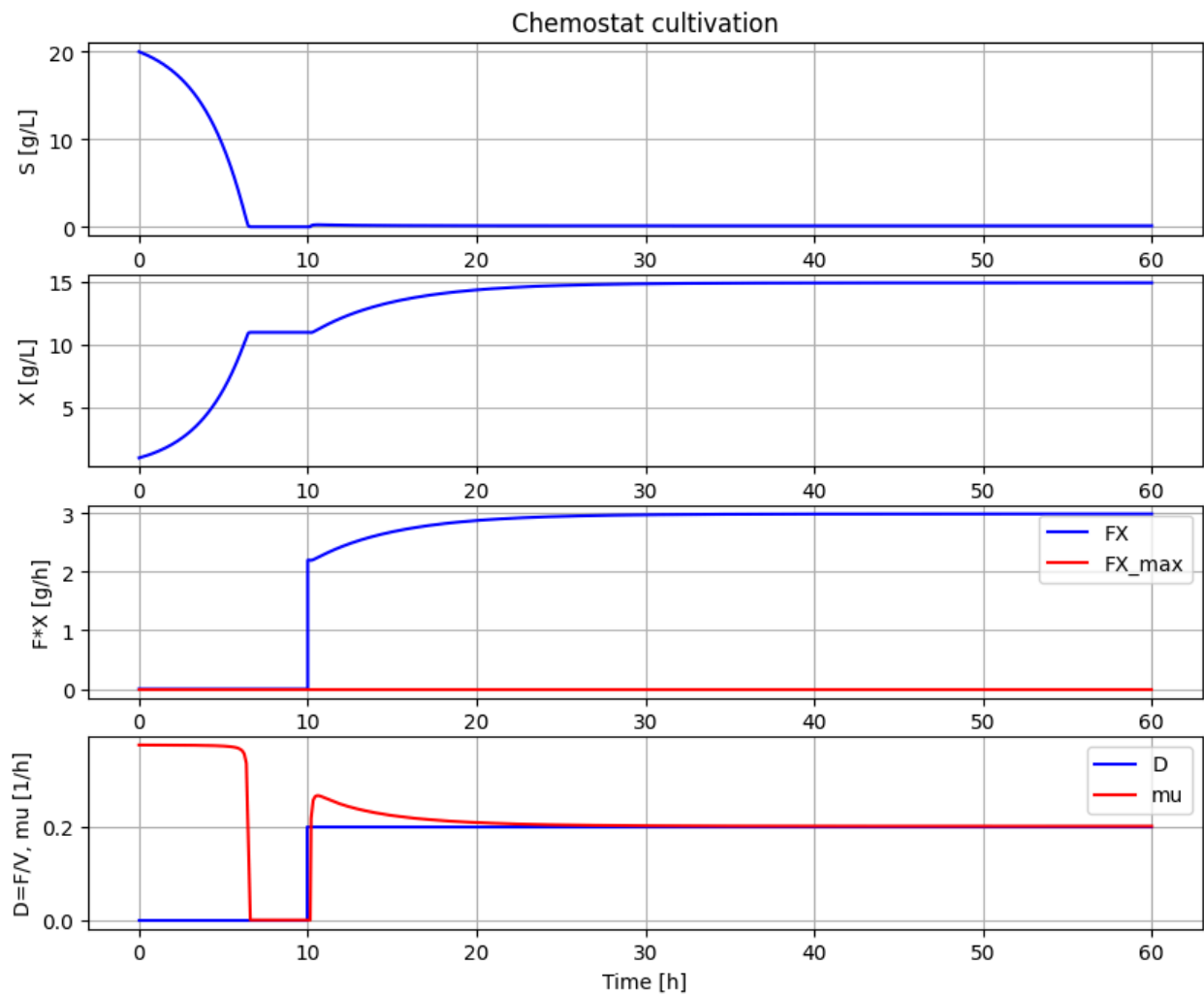
```
%matplotlib inline
plt.rcParams['figure.figsize'] = [25/2.54, 20/2.54]
```

```
process_diagram()
```

 No processDiagram.png file in the FMU, but try the file on disk.



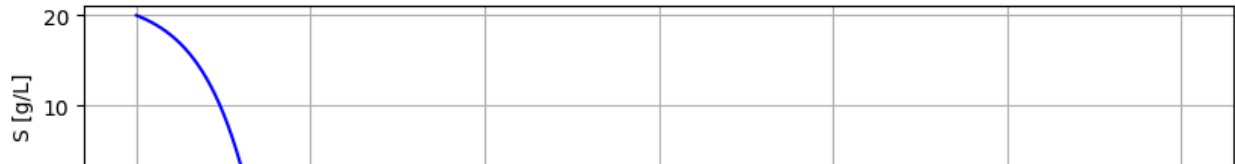
```
newplot()
par(Y=0.50, qSmax=0.75, Ks=0.1)      # Culture parameters
init(V_start=1.0, VX_start=1.0, VS_start=20)  # Bioreactor startup
par(S_in=30, t0=0, F0=0, t1=10, F1=0.2)      # Substrate feeding
simu(60)
```



```
# Test simu('cont')
newplot()
simu(20)
simu(40, 'cont')
```



### Chemostat cultivation



```
describe('parts')
```



```
['bioreactor', 'bioreactor.culture', 'D', 'feedtank', 'harvesttank', 'schemePumps']
```

```
describe('MSL')
```



```
MSL: 3.2.3 - used components: RealInput, RealOutput, CombiTimeTable, Types
```

```
system_info()
```



```
System information
```

```
-OS: Linux
```

```
-Python: 3.11.11
```

```
Colab not installed in the notebook
```