

✓ BPL_TEST2_Perfusion script with PyFMI

The key library PyFMI is installed.

After the installation a small application BPL_TEST2_Perfusion 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

```

```
%env PYTHONPATH=
```

```
env: PYTHONPATH=
```

```
!python --version
```

```
Python 3.11.11
```

```
!wget https://repo.anaconda.com/miniconda/Miniconda3-py311_24.11.1-0-Linux-x86_64.sh
```

```
!chmod +x Miniconda3-py311_24.11.1-0-Linux-x86_64.sh
```

```
!bash ./Miniconda3-py311_24.11.1-0-Linux-x86_64.sh -b -f -p /usr/local
```

```
import sys
```

```
sys.path.append('/usr/local/lib/python3.11/site-packages/')
```

```

--2025-03-25 13:02:39-- https://repo.anaconda.com/miniconda/Miniconda3-py311_24.11.1-0-Linux-x86_64.sh
Resolving repo.anaconda.com (repo.anaconda.com)... 104.16.191.158, 104.16.32.241, 2606:4700::6810:bf9e, ...
Connecting to repo.anaconda.com (repo.anaconda.com)|104.16.191.158|:443... connected.
HTTP request sent, awaiting response... 200 OK
Length: 145900576 (139M) [application/octet-stream]
Saving to: 'Miniconda3-py311_24.11.1-0-Linux-x86_64.sh'

```

```
Miniconda3-py311_24 100%[=====] 139.14M 178MB/s in 0.8s
```

```
2025-03-25 13:02:40 (178 MB/s) - 'Miniconda3-py311_24.11.1-0-Linux-x86_64.sh' saved [145900576/145900576]
```

```
PREFIX=/usr/local
```

```
Unpacking payload ...
```

```
Installing base environment...
```

```
Preparing transaction: ...working... done
```

```
Executing transaction: ...working... done
```

```
installation finished.
```

```
!conda update -n base -c defaults conda --yes
```

```

Channels:
- defaults
Platform: linux-64
Collecting package metadata (repodata.json): done
Solving environment: done

```

```
## Package Plan ##
```

```
environment location: /usr/local
```

```
added / updated specs:
```

```
- conda
```

The following packages will be downloaded:

package	build	
ca-certificates-2025.2.25	h06a4308_0	129 KB
certifi-2025.1.31	py311h06a4308_0	163 KB
openssl-3.0.16	h5eee18b_0	5.2 MB
Total:		5.5 MB

The following packages will be UPDATED:

```
ca-certificates      2024.11.26-h06a4308_0 --> 2025.2.25-h06a4308_0
certifi              2024.8.30-py311h06a4308_0 --> 2025.1.31-py311h06a4308_0
openssl              3.0.15-h5eee18b_0 --> 3.0.16-h5eee18b_0
```

Downloading and Extracting Packages:

```
openssl-3.0.16      | 5.2 MB | : 0% 0/1 [00:00<?, ?it/s]
certifi-2025.1.31   | 163 KB | : 0% 0/1 [00:00<?, ?it/s]

ca-certificates-2025 | 129 KB | : 0% 0/1 [00:00<?, ?it/s]
certifi-2025.1.31    | 163 KB | : 100% 1.0/1 [00:00<00:00, 13.29it/s]

openssl-3.0.16       | 5.2 MB | : 33% 0.328085334517827/1 [00:00<00:00, 3.22it/s]

ca-certificates-2025 | 129 KB | : 100% 1.0/1 [00:00<00:00, 8.87it/s]

ca-certificates-2025 | 129 KB | : 100% 1.0/1 [00:00<00:00, 8.87it/s]
certifi-2025.1.31    | 163 KB | : 100% 1.0/1 [00:00<00:00, 7.39it/s]
```

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

```
!conda --version
!python --version
```

```
🔄 conda 24.11.1
   Python 3.11.11
```

```
!conda config --set channel_priority strict
```

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

```
🔄
```

```

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

```

✓ Notes of BPL_TEST2_Perfusion

This notebook explore perfusion cultivation in comparison with ordinary continuous cultivation (chemostat) and use comparable settings to earlier notebook. Further you see here examples of interaction with the simplified commands `par()`, `init()`, `simu()` etc as well as direct interaction with the FMU which is called "model" here. The last simulation is always available in the workspace and called "sim_res". Note that `describe()` brings mainly up from descriptive information from the Modelica code from the FMU but is complemented by some information given in the Python setup file.

Now specific installation run a simulation and notebook for that Start with connecting to Github. Then upload the two files:

- FMU - BPL_TEST2_Perfusion_linux_om_me.fmu
- Setup-file - BPL_TEST2_Perfusion_explore.py

```

%%bash
git clone https://github.com/janpeter19/BPL_TEST2_Perfusion

```

📂 Cloning into 'BPL_TEST2_Perfusion'...

```

%cd BPL_TEST2_Perfusion

```

📂 /content/BPL_TEST2_Perfusion

```

run -i BPL_TEST2_Perfusion_explore.py

```

📂 Linux - run FMU pre-comiled OpenModelica

```

Model for the process 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()`

```

# Filter out DeprecationWarnings for 'np.float as alias' is needed - wish
import warnings
warnings.filterwarnings("ignore")

```

```

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

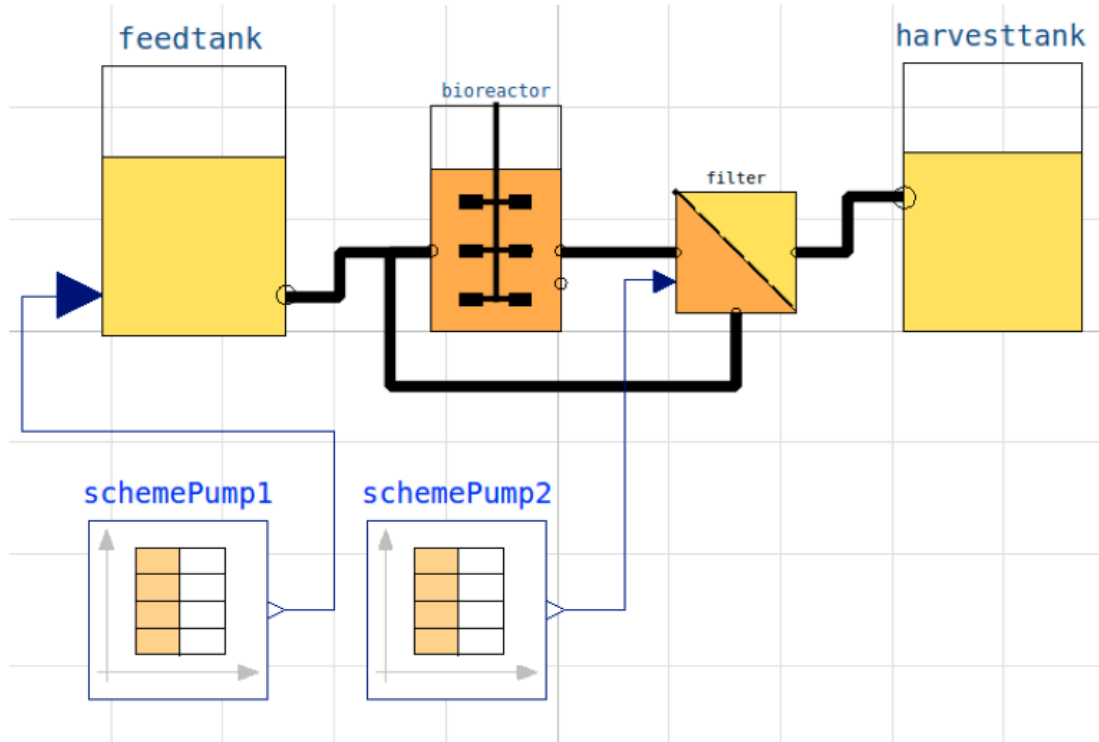
```

```

process_diagram()

```

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



```
# Process parameters used throughout
par(Y=0.5, qSmax=0.75, Ks=0.1)
par(filter_eps=0.10, filter_alpha_X=0.02, filter_alpha_S=0.10)
par(S_in=30.0)
init(V_start=1.0, VX_start=1.0)
eps = parDict['filter_eps']

# Culture
# Filter
# Inlet substrate conc
# Process initial conditions that are common
# Pump schedule parameter

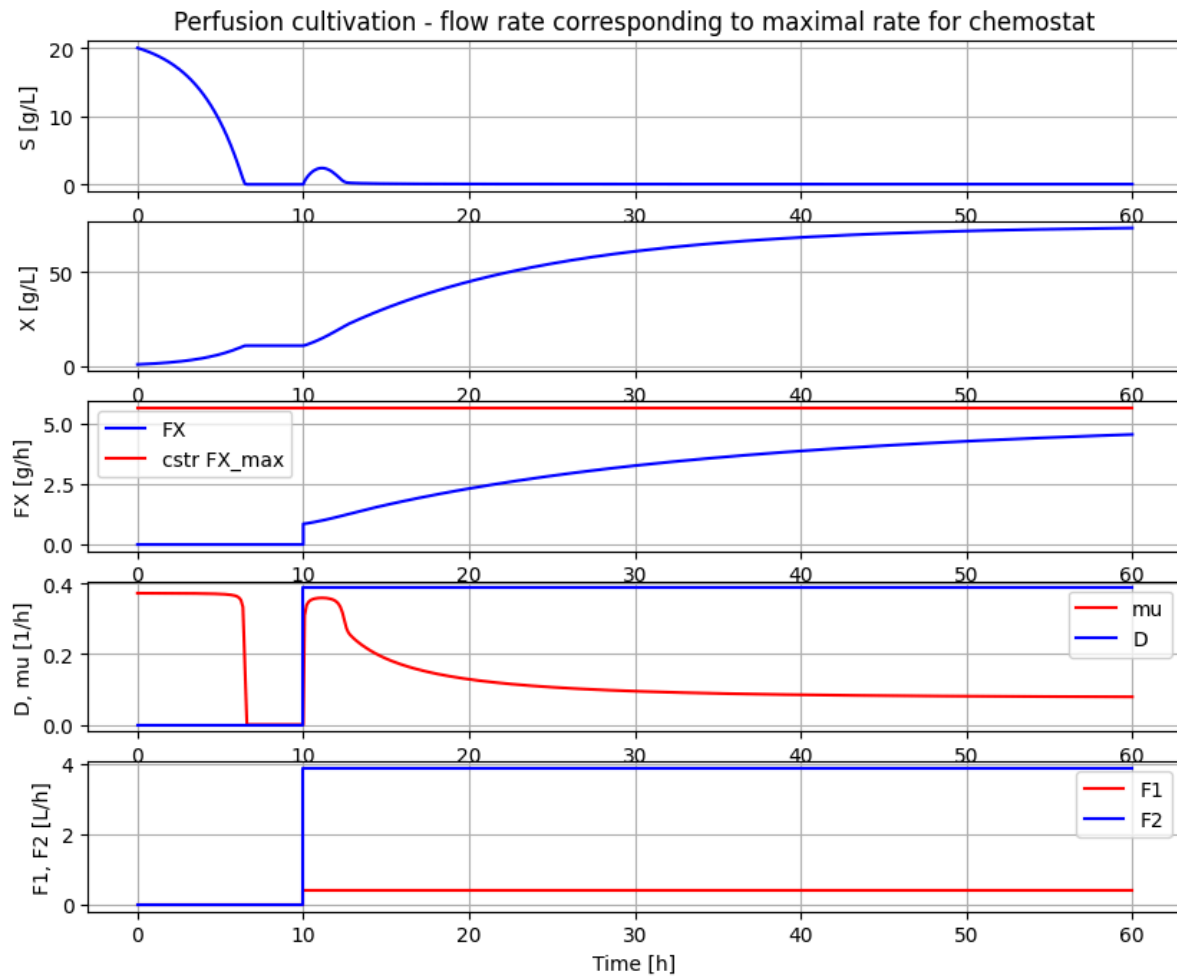
# Simulation of process with flow rate close to wash-out for chemostat

init(VS_start=20)
par(pump1_t1=10, pump2_t1=10)
par(pump1_F1=2.5*0.155, pump2_F1=2.5*0.155/eps)
par(pump1_t2=940, pump2_t2=940, pump1_t3=950, pump2_t3=950, pump1_t4=960, pump2_t4=960)

# Process initial
# Pump schedule - recycle flow 10 times perfusion flow

newplot(title='Perfusion cultivation - flow rate corresponding to maximal rate for chemostat')
simu(60)
```

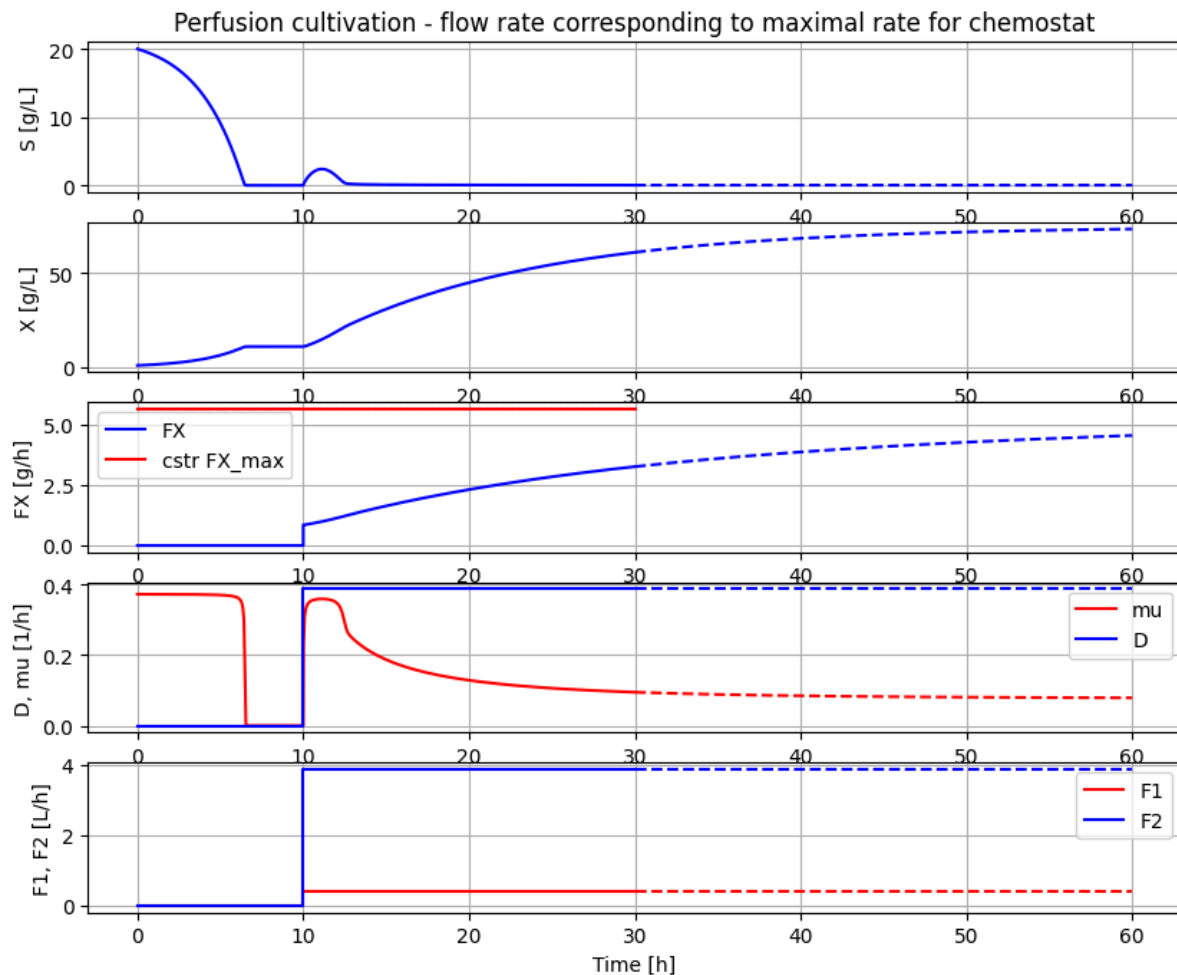
⚠ Could not find cannot import name 'dopri5' from 'assimulo.lib' (/usr/local/lib/python3.11/site-packages/assimulo)
 Could not find cannot import name 'rodas' from 'assimulo.lib' (/usr/local/lib/python3.11/site-packages/assimulo)
 Could not find cannot import name 'odassl' from 'assimulo.lib' (/usr/local/lib/python3.11/site-packages/assimulo)
 Could not find ODEPACK functions.
 Could not find RADAR5
 Could not find GLIMDA.



Simulation of process with flow rate close to wash-out for chemostat

```
init(VS_start=20)                                # Process initial
par(pump1_t1=10, pump2_t1=10)                    # Pump schedule - recycle flow 10 times perfusion flow
par(pump1_F1=2.5*0.155, pump2_F1=2.5*0.155/eps)
par(pump1_t2=940, pump2_t2=940, pump1_t3=950, pump2_t3=950, pump1_t4=960, pump2_t4=960)

newplot(title='Perfusion cultivation - flow rate corresponding to maximal rate for chemostat')
simu(30)
simu(30,'cont')
```



Note the inability of the OpenModelica FMU to handle `simu('cont')` properly.

```
# Concentration factor of the filter
c=model.get('filter.retentate.c[1]')[0]/model.get('filter.inlet.c[1]')[0]
print('Conc factor of perfusion filter =', np.round(c,3))
```

Conc factor of perfusion filter = 1.179

```
c_data=sim_res['filter.retentate.c[1]']/sim_res['filter.inlet.c[1]']
print('Conc factor variation', np.round(min(c_data[151:]), 3), np.round(max(c_data[151:]),3))
```

Conc factor variation 1.179 1.179

```
# Simulation of process with step-wise increase of perfusion rate until wash-out.
# This means that re-circulation rate change at the same time as the perfusion rate.
```

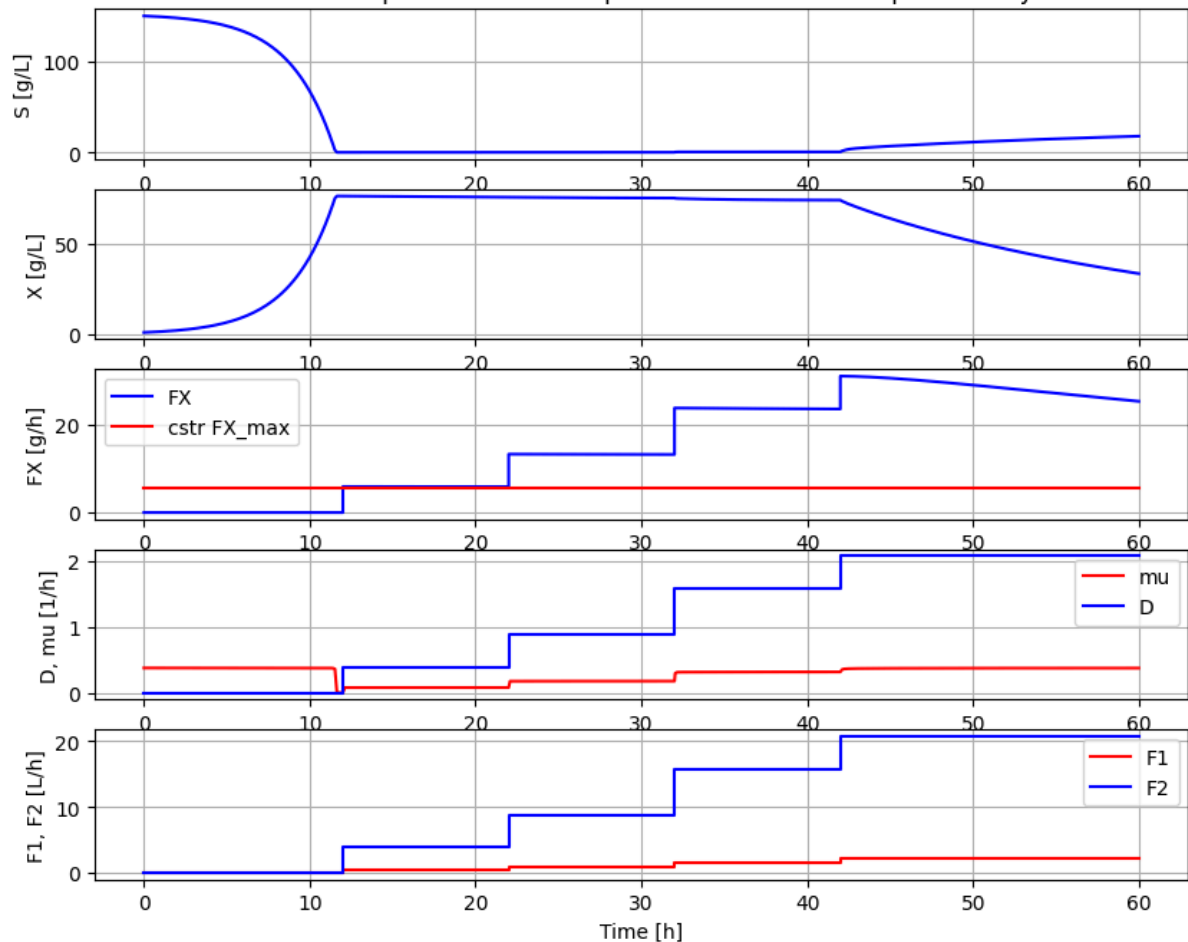
```
init(VS_start=150) # Process initial varied

par(pump1_t1=12, pump2_t1=12) # Pump schedule - recycle flow 10 times perfusion flow
par(pump1_F1=2.5*0.155, pump2_F1=2.5*0.155/eps)
par(pump1_t2=22, pump2_t2=22)
par(pump1_F2=2.5*0.35, pump2_F2=2.5*0.35/eps)
par(pump1_t3=32, pump2_t3=32)
par(pump1_F3=2.5*0.63, pump2_F3=2.5*0.63/eps)
par(pump1_t4=42, pump2_t4=42)
par(pump1_F4=2.5*0.83, pump2_F4=2.5*0.83/eps)

newplot(title='Perfusion cultivation - step wise increase of perfusion rate and max productivity 4x chemostat')
simu(60)
```



Perfusion cultivation - step wise increase of perfusion rate and max productivity 4x chemostat



```
# Simulation without a plot and just to check typical values at high production rate
simu(40)
c_data=sim_res['filter.retentate.c[1]']/sim_res['filter.inlet.c[1]']
print('Conc factor variation', np.round(min(c_data[190:]), 3), 'to', np.round(max(c_data[190:]),3))
```



Conc factor variation 1.162 to 1.179

```
#describe('cstrProdMax')
```

```
# The maximal biomass productivity before washout is obtained aroundn 40 hours
np.round(model.get('harvesttank.inlet.F')[0]*model.get('harvesttank.inlet.c[1]')[0],1)
```



23.5

```
# Thus perfusion (with this filter) brings a productivity improvement of about
np.round(23.5/5.6,1)
```



4.2

```
# Finally we check the filter flow rates at time 40 hour - note the negative sign for outflow
model.get('filter.inlet.F')[0]
```



15.749999999999998

```
model.get('filter.filtrate.F')[0]
```



-1.575

```
model.get('filter.retentate.F')[0]
```



-14.174999999999999

✓ Summary

- The perfusion filter had a concentration factor of cells around 1.08 and re-cycling flow was set to a factor 10 higher than the