

## ✓ BPL\_YEAST\_COB\_Batch script with PyFMI

The key library PyFMI is installed.

After the installation a small application BPL\_YEAST\_COB\_Batch 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 22.04.3 LTS
Release:      22.04
Codename:     jammy

```

```
%env PYTHONPATH=
```

```
env: PYTHONPATH=
```

```
!wget https://repo.anaconda.com/miniconda/Miniconda3-py310_23.1.0-1-Linux-x86_64.sh
```

```
!chmod +x Miniconda3-py310_23.1.0-1-Linux-x86_64.sh
```

```
!bash ./Miniconda3-py310_23.1.0-1-Linux-x86_64.sh -b -f -p /usr/local
```

```
import sys
```

```
sys.path.append('/usr/local/lib/python3.10/site-packages/')
```

```

--2025-01-15 07:31:14--  https://repo.anaconda.com/miniconda/Miniconda3-py310_23.1.0-1-Linux-x86_64.sh
Resolving repo.anaconda.com (repo.anaconda.com)... 104.16.32.241, 104.16.191.158, 2606:4700::6810:bf9e,
Connecting to repo.anaconda.com (repo.anaconda.com)|104.16.32.241|:443... connected.
HTTP request sent, awaiting response... 200 OK
Length: 74403966 (71M) [application/x-sh]
Saving to: 'Miniconda3-py310_23.1.0-1-Linux-x86_64.sh'

```

```
Miniconda3-py310_23 100%[=====] 70.96M 107MB/s in 0.7s
```

```
2025-01-15 07:31:15 (107 MB/s) - 'Miniconda3-py310_23.1.0-1-Linux-x86_64.sh' saved [74403966/74403966]
```

```

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 3.10.16
```

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

```
⇄
```

```

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

```

```
!pip install optlang
```

```

🔗 Collecting optlang
  Downloading optlang-1.8.3-py2.py3-none-any.whl (141 kB)
    ━━━━━━━━━━━━━━━━━━━━━━━━━━━━━━━━━ 141.8/141.8 kB 8.2 MB/s eta 0:00:00
Collecting swiglpk>=5.0.12
  Downloading swiglpk-5.0.12-cp310-cp310-manylinux_2_17_x86_64.manylinux2014_x86_64.whl (2.3 MB)
    ━━━━━━━━━━━━━━━━━━━━━━━━━━━━━━━━━ 2.3/2.3 MB 44.4 MB/s eta 0:00:00
Collecting sympy>=1.12.0
  Downloading sympy-1.13.3-py3-none-any.whl (6.2 MB)
    ━━━━━━━━━━━━━━━━━━━━━━━━━━━━━━━━━ 6.2/6.2 MB 86.6 MB/s eta 0:00:00
Collecting mpmath<1.4,>=1.1.0
  Downloading mpmath-1.3.0-py3-none-any.whl (536 kB)
    ━━━━━━━━━━━━━━━━━━━━━━━━━━━━━━━━━ 536.2/536.2 kB 43.3 MB/s eta 0:00:00
Installing collected packages: swiglpk, mpmath, sympy, optlang
Successfully installed mpmath-1.3.0 optlang-1.8.3 swiglpk-5.0.12 sympy-1.13.3

```

## ✓ Notes YEAST\_COB\_Batch constraint-based approach

Now specific installation and the run simulations. Start with connecting to Github. Then upload the two files:

- FMU - BPL\_YEAST\_AIR\_Fedbatch\_linux\_jm\_cs.fmu
- Setup-file - BPL\_YEAST\_AIR\_Fedbatch\_explore

```

%%bash
git clone https://github.com/janpeter19/CONF_2023_10_MODELICA15

```

```
🔗 Cloning into 'CONF_2023_10_MODELICA15'...
```

```
%%cd CONF_2023_10_MODELICA15
```

```
🔗 /content/CONF_2023_10_MODELICA15
```

```
run -i BPL_YEAST_COB_Batch_explore.py
```

```
🔗 Linux - run FMU pre-comiled OpenModelica 1.21.0
```

```

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

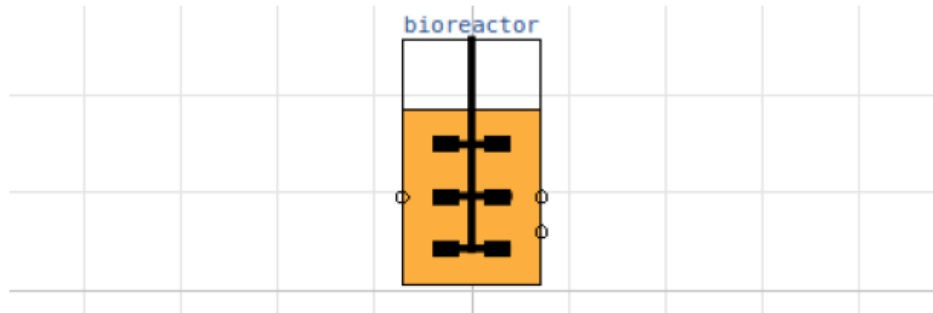
```

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()`

```
plt.rcParams['figure.figsize'] = [20/2.54, 16/2.54]
```

```
process_diagram()
```



## ✓ Try using LP in each step

```
from optlang import Model, Variable, Constraint, Objective
```

```
# Define culture constraint-based model
def culture(G, E):
```

```
    # LP calculation of the optimal qGr, qEr based on G and E values
```

```
    # - parameters
```

```
    q02max = 6.9e-3; kog = 2.3; koe = 1.6; YGr = 3.5; YEr = 1.32;
    alpha = 0.01; beta = 1.0
```

```
    # - transfer data from dynamic reactor model to static LP model
```

```
    qGr_opt = Variable('qGr_opt', lb=0)
```

```
    qEr_opt = Variable('qEr_opt', lb=0)
```

```
    # - LP model constraint and objective
```

```
    mu_max = Objective(YGr*qGr_opt + YEr*qEr_opt, direction='max')
```

```
    q02lim = Constraint(kog*qGr_opt + koe*qEr_opt, ub=q02max)
```

```
    qGlim = Constraint(qGr_opt, ub=alpha*max(0,G))
```

```
    qElim = Constraint(qEr_opt, ub=beta*max(0,E))
```

```
    # - put together the LP model
```

```
    yeast_model = Model(name='Yeast bottleneck model')
```

```
    yeast_model.objective = mu_max
```

```
    yeast_model.add(q02lim)
```

```
    yeast_model.add(qGlim)
```

```
    yeast_model.add(qElim)
```

```
    # - do LP optimization
```

```
    yeast_model.optimize()
```

```
    return (yeast_model.objective.value, yeast_model.variables.qGr_opt.primal, yeast_model.variables.qEr_opt
```

```
# Initialization
```

```
V_0=1.0
```

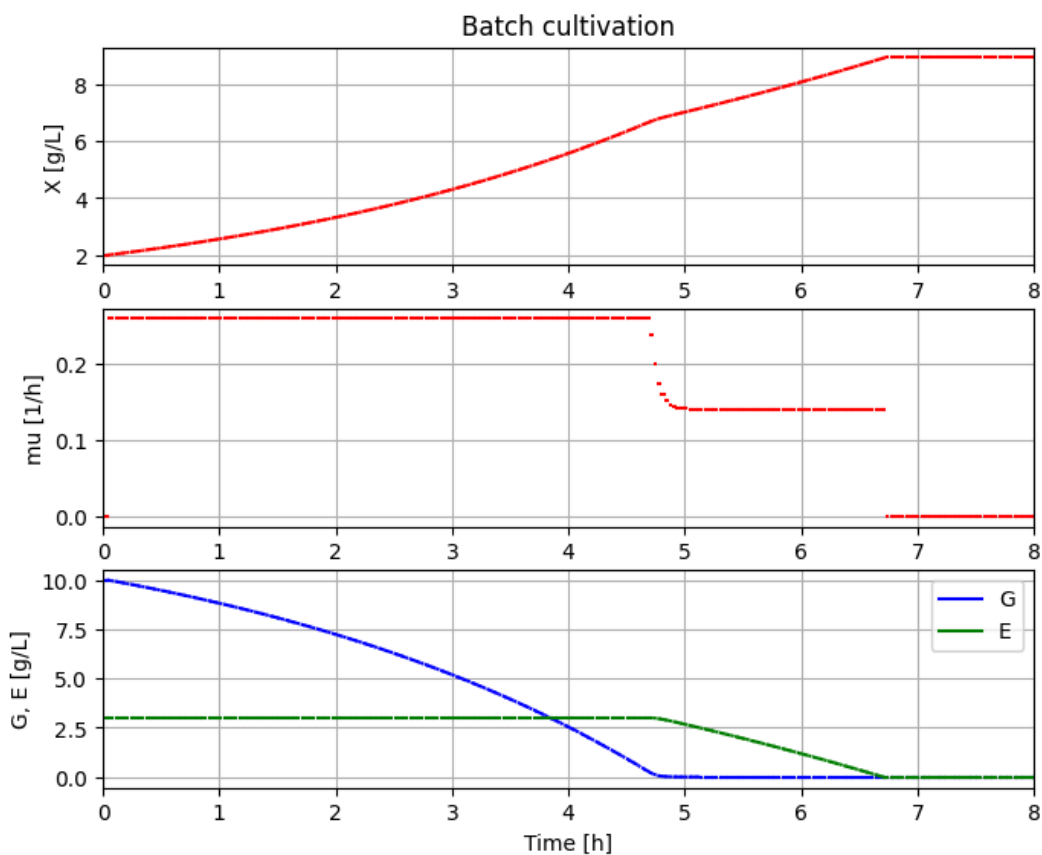
```
init(V_0=V_0, VX_0=V_0*2.0, VG_0=V_0*10, VE_0=3.0)
```

```
# Loop of simulations
t_final = 8.0
t_samp = 0.0333
n_samp = t_final/t_samp + 1

# Simulate n sample steps
newplot(title='Batch cultivation', plotType='TimeSeries2')
ax1.set_xlim([0, t_final]); ax2.set_xlim([0, t_final]); ax3.set_xlim([0, t_final])

simu(t_samp, options=opts_fast)
for i in range(int(n_samp)):
    (mum_opt, qGr_opt, qEr_opt, qO2_opt) = culture(sim_res['bioreactor.c[2]'][-1], sim_res['bioreactor.c[3]']
    par(mum=mum_opt, qGr=qGr_opt, qEr=qEr_opt, qO2=qO2_opt)
    simu(t_samp, 'cont', options=opts_fast)
```

⚠ Could not find cannot import name 'dopri5' from 'assimulo.lib' (/usr/local/lib/python3.10/site-packages/  
 Could not find cannot import name 'rodas' from 'assimulo.lib' (/usr/local/lib/python3.10/site-packages/a  
 Could not find cannot import name 'odassl' from 'assimulo.lib' (/usr/local/lib/python3.10/site-packages/  
 Could not find ODEPACK functions.  
 Could not find RADAR5  
 Could not find GLIMDA.



```
system_info()
```



```
System information
-OS: Linux
-Python: 3.10.12
-Scipy: not installed in the notebook
-PyFMI: 2.16.1
-FMU by: OpenModelica Compiler OpenModelica 1.21.0
-FMI: 2.0
-Type: FMUModelME2
-Name: BPL_YEAST_COB.Batch
-Generated: 2023-05-31T09:43:28Z
-MSL: 3.2.3
-Description: Bioprocess Library version 2.1.1
-Interaction: FMU-explore version 0.9.8
```

```
!conda list optlang
```

```
↗ # packages in environment at /usr/local:  
#  
# Name                Version             Build    Channel  
optlang                1.8.3               pypi_0   pypi
```