▼ 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 22.04.2 LTS
                       22.04
     Release:
     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/')
     --2023-09-26 06:56:18-- <a href="https://repo.anaconda.com/miniconda/Miniconda3-py310_23.1.0-1-Linux-x86_64.sh">https://repo.anaconda.com/miniconda/Miniconda3-py310_23.1.0-1-Linux-x86_64.sh</a> Resolving repo.anaconda.com (repo.anaconda.com)... 104.16.131.3, 104.16.130.3, 2606:4700::6810:8303, ...
     Connecting to repo.anaconda.com (repo.anaconda.com) | 104.16.131.3 | :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
     2023-09-26 06:56:18 (185 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 23.7.4 Python 3.10.13

!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-1.7.0-py2.py3-none-any.whl (138 kB)
                                                 - 138.3/138.3 kB 3.3 MB/s eta 0:00:00
    Requirement \ already \ satisfied: \ six>=1.9 \ in \ /usr/local/lib/python 3.10/site-packages \ (from \ optlang) \ (1.16.0)
    Collecting swiglpk>=5.0.8
      Downloading swiglpk-5.0.8-cp310-cp310-manylinux_2_17_x86_64.manylinux2014_x86_64.whl (2.3 MB)
                                                  - 2.3/2.3 MB 36.0 MB/s eta 0:00:00
    Collecting sympy>=1.12.0
      Downloading sympy-1.12-py3-none-any.whl (5.7 MB)
                                                  - 5.7/5.7 MB 78.6 MB/s eta 0:00:00
    Collecting mpmath>=0.19
      Downloading mpmath-1.3.0-py3-none-any.whl (536 kB)
                                                 - 536.2/536.2 kB 42.8 MB/s eta 0:00:00
    Installing collected packages: swiglpk, mpmath, sympy, optlang
    Successfully installed mpmath-1.3.0 optlang-1.7.0 swiglpk-5.0.8 sympy-1.12
    WARNING: Running pip as the 'root' user can result in broken permissions and conflicting behaviour with the system package
```

Notes YEAST_COB_Batch constraint-based approach

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

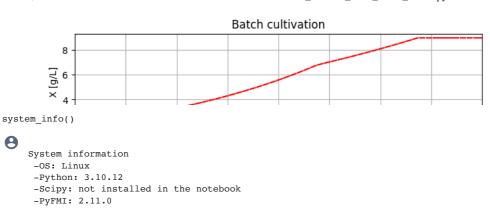
- FMU BPL_YEAST_AIR_Fedbatch_linux_jm_cs.fmu
- Setup-file BPL_YEAST_AIR_Fedbatch_explore

```
%%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
                   - simulate and plot
     - simu()
     - newplot()
                  - make a new plot
     - show()
                   - show plot from previous simulation
                   - 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()
                                     bioreactor
```

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.primal, qO2lim.pr
# 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)):
    (\texttt{mum\_opt}, \ \texttt{qGr\_opt}, \ \texttt{qGr\_opt}, \ \texttt{qO2\_opt}) = \texttt{culture}(\texttt{sim\_res['bioreactor.c[2]'][-1]}, \ \texttt{sim\_res['bioreactor.c[3]'][-1]})
    par(mum=mum_opt, qGr=qGr_opt, qEr=qEr_opt, qO2=qO2_opt)
    simu(t_samp, 'cont', options=opts_fast)
```

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-Type: FMUModelME2

-Name: BPL_YEAST_COB.Batch

-Generated: 2023-05-31T09:43:28Z

-MSL: 3.2.3

-FMI: 2.0

-Description: Bioprocess Library version 2.1.1

-FMU by: OpenModelica Compiler OpenModelica 1.21.0

-Interaction: FMU-explore version 0.9.8

<u>−</u> 5.0 +

!conda list optlang

```
# packages in environment at /usr/local:
# Name
                        Version
                                               Build Channel
optlang
                       1.7.0
                                              pypi_0 pypi
```