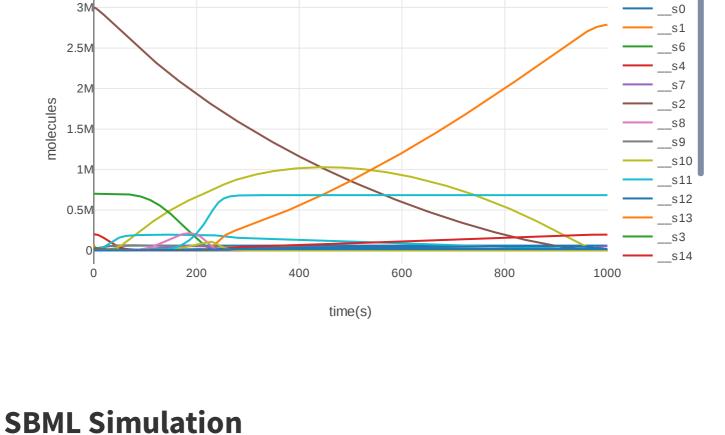
```
[7] sbml = model.toSBML()
      # print(sbml)
      from roadrunner import RoadRunner
      with open('out-sbml.xml', 'w') as f:
          f.write(sbml)
      xmod = RoadRunner(sbml)
      xmod.reset()
      xmod.integrator.relative_tolerance = 1e-12
      xmod.integrator.absolute_tolerance = 1e-20
      # xmod.oneStep()
      xmod.simulate(0,1000,1000,selections=['time']+xmod.getFloatingSpeciesIds())
      xmod.plot(xtitle='time(s)', ytitle='molecules')
      values = {s:xmod[s] for s in xmod.getFloatingSpeciesIds()}
      block_rates = [xmod['block{}_Add10'.format(k)] for k in (range(len(wiring.blocks)))]
      # pprint(block_rates)
      r_block_indices = model.getReactionBlockIndices()
```



SBML comparison import tellurium as te with open(sbfile) as f:

```
rr_model = te.loada(f.read())
rr_model.reset()
rr_model.integrator.relative_tolerance = 1e-12
rr_model.integrator.absolute_tolerance = 1e-20
# rr_model.selections =
['time','__s0','__s1','__s6','__s4','__s7','__s8','__s2','__s9','__s10','__s15']
rr_model.simulate(0,1000,2000)
for var, value in values.items():
    print('rel diff for {}: {:.3f}'.format(var,(rr_model[var]-value)/max(rr_model[var],1.)))
    rr_model[var] = value
# print('\nreactions:')
# for r_index in range(rr_model.getNumReactions()):
      r_id = rr_model.getReactionIds()[r_index]
      block_index = r_block_indices[r_index]
       print('{} rel diff: {:.3f}'.format(r_id,(rr_model[r_id]-
block_rates[block_index])/max(rr_model[r_id],1.)))
rr_model.plot(xtitle='time(s)', ytitle='molecules')
rel diff for __s15: 0.000
rel diff for __s5: 0.000
rel diff for __s16: 0.000
rel diff for __s17: -0.000
rel diff for __s18: 0.000
rel diff for __s19: -0.000
rel diff for __s20: 0.000
        ЗМ
                                                                                [_s0]
                                                                                __s1]
      2.5M
                                                                               [_s2]
                                                                               <u>[_s3]</u>
                                                                               - [_s4]
        2M
   molecules
                                                                               - [_s5]
                                                                               <u>[_s6]</u>
      1.5M
                                                                               - [_s7]
                                                                               _ [_s8]
        1M
                                                                              - [__s10]
      0.5M
                                                                               <u>[_s11]</u>
                                                                               - [_s12]
                                                                               - [_s13]
                     200
                                               600
                                                           800
                                  400
                                                                        1000
                                       time(s)
```