

---

# pyqo Documentation

*Release 0.1*

**Sebastian Krämer**

October 25, 2011



# CONTENTS

<b>1</b>	<b>pyqo Tutorial</b>	<b>3</b>
1.1	Importing <b>pyqo</b> . . . . .	3
1.2	Defining states . . . . .	3
1.3	Defining operators . . . . .	4
<b>2</b>	<b>Indices and tables</b>	<b>5</b>



Contents:



# PYQO TUTORIAL

**pyqo** is a python library similar to the *quantum optics toolbox* for Matlab. It lets the user define state vectors and operators and provides functionality to solve typical problems occurring in quantum optics, e.g. solving Schroedinger or master equations.

## 1.1 Importing pyqo

To be able to use **pyqo** from a python script the **pyqo** library has to be in your PYTHONPATH or in the same directory as the script that uses it. Then it can for example be imported as:

```
>>> import pyqo as qo
```

In all following examples it is assumed that **pyqo** was imported in that way.

## 1.2 Defining states

All objects in **pyqo** are defined as tensors of different ranks. Taking advantage of the power of numpy, they inherit from the mighty `numpy.ndarray`. Obviously their dimensionality is limited by the memory available on your system so it is impossible to calculate in infinite dimensional Hilbert spaces. For quantum states **pyqo** provides the class `pyqo.StateVector`. It can be used to directly create states from nested lists or tuples (or anything else that a numpy array can handle):

```
>>> psi = qo.StateVector([1,0])
>>> print(psi)
StateVector(2)
[ 1.+0.j  0.+0.j]
```

Alternatively there are some functions that create commonly used state vectors:

```
>>> psi = qo.basis(4,0)
>>> print(psi)
StateVector(4)
[ 1.+0.j  0.+0.j  0.+0.j  0.+0.j]
>>> psi = qo.coherent(10, 0.5)
```

Composing systems can be done with the tensor product between two states. For this the operator  $\wedge$  can be used.

```
>>> psi1 = qo.basis(2,0)
>>> psi2 = qo.basis(2,1)
>>> print(psi1 ^ psi2)
StateVector(2 x 2)
```

```
[[ 0.+0.j  1.+0.j]
 [ 0.+0.j  0.+0.j]]
```

---

**Note:** The `^` operator follows the built-in operator precedence. That means “`*`” and “`+`” have higher precedence!

---

## 1.3 Defining operators

Operators are represented by the `pyqo.Operator`. Like in the case of state vectors operators can be constructed directly from a list or tuple:

```
>>> A = qo.Operator([[1,0], [0,-1]])
>>> print(A)
Operator
2 -> 2
[[ 1.+0.j  0.+0.j]
 [ 0.+0.j -1.+0.j]]
```

Operators have some constraint on their shape - it has to be of the form  $(n_1, n_2, \dots, n_N, n_1, n_2, \dots, n_N)$ .

Many commonly used operators are already defined:

```
>>> print(qo.sigmax)
Operator
2 -> 2
[[ 0.+0.j  1.+0.j]
 [ 1.+0.j  0.+0.j]]
>>> print(qo.create(3))
Operator
3 -> 3
[[ 0.00000000+0.j  0.00000000+0.j  0.00000000+0.j]
 [ 1.00000000+0.j  0.00000000+0.j  0.00000000+0.j]
 [ 0.00000000+0.j  1.41421356+0.j  0.00000000+0.j]]
```

Composing operators of different systems can be done in the following way:

```
>>> s_z = qo.sigmaz
>>> s_p = qo.sigmamp
>>> print(s_z^s_p)
Operator
2 x 2 -> 2 x 2
[[[ 0.+0.j  0.+0.j]
  [ 0.+0.j  0.+0.j]]

 [ 1.+0.j  0.+0.j]
 [ 0.+0.j  0.+0.j]]]

[[[ 0.+0.j  0.+0.j]
  [-0.+0.j -0.+0.j]]

 [ 0.+0.j  0.+0.j]
 [-1.+0.j -0.+0.j]]]]
```



# EXAMPLES

## 2.1 Rabi oscillation

### 2.1.1 Code

```
import imp
qo = imp.load_module("pyqo", *imp.find_module("pyqo", [".."]))
import numpy as np

Delta = 2
Omega = 1
phi = np.pi

H = 1./2*(- Delta * qo.sigmaz\
          + Omega * np.exp(1j*phi) * qo.sigmam\
          + Omega * np.exp(-1j*phi) * qo.sigmam)

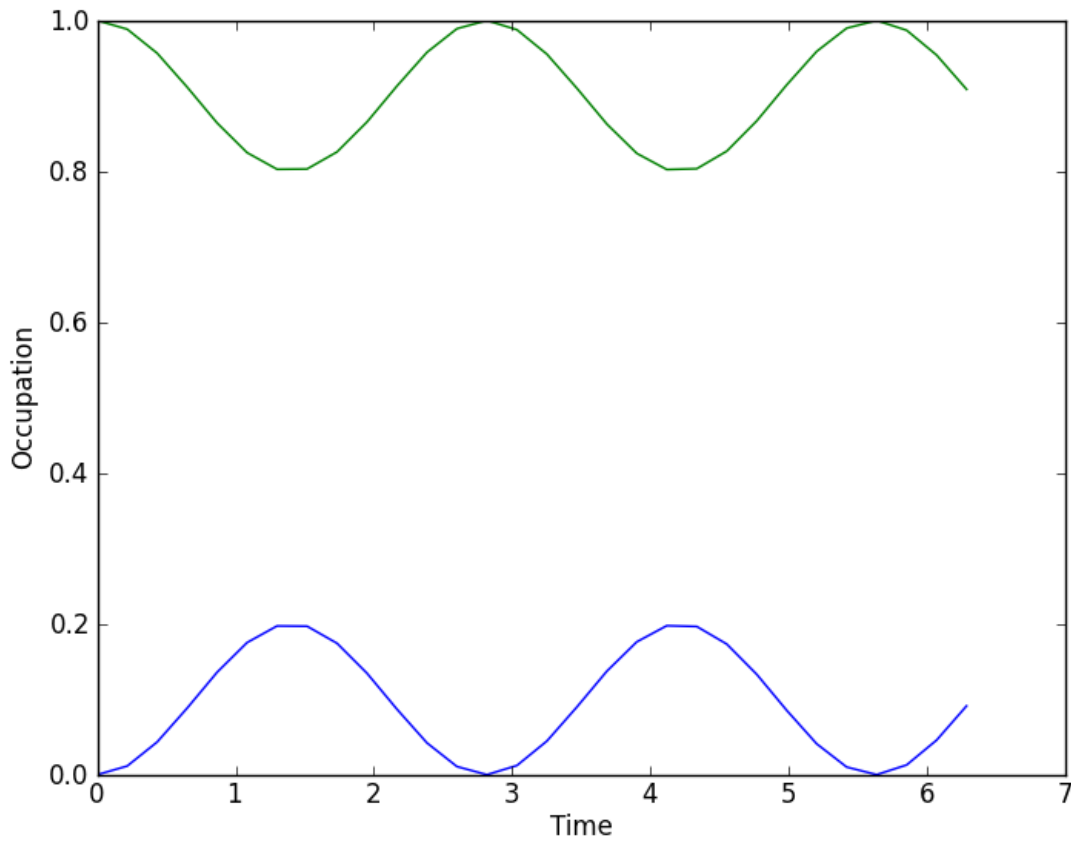
psi_0 = qo.basis(2,0)

T = np.linspace(0,2*np.pi,30)
psi = qo.solve_ode(H, psi_0, T)

e0 = qo.expect(qo.sigmam*qo.sigmam, psi)
e1 = qo.expect(qo.sigmam*qo.sigmam, psi)

import pylab
pylab.plot(T, e0, T, e1)
pylab.xlabel("Time")
pylab.ylabel("Occupation")
pylab.show()
```

## 2.1.2 Output



## 2.2 Jaynes-Cummings model

### 2.2.1 Code

```
import imp
qo = imp.load_module("pyqo", *imp.find_module("pyqo", [".."]))
import numpy as np

N = 10 # dimension of field Hilbert space
delta_c = 1
delta_a = 2
g = 1
gamma = 0.1
kappa = 0.1

# Field
id_f = qo.identity(N)
a = qo.destroy(N)
at = qo.create(N)
n = qo.number(N)

# Atom
```

```

id_a = qo.identity(2)

# Initial state
psi_0 = qo.basis(N,0) ^ qo.basis(2,1)

# Hamiltonian
H = delta_c*(at*a^id_a)\
    + delta_a*(id_f^qo.sigmam*qo.sigmam)\
    + g*(a^qo.sigmam) + g*(at^qo.sigmam)

# Solve Master equation
T = np.linspace(0, 2*np.pi, 30)
rho = qo.solve_ode(H, psi_0, T,
                  [gamma**(1/2)*(id_f^qo.sigmam), kappa**(1/2)*(a^id_a)])

# Expectation values
n_exp = qo.expect(n^id_a, rho)
e_exp = qo.expect(id_f^qo.sigmam*qo.sigmam, rho)

# Q-function
x = np.linspace(-4,4,40)
y = np.linspace(-4,4,40)
X, Y = np.meshgrid(x,y)

# Visualization
import pylab
pylab.figure(1)
pylab.subplot(211)
pylab.xlabel("time")
pylab.ylabel(r"$\langle n \rangle$")
pylab.plot(T, n_exp)
pylab.subplot(212)
pylab.xlabel("time")
pylab.ylabel(r"$\langle P_1 \rangle$")
pylab.plot(T, e_exp)
pylab.show()

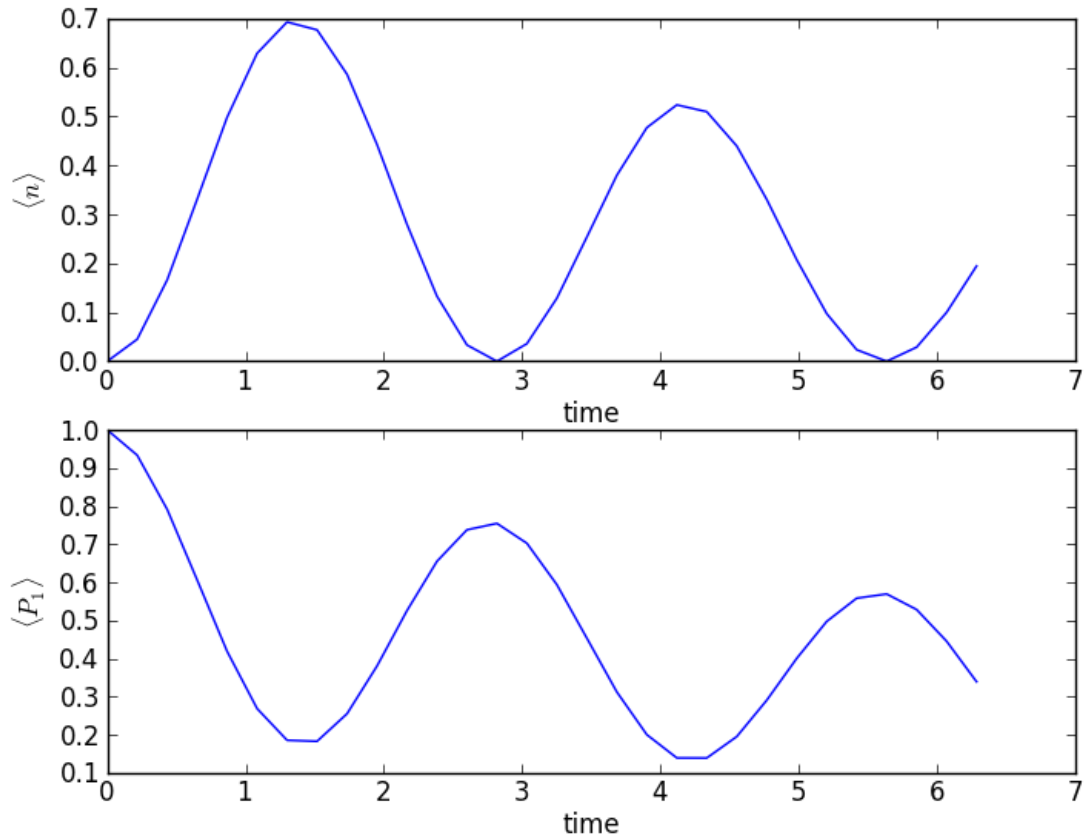
Q = []
for rho_t in rho:
    rho_f = qo.ptrace(rho_t,1)
    Q.append(np.abs(qo.qfunc(rho_f,X,Y)))

def qplot(fig,step):
    axes = fig.add_subplot(111)
    axes.clear()
    axes.imshow(Q[step])

qo.animate(len(rho), qplot)

```

## 2.2.2 Output



## 2.3 Single atom laser

### 2.3.1 Code

```
import imp
qo = imp.load_module("pyqo", *imp.find_module("pyqo", [".."]))
import numpy as np

N = 10 # dimension of field Hilbert space
delta_c = 1
delta_a = 1
g = 1
gamma = 0.2
kappa = 0.1
R = 0.5

# Field
id_f = qo.identity(N)
a = qo.destroy(N)
at = qo.create(N)
n = qo.number(N)
```

---

```

# Atom
id_a = qo.identity(2)

# Initial state
psi_0 = qo.basis(N,0) ^ qo.basis(2,1)

# Hamiltonian
H = delta_c*(a^id_a) + delta_a*(id_f^qo.sigmap*qo.sigmam) + g*(a^qo.sigmap) + g*(a^qo.sigmam)

# Jump operators
j1 = gamma**(1./2)*(id_f^qo.sigmam)
j2 = kappa**(1./2)*(a^id_a)
j3 = R**(1./2)*(id_f^qo.sigmap)
J = [j1, j2, j3]

# Solve Master equation
T = np.linspace(0, 12*np.pi, 80)
rho = qo.solve_ode(H, psi_0, T, J)
#rho = qo.solve_es(H, psi_0, T, J)

# Expectation values
n_exp, e_exp = qo.expect((n^id_a, id_f^qo.sigmap*qo.sigmam), rho)

# Calculate Q-function and photon number distribution
x_min, x_max = -5, 5
y_min, y_max = -5, 5
x = np.linspace(x_min, x_max, 30)
y = np.linspace(y_min, y_max, 30)
X, Y = np.meshgrid(x,y)

Q = []
F = []
for rho_t in rho:
    rho_f = qo.ptrace(rho_t,1)
    Q.append(np.abs(qo.qfunc(rho_f,X,Y)))
    F.append(np.abs(np.diag(rho_f)))

# Visualization
import pylab
pylab.figure(1)
pylab.subplot(211)
pylab.xlabel("time")
pylab.ylabel(r"$\langle n \rangle$")
pylab.plot(T, n_exp)
pylab.ylim(ymin=0)
pylab.subplot(212)
pylab.xlabel("time")
pylab.ylabel(r"$\langle P_1 \rangle$")
pylab.ylim((0,1))
pylab.plot(T, e_exp)
pylab.show()

from scipy.misc import factorial

F_x = np.arange(0, N)
def fplot(fig, step):
    axes = fig.add_subplot(111)
    axes.clear()

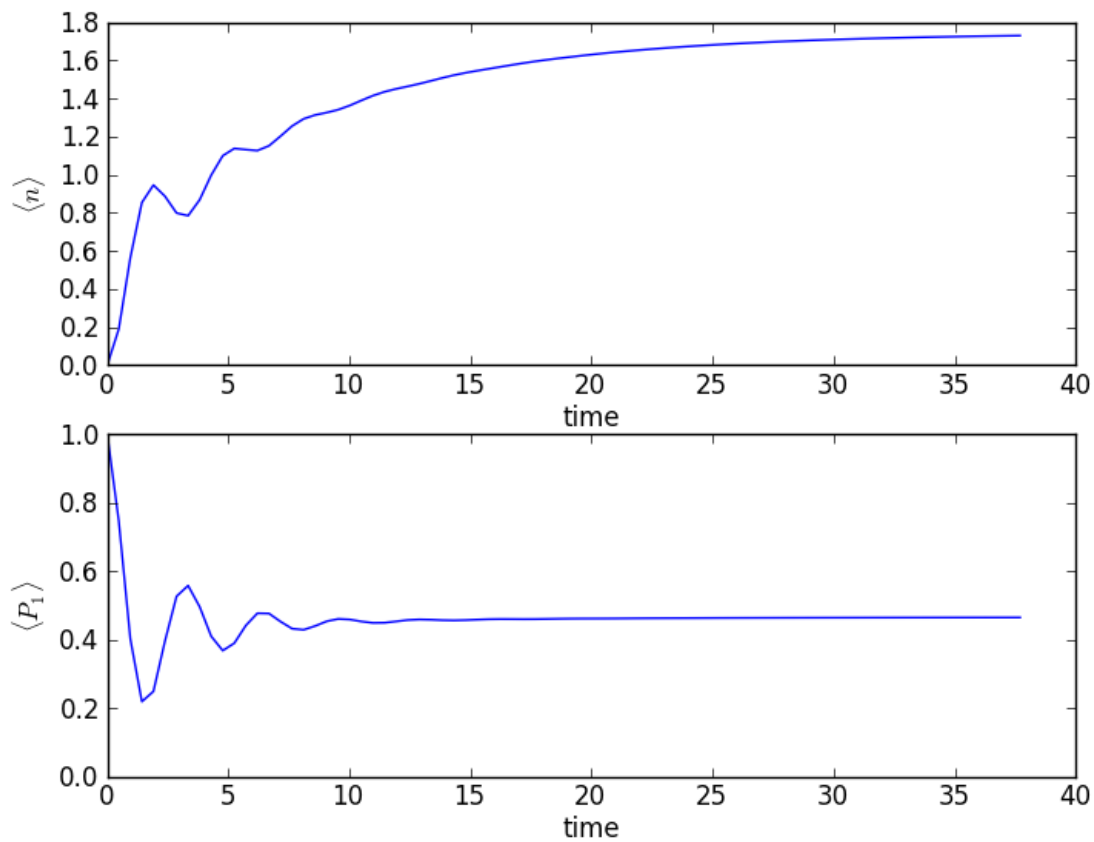
```

```
pylab.plot(F_x, F[step], "o")
n_ = np.abs(n_exp[step])
pylab.plot(F_x, np.exp(-n_) * n_**F_x / factorial(F_x))
qo.animate(len(rho), fplot)

def qplot(fig, step):
    axes = fig.add_subplot(111)
    axes.clear()
    axes.imshow(Q[step], interpolation='bilinear', origin='lower',
                extent=(x_min, x_max, y_min, y_max))

qo.animate(len(rho), qplot)
```

## 2.3.2 Output



1. *Rabi oscillation*
2. *Jaynes-Cummings model*
3. *Single atom laser*

# INDICES AND TABLES

- *genindex*
- *modindex*
- *search*