example-rabi-oscillations

June 25, 2014

1 QuTiP example: Vacuum Rabi oscillations in the Jaynes-Cummings model

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This ipython notebook demonstrates how to simulate the quantum vacuum rabi oscillations in the Jaynes-Cumming model, using QuTiP: The Quantum Toolbox in Python.

For more information about QuTiP see project web page: http://code.google.com/p/qutip/

Populating the interactive namespace from numpy and matplotlib

2 Introduction

The Jaynes-Cumming model is the simplest possible model of quantum mechanical light-matter interaction, describing a single two-level atom interacting with a single electromagnetic cavity mode. The Hamiltonian for this system is (in dipole interaction form)

```
H = \hbar \omega_c a^{\dagger} a + \frac{1}{2} \hbar \omega_a \sigma_z + \hbar g(a^{\dagger} + a)(\sigma_- + \sigma_+) or with the rotating-wave approximation H_{\text{RWA}} = \hbar \omega_c a^{\dagger} a + \frac{1}{2} \hbar \omega_a \sigma_z + \hbar g(a^{\dagger} \sigma_- + a \sigma_+) where \omega_c and \omega_a are the frequencies of the cavity and atom, respectively, and g is the interaction strength.
```

2.0.1 Problem parameters

Here we use units where $\hbar = 1$:

```
In [2]: wc = 1.0 * 2 * pi # cavity frequency
    wa = 1.0 * 2 * pi # atom frequency
    g = 0.05 * 2 * pi # coupling strength
    kappa = 0.005 # cavity dissipation rate
    gamma = 0.05 # atom dissipation rate
    N = 15 # number of cavity fock states
    n_th_a = 0.0 # temperature in frequency units
    use_rwa = True

tlist = linspace(0,25,100)
```

2.0.2 Setup the operators, the Hamiltonian and initial state

```
In [3]: # intial state
    psi0 = tensor(basis(N,0), basis(2,1)) # start with an excited atom

# operators
a = tensor(destroy(N), qeye(2))
sm = tensor(qeye(N), destroy(2))

# Hamiltonian
if use_rwa:
    H = wc * a.dag() * a + wa * sm.dag() * sm + g * (a.dag() * sm + a * sm.dag())
else:
    H = wc * a.dag() * a + wa * sm.dag() * sm + g * (a.dag() + a) * (sm + sm.dag())
```

2.0.3 Create a list of collapse operators that describe the dissipation

```
In [4]: c_op_list = []

    rate = kappa * (1 + n_th_a)
    if rate > 0.0:
        c_op_list.append(sqrt(rate) * a)

    rate = kappa * n_th_a
    if rate > 0.0:
        c_op_list.append(sqrt(rate) * a.dag())

    rate = gamma
    if rate > 0.0:
        c_op_list.append(sqrt(rate) * sm)
```

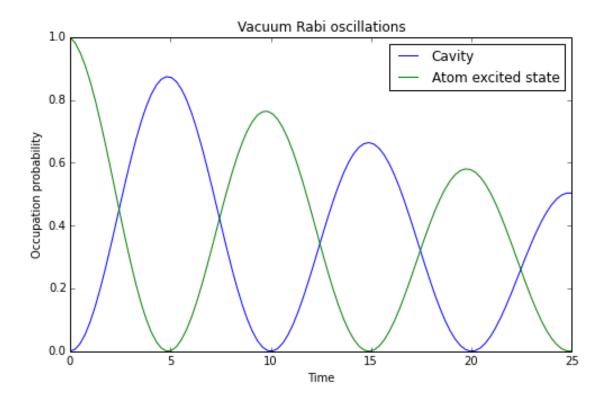
2.0.4 Evolve the system

Here we evolve the system with the Lindblad master equation solver, and we request that the expectation values of the operators $a^{\dagger}a$ and $\sigma_{+}\sigma_{-}$ are returned by the solver by passing the list [a.dag()*a, sm.dag()*sm] as the fifth argument to the solver.

```
In [5]: output = mesolve(H, psi0, tlist, c_op_list, [a.dag() * a, sm.dag() * sm])
```

2.1 Visualize the results

Here we plot the excitation probabilities of the cavity and the atom (these expectation values were calculated by the mesolve above). We can clearly see how energy is being coherently transferred back and forth between the cavity and the atom.



2.1.1 Software version:

Out[7]: <IPython.core.display.HTML at 0x7f2358c26ba8>