# **QuTiP lecture: The Dicke model**

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Latest version of this ipython notebook lecture is available at: http://github.com/jrjohansson/qutip-lectures

### Introduction

The Dicke Hamiltonian consists of a cavity mode and N spin-1/2 coupled to the cavity:

$$\begin{split} H_D &= \omega_0 \sum_{i=1}^{N} \sigma_z^{(i)} + \omega a^{\dagger} a + \sum_{i}^{N} \frac{\lambda}{\sqrt{N}} \left( a + a^{\dagger} \right) \left( \sigma_{+}^{(i)} + \sigma_{-}^{(i)} \right) \\ H_D &= \omega_0 J_z + \omega a^{\dagger} a + \frac{\lambda}{\sqrt{N}} \left( a + a^{\dagger} \right) (J_+ + J_-) \end{split}$$

where  $J_z$  and  $J_\pm$  are the collective angular momentum operators fpr a pseudospin of length j=N/2:

$$J_z = \sum_{i=1}^{N} \sigma_z^{(i)}$$
$$J_{\pm} = \sum_{i=1}^{N} \sigma_{\pm}^{(i)}$$

### References

- R.H. Dicke, Phys. Rev. 93, 99-110 (1954)
- Lambert et al., Phys. Rev. B 80, 165308 (2009)
- Lambert et al., Phys. Rev. Lett. 92, 073602 (2004)

## Setup problem in QuTiP

```
In [52]: M = 20
N = 8
j = N/2.0
n = 2*j + 1

a = tensor(destroy(M), qeye(n))
Jp = tensor(qeye(M), jmat(j, '+'))
Jm = tensor(qeye(M), jmat(j, '-'))
Jz = tensor(qeye(M), jmat(j, 'z'))

H0 = w * a.dag() * a + w0 * Jz
H1 = 1.0 / sqrt(N) * (a + a.dag()) * (Jp + Jm)
H = H0 + g * H1
```

Out [52]: Quantum object: dims = [[20, 9], [20, 9]], shape = [180, 180], type = oper, isHerm = True

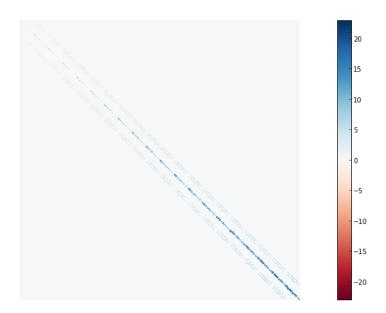
```
0.0 0.0 0.0 0.0 ...
                             0.0
                                   0.0
                                         0.0
                                               0.0
                                                     0.0
    3.0 0.0
             0.0 0.0 ...
                                   0.0
                                         0.0
                                               0.0
                                                     0.0
0.0
    0.0 2.0
              0.0
                   0.0 ...
                             0.0
                                   0.0
                                         0.0
                                               0.0
                                                     0.0
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    0.0 0.0
              1.0
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                                         ÷
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                             19.0
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    0.0
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              0.0
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                                                     0.0
                                               0.0
0.0
    0.0 0.0
              0.0
                   0.0 ...
                             0.0
                                   18.0
                                         0.0
                                                     0.0
0.0
    0.0 0.0
             0.0 0.0 ...
                             0.0
                                   0.0
                                        17.0
                                               0.0
                                                     0.0
```

```
 \begin{bmatrix} 0.0 & 0.0 & 0.0 & 0.0 & 0.0 & \cdots & 0.0 & 0.0 & 0.0 & 16.0 & 0.0 \\ 0.0 & 0.0 & 0.0 & 0.0 & \cdots & 0.0 & 0.0 & 0.0 & 0.0 & 15.0 \end{bmatrix}
```

#### Structure of the Hamiltonian

```
In [53]: fig, ax = subplots(1, 1, figsize=(10,10))
hinton(H, ax=ax)

Out [53]: <matplotlib.axes.AxesSubplot at 0x2be01b90>
```



# Find the ground state as a function of cavity-spin interaction strength

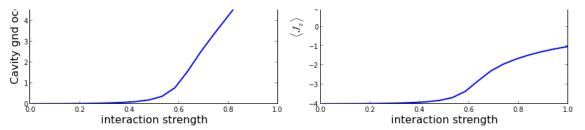
```
In [54]: g_vec = linspace(0.01, 1.0, 20)
# Ground state and steady state for the Hamiltonian: H = H0 + g * H1
psi_gnd_list = [(H0 + g * H1).groundstate()[1] for g in g_vec]
```

### Cavity ground state occupation probability

```
In [55]: n_gnd_vec = expect(a.dag() * a, psi_gnd_list)
Jz_gnd_vec = expect(Jz, psi_gnd_list)

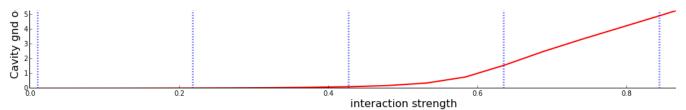
In [56]: fig, axes = subplots(1, 2, sharex=True, figsize=(12,4))
    axes[0].plot(g_vec, n_gnd_vec, 'b', linewidth=2, label="cavity occupation")
    axes[0].set_ylim(0, max(n_gnd_vec))
    axes[0].set_ylabel("Cavity gnd occ. prob.", fontsize=16)
    axes[0].set_xlabel("interaction strength", fontsize=16)

    axes[1].plot(g_vec, Jz_gnd_vec, 'b', linewidth=2, label="cavity occupation")
    axes[1].set_ylim(-j, j)
    axes[1].set_ylabel(r"s\langle J_z\rangle\sigma", fontsize=16)
    axes[1].set_xlabel("interaction strength", fontsize=16)
    fig.tight_layout()
```



### Cavity Wigner function and Fock distribution as a function of coupling strength

```
In [57]: psi_gnd_sublist = psi_gnd_list[::4]
          xvec = linspace(-7,7,200)
          fig_grid = (3, len(psi_gnd_sublist))
          fig = figure(figsize=(4*len(psi_gnd_sublist),12))
          for idx, psi_gnd in enumerate(psi_gnd_sublist):
               # trace out the cavity density matrix
               rho\_gnd\_cavity = ptrace(psi\_gnd, \theta)
               # calculate its wigner function
               W = wigner(rho_gnd_cavity, xvec, xvec)
               # plot its wigner function
               ax = subplot2grid(fig_grid, (0, idx))
               ax.contourf(xvec, xvec, W, 100)
               # plot its fock-state distribution
               ax = subplot2grid(fig_grid, (1, idx))
               ax.bar(arange(0, M), real(rho_gnd_cavity.diag()), color="blue", alpha=0.6)
               ax.set_ylim(0, 1)
               ax.set_xlim(0, M)
          # plot the cavity occupation probability in the ground state
          ax = subplot2grid(fig_grid, (2, 0), colspan=fig_grid[1])
          ax.plot(g_vec, n_gnd_vec, 'r', linewidth=2, label="cavity occupation")
          ax.set_xlim(0, max(g_vec))
ax.set_ylim(0, max(n_gnd_vec)*1.2)
          ax.set_ylabel("Cavity gnd occ. prob.", fontsize=16)
ax.set_xlabel("interaction strength", fontsize=16)
          for g in g_vec[::4]:
               \overline{ax.plot([g,g],[0,max(n_gnd_vec)*1.2], b:', linewidth=2.5)}
             1.0
                                              1.0
                                                                               1.0
                                                                                                                 1.0
                                                                                                                                                  1.0
             0.8
                                              0.8
                                                                               0.8
                                                                                                                0.8
                                                                                                                                                  0.8
             0.6
                                              0.6
                                                                               0.6
                                                                                                                0.6
                                                                                                                                                  0.6
             0.4
                                              0.4
                                                                               0.4
                                                                                                                0.4
                                                                                                                                                  0.4
             0.2
                                               0.2
                                                                               0.2
                                                                                                                 0.2
                                                                                                                                                  0.2
                                                                               0.0
                                              0.0
                                                                                                                0.0
             0.0
                                                                                                                                                  0.0
                             10
                                    15
                                                                     15
                                                                                                      15
                                                                                                                                       15
                                                                                                                                              20
```



### Entropy/Entanglement between spins and cavity

# Entropy as a function interaction strength for increasing N

0.2

See Lambert et al., Phys. Rev. Lett. 92, 073602 (2004).

0.2

0.0

interaction strength

0.8

```
entropy_cavity = zeros(shape(g_vec))
entropy_spin = zeros(shape(g_vec))

for idx, psi_gnd in enumerate(psi_gnd_list):
    rho_gnd_cavity = ptrace(psi_gnd, 0)
    rho_gnd_spin = ptrace(psi_gnd, 1)

    entropy_cavity[idx] = entropy_vn(rho_gnd_cavity, 2)
    entropy_spin[idx] = entropy_vn(rho_gnd_spin, 2)

return entropy_cavity, entropy_spin
```

```
In [61]: g_vec = linspace(0.2, 0.8, 60)
N_vec = [4, 8, 12, 16, 24, 32]
MM = 25

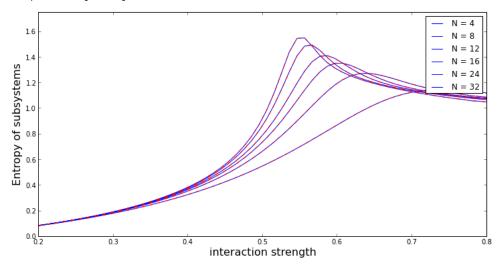
fig, axes = subplots(1, 1, figsize=(12,6))

for NN in N_vec:
    entropy_cavity, entropy_spin = calulcate_entropy(MM, NN, g_vec)

    axes.plot(g_vec, entropy_cavity, 'b', label="N = %d" % NN)
    axes.plot(g_vec, entropy_spin, 'r--')

axes.set_ylim(0, 1.75)
axes.set_ylabel("Entropy of subsystems", fontsize=16)
axes.set_xlabel("interaction strength", fontsize=16)
axes.legend()
```

Out [61]: <matplotlib.legend.Legend at 0x4ec67790>



# Dissipative cavity: steady state instead of the ground state

```
In [62]: # average number thermal photons in the bath coupling to the resonator
n_th = 0.25

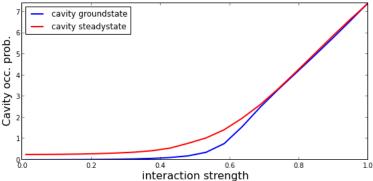
c_ops = [sqrt(kappa * (n_th + 1)) * a, sqrt(kappa * n_th) * a.dag()]
#c_ops = [sqrt(kappa) * a, sqrt(gamma) * Jm]
```

### Find the ground state as a function of cavity-spin interaction strength

```
In [63]: g_vec = linspace(0.01, 1.0, 20)
# Ground state for the Hamiltonian: H = H0 + g * H1
rho_ss_list = [steadystate(H0 + g * H1, c_ops) for g in g_vec]
```

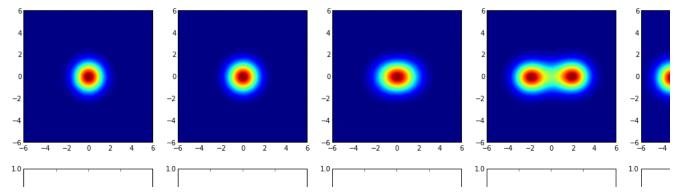
### Cavity ground state occupation probability

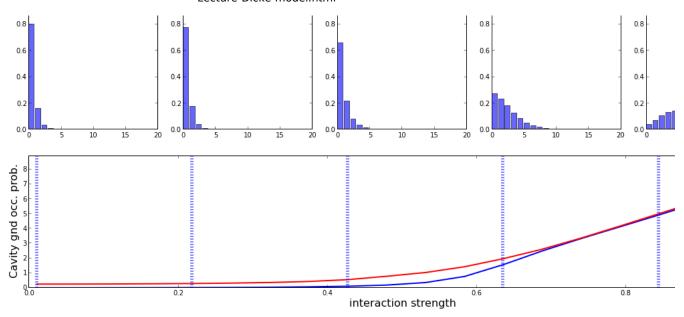
```
In [64]: # calculate the expectation value of the number of photons in the cavity
    n_ss_vec = expect(a.dag() * a, rho_ss_list)
```



### Cavity Wigner function and Fock distribution as a function of coupling strength

```
In [66]: rho_ss_sublist = rho_ss_list[::4]
            xvec = linspace(-6,6,200)
            fig_grid = (3, len(rho_ss_sublist))
            fig = figure(figsize=(4*len(rho_ss_sublist),12))
            for idx, rho_ss in enumerate(rho_ss_sublist):
                 # trace out the cavity density matrix
                 rho_ss_cavity = ptrace(rho_ss, 0)
                 # calculate its wigner function
                 W = wigner(rho_ss_cavity, xvec, xvec)
                 # plot its wigner function
                 ax = subplot2grid(fig_grid, (0, idx))
                 ax.contourf(xvec, xvec, W, 100)
                 # plot its fock-state distribution
                 ax = subplot2grid(fig_grid, (1, idx))
                 ax.bar(arange(0, M), real(rho_ss_cavity.diag()), color="blue", alpha=0.6)
                 ax.set_ylim(0, 1)
            # plot the cavity occupation probability in the ground state
ax = subplot2grid(fig_grid, (2, 0), colspan=fig_grid[1])
ax.plot(g_vec, n_gnd_vec, 'b', linewidth=2, label="cavity groundstate")
ax.plot(g_vec, n_ss_vec, 'r', linewidth=2, label="cavity steadystate")
            ax.set_xlim(0, max(g_vec))
            ax.set_ylim(0, max(n_ss_vec)*1.2)
            ax.set_ylabel("Cavity gnd occ. prob.", fontsize=16)
ax.set_xlabel("interaction strength", fontsize=16)
            for g in g_vec[::4]:
                 ax.plot([g,g],[0,max(n_ss_vec)*1.2], 'b:', linewidth=5)
```





# **Entropy**

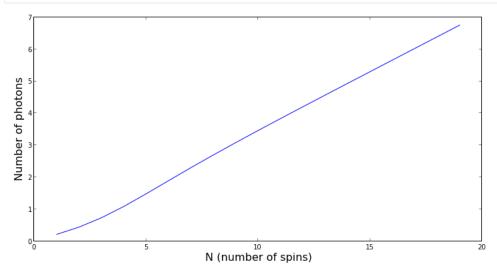
```
In [67]: entropy_tot
                                     = zeros(shape(g_vec))
               entropy_cavity = zeros(shape(g_vec))
               entropy_spin
                                     = zeros(shape(g_vec))
               for idx, rho_ss in enumerate(rho_ss_list):
                      rho_gnd_cavity = ptrace(rho_ss, 0)
                      rho_gnd_spin = ptrace(rho_ss, 1)
                      entropy_tot[idx]
                                                     = entropy_vn(rho_ss, 2)
                     entropy_cavity[idx] = entropy_vn(rho_gnd_cavity, 2)
entropy_spin[idx] = entropy_vn(rho_gnd_spin, 2)
In [68]: fig, axes = subplots(1, 1, figsize=(12,6))
    axes.plot(g_vec, entropy_tot, 'k', label="total")
    axes.plot(g_vec, entropy_cavity, 'b', label="cavity")
    axes.plot(g_vec, entropy_spin, 'r--', label="spin")
              #axes.set_ylim(0, 1.5)
axes.set_ylabel("Entropy of subsystems", fontsize=16)
axes.set_xlabel("interaction strength", fontsize=16)
               axes.legend(loc=0)
               fig.tight_layout()
                                                                                                                                                                              total
                                                                                                                                                                              cavity
                                                                                                                                                                              spin
                     2.5
                Entropy of subsystems
                     1.0
                    0.5 L
0.0
                                                       0.2
                                                                                                                                                     0.8
                                                                                                                                                                                     1.0
                                                                                        interaction strength
```

# Superradiance?

start with all atoms in the excited state

```
In [71]: N_vec = arange(1, 20, 1)
    MM = 25
    a

    fig, axes = subplots(1, 1, figsize=(12,6))
    n_cavity = [calulcate_intensity(MM, NN, 0.7) for NN in N_vec]
    axes.plot(N_vec, n_cavity, 'b')
    axes.set_ylabel("Number of photons", fontsize=16)
    axes.set_xlabel("N (number of spins)", fontsize=16)
    axes.legend()
```



```
\textbf{return} \ \texttt{res.expect[0]}
```

## Out [73]: [<matplotlib.lines.Line2D at 0x53937ad0>]

