# scratchpad1

February 2, 2019

# 1 Imports

```
In [71]: # # Imports
         # ## Qutip imports 1
         # In [17:
         from qutip.operators import sigmax, sigmay, sigmaz, identity, qeye
         from qutip.operators import position, momentum, num, create, destroy
         from qutip.operators import commutator, qdiags
         from qutip.simdiag import simdiag
         from qutip.states import basis, ket2dm
         from qutip.tensor import tensor
         from qutip.qip.gates import swap, rx, ry, rz, cnot
         from qutip.qobj import Qobj
         from qutip.visualization import hinton
         from qutip.visualization import matrix_histogram_complex, matrix_histogram
         from qutip.random_objects import rand_herm, rand_unitary, rand_dm
         # ## Qutip imports 2
         # In[3]:
         #from
         # ## Numpy imports
         # In[2]:
         from numpy import sin, cos, tan, real, imag, log, conj
         from numpy import array, append, linspace, arange
```

```
from numpy import add, sqrt, abs, dot
from numpy.random import random, random_sample, rand, seed, RandomState
from numpy import concatenate, hstack, vstack, block, dstack, vsplit
from numpy import trace, diag
from numpy import ones, zeros, ones like, zeros like
from numpy import amax, amin, nanmax, nanmin
from numpy import outer, multiply
# from numpy import pi
# ## Scipy imports
# In[17]:
from scipy.integrate import ode, odeint, complex_ode
from scipy.optimize import minimize
from scipy.linalg import eigh, inv, norm, expm
# from scipy.linalg import
# from scipy import
# ## Matplotlib imports
# In[18]:
from matplotlib.pyplot import plot, figure, show, savefig, axes
from matplotlib.pyplot import xlabel, ylabel, title, legend
from matplotlib import rcParams
from matplotlib.pyplot import style
from matplotlib.pyplot import xlim, ylim, axis
# beware not same as axes
from matplotlib.pyplot import subplot, subplots, text
from matplotlib.pyplot import GridSpec
from matplotlib.pyplot import scatter, colorbar
pgf_with_rc_fonts = {"pgf.texsystem": "pdflatex"}
rcParams.update(pgf_with_rc_fonts)
style.use('seaborn-whitegrid')
# ## Math imports
# In[19]:
```

```
from math import pi
from math import exp
# ## Cmath imports
# ## Date and datetime imports
# In[20]:
from datetime import date
from datetime import datetime# now
# ## Os imports
# In[21]:
from os import getcwd, mkdir, chdir
from os.path import abspath, join
# ## Sympy imports
# In[22]:
from sympy import Function, dsolve, Eq, Derivative, symbols
\# x, y, z, t = symbols('x y z t')
\# k, m, n = symbols('k m n', integer=True)
# f, g, h = symbols('f g h', cls=Function)
# ## Miscellaneous imports
# ## Extra useful functions
# In[23]:
def rint(x):
   print("x = ", x)
    return None
# # Next chapter
```

```
# ## sub topic 1
# ## sub topic 2
# ## sub topic 3
# ### sub sub topic 1
```

# 2 Next chapter

## 2.1 memory clear (uses regex, so be careful)

In [30]: %reset\_selective -f var1, var2 # replace var1, var2 with your defined ones

### 2.2 sub topic 2

In [31]: cnot()

### Out[31]:

Quantum object: dims = [[2, 2], [2, 2]], shape = (4, 4), type = oper, isherm = True

$$\left(\begin{array}{ccccc}
1.0 & 0.0 & 0.0 & 0.0 \\
0.0 & 1.0 & 0.0 & 0.0 \\
0.0 & 0.0 & 0.0 & 1.0 \\
0.0 & 0.0 & 1.0 & 0.0
\end{array}\right)$$

In [32]: (qeye(2) - sigmaz())/2

#### Out[32]:

Quantum object: dims = [[2], [2]], shape = (2, 2), type = oper, isherm = True

$$\left(\begin{array}{cc} 0.0 & 0.0 \\ 0.0 & 1.0 \end{array}\right)$$

In [33]: (-qeye(2) + sigmax())

### Out[33]:

Quantum object: dims = [[2], [2]], shape = (2, 2), type = oper, isherm = True

$$\left(\begin{array}{cc} -1.0 & 1.0 \\ 1.0 & -1.0 \end{array}\right)$$

In [34]: tensor((qeye(2) - sigmaz())/2,(-qeye(2) + sigmax())

#### Out[34]:

Quantum object: dims = [[2, 2], [2, 2]], shape = (4, 4), type = oper, isherm = True

$$\begin{pmatrix}
0.0 & 0.0 & 0.0 & 0.0 \\
0.0 & 0.0 & 0.0 & 0.0 \\
0.0 & 0.0 & -1.0 & 1.0 \\
0.0 & 0.0 & 1.0 & -1.0
\end{pmatrix}$$

In [35]: (1j\*(pi/4)\*tensor((qeye(2)-sigmaz())/2,(-qeye(2)+sigmax())).expm()

### Out[35]:

Quantum object: dims = [[2, 2], [2, 2]], shape = (4, 4), type = oper, isherm = False

$$\begin{pmatrix} 1.0 & 0.0 & 0.0 & 0.0 \\ 0.0 & 1.0 & 0.0 & 0.0 \\ 0.0 & 0.0 & (0.500 - 0.500j) & (0.500 + 0.500j) \\ 0.0 & 0.0 & (0.500 + 0.500j) & (0.500 - 0.500j) \end{pmatrix}$$

In [10]: (1\*(pi/4)\*tensor((qeye(2)-sigmaz())/2,(-qeye(2)+sigmax())).expm()

### Out[10]:

Quantum object: dims = [[2, 2], [2, 2]], shape = (4, 4), type = oper, isherm = True

$$\begin{pmatrix}
1.0 & 0.0 & 0.0 & 0.0 \\
0.0 & 1.0 & 0.0 & 0.0 \\
0.0 & 0.0 & 0.604 & 0.396 \\
0.0 & 0.0 & 0.396 & 0.604
\end{pmatrix}$$

In [12]: (1j\*(pi/4)\*(1j)\*tensor((qeye(2)-sigmaz())/2,(-qeye(2)+sigmax()))).expm()

### Out[12]:

Quantum object: dims = [[2, 2], [2, 2]], shape = (4, 4), type = oper, isherm = True

$$\begin{pmatrix}
1.0 & 0.0 & 0.0 & 0.0 \\
0.0 & 1.0 & 0.0 & 0.0 \\
0.0 & 0.0 & 2.905 & -1.905 \\
0.0 & 0.0 & -1.905 & 2.905
\end{pmatrix}$$

### 2.3 sub topic 3

### 2.3.1 sub sub topic 1

In [13]: cnot()

### Out[13]:

Quantum object: dims = [[2, 2], [2, 2]], shape = (4, 4), type = oper, isherm = True

$$\left(\begin{array}{ccccc}
1.0 & 0.0 & 0.0 & 0.0 \\
0.0 & 1.0 & 0.0 & 0.0 \\
0.0 & 0.0 & 0.0 & 1.0 \\
0.0 & 0.0 & 1.0 & 0.0
\end{array}\right)$$

In [14]: cnot().eigenenergies()

Out[14]: array([-1., 1., 1., 1.])

In [15]: cnot().eigenstates()

```
Out[15]: (array([-1., 1., 1., 1.]),
          array([Quantum object: dims = [[2, 2], [1, 1]], shape = (4, 1), type = ket
          Qobj data =
          [[ 0.
           ΓО.
                       ]
           [ 0.70710678]
           [-0.70710678],
                 Quantum object: dims = [[2, 2], [1, 1]], shape = (4, 1), type = ket
          Qobj data =
          [[0.]
           [1.]
           [0.]
           [0.]],
                 Quantum object: dims = [[2, 2], [1, 1]], shape = (4, 1), type = ket
          Qobj data =
          [[1.]]
           [0.]
           [0.]
           [0.]],
                 Quantum object: dims = [[2, 2], [1, 1]], shape = (4, 1), type = ket
          Qobj data =
          [[0.
                      ]
           ΓΟ.
                      1
           [0.70710678]
           [0.70710678]]], dtype=object))
In [16]: cnot().diag()
Out[16]: array([1., 1., 0., 0.])
In [20]: c_eigenvalue, c_eigenvec_arr = cnot().eigenstates()
In [23]: c_eigenvec
Out[23]: array([Quantum object: dims = [[2, 2], [1, 1]], shape = (4, 1), type = ket
         Qobj data =
         [[ 0.
                      ]
                      ]
          [ 0.
          [ 0.70710678]
          [-0.70710678],
                Quantum object: dims = [[2, 2], [1, 1]], shape = (4, 1), type = ket
         Qobj data =
         [[0.]
          [1.]
          [0.]
          [0.]],
                Quantum object: dims = [[2, 2], [1, 1]], shape = (4, 1), type = ket
         Qobj data =
         [[1.]]
```

```
[0.]
          [0.]
          [0.]],
                Quantum object: dims = [[2, 2], [1, 1]], shape = (4, 1), type = ket
         Qobj data =
         [[0.
          [0.
          [0.70710678]
          [0.70710678]]], dtype=object)
In [24]: c_eigenvalue
Out[24]: array([-1., 1., 1., 1.])
In [27]: qdiags(2)
       TypeError
                                                  Traceback (most recent call last)
        <ipython-input-27-61a25a066823> in <module>()
   ----> 1 qdiags(2)
        TypeError: qdiags() missing 1 required positional argument: 'offsets'
In [37]: simdiag(array([sigmax(), qeye(2)]))
       AttributeError
                                                  Traceback (most recent call last)
        <ipython-input-37-54d465f50d4b> in <module>()
   ---> 1 simdiag(array([sigmax(), qeye(2)]))
        /anaconda3/envs/qutip-env/lib/python3.6/site-packages/qutip/simdiag.py in simdiag(ops,
         82
                eigvals, eigvecs = la.eig(A.full())
         83
                zipped = zip(-eigvals, range(len(eigvals)))
                zipped.sort()
    ---> 84
         85
               ds, perm = zip(*zipped)
                ds = -np.real(np.array(ds))
         86
        AttributeError: 'zip' object has no attribute 'sort'
```

```
In [26]: d = qdiags(c_eigenvalue)
                                                                  Traceback (most recent call last)
          TypeError
          <ipython-input-26-c2334747d78d> in <module>()
     ----> 1 d = qdiags(c_eigenvalue)
             2 d
          TypeError: qdiags() missing 1 required positional argument: 'offsets'
In [38]: Qobj(diag(c_eigenvalue.full()))
                                                                  Traceback (most recent call last)
          AttributeError
          <ipython-input-38-66948a1e6342> in <module>()
     ---> 1 Qobj(diag(c_eigenvalue.full()))
          AttributeError: 'numpy.ndarray' object has no attribute 'full'
In [40]: ceq = Qobj(diag(c_eigenvalue))
   Out[40]:
   Quantum object: dims = [[4], [4]], shape = (4, 4), type = oper, isherm = True
                                       \left(\begin{array}{ccccc} -1.000 & 0.0 & 0.0 & 0.0 \\ 0.0 & 1.0 & 0.0 & 0.0 \\ 0.0 & 0.0 & 1.0 & 0.0 \\ 0.0 & 0.0 & 0.0 & 1.0 \end{array}\right)
In [42]: Qobj(arange(9).reshape((3,3)))
   Out [42]:
   Quantum object: dims = [[3], [3]], shape = (3, 3), type = oper, isherm = False

\left(\begin{array}{cccc}
0.0 & 1.0 & 2.0 \\
3.0 & 4.0 & 5.0 \\
6.0 & 7.0 & 8.0
\end{array}\right)
```

```
In [44]: inter = [a.full() for a in c_eigenvec]
In [45]: inter
Out[45]: [array([[ 0.
                              +0.j],
                              +0.j],
                 [ 0.
                 [ 0.70710678+0.j],
                 [-0.70710678+0.j]]), array([[0.+0.j],
                 [1.+0.j],
                 [0.+0.j],
                 [0.+0.j]]), array([[1.+0.j],
                 [0.+0.j],
                 [0.+0.j],
                 [0.+0.j]]), array([[0.
                                              +0.j],
                            +0.j],
                 [0.70710678+0.j],
                 [0.70710678+0.j]])]
In [46]: array(c_eigenvec)
Out [46]: array([Quantum object: dims = [[2, 2], [1, 1]], shape = (4, 1), type = ket
         Qobj data =
         [[ 0.
                      ]
          [ 0.
                      ]
          [ 0.70710678]
          [-0.70710678],
                Quantum object: dims = [[2, 2], [1, 1]], shape = (4, 1), type = ket
         Qobj data =
         [[0.]
          [1.]
          [0.]
          [0.]],
                Quantum object: dims = [[2, 2], [1, 1]], shape = (4, 1), type = ket
         Qobj data =
         [[1.]]
          [0.]
          [0.]
          [0.]],
                Quantum object: dims = [[2, 2], [1, 1]], shape = (4, 1), type = ket
         Qobj data =
                     ]
         [[0.
          [0.
                     ]
          [0.70710678]
          [0.70710678]]], dtype=object)
In [47]: array(inter)
Out[47]: array([[[ 0.
                              +0.j],
                 [ 0.
                             +0.j],
```

```
[0.70710678+0.j],
                 [-0.70710678+0.j],
                [[ 0.
                            +0.j],
                 [ 1.
                           +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],
                 [ 0.70710678+0.j],
                 [ 0.70710678+0.j]])
In [48]: crq = Qobj(array(inter))
         crq
       TypeError
                                                  Traceback (most recent call last)
        <ipython-input-48-fb8a58b562e3> in <module>()
   ----> 1 crq = Qobj(array(inter))
          2 crq
        /anaconda3/envs/qutip-env/lib/python3.6/site-packages/qutip/qobj.py in __init__(self,
        277
                       do\_copy = copy
        278
                       if not isinstance(inpt, fast_csr_matrix):
    --> 279
                            _tmp = sp.csr_matrix(inpt, dtype=complex, copy=do_copy)
        280
                            _tmp.sort_indices() #Make sure indices are sorted.
                            do_{copy} = 0
        281
        /anaconda3/envs/qutip-env/lib/python3.6/site-packages/scipy/sparse/compressed.py in ___
        77
                                    self.format)
        78
                        from .coo import coo_matrix
    ---> 79
                        self._set_self(self.__class__(coo_matrix(arg1, dtype=dtype)))
        80
        81
                    # Read matrix dimensions given, if any
        /anaconda3/envs/qutip-env/lib/python3.6/site-packages/scipy/sparse/coo.py in __init__(
```

```
179
180
if M.ndim != 2:
--> 181
raise TypeError('expected dimension <= 2 array or matrix')
182
else:
self._shape = check_shape(M.shape)
```

TypeError: expected dimension <= 2 array or matrix</pre>

In [56]: crq

#### Out [56]:

Quantum object: dims = [[4], [4]], shape = (4, 4), type = oper, isherm = False

$$\begin{pmatrix}
0.0 & 0.0 & 1.0 & 0.0 \\
0.0 & 1.0 & 0.0 & 0.0 \\
0.707 & 0.0 & 0.0 & 0.707 \\
-0.707 & 0.0 & 0.0 & 0.707
\end{pmatrix}$$

In [57]: ceq

#### Out [57]:

Quantum object: dims = [[4], [4]], shape = (4, 4), type = oper, isherm = True

$$\begin{pmatrix}
-1.000 & 0.0 & 0.0 & 0.0 \\
0.0 & 1.0 & 0.0 & 0.0 \\
0.0 & 0.0 & 1.0 & 0.0 \\
0.0 & 0.0 & 0.0 & 1.0
\end{pmatrix}$$

In [58]: crq\*ceq\*(crq.dag())

### Out [58]:

Quantum object: dims = [[4], [4]], shape = (4, 4), type = oper, isherm = True

$$\begin{pmatrix}
1.0 & 0.0 & 0.0 & 0.0 \\
0.0 & 1.0 & 0.0 & 0.0 \\
0.0 & 0.0 & 0.0 & 1.000 \\
0.0 & 0.0 & 1.000 & 0.0
\end{pmatrix}$$

In [60]: cnot()

### Out[60]:

Quantum object: dims = [[2, 2], [2, 2]], shape = (4, 4), type = oper, isherm = True

$$\left(\begin{array}{ccccc}
1.0 & 0.0 & 0.0 & 0.0 \\
0.0 & 1.0 & 0.0 & 0.0 \\
0.0 & 0.0 & 0.0 & 1.0 \\
0.0 & 0.0 & 1.0 & 0.0
\end{array}\right)$$

In [59]: (crq.dag())\*ceq\*crq

### Out[59]:

Quantum object: dims = [[4], [4]], shape = (4, 4), type = oper, isherm = True

$$\begin{pmatrix}
1.0 & 0.0 & 0.0 & 0.0 \\
0.0 & 1.0 & 0.0 & 0.0 \\
0.0 & 0.0 & -1.000 & 0.0 \\
0.0 & 0.0 & 0.0 & 1.0
\end{pmatrix}$$

In [62]: = tensor(sigmax(), sigmax()) + tensor(sigmay(), sigmay())

### Out[62]:

Quantum object: dims = [[2, 2], [2, 2]], shape = (4, 4), type = oper, isherm = True

$$\left(\begin{array}{ccccc}
0.0 & 0.0 & 0.0 & 0.0 \\
0.0 & 0.0 & 2.0 & 0.0 \\
0.0 & 2.0 & 0.0 & 0.0 \\
0.0 & 0.0 & 0.0 & 0.0
\end{array}\right)$$

In [63]: (-1j\*(1/8)\*).expm()

### Out[63]:

Quantum object: dims = [[2, 2], [2, 2]], shape = (4, 4), type = oper, isherm = False

$$\begin{pmatrix}
1.0 & 0.0 & 0.0 & 0.0 \\
0.0 & 0.969 & -0.247j & 0.0 \\
0.0 & -0.247j & 0.969 & 0.0 \\
0.0 & 0.0 & 0.0 & 1.0
\end{pmatrix}$$

In [64]: (1j\*(1/8)\*).expm()

#### Out [64]:

Quantum object: dims = [[2, 2], [2, 2]], shape = (4, 4), type = oper, isherm = False

$$\begin{pmatrix}
1.0 & 0.0 & 0.0 & 0.0 \\
0.0 & 0.969 & 0.247j & 0.0 \\
0.0 & 0.247j & 0.969 & 0.0 \\
0.0 & 0.0 & 0.0 & 1.0
\end{pmatrix}$$

In [65]: (1j\*(1/4)\*).expm()

### Out [65]:

Quantum object: dims = [[2, 2], [2, 2]], shape = (4, 4), type = oper, isherm = False

$$\begin{pmatrix}
1.0 & 0.0 & 0.0 & 0.0 \\
0.0 & 0.878 & 0.479j & 0.0 \\
0.0 & 0.479j & 0.878 & 0.0 \\
0.0 & 0.0 & 0.0 & 1.0
\end{pmatrix}$$

In [66]: (1j\*(pi/4)\*).expm()

#### Out [66]:

Quantum object: dims = [[2, 2], [2, 2]], shape = (4, 4), type = oper, isherm = False

$$\left(\begin{array}{ccccc}
1.0 & 0.0 & 0.0 & 0.0 \\
0.0 & 0.0 & 1.0j & 0.0 \\
0.0 & 1.0j & 0.0 & 0.0 \\
0.0 & 0.0 & 0.0 & 1.0
\end{array}\right)$$

In [67]: (-1j\*(pi/8)\*).expm()

#### Out [67]:

Quantum object: dims = [[2, 2], [2, 2]], shape = (4, 4), type = oper, isherm = False

$$\begin{pmatrix}
1.0 & 0.0 & 0.0 & 0.0 \\
0.0 & 0.707 & -0.707j & 0.0 \\
0.0 & -0.707j & 0.707 & 0.0 \\
0.0 & 0.0 & 0.0 & 1.0
\end{pmatrix}$$

In [69]: = tensor(sigmax(), sigmax())

### Out[69]:

Quantum object: dims = [[2, 2], [2, 2]], shape = (4, 4), type = oper, isherm = True

$$\left(\begin{array}{ccccc}
0.0 & 0.0 & 0.0 & 1.0 \\
0.0 & 0.0 & 1.0 & 0.0 \\
0.0 & 1.0 & 0.0 & 0.0 \\
1.0 & 0.0 & 0.0 & 0.0
\end{array}\right)$$

In [70]: (1j\*(pi/4)\*).expm()

### Out[70]:

Quantum object: dims = [[2, 2], [2, 2]], shape = (4, 4), type = oper, isherm = False

$$\begin{pmatrix}
0.707 & 0.0 & 0.0 & 0.707j \\
0.0 & 0.707 & 0.707j & 0.0 \\
0.0 & 0.707j & 0.707 & 0.0 \\
0.707j & 0.0 & 0.0 & 0.707
\end{pmatrix}$$

#### Out [75]:

Quantum object: dims = [[2], [1]], shape = (2, 1), type = ket

$$\begin{pmatrix} 1.0\\ 0.0 \end{pmatrix}$$

In [76]: zero = basis(2,1)
 zero

### Out [76]:

Quantum object: dims = [[2], [1]], shape = (2, 1), type = ket

$$\begin{pmatrix} 0.0\\ 1.0 \end{pmatrix}$$

In [78]: aas = (1/sqrt(2))\*(tensor(one, zero) + 1j\*tensor(zero, one))

In [79]: aas

#### Out [79]:

Quantum object: dims = [[2, 2], [1, 1]], shape = (4, 1), type = ket

$$\begin{pmatrix}
0.0 \\
0.707 \\
0.707j \\
0.0
\end{pmatrix}$$

In [80]: ket2dm(aas)

#### Out[80]:

Quantum object: dims = [[2, 2], [2, 2]], shape = (4, 4), type = oper, isherm = True

$$\begin{pmatrix}
0.0 & 0.0 & 0.0 & 0.0 \\
0.0 & 0.500 & -0.500j & 0.0 \\
0.0 & 0.500j & 0.500 & 0.0 \\
0.0 & 0.0 & 0.0 & 0.0
\end{pmatrix}$$

In [70]: (1j\*(pi/4)\*).expm()\*

### Out[70]:

Quantum object: dims = [[2, 2], [2, 2]], shape = (4, 4), type = oper, isherm = False

$$\begin{pmatrix}
0.707 & 0.0 & 0.0 & 0.707j \\
0.0 & 0.707 & 0.707j & 0.0 \\
0.0 & 0.707j & 0.707 & 0.0 \\
0.707j & 0.0 & 0.0 & 0.707
\end{pmatrix}$$

In []: 
$$\#g = pi/(4 * T) T = 2 * np.pi$$
  
 $\#H = g * (tensor(sigmax(), sigmax()) + tensor(sigmay(), sigmay()))$ 

In [84]: H = (1/(4 \* 2))\* (tensor(sigmax(), sigmax()) + tensor(sigmay(), sigmay()))

In [85]: (-1j\*H\*2\*pi).expm()

#### Out[85]:

Quantum object: dims = [[2, 2], [2, 2]], shape = (4, 4), type = oper, isherm = False

$$\left(\begin{array}{cccc}
1.0 & 0.0 & 0.0 & 0.0 \\
0.0 & 0.0 & -1.0j & 0.0 \\
0.0 & -1.0j & 0.0 & 0.0 \\
0.0 & 0.0 & 0.0 & 1.0
\end{array}\right)$$

### Out[86]:

Quantum object: dims = [[2, 2], [2, 2]], shape = (4, 4), type = oper, isherm = False

$$\left(\begin{array}{ccccc}
1.0 & 0.0 & 0.0 & 0.0 \\
0.0 & 0.0 & 1.0j & 0.0 \\
0.0 & 1.0j & 0.0 & 0.0 \\
0.0 & 0.0 & 0.0 & 1.0
\end{array}\right)$$

### Out[87]:

Quantum object: dims = [[2, 2], [2, 2]], shape = (4, 4), type = oper, isherm = True

$$\left(\begin{array}{ccccc}
1.0 & 0.0 & 0.0 & 0.0 \\
0.0 & 2.509 & 2.301 & 0.0 \\
0.0 & 2.301 & 2.509 & 0.0 \\
0.0 & 0.0 & 0.0 & 1.0
\end{array}\right)$$