

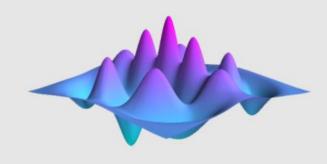


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QuTiP

Quantum Toolbox in Python

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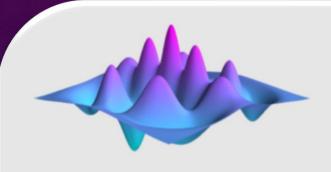
A Technical Staff position for QuTiP is available, check it out under the jobs page.

QuTiP is open-source software for simulating the dynamics of open quantum systems. The QuTiP library depends on the excellent Numpy, Scipy, and Cython numerical packages. In addition, graphical output is provided by Matplotlib. QuTiP aims to provide user-friendly and efficient numerical simulations of a wide variety of Hamiltonians, including those with arbitrary time-dependence, commonly found in a wide range of physics applications such as quantum optics, trapped ions, superconducting circuits, and quantum nanomechanical resonators. QuTiP is freely available for use and/or modification on all major platforms such as Linux, Mac OSX, and Windows*. Being free of any licensing fees, QuTiP is ideal for exploring quantum mechanics and dynamics in the classroom.

*QuTiP is developed on Unix platforms only, and some features may not be available under Windows.







QuTiP

Quantum Toolbox in Python

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Users guide

Latest release

Version 4.5



Online HTML documentation



PDF documentation



BASIC OPERATIONS ON QUANTUM OBJECTS

The key difference between classical and quantum mechanics

in computing the dynamics of quantum systems we need a data structure that is capable of encapsulating the properties of a quantum operator and ket/bra vectors. The quantum object class, qutip.Qobj, accomplishes this using matrix representation.



lies in the use of operators instead of numbers as variables.

```
In [3]: Qobj()
Out[3]:
Quantum object: dims = [[1], [1]], shape = [1, 1], type = oper, isherm = True
Qobi data =
[[ 0.]]
      In [4]: Qobj([[1],[2],[3],[4],[5]])
     Out[4]:
      Quantum object: dims = [[5], [1]], shape = [5, 1], type = ket
      Qobj data =
      [[ 1.]
      [ 2.]
      [ 3.]
      [ 4.]
      [5.]]
      In [5]: x = np.array([[1, 2, 3, 4, 5]])
      In [6]: Qobj(x)
      Out[6]:
      Quantum object: dims = [[1], [5]], shape = [1, 5], type = bra
      Qobj data =
```

[[1. 2. 3. 4. 5.]]





States	Command (# means optional)	Inputs
Fock state ket vector	basis(N,#m) / fock(N,#m)	N = number of levels in Hilbert space, m = level containing excitation (0 if no m given)
Fock density matrix (outer product of basis)	<pre>fock_dm(N, #p)</pre>	same as basis(N,m) / fock(N,m)
Coherent state	<pre>coherent(N,alpha)</pre>	alpha = complex number (eigenvalue) for requested coherent state
Coherent density matrix (outer product)	<pre>coherent_dm(N,alpha)</pre>	same as coherent(N,alpha)
Thermal density matrix (for n particles)	thermal_dm(N,n)	n = particle number expectation value

Operators	Command (# means optional)	Inputs
Identity	qeye(N)	N = number of levels in Hilbert space.
Lowering (destruction) operator	destroy(N)	same as above
Raising (creation) operator	<pre>create(N)</pre>	same as above
Number operator	num(N)	same as above
Single-mode displacement operator	displace(N,alpha)	N=number of levels in Hilbert space, alpha = complex displacement amplitude.
Single-mode squeezing operator	squeeze(N,sp)	N=number of levels in Hilbert space, sp = squeezing parameter.
Sigma-X	sigmax()	
Sigma-Y	sigmay()	
Sigma-Z	sigmaz()	
Sigma plus	sigmap()	
Sigma minus	sigmam()	





EXPECTATION VALUES

$$|0\rangle = {1 \choose 0}$$

$$|1\rangle = {0 \choose 1}$$

$$s_z = {1 \quad 0 \choose 0 \quad -1}$$

$$\langle s_z \rangle_0 = (1 \quad 0) \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix} \begin{pmatrix} 1 \\ 0 \end{pmatrix} = 1$$

$$\langle s_z \rangle_1 = (0 \ 1) \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix} \begin{pmatrix} 0 \\ 1 \end{pmatrix} = -1$$

```
In [72]: up = basis(2, 0)
In [73]: down = basis(2, 1)
In [74]: expect(sigmaz(), up)
Out[74]: 1.0
In [75]: expect(sigmaz(), down)
Out[75]: -1.0
```



TENSOR PRODUCTS



To describe the states of multipartite quantum systems - such as two coupled qubits, a qubit coupled to an oscillator, etc. - we need to expand the Hilbert space by taking the tensor product of the state vectors for each of the system components.

For example, the state vector describing two qubits in their ground states is formed by taking the tensor product of the two single-qubit ground state vectors:

```
In [1]: tensor(basis(2, 0), basis(2, 0))
Out[1]:
Quantum object: dims = [[2, 2], [1, 1]], shape = (4, 1), type = ket
Qobj data =
[[1.]
       [0.]
       [0.]
       [0.]]
```

```
In [4]: tensor(sigmax(), sigmax())  
Out[4]:  
Quantum object: dims = [[2, 2], [2, 2]], shape = (4, 4), type = oper, isherm = True  
Qobj data =  
[[0. 0. 0. 1.]  
[0. 0. 1. 0.]  
[0. 1. 0. 0.]  
[1. 0. 0. 0.]]  
\sigma_{\chi} \otimes \sigma_{\chi}
```

Partial trace





The partial trace is an operation that reduces the dimension of a Hilbert space by eliminating some degrees of freedom by averaging (tracing).

For example, the density matrix describing a single qubit obtained from a coupled two-qubit system is obtained via:

Example:

The two-qubit spin singlet $|\psi^{-}\rangle = \frac{1}{\sqrt{2}}(|01\rangle - |10\rangle)$ corresponds to the density matrix

$$|\psi^{-}\rangle\langle\psi^{-}| = \frac{1}{2}(|01\rangle\langle01| - |01\rangle\langle10| - |10\rangle\langle01| + |10\rangle\langle10|),$$

on which we act with the partial trace acording to the definition (1)

$$\operatorname{tr}_{B} |\psi^{-}\rangle\langle\psi^{-}| = \frac{1}{2} (|0\rangle\langle0|\langle1|1\rangle - |0\rangle\langle1|\langle0|1\rangle - |1\rangle\langle0|\langle1|0\rangle + |1\rangle\langle1|\langle0|0\rangle)$$
$$= \frac{1}{2} (|0\rangle\langle0| + |1\rangle\langle1|),$$

The partial trace always results in a density matrix (mixed state), regardless of whether the composite system is a pure state (described by a state vector) or a mixed state (described by a density matrix)

```
In [70]: psi
Out[70]:
Quantum object: dims = [[2, 2], [1, 1]], shape = (4, 1), type = ket
Qobj data =
 [-0.70710678]
 [ 0.70710678]
In [71]: psi.ptrace(0)
Out[71]:
Quantum object: dims = [[2], [2]], shape = (2, 2), type = oper, isherm = True
Qobi data =
[[0.5 0. ]
 [0. 0.5]]
In [72]: psi.ptrace(1)
Out[72]:
Quantum object: dims = [[2], [2]], shape = (2, 2), type = oper, isherm = True
Qobj data =
[[0.5 0. ]
 [0. 0.5]]
```





QUANTUM INFORMATION TOOLS IN QUTIP

Name	Equation	Qutip Command
Von Neumann entropy	$H = -tr(\rho ln\rho)$ $= -\Sigma \lambda_i ln \lambda_i$	entropy_vn(rho)
Linear entropy of a density matrix	$S_L = 1 - tr(\rho^2)$	*entropy_linear (rho)
Mutual information	I(A:B) = H(A) + H(B) - H(AB)	entropy_mutual(rho, selA, selB)
Concurrence entanglement	$C(\rho) = \max\{0, \lambda_1 - \lambda_2 - \lambda_3 - \lambda_4\}$	concurrence(rho)
Negativity	$N(\rho) = \frac{\left \left \rho^{\Gamma_A}\right \right _1 - 1}{2}$	negativity(rho, subsys)

^{*}The linear entropy can range between zero, corresponding to a completely pure state, and (1-1/d), corresponding to a completely mixed state. (Here, d is the dimension of the density matrix.)

TWO-QUBIT SYSTEM





$$\psi_1 = |00>$$

$$\psi_2 = \frac{1}{\sqrt{2}}(|00> + |11>)$$

$$\rho_3 = \frac{1}{\sqrt{2}}(|0> < 0| \otimes |0> < 0| + |+> < + |\otimes|1> < 1|)$$

Quantum State	ψ_1	ψ_2	P3 *** *** *** *** *** *** *** *** *** *	
Von Neumann entropy	0	0	0	
Linear entropy of a density matrix	0	0	0.5	
Concurrence entanglement	0	0.999	0	
	Pure separable state	Pure entangled state	Mixed state	



QUANTUM DISCORD



In quantum information theory, **quantum discord** is a measure of nonclassical correlations between two subsystems of a quantum system. It includes correlations that are due to quantum physical effects but do not necessarily involve quantum entanglement.

The quantum discord, a measure of the quantumness of correlations, of a state ρ_{AB} under a von Neumann measurement $\{\Pi_a\}$ is defined as a difference between the quantum forms of mutual information,

$$\mathcal{D} = I(A:B) - J(A:B) = S(\rho_A) - S(\rho_{AB}) + \min_{\{\Pi_a\}} \sum_a p_a S(\rho_{B|a}),$$
(0.1)

where the minimization is over all local projectors.

Example: Consider the following state

$$\rho_{AB} = p |0\rangle \langle 0| \otimes |0\rangle \langle 0| + (1-p) |+\rangle \langle +| \otimes |1\rangle \langle 1|$$

where $|+\rangle = \frac{1}{\sqrt{2}}(|0\rangle + |1\rangle)$.

$$\rho_{B|a} = \operatorname{tr}_A[(\Pi_a \otimes I)\rho_{AB}(\Pi_a^{\dagger} \otimes I)]/p_a, \quad \left(p_a = \operatorname{tr}[(\Pi_a \otimes I)\rho_{AB}(\Pi_a^{\dagger} \otimes I)]\right)$$

where von Neumann measurements are given by $\Pi_1 = |0\rangle \langle 0|$ and $\Pi_2 = |1\rangle \langle 1|$. Then we have

$$p_1 = \operatorname{tr}[(\Pi_1 \otimes I)\rho_{AB}(\Pi_1^{\dagger} \otimes I)] = \operatorname{tr}[p|0\rangle\langle 0| \otimes |0\rangle\langle 0| + \frac{(1-p)}{2}|0\rangle\langle 0| \otimes |1\rangle\langle 1|] = \frac{(1+p)}{2}$$

$$p_2 = \operatorname{tr}[(\Pi_2 \otimes I)\rho_{AB}(\Pi_2^{\dagger} \otimes I)] = \operatorname{tr}[\frac{(1-p)}{2} |1\rangle \langle 1| \otimes |1\rangle \langle 1|] = \frac{(1-p)}{2}$$

$$\rho_{B|a=1} = \operatorname{tr}_{A}[(\Pi_{1} \otimes I)\rho_{AB}(\Pi_{1}^{\dagger} \otimes I)]/p_{1}$$

$$= \operatorname{tr}_{A}[p|0\rangle\langle 0| \otimes |0\rangle\langle 0| + \frac{(1-p)}{2}|0\rangle\langle 0| \otimes |1\rangle\langle 1|]/p_{1}$$

$$= [p|0\rangle\langle 0| + \frac{(1-p)}{2} \otimes |1\rangle\langle 1|]/(\frac{1+p}{2})$$
(0.2)

$$\rho_{B|a=2} = \operatorname{tr}_{A}[(\Pi_{2} \otimes I)\rho_{AB}(\Pi_{2}^{\dagger} \otimes I)]/p_{2}$$

$$= \operatorname{tr}_{A}[\frac{(1-p)}{2} |1\rangle \langle 1| \otimes |1\rangle \langle 1|]/p_{2}$$

$$= |1\rangle \langle 1| \qquad (0.3)$$









$$\rho_{A} = \operatorname{tr}_{B}(\rho_{AB})$$

$$= \operatorname{tr}_{B}(p|0)\langle 0| \otimes |0\rangle\langle 0| + \frac{(1-p)}{2} \left((|0\rangle\langle 0| + |0\rangle\langle 1| + |1\rangle\langle 0| + |1\rangle\langle 1|) \otimes |1\rangle\langle 1| \right))$$

$$= p|0\rangle\langle 0| + \frac{(1-p)}{2} (|0\rangle\langle 0| + |0\rangle\langle 1| + |1\rangle\langle 0| + |1\rangle\langle 1|)$$

$$= \frac{(1+p)}{2} |0\rangle\langle 0| + \frac{(1-p)}{2} (|0\rangle\langle 1| + |1\rangle\langle 0| + |1\rangle\langle 1|)$$
(0.5)

Then we have

Eigenvalues of
$$\rho_A$$
: $\left\{ \frac{1}{2} \left(1 - \sqrt{2p^2 - 2p + 1} \right), \frac{1}{2} \left(\sqrt{2p^2 - 2p + 1} + 1 \right) \right\}$

Eigenvalues of ρ_{AB} : $\{0, 0, 1-p, p\}$

Eigenvalues of
$$\rho_{B|a=1}: \left\{ \frac{1-p}{p+1}, \frac{2p}{p+1} \right\}$$

Eigenvalues of $\rho_{B|a=2}:\{1,0\}$





$$\mathcal{D} = I(A:B) - J(A:B)$$

$$= S(\rho_A) - S(\rho_{AB}) + \min_{\{\Pi_a\}} \sum_{a} p_a S(\rho_{B|a})$$

$$= -\left[\frac{1}{2}\left(1 - \sqrt{2p^2 - 2p + 1}\right) \log \frac{1}{2}\left(1 - \sqrt{2p^2 - 2p + 1}\right) + \frac{1}{2}\left(1 + \sqrt{2p^2 - 2p + 1}\right) \log \frac{1}{2}\left(1 + \sqrt{2p^2 - 2p + 1}\right)\right]$$

$$+ \left[(1 - p) \log (1 - p) + p \log p\right]$$

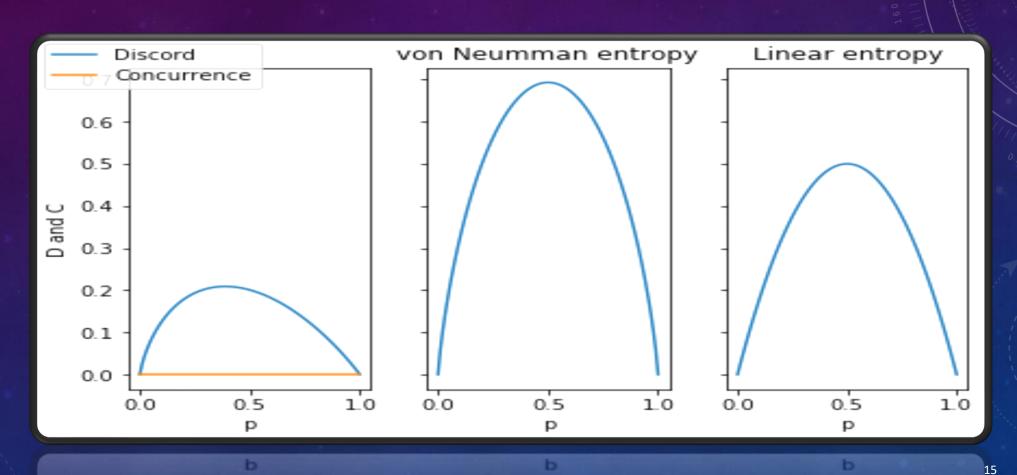
$$+ p_1 \left(-\frac{1 - p}{p + 1} \log(\frac{1 - p}{p + 1}) - \frac{2p}{p + 1} \log(\frac{2p}{p + 1})\right) + p_2 \left(-1 \log 1\right)$$

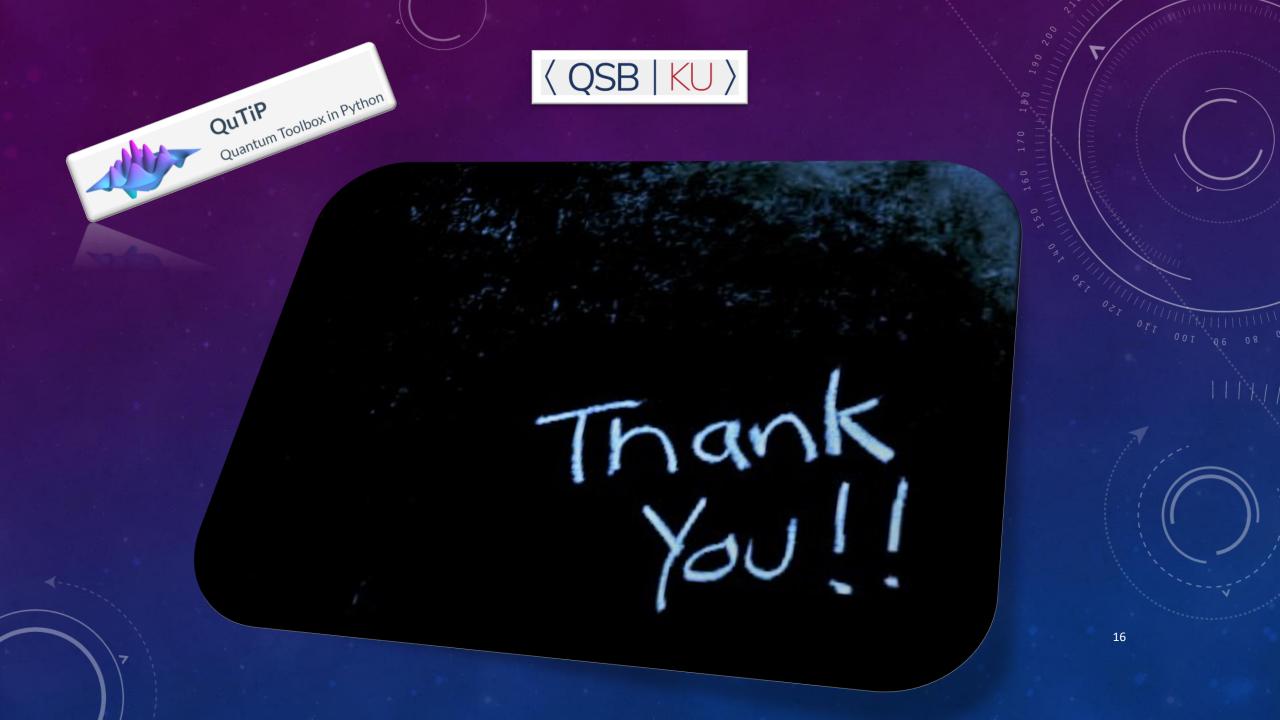
$$(0.6)$$





$\rho_{AB} = p|0><0|\otimes|0><0|+(1-p)|+><+|\otimes|1><1|$





KUANTUM TEKNOLOJİLERİ & KAYNAK TEORİLERİ

10 Nisan KUANTUM KAYNAK TEORİLERİNE GİRİŞ	
14.00-14.50 Matematiksel Formaliz	zm
14.50-15.40 Kuantum Üst Üste Bin Kuantum Eşevrelilik	me
15.50-16.40 Kuantum Dolaşıklık	
16.40-17.30 Kuantum Uyumsuzluk	
17.30-18.15 QuTiP'e Giriş	

11 Nisan	KUANTUM TEKNOLOJİLERİNE ÖRNEKLER
14.00-14.50	Kuantum Enformasyon ve Hesaplama
14.50-15.40	Kuantum Termodinamik
15.50-16.40	Kuantum Metroloji
16.40-17.30	Kuantum Biyoloji

Eğitimler Dr.Onur Pusuluk, Dr. Gökhan Torun ve Mohsen Izadyari tarafından verilecektir.

 $\langle \mathrm{Q} | \mathrm{Turkey} \rangle$