

A quantum computer can be simulated by applying rotations to a unit vector  $u \in \mathbb{C}^{2^n}$  where  $n$  is the number of qubits. The dimension is  $2^n$  because a register with  $n$  qubits has  $2^n$  eigenstates. Quantum operations are “rotations” because they preserve  $|u| = 1$ .

The Eigenmath function  $rotate(u, s, k, \dots)$  rotates vector  $u$  and returns the result. Vector  $u$  is required to have  $2^n$  elements where  $n$  is an integer from 1 to 15. Arguments  $s, k, \dots$  are a sequence of rotation codes where  $s$  is an upper case letter and  $k$  is a qubit number from 0 to  $n-1$ . Rotations are evaluated from left to right. The available rotation codes are

$C, k$	Control prefix
$H, k$	Hadamard
$P, k, \phi$	Phase modifier
$Q, k$	Quantum Fourier transform
$V, k$	Inverse quantum Fourier transform
$W, k, j$	Swap bits
$X, k$	Pauli X
$Y, k$	Pauli Y
$Z, k$	Pauli Z

Control prefix  $C, k$  modifies the next rotation code so that it is a controlled rotation with  $k$  as the control qubit. Fourier rotations  $Q, k$  and  $V, k$  are applied to qubits 0 through  $k$ . ( $Q$  and  $V$  ignore any control prefix.)

Error codes

- 1 Argument  $u$  is not a vector or does not have  $2^n$  elements where  $n = 1, 2, \dots, 15$ .
- 2 Unexpected end of argument list (i.e., missing argument).
- 3 Bit number format error or range error.
- 4 Unknown rotation code.

Eigenstates  $|j\rangle$  are represented by the following vectors. (Each vector has  $2^n$  elements.)

$$\begin{aligned}
|0\rangle &= (1, 0, 0, \dots, 0) \\
|1\rangle &= (0, 1, 0, \dots, 0) \\
|2\rangle &= (0, 0, 1, \dots, 0) \\
&\vdots \\
|2^n - 1\rangle &= (0, 0, 0, \dots, 1)
\end{aligned}$$

A quantum computing algorithm is a sequence of rotations applied to the initial state  $|0\rangle$ . (Mathematically, the sequence can be collapsed into a single rotation.) Let  $\psi$  be the final state of the quantum computer after all the rotations have been applied. Like any other state,  $\psi$  is a linear combination of eigenstates.

$$\psi = \sum_{j=0}^{2^n-1} c_j |j\rangle, \quad |\psi| = 1$$

The final step is to measure  $\psi$  and get a result. Measurement rotates  $\psi$  to an eigenstate  $|j\rangle$ . The measurement result is  $|j\rangle$ . The probability of obtaining a specific  $|j\rangle$  is

$$P_j = c_j c_j^*$$

Note that if  $\psi$  is already an eigenstate then no rotation occurs. (The probability of rotating to a different eigenstate is zero.) Since the measurement result is always an eigenstate, the coefficients  $c_j$  cannot be observed. However, the same calculation can be run multiple times to obtain a probability distribution. The probability distribution is an estimate of  $c_j c_j^*$  for each  $|j\rangle$ .

Eigenmath code snippets for before and after *rotate*:

Initialize  $\psi = |0\rangle$ .

```
n = 4          -- number of qubits (example)
N = 2^n        -- number of eigenstates
psi = zero(N)
psi[1] = 1
```

Compute the probability distribution for state  $\psi$ .

```
P = psi conj(psi)
```

Hence

```
P[1] = probability that |0> will be the result
P[2] = probability that |1> will be the result
P[3] = probability that |2> will be the result
:
P[N] = probability that |N - 1> will be the result
```

Draw the probability distribution.

```
xrange = (0,N)
yrange = (0,1)
draw(P[ceiling(x)],x)
```

Compute an expectation value.

```
sum(k,1,N, (k - 1) P[k])
```

Make the high order bit “don’t care.”

```
for(k,1,N/2, P[k] = P[k] + P[k + N/2])
```

Hence for  $N = 16$

```
P[1] = probability that the result will be |0> or |8>
P[2] = probability that the result will be |1> or |9>
P[3] = probability that the result will be |2> or |10>
:
P[8] = probability that the result will be |7> or |15>
```