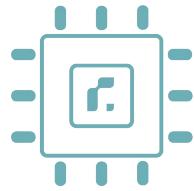


# rigetti

Unsupervised Machine Learning on a Hybrid Quantum Computer  
Johannes Otterbach

Bay Area Quantum Computing Meetup - YCombinator  
February 1, 2018

# Full Stack Quantum Computing



Quantum  
Processor



Hardware



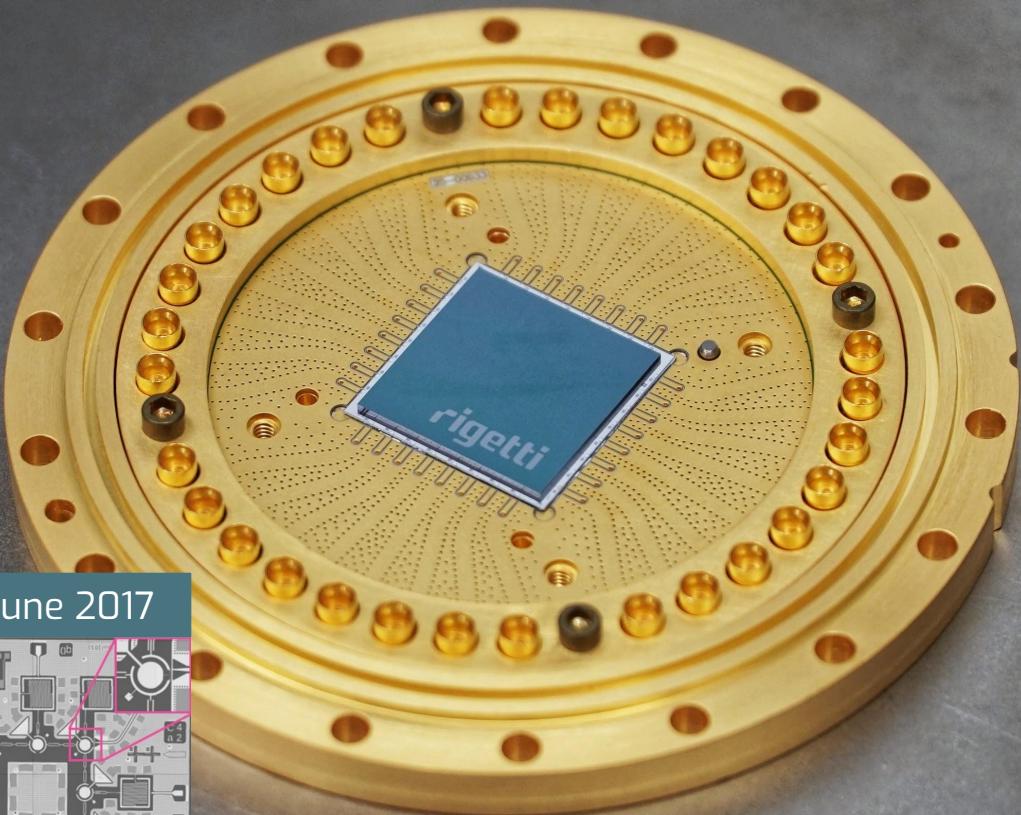
Cloud based Quantum  
Operating System



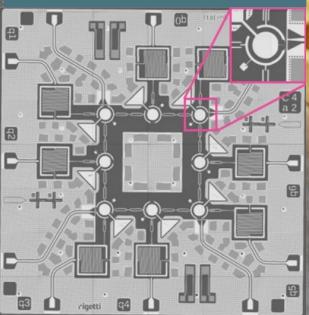
Applications



# 19Q-Acorn available now

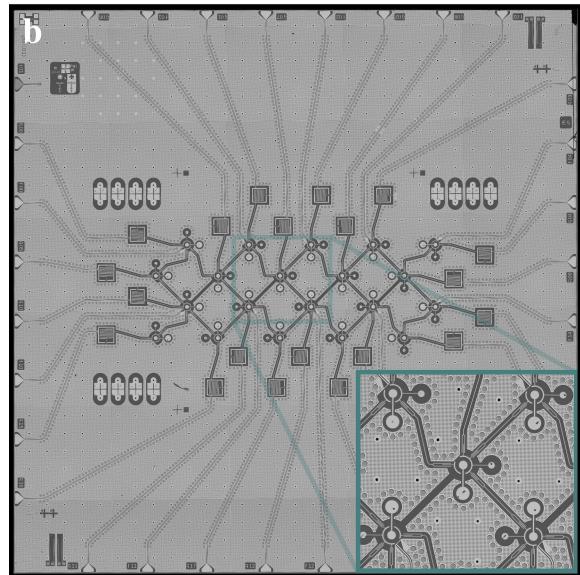
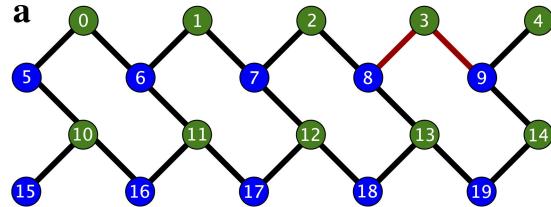


8Q June 2017

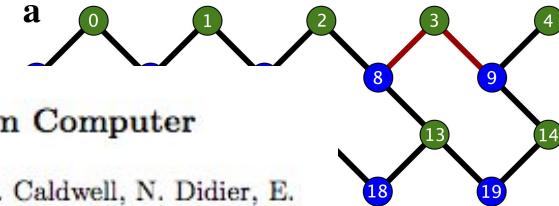


[pyquil.readthedocs.io](http://pyquil.readthedocs.io)

rigetti



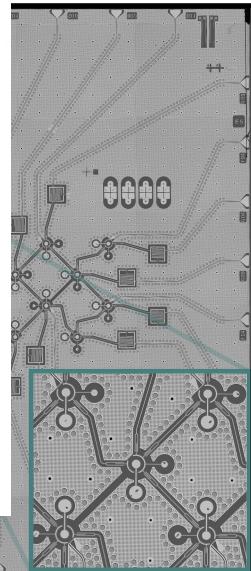
arXiv:1712.05771



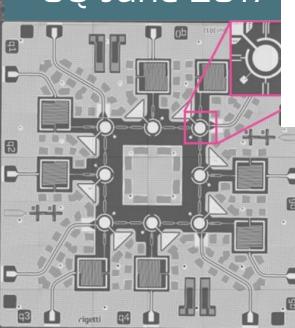
## Unsupervised Machine Learning on a Hybrid Quantum Computer

J. S. Otterbach, R. Manenti, N. Alidoust, A. Bestwick, M. Block, B. Bloom, S. Caldwell, N. Didier, E. Schuyler Fried, S. Hong, P. Karalekas, C. B. Osborn, A. Papageorge, E. C. Peterson, G. Prawiroatmodjo, N. Rubin, Colm A. Ryan, D. Scarabelli, M. Scheer, E. A. Sete, P. Sivarajah, Robert S. Smith, A. Staley, N. Tezak, W. J. Zeng, A. Hudson, Blake R. Johnson, M. Reagor, M. P. da Silva, and C. Rigetti  
*Rigetti Computing, Inc., Berkeley, CA*  
(Dated: December 18, 2017)

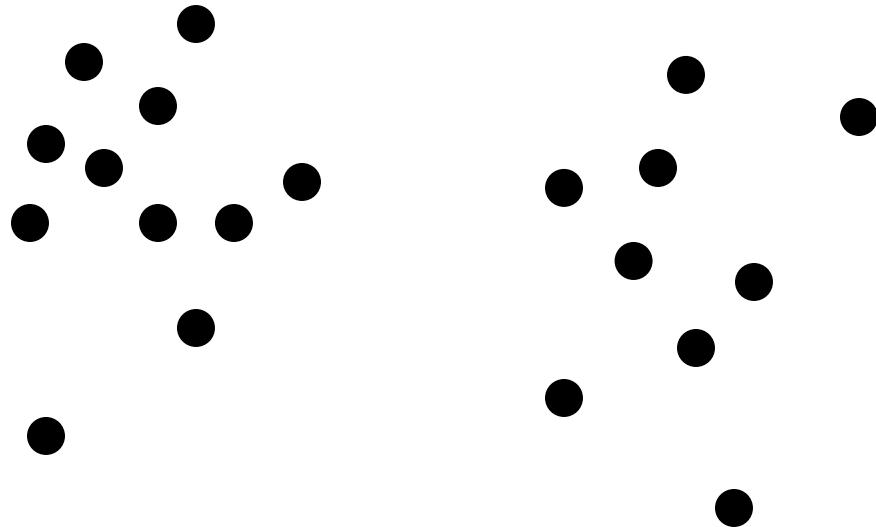
Machine learning techniques have led to broad adoption of a statistical model of computing. The statistical distributions natively available on quantum processors are a superset of those available classically. Harnessing this attribute has the potential to accelerate or otherwise improve machine learning relative to purely classical performance. A key challenge toward that goal is learning to hybridize classical computing resources and traditional learning techniques with the emerging capabilities of general purpose quantum processors. Here, we demonstrate such hybridization by training a 19-qubit gate model processor to solve a clustering problem, a foundational challenge in unsupervised learning. We use the quantum approximate optimization algorithm in conjunction with a gradient-free Bayesian optimization to train the quantum machine. This quantum/classical hybrid algorithm shows robustness to realistic noise, and we find evidence that classical optimization can be used to train around both coherent and incoherent imperfections.



8Q June 2017



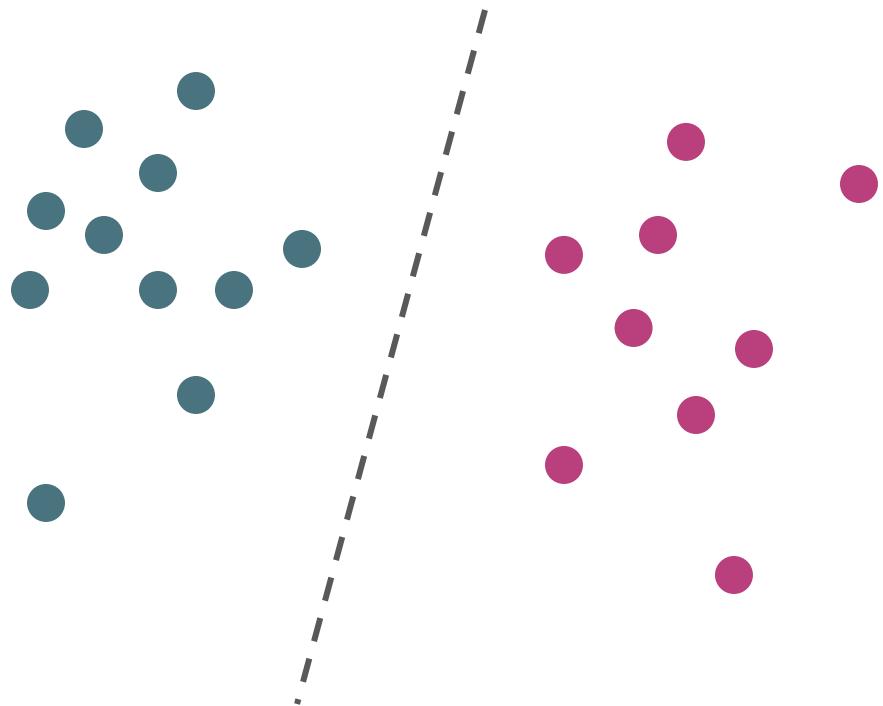
# Clustering



Given an unlabeled set of points



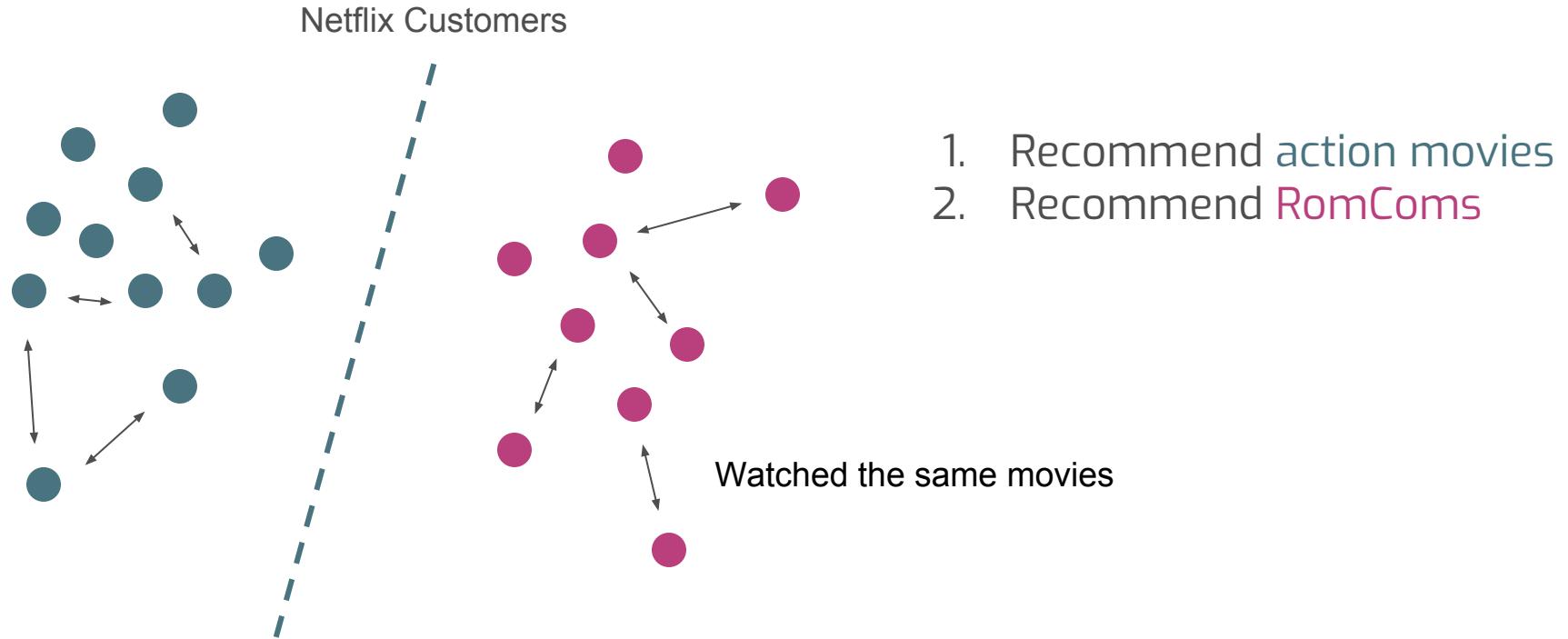
# Clustering



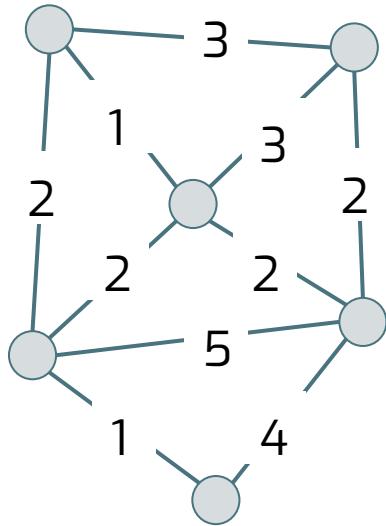
Given an unlabeled set of points, find labels based upon ***similarity*** metric (e.g. Euclidean distance).



# Clustering Example - Recommender Systems



# Clustering as MAXCUT

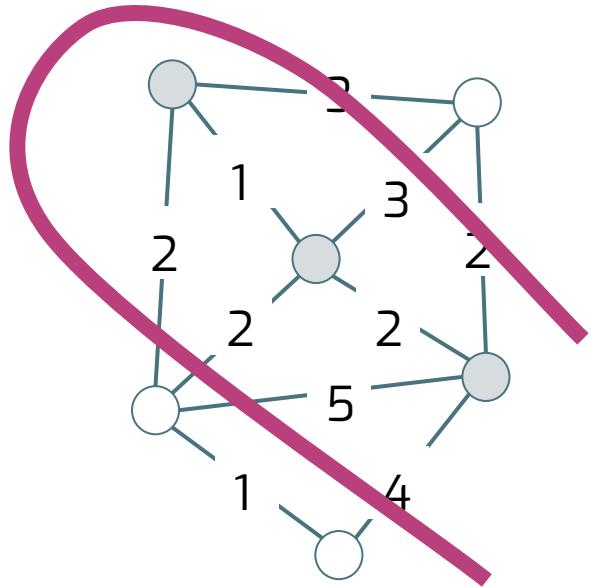


Construct a graph  $G=(V,E)$  where the edge weights  $w_{i,j}$  are determined by the distance metric. Then, MAXCUT is a clustering algorithm for the original points.

$$\text{MAXCUT} = \max_{\text{cut } S \subset E} \sum_{(i,j) \in S} w_{ij}$$



# Clustering as MAXCUT



Construct a graph  $G=(V,E)$  where the edge weights  $w_{i,j}$  are determined by the distance metric. Then, MAXCUT is a clustering algorithm for the original points.

Clustering transformed into an ***optimization*** problem.

$$\text{MAXCUT} = \max_{\text{cut } S \subset E} \sum_{(i,j) \in S} w_{ij}$$

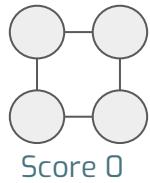


# MAXCUT as Energy Functional

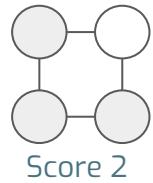
*“Maximize disagreement on a colored graph”*



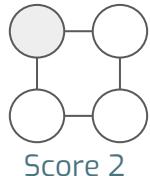
4-node “ring of disagrees”



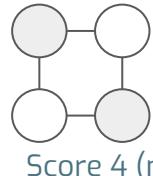
Score 0



Score 2



Score 2



Score 4 (max)

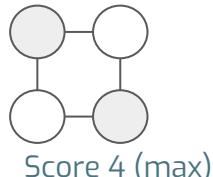
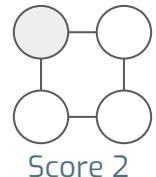
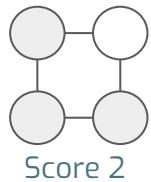
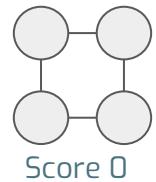


# MAXCUT as Energy Functional

*“Maximize disagreement on a colored graph”*



4-node “ring of disagrees”



Binary variable

$$\left. \begin{array}{l} \textcircled{1} = 1 \\ \textcircled{0} = 0 \end{array} \right\} \sigma^z \in \{0, 1\}$$

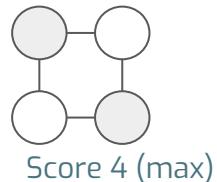
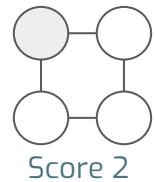
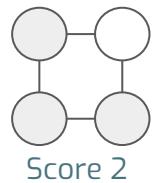
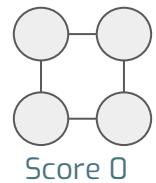


# MAXCUT as Energy Functional

“Maximize disagreement on a colored graph”



4-node “ring of disagrees”

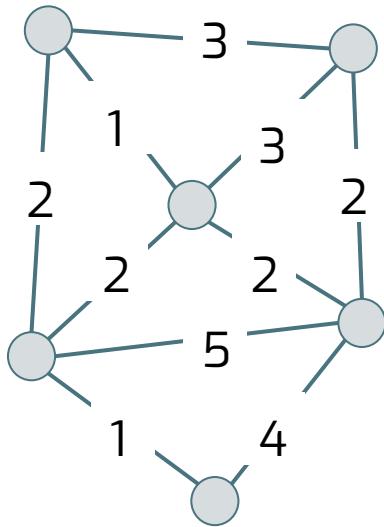


Binary variable

$$\left. \begin{array}{l} \textcircled{1} = 1 \\ \textcircled{0} = 0 \end{array} \right\} \sigma^z \in \{0, 1\}$$



# MAXCUT as Energy Functional

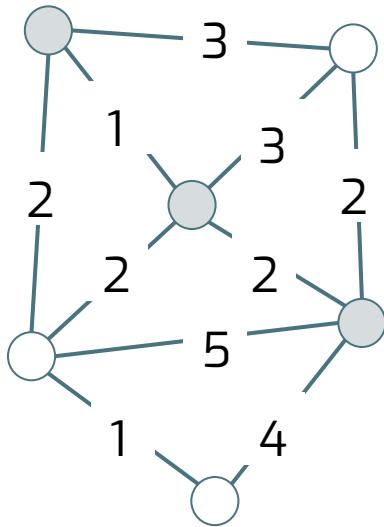


$$\begin{array}{l} \textcircled{\text{1}} = 1 \\ \textcircled{\text{0}} = 0 \end{array} \quad \left. \begin{array}{l} \sigma^z \in \{0, 1\} \end{array} \right\}$$

$$\begin{aligned} & \max_{\text{cut } S \subset E} \sum_{(i,j) \in S} w_{ij} \\ &= \max_{\sigma_i^z \in \{0,1\}} \sum_{i,j \in V} w_{ij} \sigma_i^z \sigma_j^z \end{aligned}$$



# MAXCUT as Energy Functional



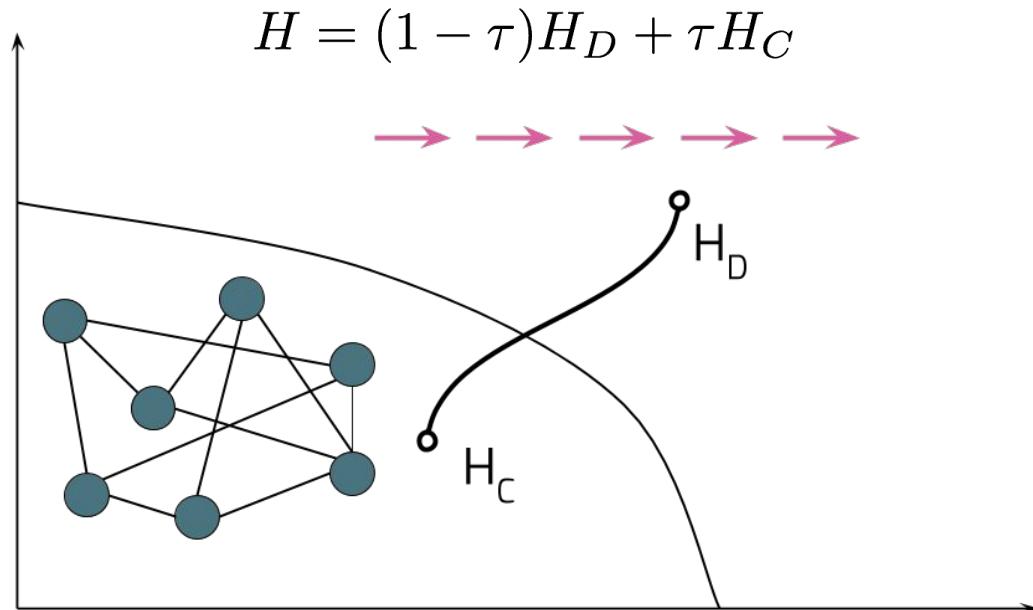
$$\begin{array}{l} \textcircled{\text{1}} = 1 \\ \textcircled{\text{0}} = 0 \end{array} \quad \left. \right\} \sigma^z \in \{0, 1\}$$

$$\begin{aligned} & \max_{\text{cut } S \subset E} \sum_{(i,j) \in S} w_{ij} \\ &= \max_{\sigma_i^z \in \{0,1\}} \sum_{i,j \in V} w_{ij} \sigma_i^z \sigma_j^z \end{aligned}$$

Find the right bit-string assignment that maximizes the energy



# QAOA - Quantum Approximate Optimization Algorithm



IDEA

Start at easy to prepare initial state of energy functional  $H_D$

“Cool” the system until it freezes in the low energy state of  $H_C$



# Discretize the Cooling Protocol

- Gate model of Optimization (Farhi, Goldstone, Gutman, arxiv:1411.4028)

$$V(\beta) = e^{-i\beta H_C}$$

$$U(\gamma) = e^{-i\gamma H_D}$$

- Angles  $\beta, \gamma$  need not be small
- How to find optimal  $\beta, \gamma$  ?



# Back to MAXCUT

- Initial state is ground state of  $H_D$ :  $| \rightarrow \rangle = H^{\otimes n} |0\rangle$
- Run the QAOA prescription:

$$| \beta, \gamma \rangle = U(\gamma) V(\beta) H^{\otimes n} |0\rangle$$

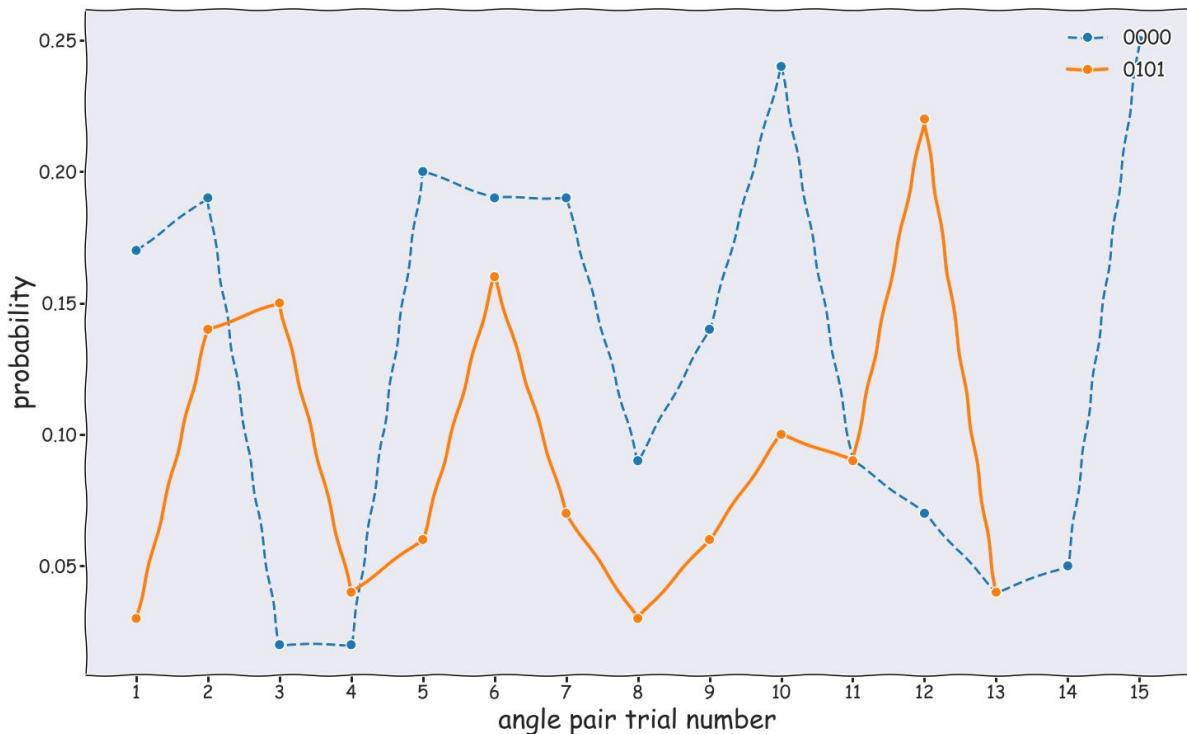
- Intuitively: Superposition of bitstring configurations

$$| \beta, \gamma \rangle \equiv \sqrt{p_1} \begin{array}{c} \text{---} \\ | \quad | \\ \text{---} \end{array} + \sqrt{p_2} \begin{array}{c} \text{---} \\ | \quad | \\ \text{---} \\ | \quad | \\ \text{---} \end{array} + \dots + \sqrt{p_{16}} \begin{array}{c} \text{---} \\ | \quad | \\ \text{---} \\ | \quad | \\ \text{---} \\ | \quad | \\ \text{---} \end{array}$$



# Effects of the different Angles

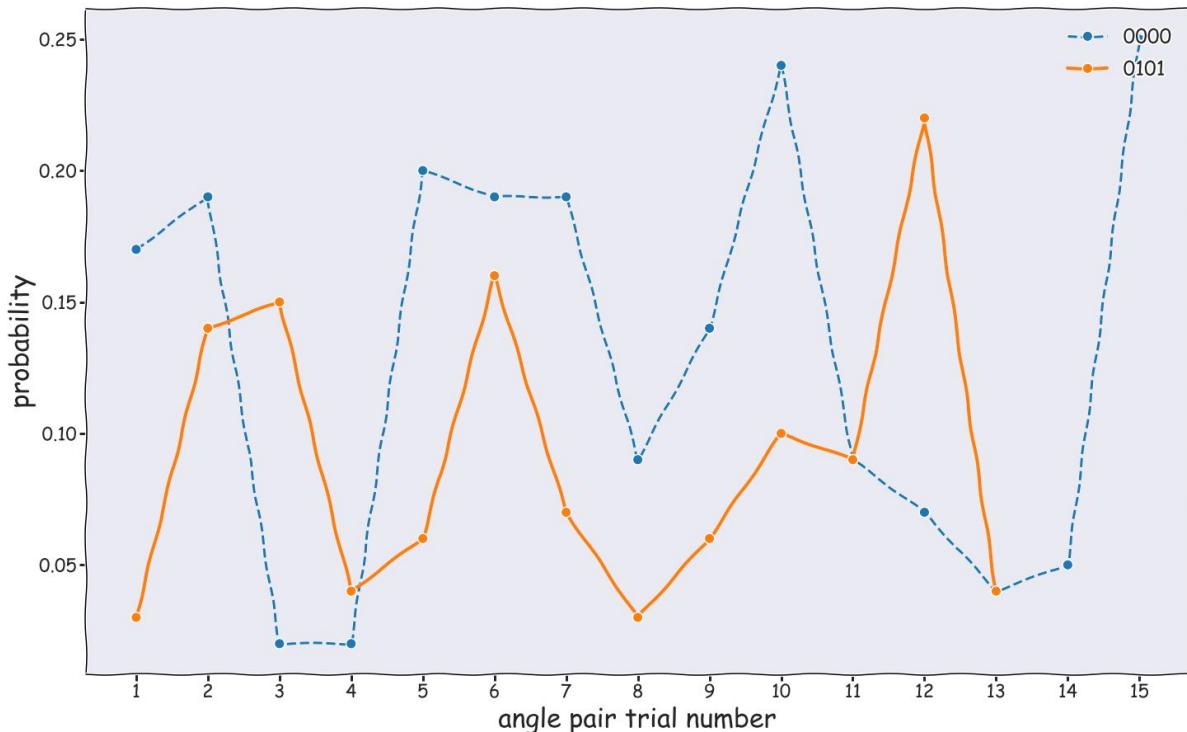
Angles change the probability to sample different bit strings



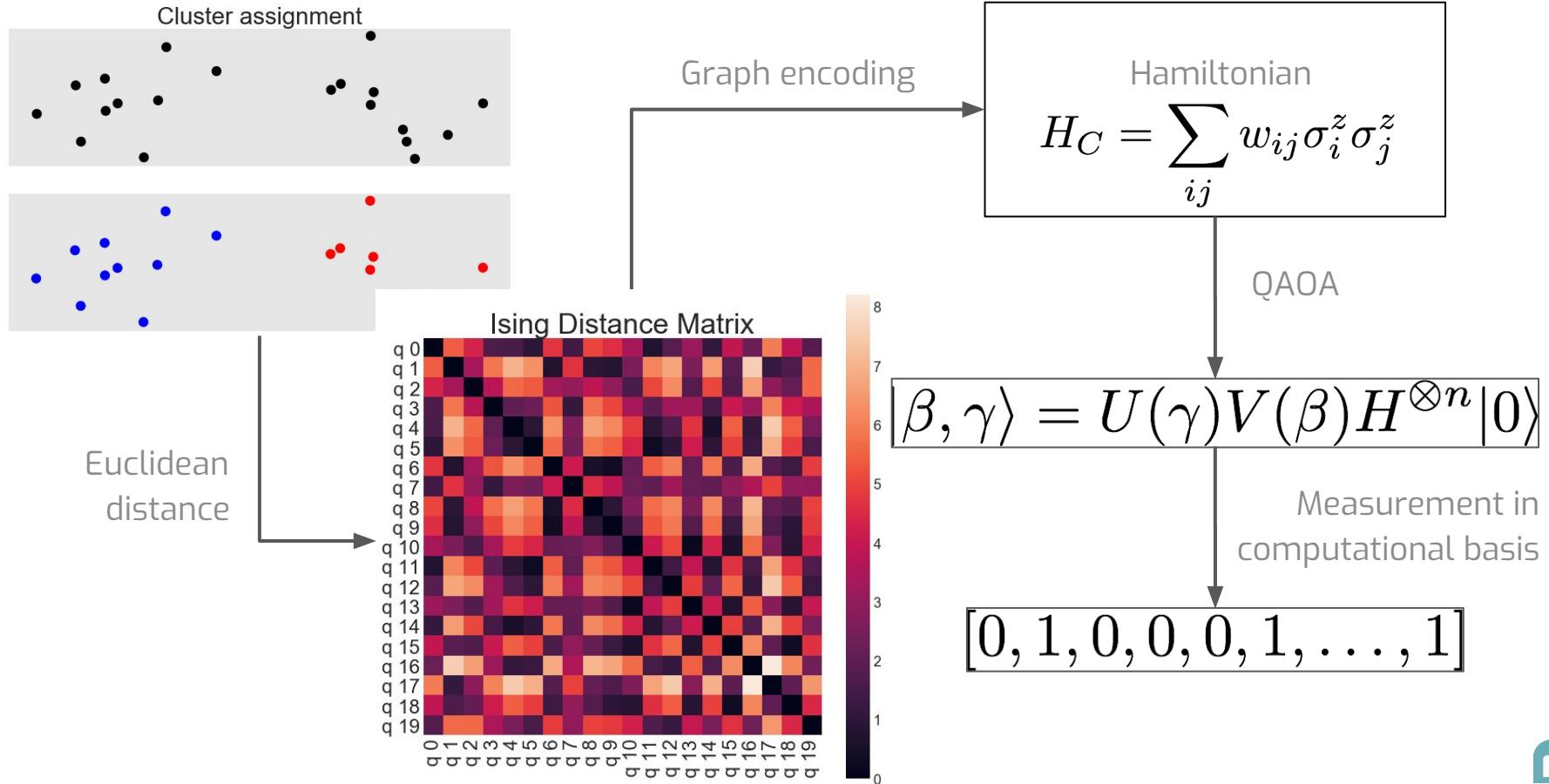
# Effects of the different Angles

Angles change the probability to sample different bit strings

Want to maximize the probability to sample the “correct” bitstring



# Clustering Procedure



# Forest - Quil

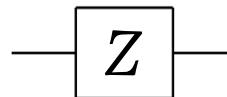
## Quantum instructions



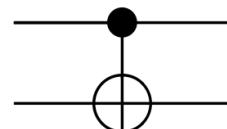
x



H



z



CNOT



MEASURE

## Quil [01⟩

Quantum Instruction Language

```
X q
H q
Z q
CNOT p q
MEASURE q [0]
MEASURE p [1]
```



# Forest - pyQuil

## QAOA

In **14** lines of code

```
from pyquil.quil import Program
from pyquil.gates import H
from pyquil.paulis import sI, sx, sz, exponentiate_commuting_pauli_sum
from pyquil.api import QVMConnection

graph = [(0, 1), (1, 2), (2, 3)]
nodes = range(4)

init_state_prog = sum([H(i) for i in nodes], Program())
h_cost = -0.5 * sum(sI(nodes[0]) - sz(i) * sz(j) for i, j in graph)
h_driver = -1. * sum(sx(i) for i in nodes)

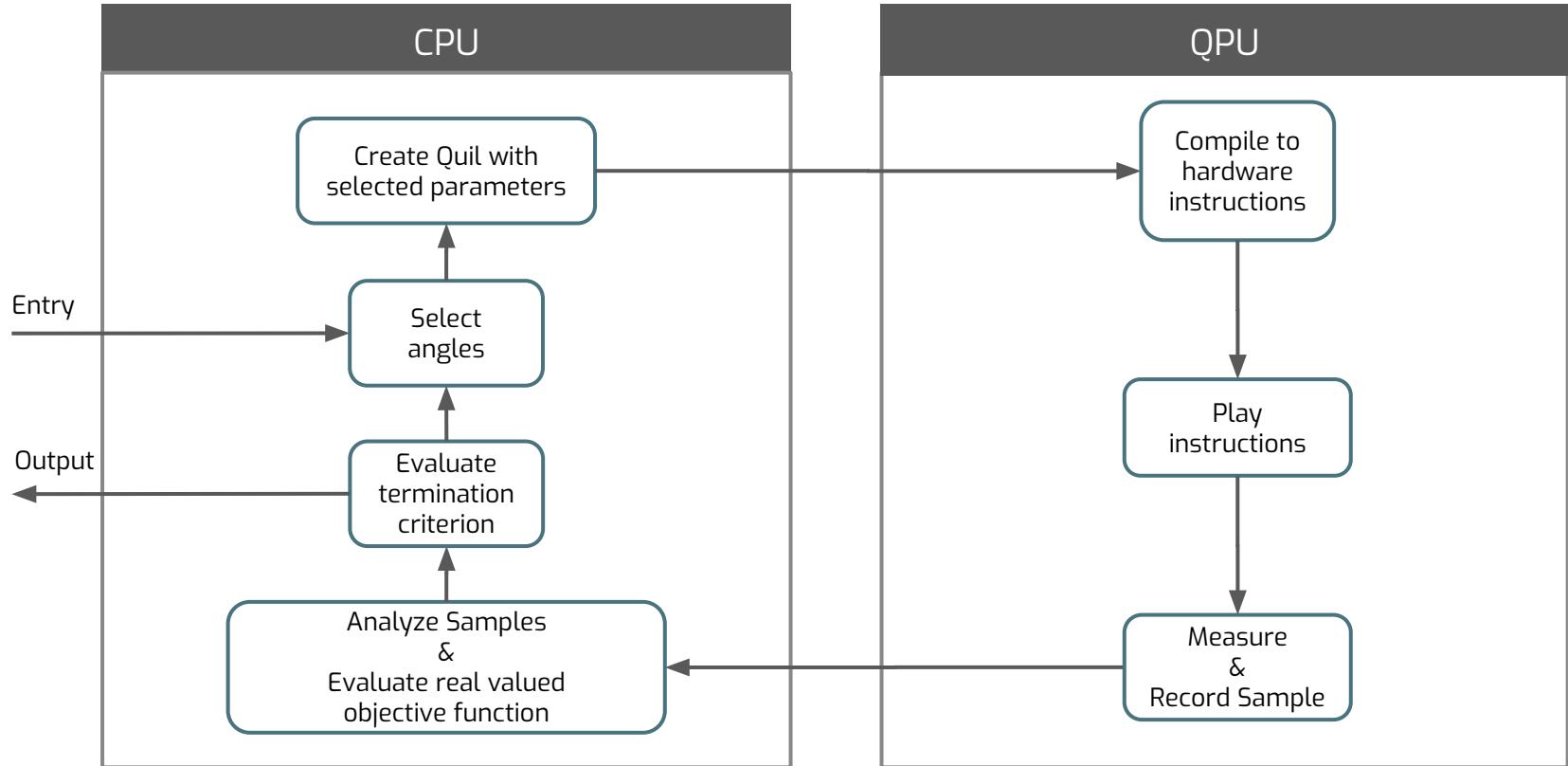
def qaoa_ansatz(betas, gammas):
    return sum([exponentiate_commuting_pauli_sum(h_cost)(g) + exponentiate_commuting_pauli_sum(h_driver)(b) \
               for g, b in zip(gammas, betas)]), Program()

program = init_state_prog + qaoa_ansatz([0., 0.5], [0.75, 1.])

qvm = QVMConnection()
qvm.run_and_measure(program, qubits=nodes, trials=10)
```



# When are we done?

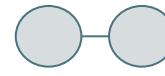


# Objective Function

- Loss/Reward Function:

$$c_{\beta, \gamma} : \{0, 1\}^n \mapsto \mathbb{R}$$

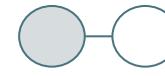
*"quality of a sampled bit-string"*



Score 0



Score 0



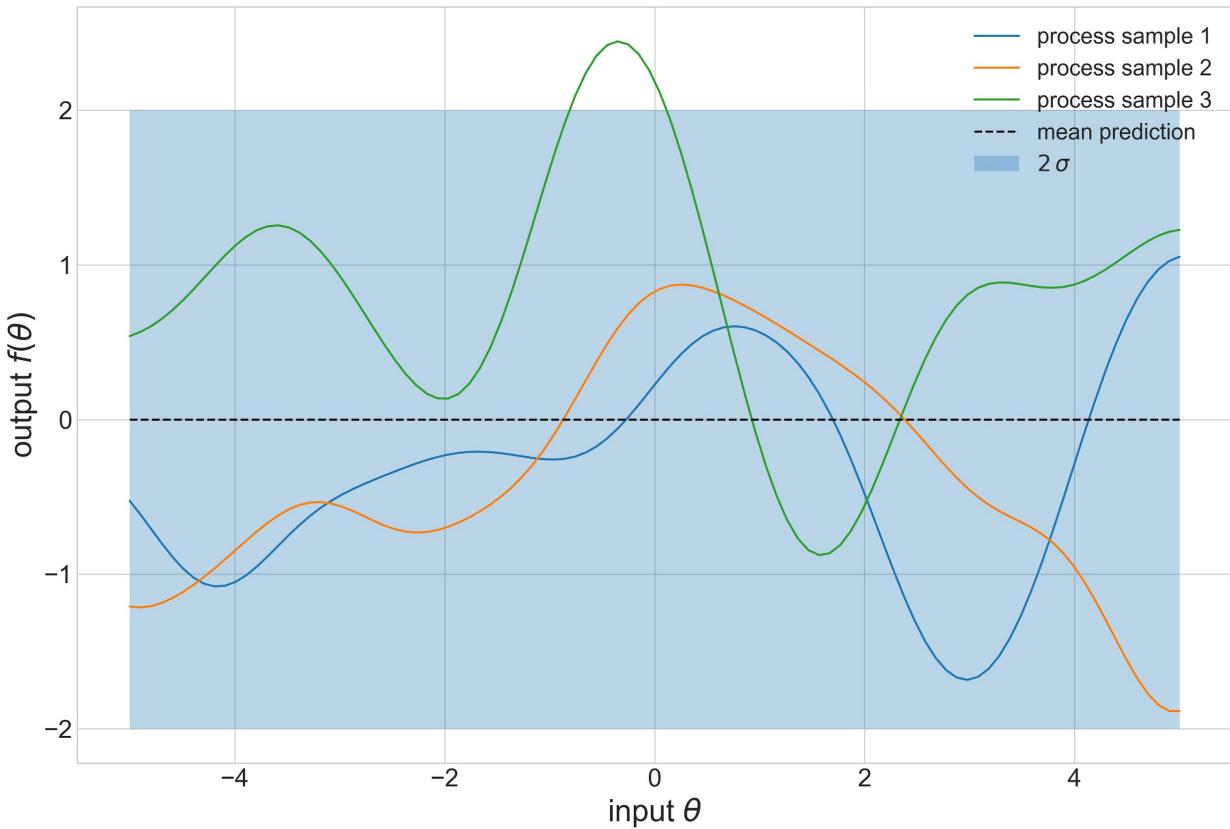
Score +1

- Find the optimal value of the Reward function:
  - No easy access to gradients, need derivative free methods
  - E.g. Bayesian Methods



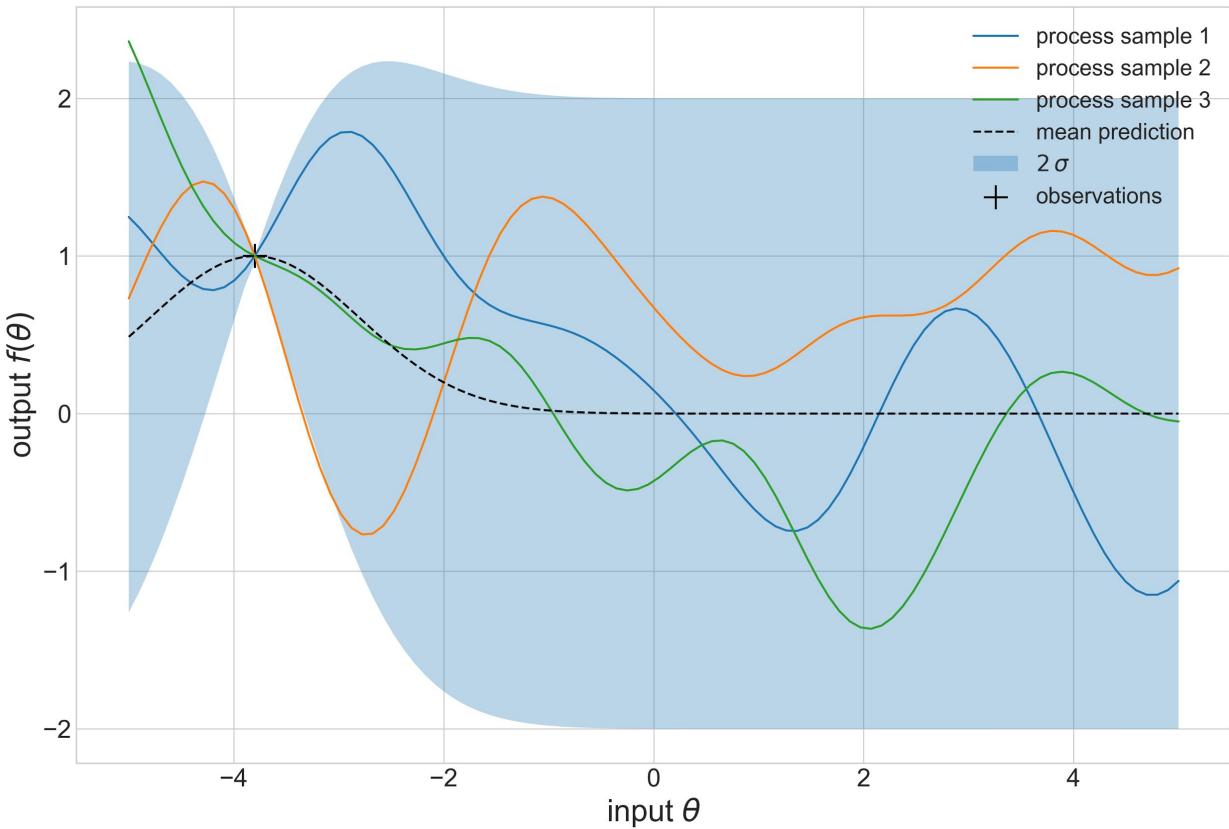
# Bayesian Optimization

- Assume objective function is Gaussian



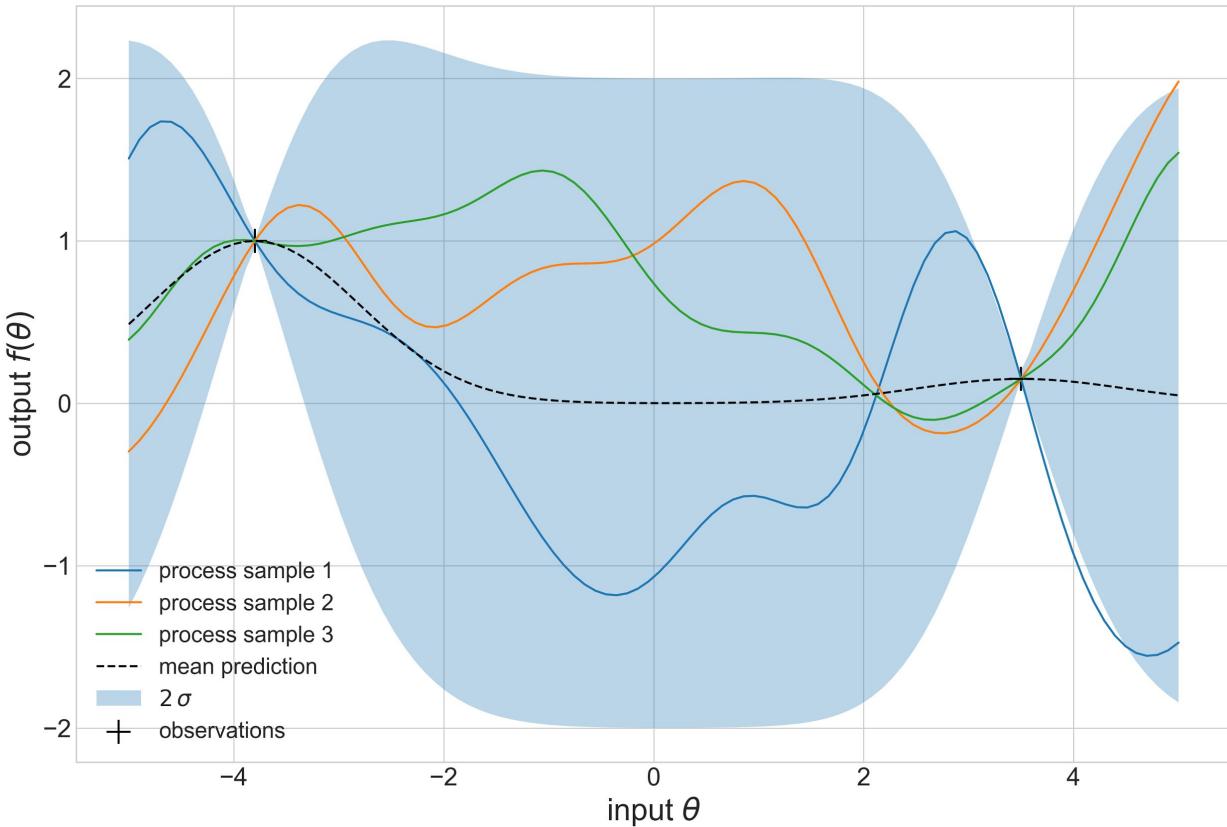
# Bayesian Optimization

- Assume objective function is Gaussian
- Measure and update Prior



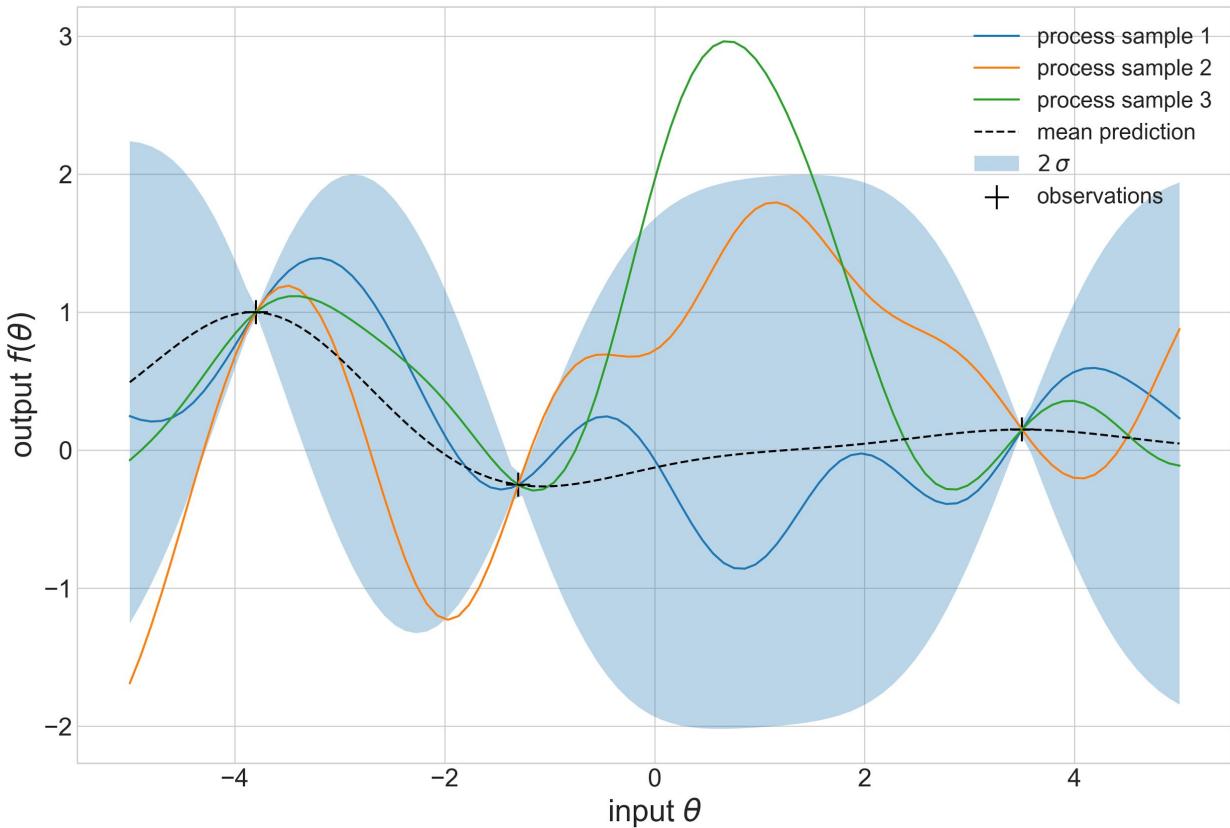
# Bayesian Optimization

- Assume objective function is Gaussian
- Measure and update Prior
- Choose next point to measure and update



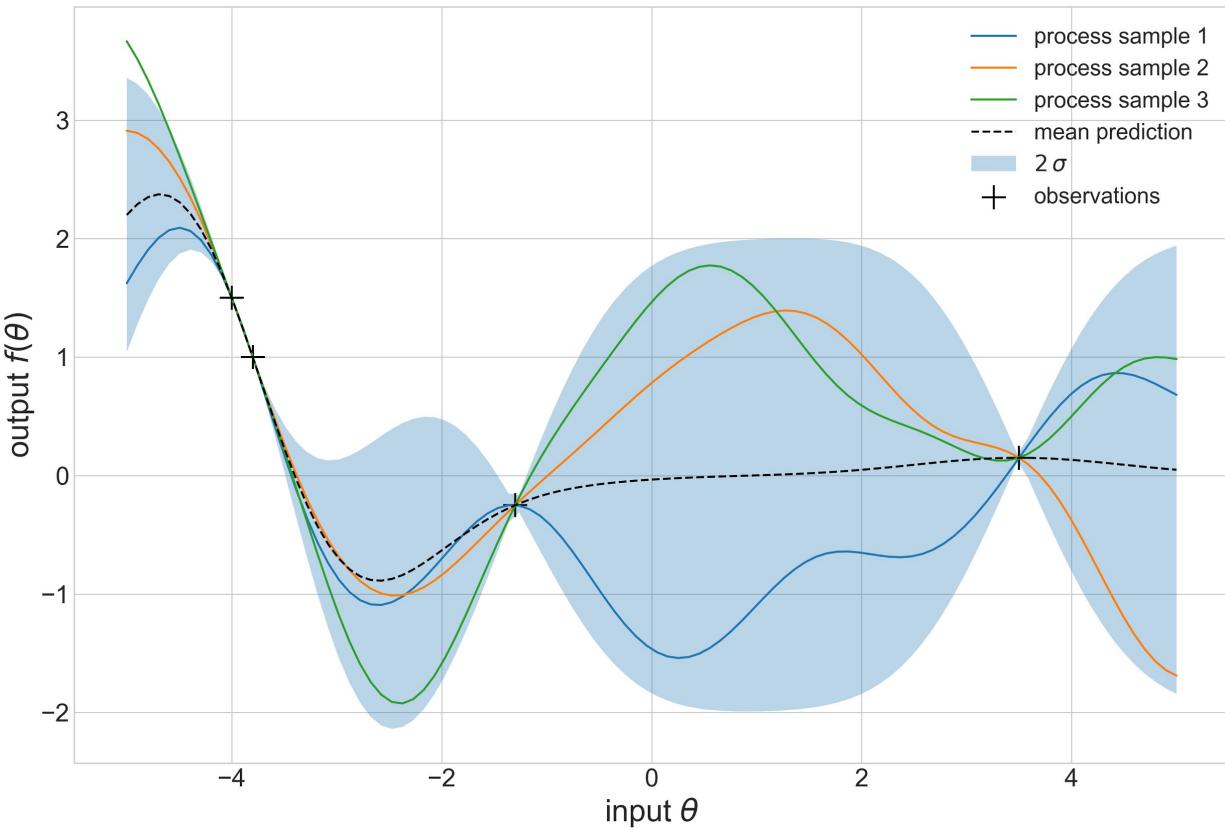
# Bayesian Optimization

- Assume objective function is Gaussian
- Measure and update Prior
- Choose next point to measure and update
- Again



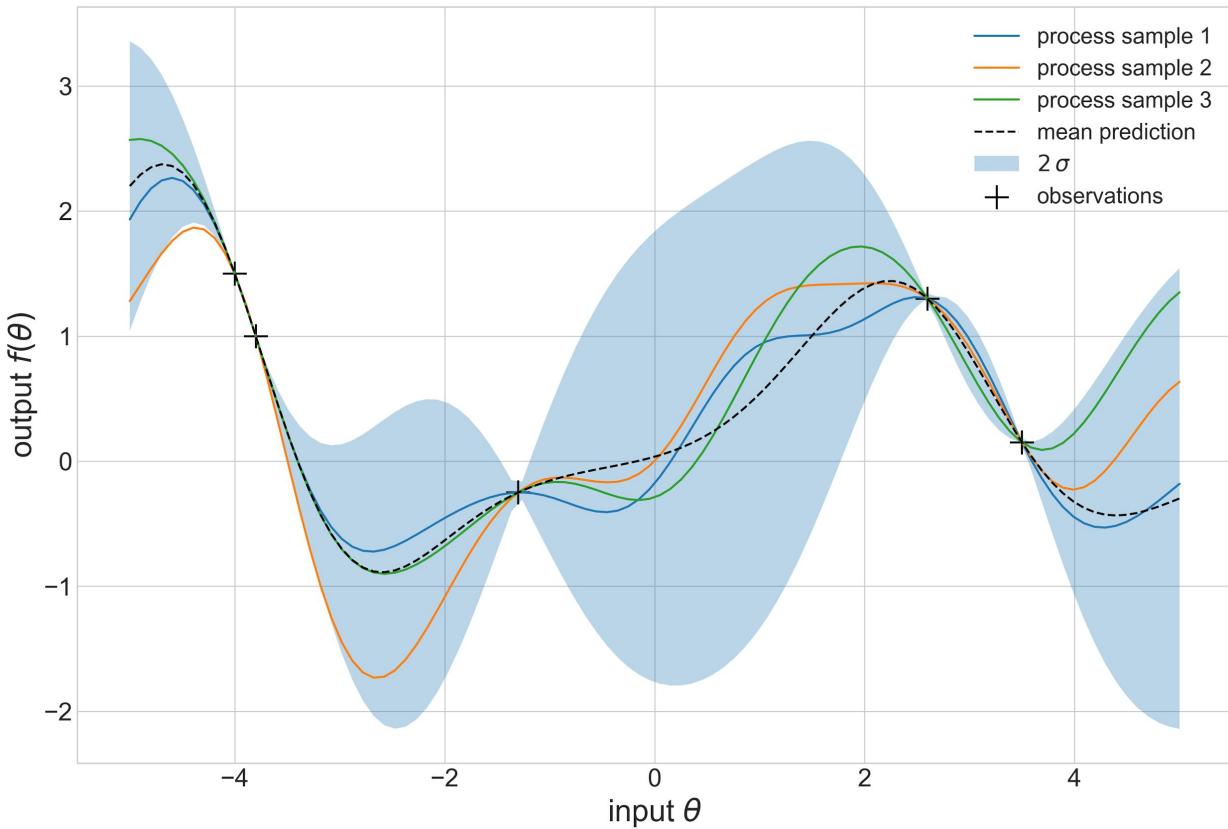
# Bayesian Optimization

- Assume objective function is Gaussian
- Measure and update Prior
- Choose next point to measure and update
- Again
- ...



# Bayesian Optimization

- Assume objective function is Gaussian
- Measure and update Prior
- Choose next point to measure and update
- Again
- ...
- ...



# That's how it looks in practice

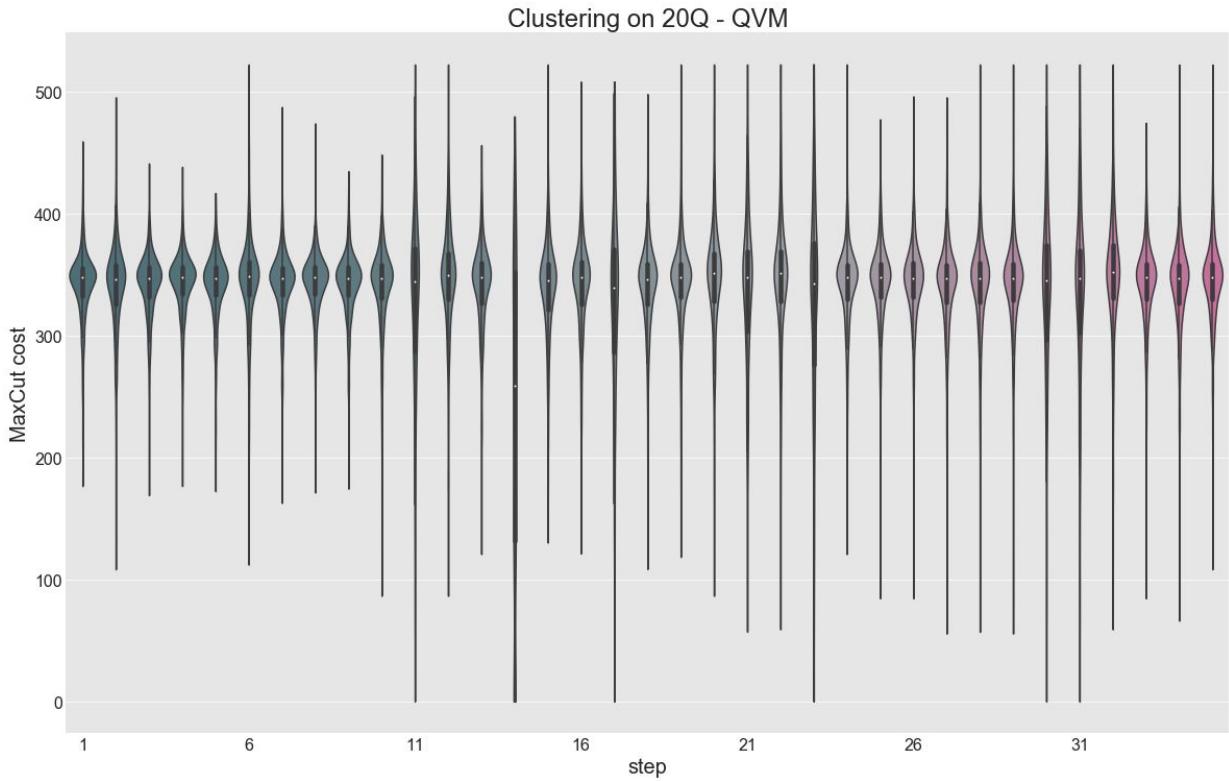
At each step

Sample a new angle pair from  
the Gaussian prior

Run the Quil program and  
sample several bitstrings

Evaluate the MAXCUT cost and  
return maximum as the value

Update Prior and repeat the  
process

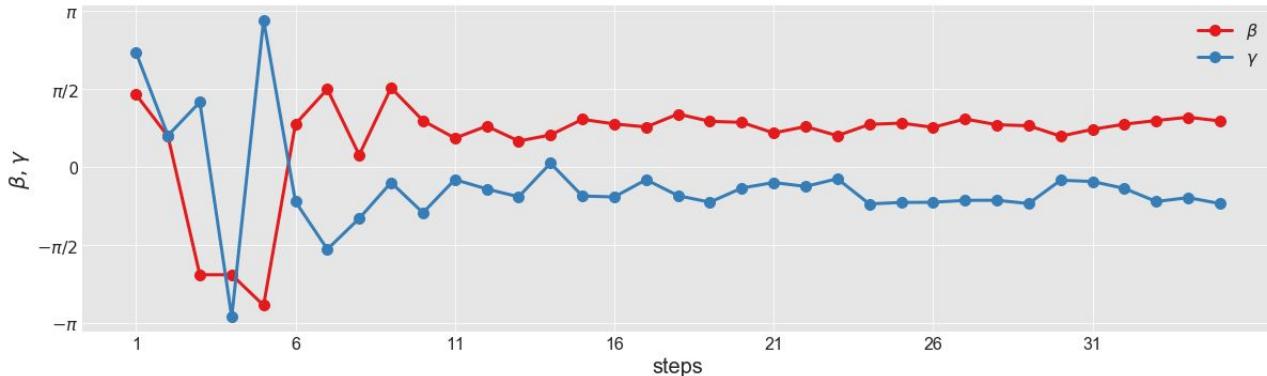
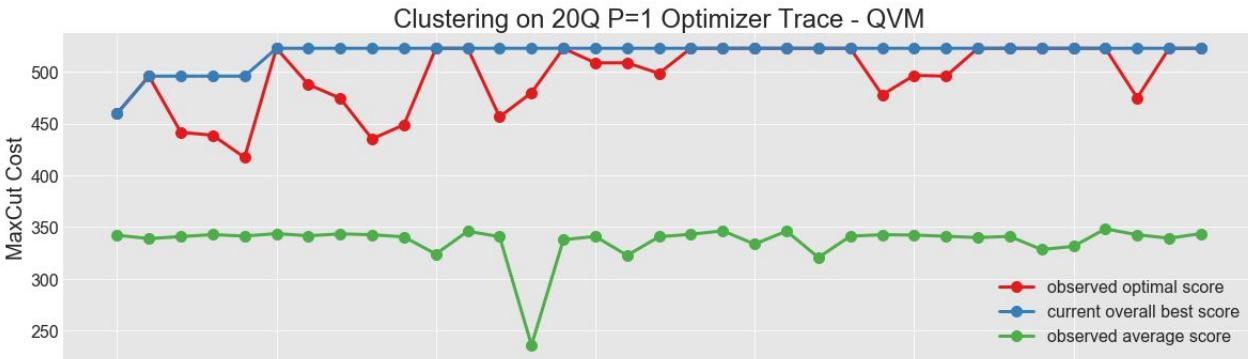


# That's how it looks in practice

At each step

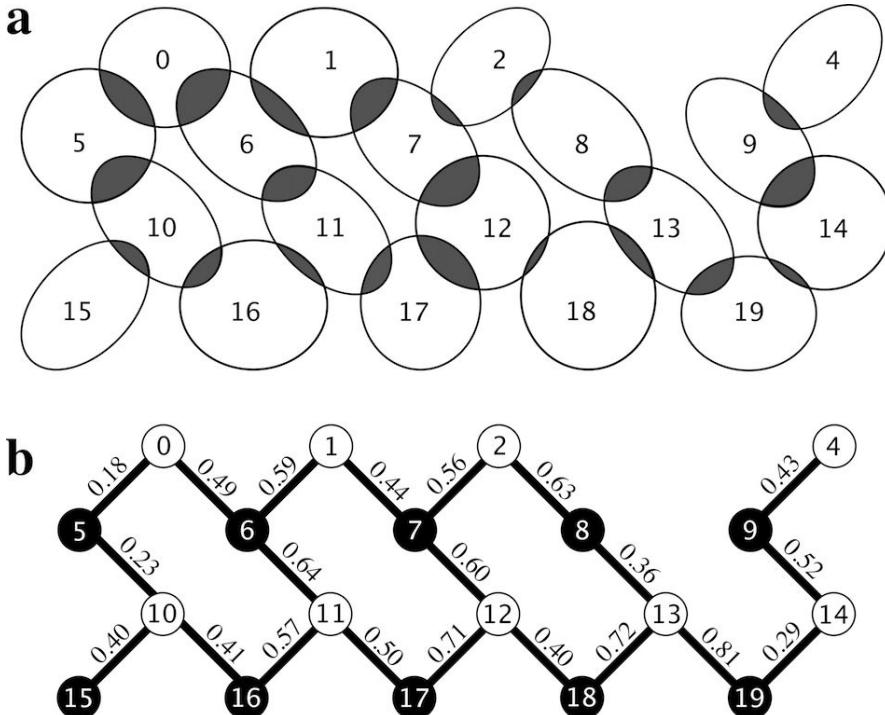
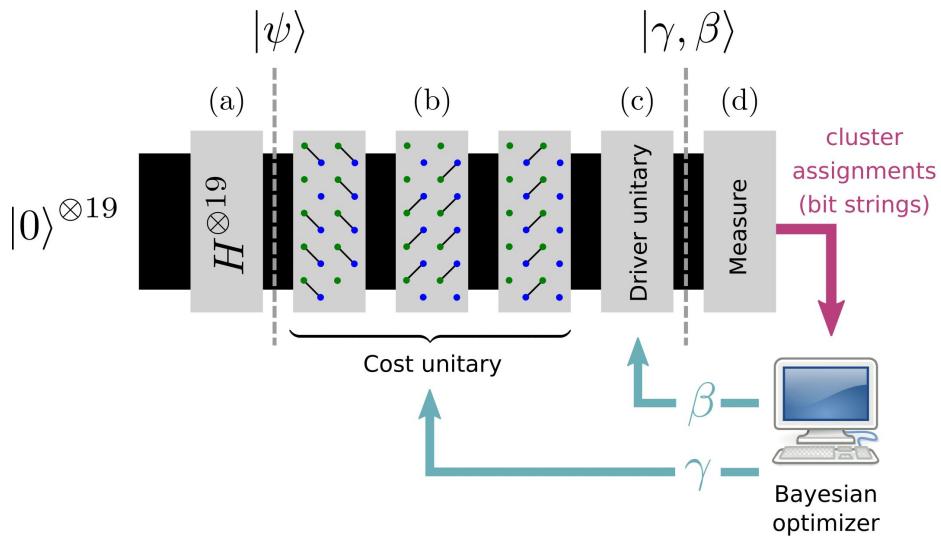
Keep the historic best value

Angles start to converge for large steps numbers

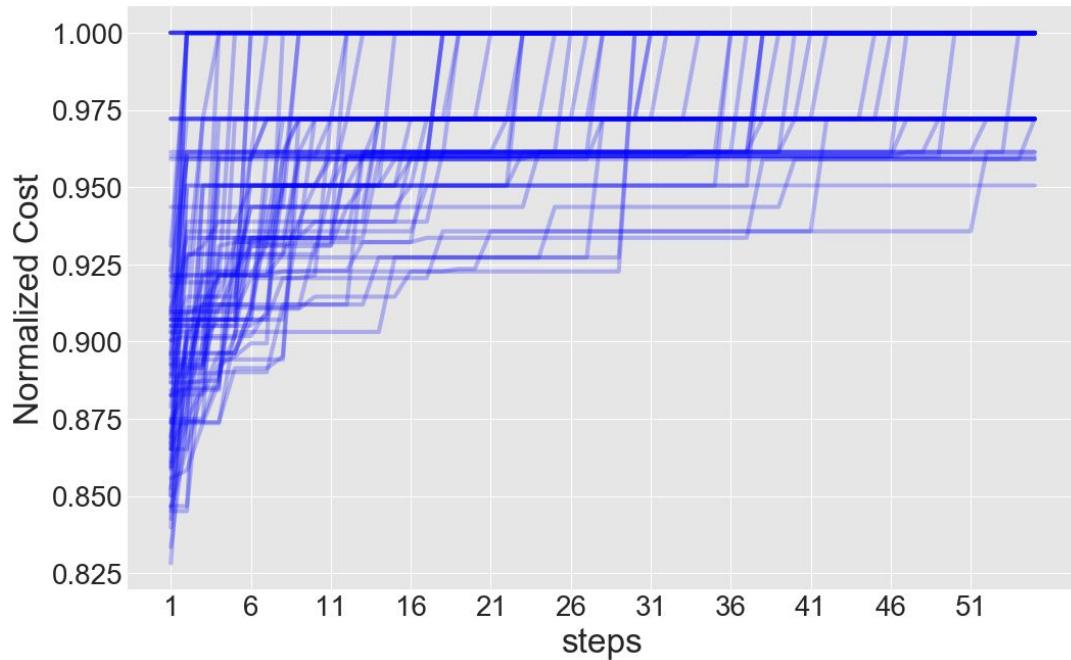


# Clustering on a 19-Q Chip

For demonstration purposes chose problem instance to match chip topology



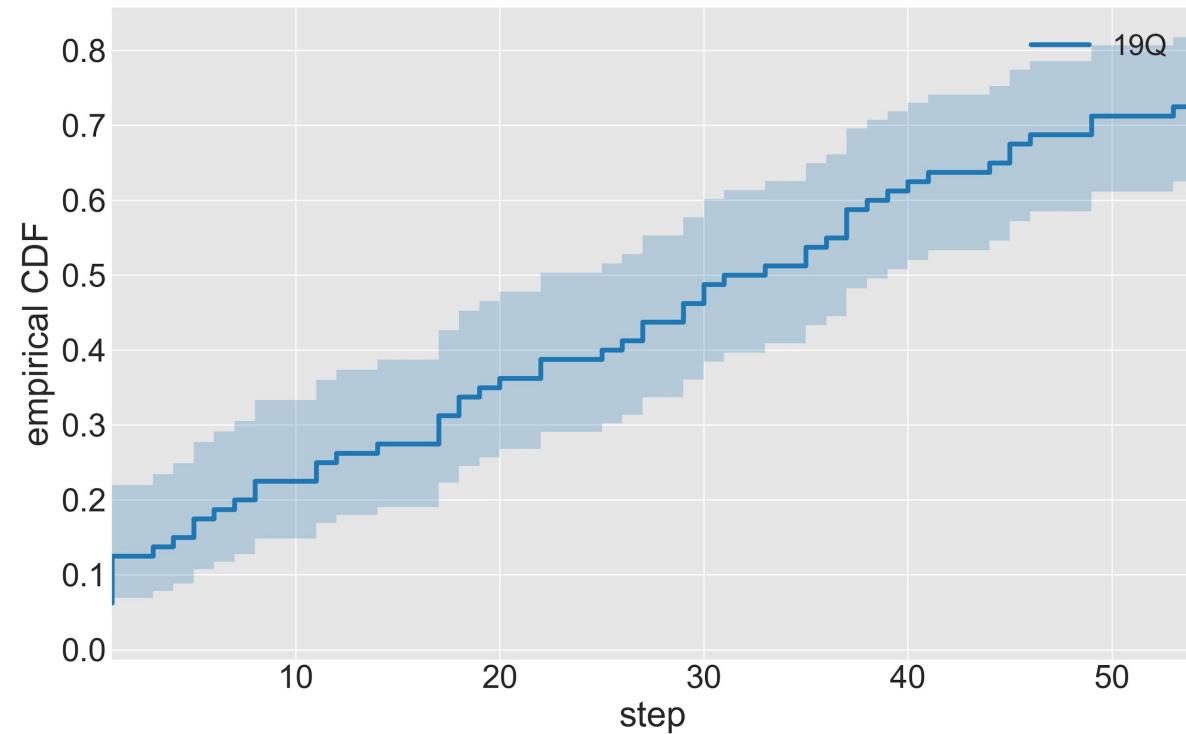
# Putting it all together



- 83 trial runs on the QPU
- Algorithm finds the optimum most of the time
- Calculate success probability from the traces



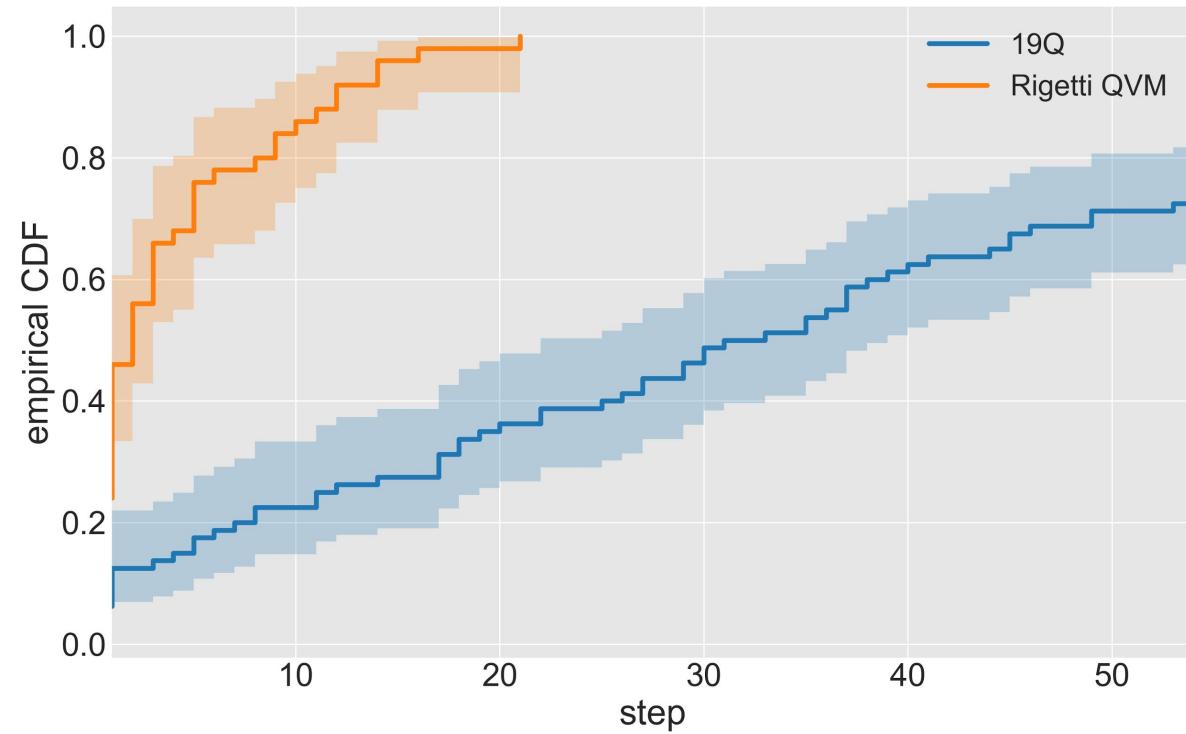
# Empirical performance



- Success probability monotonically increases with number of steps.



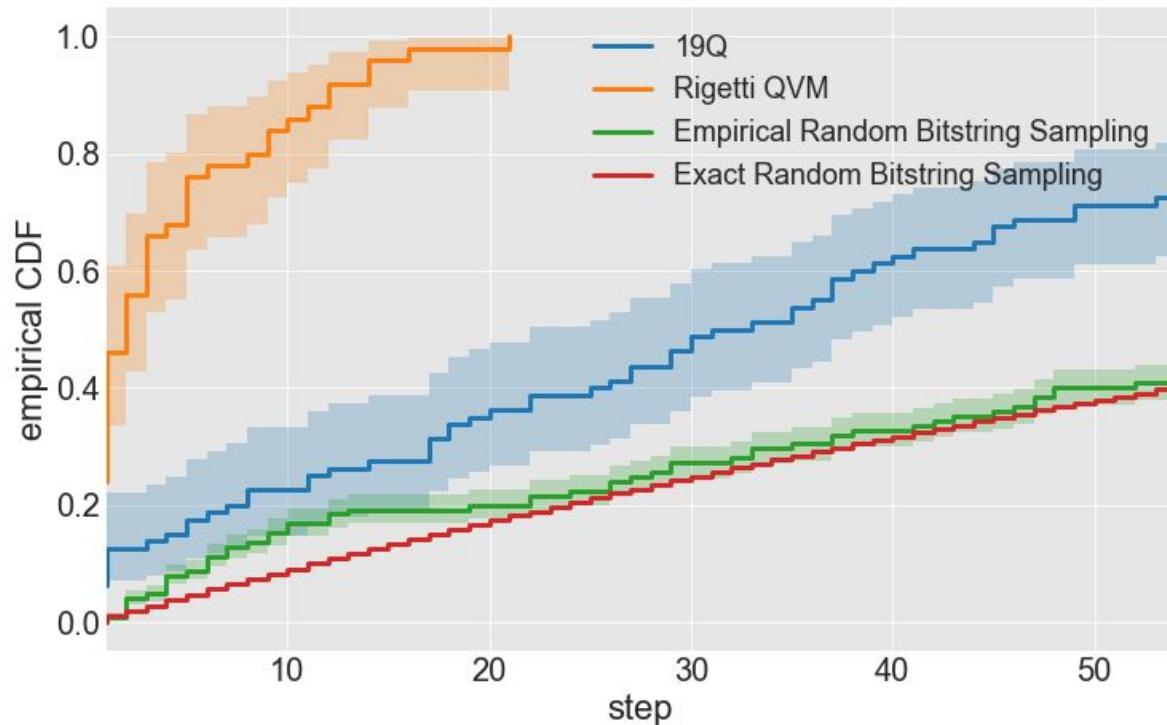
# Empirical performance



- Success probability monotonically increases with number of steps.
- Noise in 19Q has a significant impact on performance.



# Empirical performance



- Success probability monotonically increases with number of steps.
- Noise in 19Q has a significant impact on performance.
- Approach clearly outperforms random sampling.



# Forest



Join our community Slack:

[slack.rigetti.com](https://slack.rigetti.com)



Find us on Github:

[github.com/rigetticomputing](https://github.com/rigetticomputing)

Sign-Up @ [rigetti.com/forest](https://rigetti.com/forest)

QPU access @ [rigetti.com/qpu-request](https://rigetti.com/qpu-request)



# Thank you

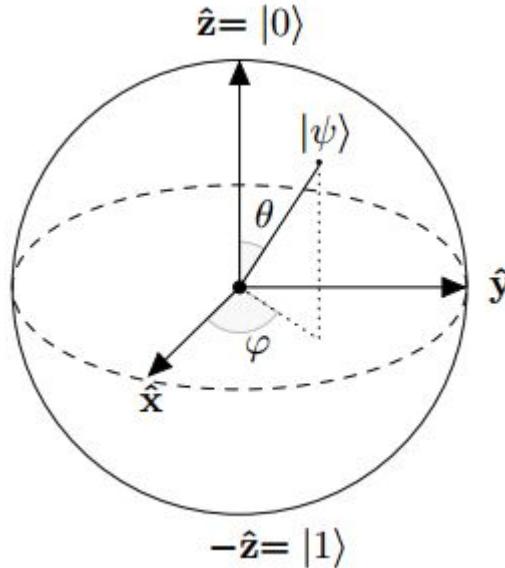
More details in our pre-print arXiv: 1712.05771

# Spare slides



# Bloch Sphere

Lives on the surface of the Bloch sphere



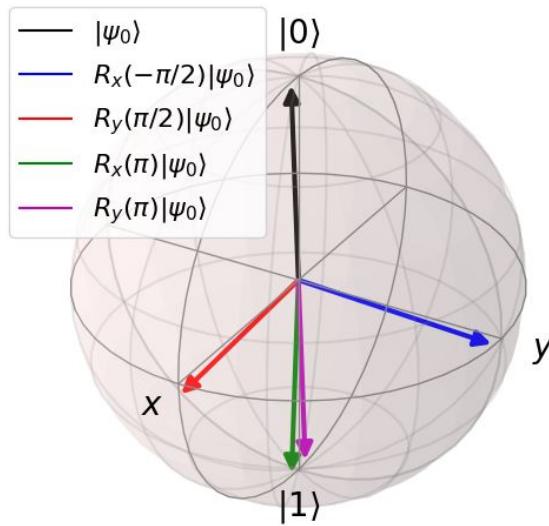
$$|\psi\rangle = \cos(\theta/2)|0\rangle + e^{i\varphi}\sin(\theta/2)|1\rangle$$



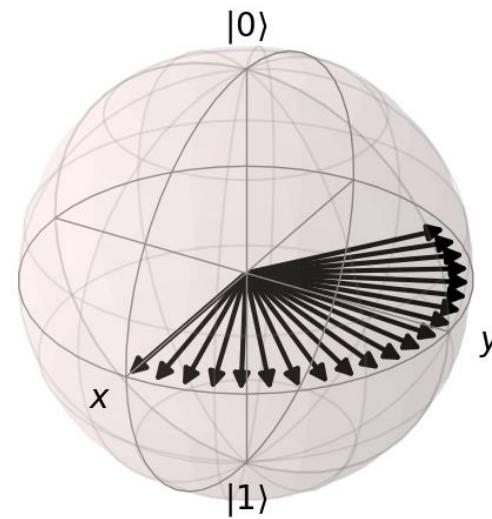
# Quantum Control on the Bloch Sphere

"Machine that natively executes unitary operations on quantum systems"

Unitaries are rotations



X and Y rotations by driving



Z rotations by waiting

