

# Quantum Computing

## Lecture $|16\rangle$ : **Variational Quantum Algorithms**

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# Agenda

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- Variational Algorithms and the NISQ Era
- Variational Quantum Eigensolver and Optimization

# Variational Algorithms

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The quantum circuits seen so far (Grover's, Shor's, *etc.*) depended on the input size:

- the same circuit is used for inputs of some maximum length;
- if the input gets larger, one needs a different, larger circuit;

## Variational Quantum Algorithms:

- quantum circuits are **updated** as a way to solve an optimization problem;
- the circuits are usually small and not exceedingly deep;
- the circuits can be run on NISQ (Noisy Intermediate-Scale Quantum) computers.

# Extremal Eigenvalue Problem

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## Definition (Extremal Eigenvalue Problem)

Given an Hermitian matrix, find its **extremal** eigenvalues.

Given a generic optimization problem where  $C(\cdot)$  is a real cost function and  $S$  is a set representing some constraints

$$\begin{aligned} & \max / \min C(x) \\ & \text{subject to } x \in S \end{aligned}$$

can be reduced to an extremal eigenvalue problem.

# Optimization as an Extremal Eigenvalue Problem

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Assume that the set of solutions  $Y$  is finite, so can use finite bit-strings to label the solutions.

We build the Hermitian operator

$$H_C = \sum_{y \in Y} C(y) |y\rangle\langle y|$$

which implies that  $H_C |a\rangle = C(a) |a\rangle$  for any  $a \in Y$ .

Therefore:

- $\min_x C(x) = \text{minimal eigenvalue of } H_C$ ;
- $\max_x C(x) = \text{maximal eigenvalue of } H_C = \text{minimal eigenvalue of } -H_C$

# Variational Quantum Eigensolver

How to find the minimal eigenvalue of an Hermitian operator?

## Theorem

*Let  $A$  be an Hermitian operator/matrix on an Hilbert space  $\mathcal{H}$  and  $\lambda_{\min}$  its least eigenvalue. Then:*

$$\forall |\psi\rangle \in \mathcal{H} \quad \langle \psi | A \psi \rangle \geq \lambda_{\min}$$

*with equality iff  $|\psi\rangle = |\psi\rangle_{\min}$  (an eigenvector associated to  $\lambda_{\min}$ ).*

Therefore, we can solve our extremal eigenvalue problem by **minimizing** the function  $f : \mathcal{H} \rightarrow \mathbb{R}$  defined as:

$$f(|\psi\rangle) = \langle \psi | A \psi \rangle .$$

Note that  $f$  is a well-behaved function. Problem: the state space is usually huge!

# Variational Quantum Eigensolver: The Main Idea

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To generate a **sequence** of quantum circuits whose output is **close** to  $\langle \psi_{\min} | A | \psi_{\min} \rangle$ .

The quantum circuits are parameterized by a number (say  $p$ ) of **real parameters**.

$$\text{parameters } \theta \in \mathbb{R}^p \xrightarrow{\text{circuit } Q_\theta} \text{evaluate circuit} \xrightarrow{|\psi_\theta\rangle} \text{estimate } \langle \psi_\theta | A | \psi_\theta \rangle$$

How to estimate  $\langle \psi_\theta | A | \psi_\theta \rangle$ ? (Remember: huge state spaces!)

We “simply” measure  $A$  (it is an Hermitian operator, hence a valid observable) multiple times in order to estimate the amplitudes of the basis states.

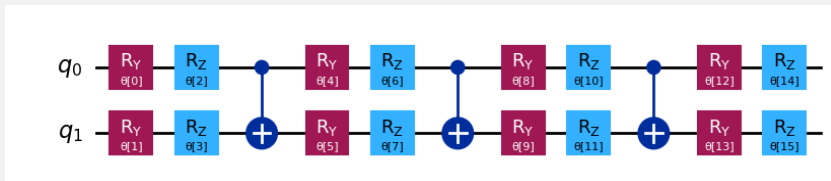
Recall that given a cost function  $C(\cdot)$  we have

$$\langle \psi | H_C | \psi \rangle = \sum_y |\alpha_y|^2 C(y) \quad \text{where } |\psi\rangle = \sum_y \alpha_y |y\rangle$$

# Variational Quantum Eigensolver: The Main Idea

The quantum circuits are built from an **ansatz** (a 'template' or 'educated guess') circuit containing single-qubit parameterized rotations and 2-qubit gates.

An example of ansatz on two qubits:



from <https://learning.quantum.ibm.com/tutorial/variational-quantum-eigensolver>

The ansatz should be able to reach much of the Hilbert space by an appropriate choice of parameters. Choosing the right ansatz is quite an art.



# Variational Quantum Optimization

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We can estimate  $f(|\psi_\theta\rangle) = \langle\psi_\theta|A|\psi_\theta\rangle$ , and since  $f : \mathbb{R}^p \rightarrow \mathbb{R}$  is a classical function, we can use any standard (classical) optimization technique to minimize  $f$ , e.g.:

- stochastic gradient descent;
- simulated annealing;
- particle swarm;
- genetic algorithms;
- etc.

# Variational Quantum Optimization: The Algorithm

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## Algorithm 1: Optimization by VQE

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**Input:** Cost function  $C$ , number of circuit evaluations  $N$

**Output:** An approximation of  $\min C$

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1  $\theta = \theta_0$ ; done = false;  
2 while not done do  
3     generate circuit  $Q_\theta$  from ansatz with parameters  $\theta$   
4     for  $N$  times do  
5          $|\psi_\theta\rangle = Q_\theta |00\dots 0\rangle$   
6         measure  $H_C$  on  $|\psi_\theta\rangle$   
7      $l_\theta = \text{estimate } \langle \psi_\theta | A | \psi_\theta \rangle$  from measurements  
8     if classical optimization algorithm decides  $l_\theta$  is OK then  
9         done = true  
10    else  
11        update  $\theta$ 
```