

# **CPEN 400Q / EECE 571Q Lecture 18**

## **Error mitigation, and introducing QAOA**

Thursday 17 March 2022

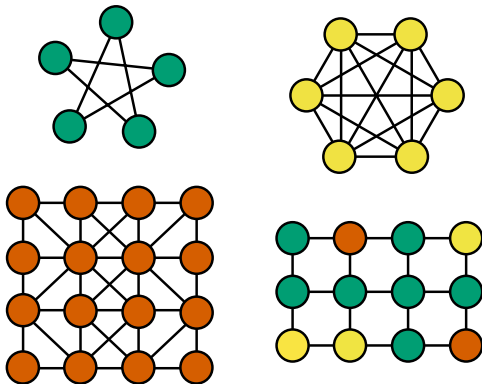
- Assignment 4 coming (will be due at end of term)

## Last time

We learned how to use trace distance ( $T$ ) and fidelity ( $F$ ) to compare two density matrices:

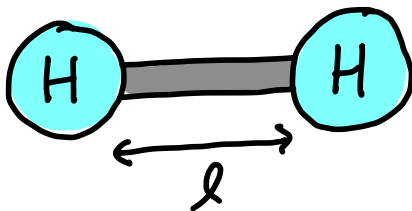
## Last time

We discussed some ways by which quantum computers are compared.



## Last time

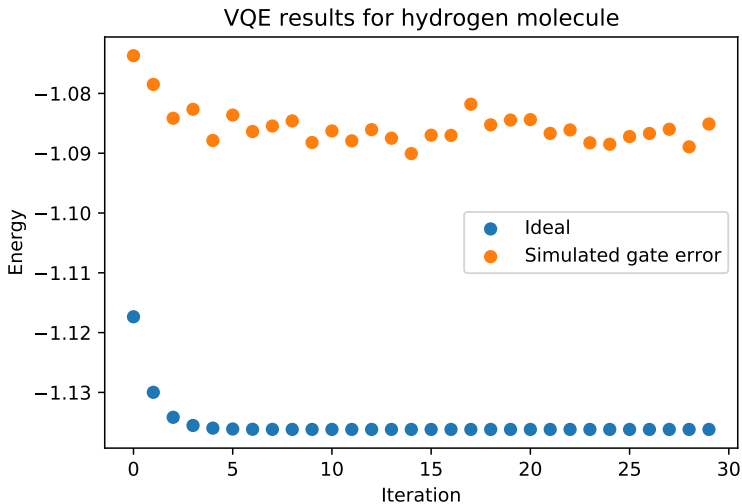
We used VQE to solve a small quantum chemistry problem:



The Hamiltonian uses four qubits, and we needed only a single variational parameter to produce its ground state,

## Last time

Then we tried running VQE on a simulated noisy device...



- Mitigate noise using zero-noise extrapolation
- Perform basic quantum state tomography
- Describe the underlying ideas of adiabatic quantum computation, and the quantum approximate optimization algorithm

Current quantum hardware is noisy. Noise comes from a variety of sources, and depends on the qubit technology and architecture.

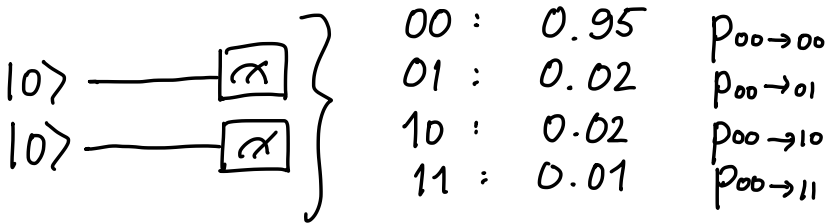
We need to do a combination of:

- Better-characterizing the behaviour of devices to learn how to improve their operation
- Processing the results to mitigate the effects of noise as much as possible



# Measurement error mitigation

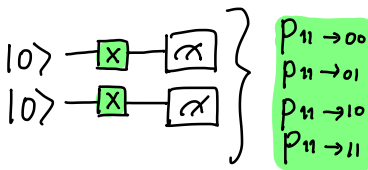
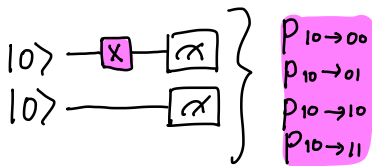
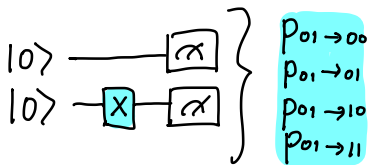
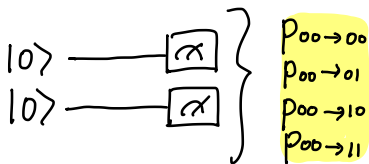
Errors can occur in the measurement process where states are read out incorrectly.



These kinds of errors are quite straightforward to mitigate.

# Measurement error mitigation

Check what happens with all possible input states:



## Measurement error mitigation

We can put the results from our calibration circuits into a matrix:

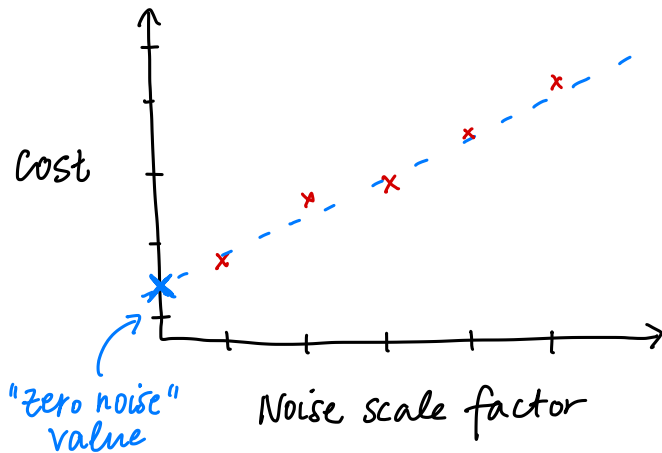
$$M = \begin{bmatrix} p_{00 \rightarrow 00} & p_{01 \rightarrow 00} & p_{10 \rightarrow 00} & p_{11 \rightarrow 00} \\ p_{00 \rightarrow 01} & p_{01 \rightarrow 01} & p_{10 \rightarrow 01} & p_{11 \rightarrow 01} \\ p_{00 \rightarrow 10} & p_{01 \rightarrow 10} & p_{10 \rightarrow 10} & p_{11 \rightarrow 10} \\ p_{00 \rightarrow 11} & p_{01 \rightarrow 11} & p_{10 \rightarrow 11} & p_{11 \rightarrow 11} \end{bmatrix}$$

Can suppose that the probability vector  $P_{noisy}$  we get at the end of a quantum algorithm is related to the ideal one,  $P_{ideal}$ , under multiplication by  $M$  since that's what we see:

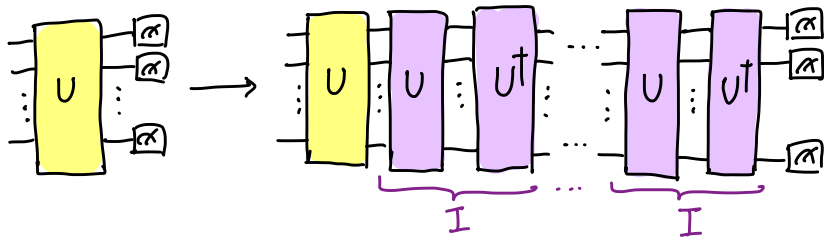
So to get the ideal results:

## Gate error mitigation: zero-noise extrapolation

Modify the circuits to systematically *increase* scale of the noise, then extrapolate down to the zero-noise limit.



## Unitary folding

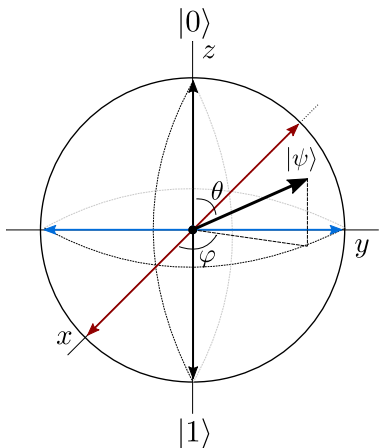


Let's code up a very basic version of this in PennyLane.

More sophisticated version: Python package `mitiq`  
<https://github.com/unitaryfund/mitiq>

# Quantum state tomography

In order to quantify how close the state obtained from a noisy process is from the true state, we need a way of determining what that state is (i.e., its density matrix). We can reconstruct a state by taking an *informationally complete* set of measurements.



Measure:

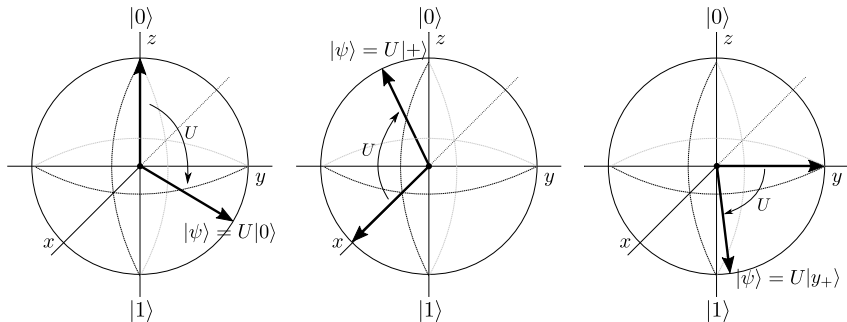
$$X = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}$$

$$Y = \begin{pmatrix} 0 & -i \\ i & 0 \end{pmatrix}$$

$$Z = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}$$

# Quantum process tomography

Similar method for learning about quantum processes: reconstruct an operation based on how it acts on known states.



(Example: a unitary operation on a single qubit.)

## Quantum tomography: MUBs

Tomography is, in some sense, a “solved problem”: we know what **optimal sets of measurements** look like for most cases. (The amount of them scales exponentially in the number of qubits...)

In dimension  $d$ , two bases,  $A = \{|a_i\rangle\}$  and  $B = \{|b_i\rangle\}$  are *mutually unbiased* if for all  $|a_i\rangle, |b_j\rangle$ ,

A complete set of  $d + 1$  mutually unbiased bases (**MUBs**) comprises an optimal set of measurement bases.



## Quantum tomography: MUBs

**Pro:** systematic method of construction using finite fields and the Pauli group

**Con:** complete sets are only known in systems with  $d$  prime or power-of-prime.

Example:  $d = 2^2$  (2-qubit case). Partition Paulis into  $d + 1$  sets of  $d - 1$  commuting operators; their shared eigenbases are the MUBs.

Set ID	Paulis		
1	$ZZ$	$IZ$	$ZI$
2	$XX$	$IX$	$XI$
3	$YY$	$IY$	$YI$
4	$XY$	$YZ$	$ZX$
5	$YX$	$XZ$	$ZY$

## Quantum tomography: SIC-POVMs

Alternative: symmetric, informationally complete positive operator-valued measures (**SIC-POVMs**).

In dimension  $d$ , this is a POVM with  $d^2$  elements  $\{\Pi_i\}^1$  and some additional special properties:

- all  $\Pi_i$  are *rank-1 projectors*, i.e.,  $\Pi_i = |\psi_i\rangle\langle\psi_i|$
- all pairwise inner products are equal:

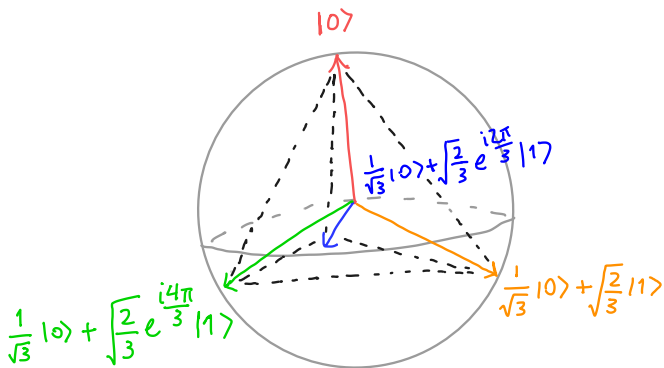
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<sup>1</sup>Recall that for a POVM,  $\sum_i \Pi_i = I$ .

# Quantum tomography: SIC-POVMs

**Pro:** unlike MUBs, conjectured to exist in every dimension

**Con:** no one has been able to *prove* this (if you can do so, you would be famous)



Let's carry out a very simple state reconstruction on our noisy state.

The  $d$ -dimensional Pauli operators are a basis, and are orthogonal w.r.t. the inner product  $\text{Tr}(P_i P_j) = d\delta_{ij}$ .

## Quantum tomography

What happens when we multiply  $\rho$  by a particular Pauli operator  $P_j$  and take the trace?

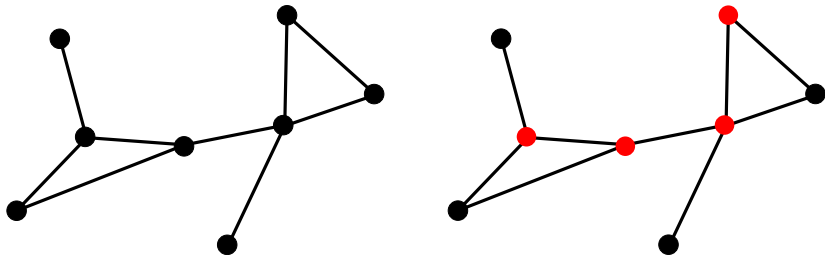
A simple way to reconstruct the state is to compute the expectation value of every Pauli operator for the final state.

Let's try it.

# Quantum approximate optimization algorithm

# Combinatorial optimization

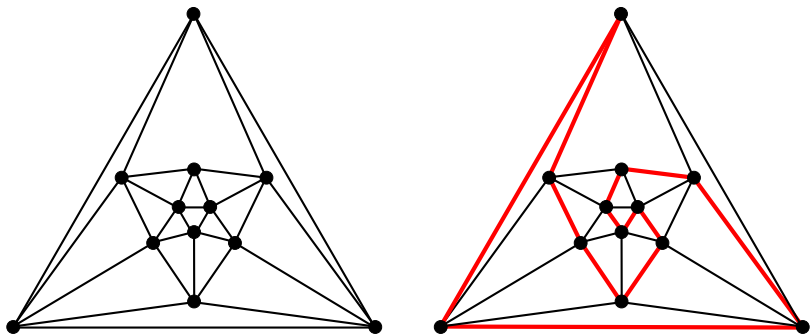
Example: Given a graph  $G = (V, E)$ , what is the *smallest number of vertices* you can colour such that every edge in the graph is attached to at least one coloured vertex?





# Combinatorial optimization

Example: Given a graph, can we find a path through it that visits every node *exactly once* and returns to the starting point?



(In graph theory terms, can we find a *Hamiltonian cycle*?)

# Combinatorial optimization

Example: You have the opportunity to purchase 100 units of stocks from a fixed list of assets. You know the average returns of each stock, and their covariances.

Stock	Avg. return
AAA	3.44 %
BBB	2.21 %
CCC	-0.28 %
$\vdots$	$\vdots$

Cov.	AAA	BBB	...
AAA	0.0038	0.002	...
BBB	0.002	-0.006	...
CCC	0.014	-0.0008	...
$\vdots$	$\vdots$	$\vdots$	$\ddots$

Suppose you're restricted to buying no more than 5 of any stock.

*Which stocks, and how many of each, should you purchase, to maximize your profits?*

# Adiabatic quantum computing (AQC)

The structure of a classical optimization problem is something like:

$$\min_{\vec{x}} \text{cost}(\vec{x}) \quad \text{subject to constraints}(\vec{x})$$

where  $\vec{x}$  is a multi-dimensional vector of parameters in the problem space.

# Adiabatic quantum computing (AQC)

The structure of a classical optimization problem is something like:

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where  $\vec{x}$  is a multi-dimensional vector of parameters in the problem space.

In a physical context, optimization can be interpreted as an energy minimization problem.

Optimization	Physical system
$\vec{x}$	State of the system
$\text{cost}(\vec{x})$	Hamiltonian
Optimum $\vec{x}^*$	Ground state
$\text{cost}(\vec{x}^*)$	Ground state energy

# Adiabatic quantum computing (AQC)

Recall that every unitary  $U$  is directly related to a Hermitian Hamiltonian  $H$  under the correspondence

We know that we can use gate model QC to *simulate* the evolution of a Hamiltonian.

Instead of simulating the Hamiltonians, **adiabatic quantum computing** works with them directly to perform computations. It is generally used to solve **optimization problems**.

# Adiabatic quantum computing (AQC)

1. Design a cost Hamiltonian whose ground state represents the solution to our optimization problem
2. Prepare a system in the ground state of an easy-to-prepare mixer Hamiltonian
3. Perform adiabatic evolution to transform the system from the ground state of mixer Hamiltonian to the ground state of the cost Hamiltonian, which is our solution

# The adiabatic theorem

Why would we want to do this?

## Theorem:

*"A physical system remains in its instantaneous eigenstate if a given perturbation is acting on it slowly enough and if there is a gap between the eigenvalue and the rest of the Hamiltonian's spectrum."*

## What we can take from this:

If we initialize a system in the lowest energy state and perturb it slowly enough, it will remain in the lowest energy state (with respect to the changed system)

# Adiabatic quantum computing (AQC)

Let  $H_m$  be a **mixer Hamiltonian** whose ground state can be easily prepared.

Let  $H_c$  be a **cost Hamiltonian** whose ground state represents the solution to a problem of interest.

Adiabatic evolution is expressed mathematically as the function

The parameter  $s$  is representative of time;  $s$  goes from 0 to 1;  $A(s)$  decreases to 0 with time and  $B(s)$  increases from 0.



# Quantum annealing

D-Wave makes **quantum annealers**: these are a physical implementation of AQC for a limited set of Hamiltonians.



Image credit:

[www.dwavesys.com/tutorials/background-reading-series/introduction-d-wave-quantum-hardware](http://www.dwavesys.com/tutorials/background-reading-series/introduction-d-wave-quantum-hardware)

# Quantum approximate optimization algorithm (QAOA)

QAOA is a gate-model algorithm that can obtain approximate solutions to combinatorial optimization problems.

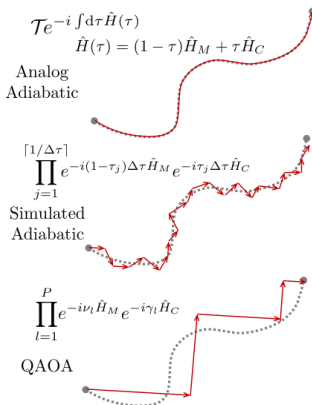
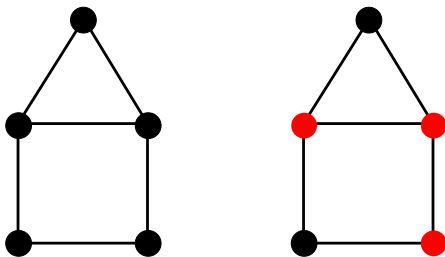


Image credit: G. Verdon, M. Broughton, J. Biamonte. *A quantum algorithm to train neural networks using low-depth circuits*. <https://arxiv.org/abs/1712.05304>

## Motivating example: vertex cover

How do we turn an optimization problem for some graph into a Hamiltonian?

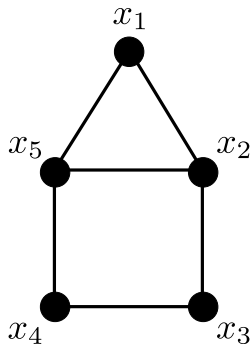
Let's start small, and consider the problem of vertex cover of a graph  $G = (V, E)$ .



First, we will define a cost function, whose minimum cost will correspond to the optimal set of vertices to colour. Then, we will turn it into a Hamiltonian.

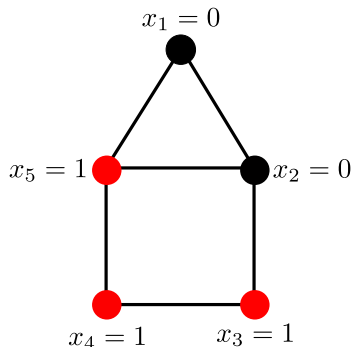
## Motivating example: vertex cover

Whether or not a vertex is coloured is a *binary variable*.



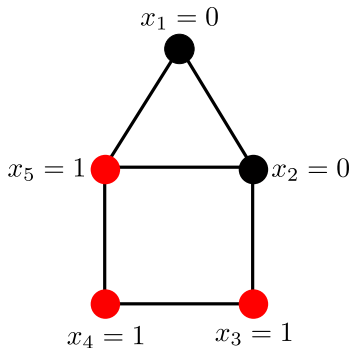
## Motivating example: vertex cover

Let's assign coloured vertices to have value 1, and un-coloured 0.



Now that we have our variables, how do we come up with a minimizable cost function that represents the problem?

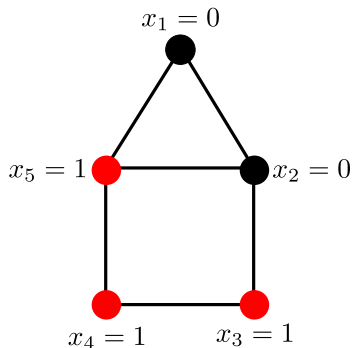
## Motivating example: vertex cover



We need every edge to be next to a coloured vertex. Design a cost function that penalizes edges that are not, but favours ones that are.

Intuitively, find a function of two vertices that is 0 if the colouring is valid, and 1 if it is not.

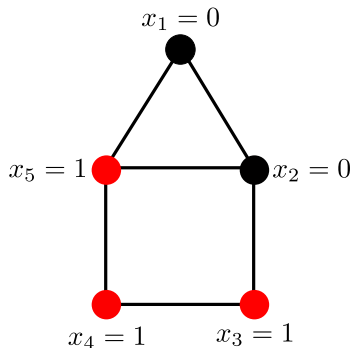
## Motivating example: vertex cover



Consider for each edge  $ij$  the function

The possible values are:

## Motivating example: vertex cover



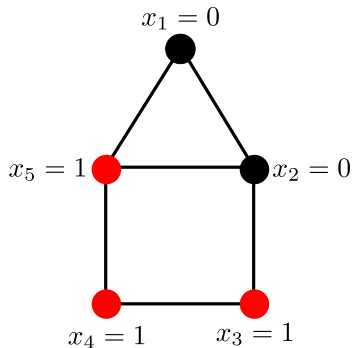
Then in an optimal colouring,

for all edges  $ij \in E$ .

So we can write



## Motivating example: vertex cover



However recall that we also want to colour the fewest vertices. The cost should also depend on the number of coloured vertices.

Solution: add to our cost

## Motivating example: vertex cover

The full cost function is then

1. How do we turn this into a Hamiltonian?
2. How do we find its minimum energy / configuration on a quantum computer?

## Hamiltonian translation

$$\min_{\vec{x}} \left( \sum_{ij \in E} (1 - x_i)(1 - x_j) + \sum_{i \in V} x_i \right)$$

First thing to consider is the problem domain:  $x_i$  are binary variables. We have qubits, which can be  $|0\rangle$  and  $|1\rangle$ .

But since we want to turn this into a Hamiltonian and compute a cost (i.e., measure its expectation value), it's more straightforward to map 0 and 1 to *expectation values* associated to  $|0\rangle$  and  $|1\rangle$ .

# Hamiltonian translation

Usually we consider expectation values of Pauli  $Z$ .

We will make the mapping

This associates  $x_i = 0$  to  $z_i = 1$  (corresponds to  $|0\rangle$ ), and  $x_i = 1$  to  $z_i = -1$  (corresponds to  $|1\rangle$ ).

## Hamiltonian translation

Let's expand our cost function and make this substitution.

# Hamiltonian translation

Substitute:

Expand:

Collect:

## Hamiltonian translation

Consider now that: the total number of edges and vertices are constant - they will provide only an “offset” to the cost, and the values of the variables don’t matter.

And finally, the absolute value doesn’t matter, so we can rescale:

# Hamiltonian translation

Can also weight the terms differently depending on which constraint is more important (i.e., if you care more about just getting a valid colouring, weight the first one more).

To turn this into a Hamiltonian, recall that

- Each  $z_i$  represents an expectation value of  $Z_i$
- Computing expectation values is linear



Next time: we will look at the actual QAOA that can find the optimal configuration / minimum energy.

# Next time

## Content:

- Continue with QAOA

## Action items:

1. Final project

## Recommended reading:

- Qiskit tutorial on measurement error mitigation:  
<https://qiskit.org/textbook/ch-quantum-hardware/measurement-error-mitigation.html>
- mitiq documentation, for more fun error mitigation: <https://mitiq.readthedocs.io/en/stable/guide/guide.html>
- New preprint, *Error mitigation increases the effective quantum volume of quantum computers*.  
<https://arxiv.org/abs/2203.05489>