

# Quadratic optimization with quantum computing

Biweekly Presentation IV

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# Mathematical background

- Expected value of a matrix:  $\langle \mathbf{b} | \mathbf{Q} | \mathbf{b} \rangle$

$$\langle 10010 | \mathbf{Q} | 10010 \rangle = \text{scalar}$$

- $\mathbf{Q}$  is a symmetric matrix (n x n)
- $\mathbf{b}$  is a **binary** vector (n)

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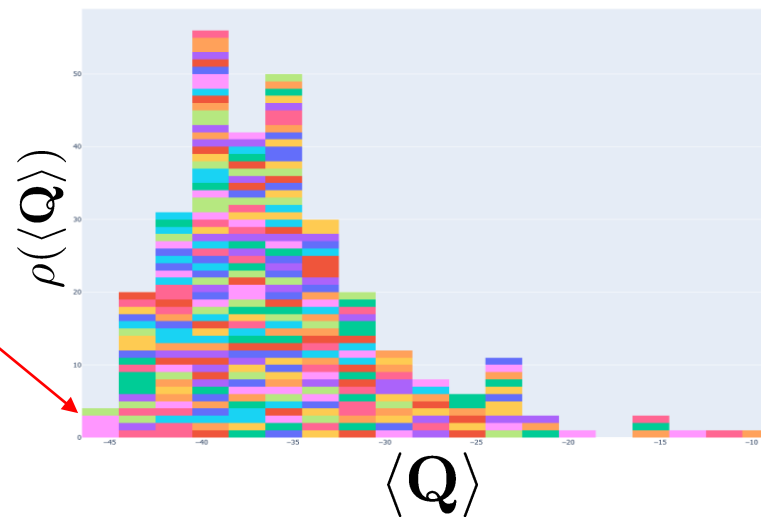
$$\langle 10010 | \mathbf{Q} | 10010 \rangle = \text{scalar}$$

- $\mathbf{Q}$  is a symmetric matrix ( $n \times n$ )
- $\mathbf{b}$  is a **binary** vector ( $n$ )

- Find  $|\mathbf{b}\rangle$  such that  $\min(\langle \mathbf{b} | \mathbf{Q} | \mathbf{b} \rangle)$

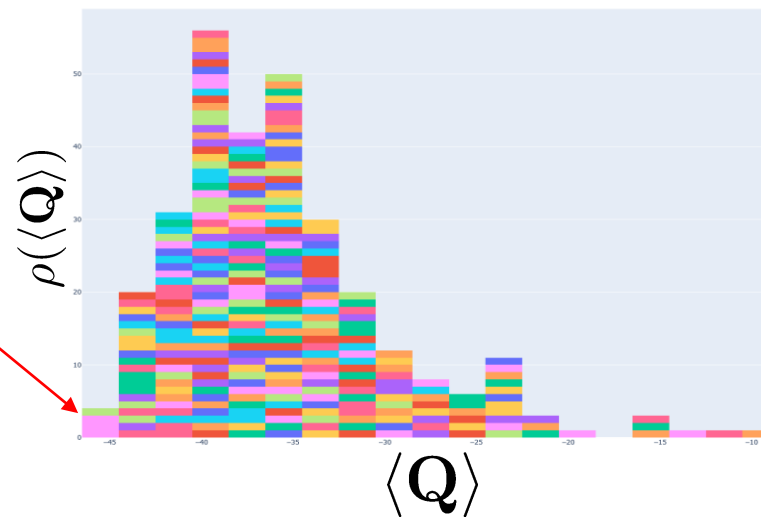
- Explicitly (calculate **all**  $\langle \mathbf{Q} \rangle$ )

- ★ • By **sampling**  $\rho(\langle \mathbf{Q} \rangle)$



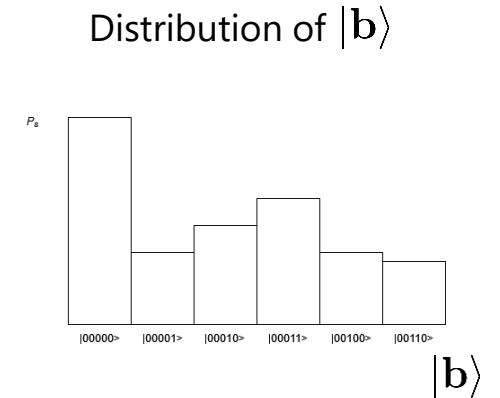
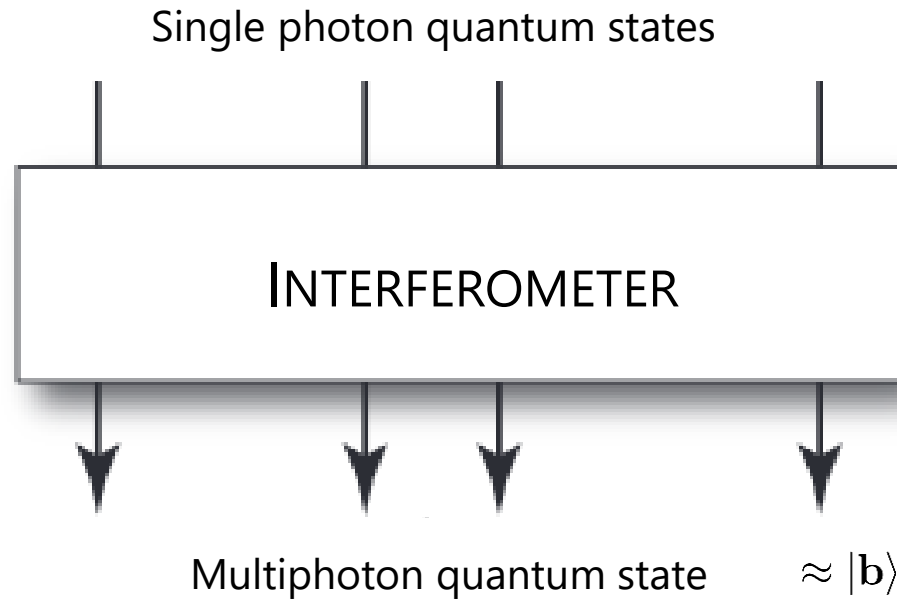
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- Find  $|\mathbf{b}\rangle$  such that  $\min(\langle \mathbf{b} | \mathbf{Q} | \mathbf{b} \rangle)$ 
  - Explicitly (calculate **all**  $\langle \mathbf{Q} \rangle$ )
  - ★ • By **sampling**  $\rho(\langle \mathbf{Q} \rangle)$
- Better than random sampling?
  - Optimizing is hard, because  $|\mathbf{b}\rangle$  is discrete
  - Need to find **continuous parameters** to repr  $|\mathbf{b}\rangle$



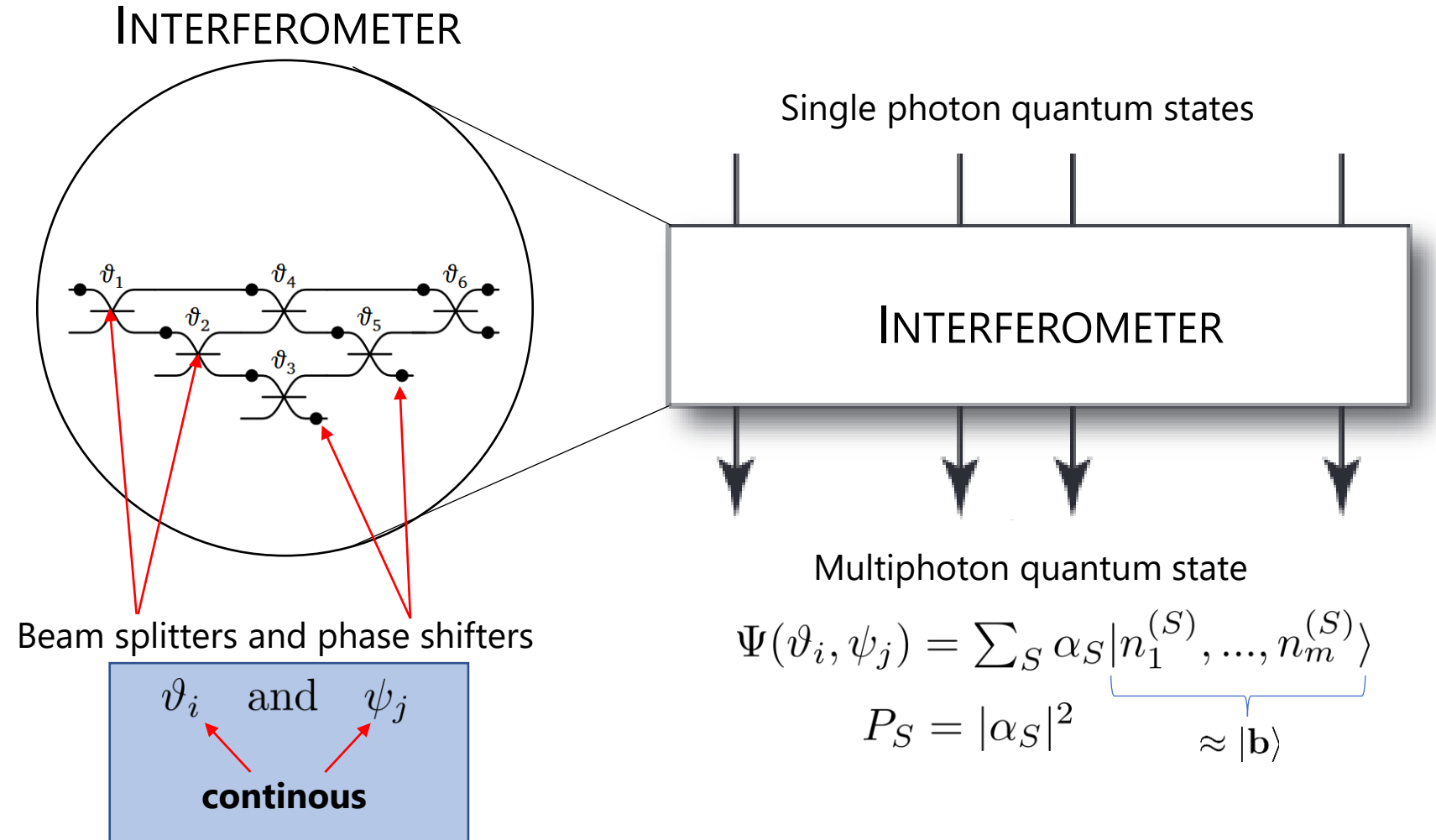
# Boson sampling

- Need to find **continuous parameters** to repr  $|b\rangle$



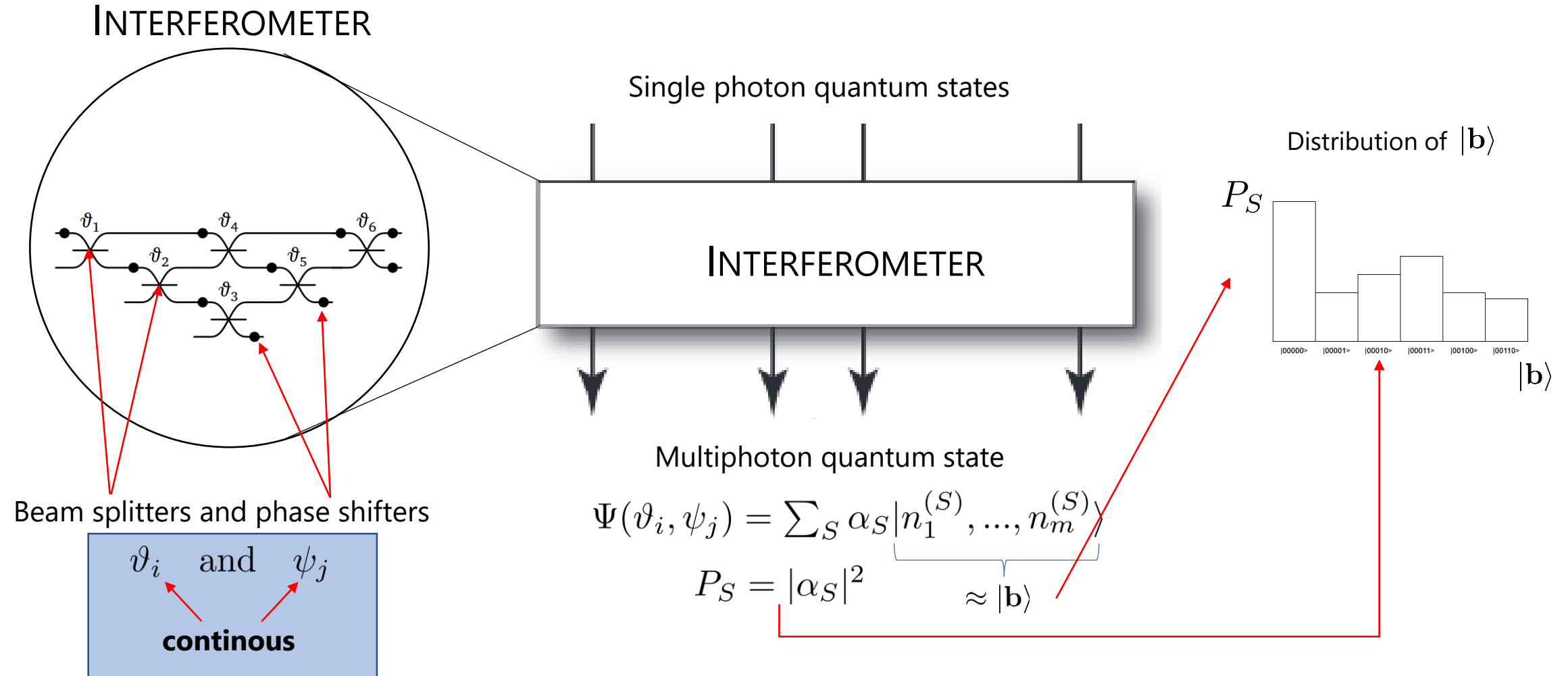
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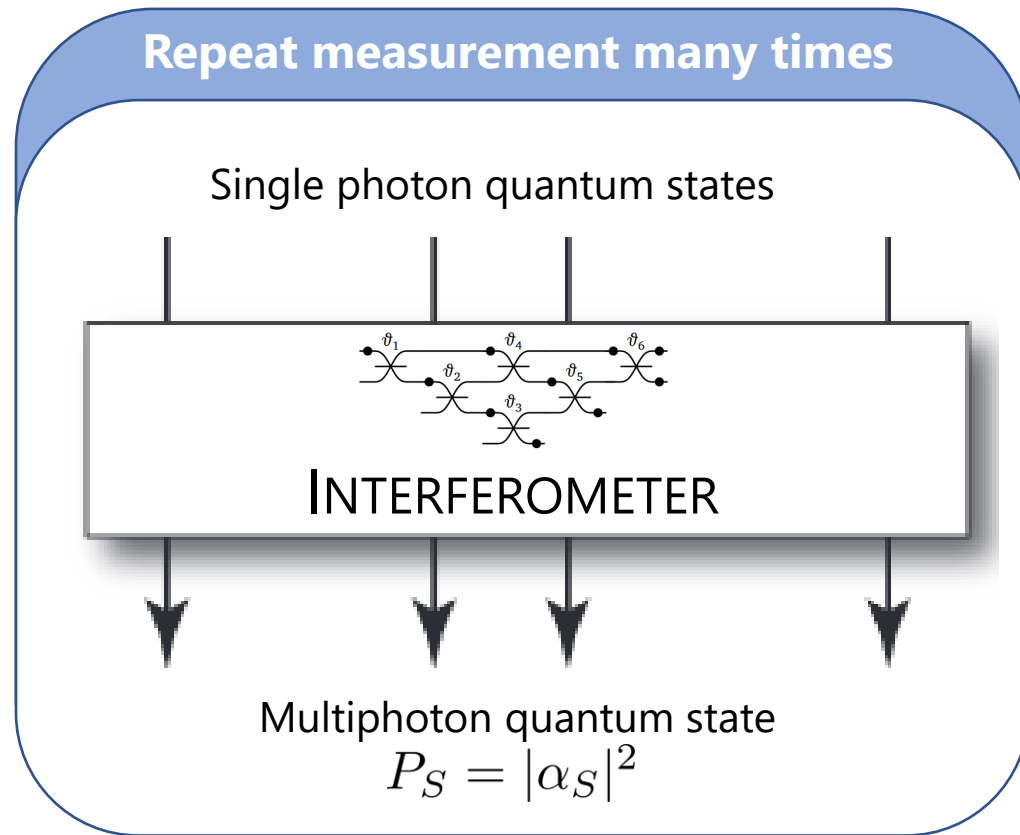
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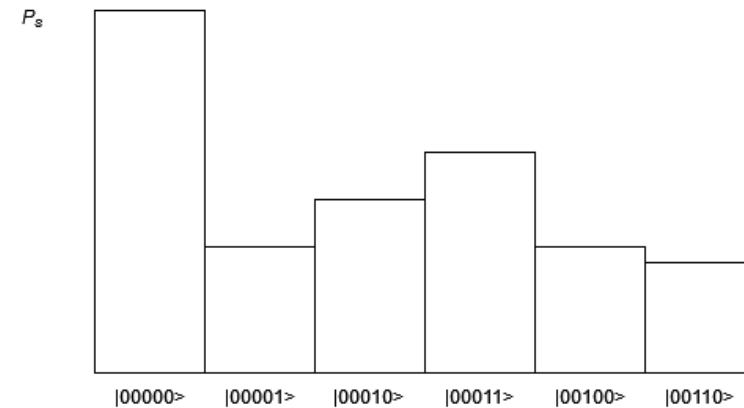


Continuous variables  $\vartheta_i$  and  $\psi_j$



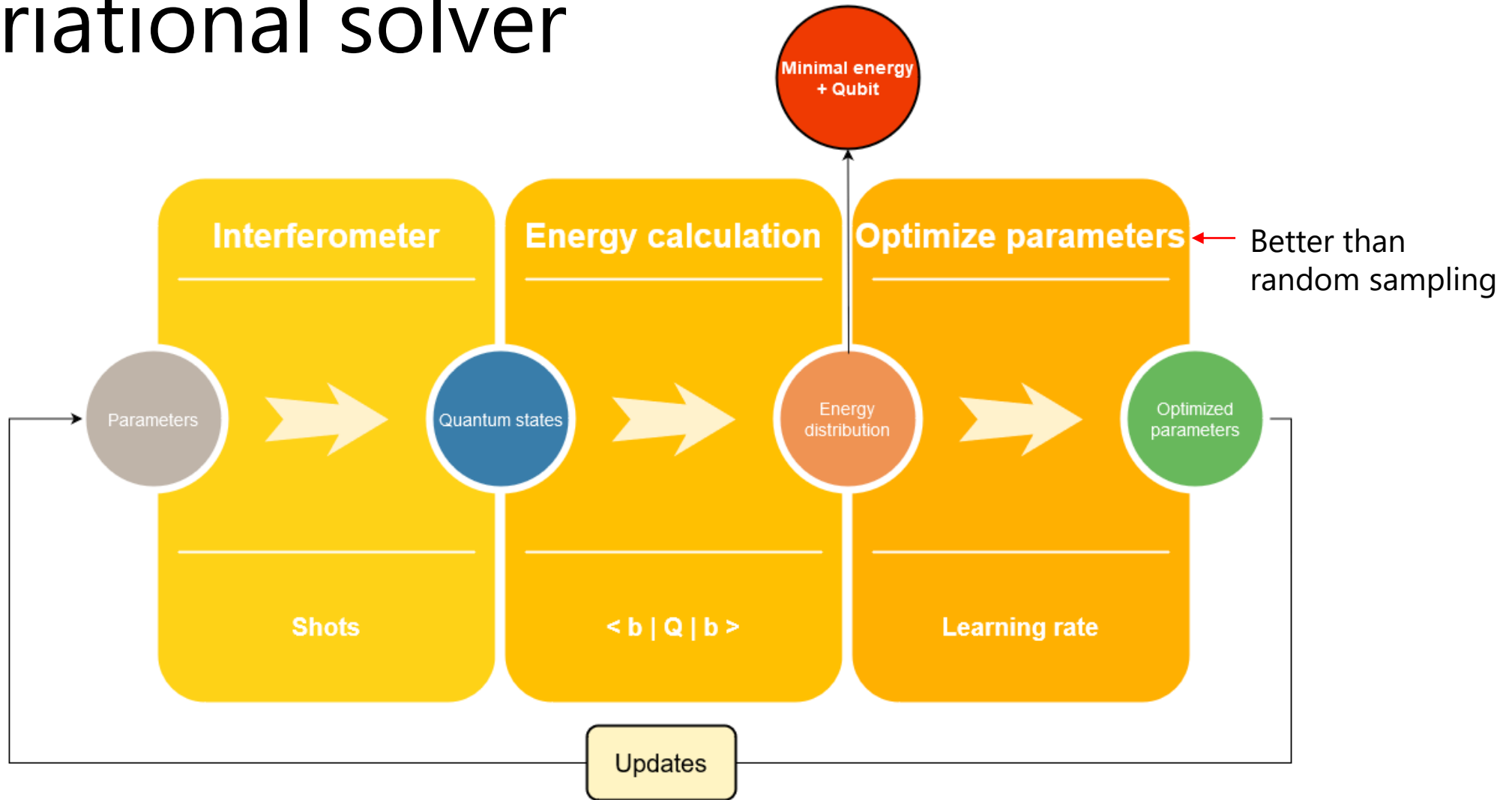
Distribution of quantum states:

$$P_S(\vartheta, \psi) = |\alpha_S|^2$$





# Variational solver



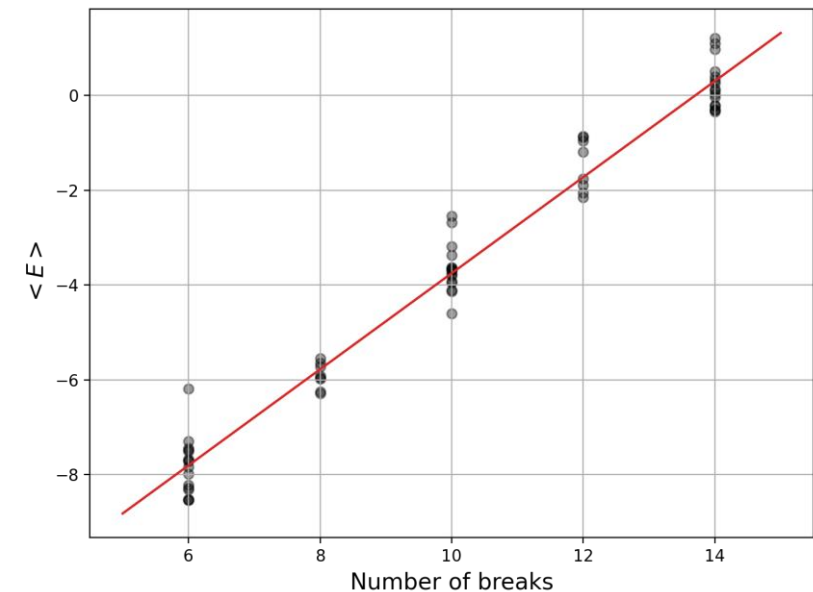
# Setup

- **Piquasso** – simulator for photonic quantum computations
- **Piquassoboost** – for performance improvement
- Personal use:
  - **Ubuntu on Docker** for a separated application environment
- Simulating a Boson Sampler is computationally expensive
- **Budapest Quantum Computing Group** server:
  - 64 Core CPU
  - FPGA server

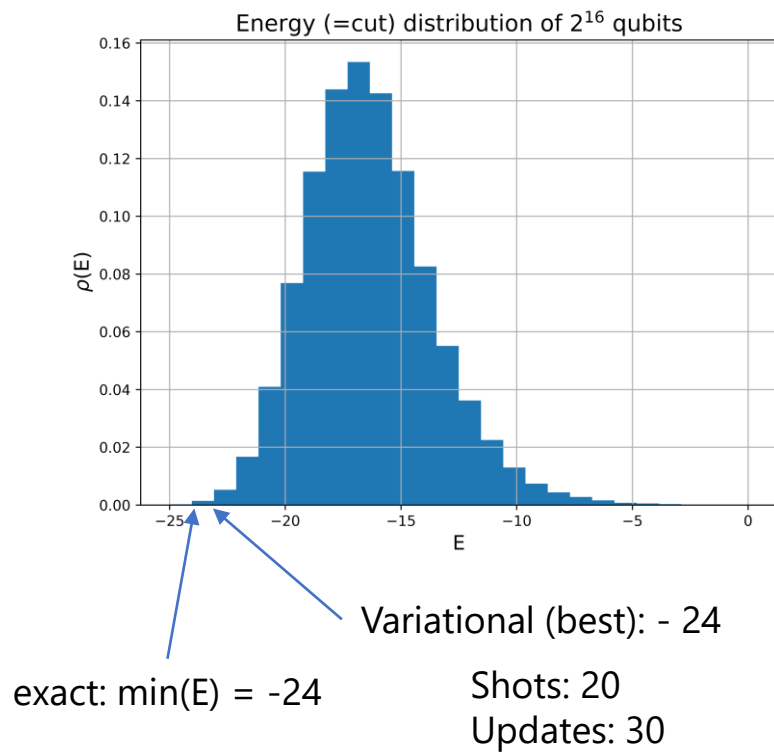


# Results: break minimization

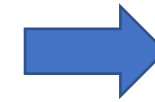
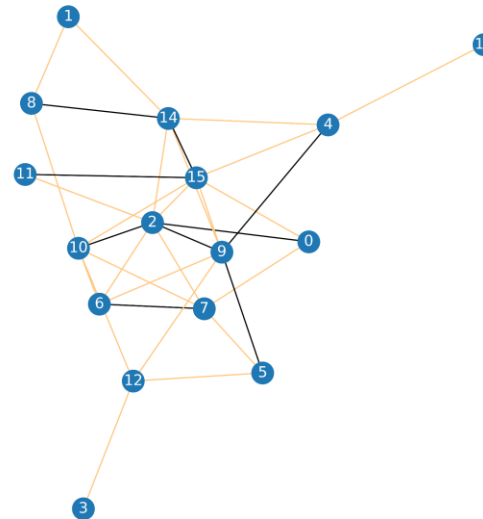
- **Break:** team plays at the same place two times in a row
- Compared **energy** and **number of breaks**
- Expected a linear connection between breaks and energy ( $\langle \mathbf{Q} \rangle = \langle \mathbf{E} \rangle$ )
- **The lowest energy** configuration paired with a min-break → **quantum annealing** is viable
- The global minimum (6) was successfully **found** with the **Piquasso** model



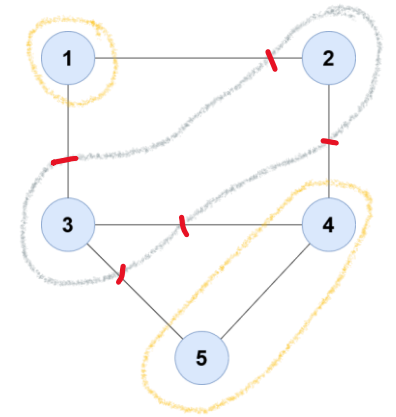
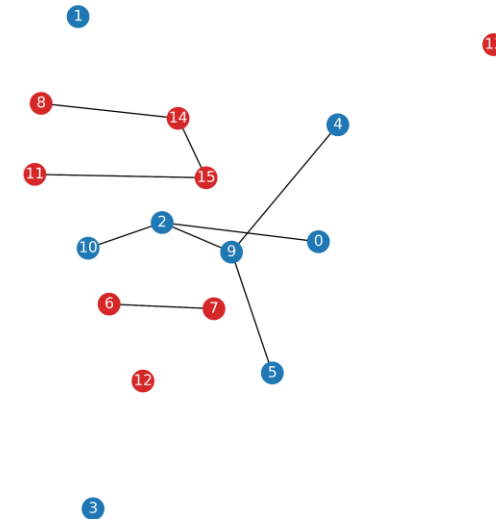
# Results: Max-cut



Network without  
the cut  
 $N = 16$



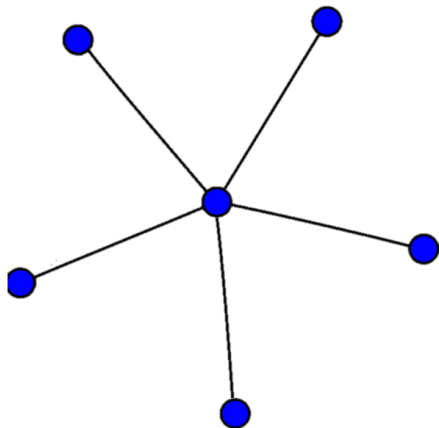
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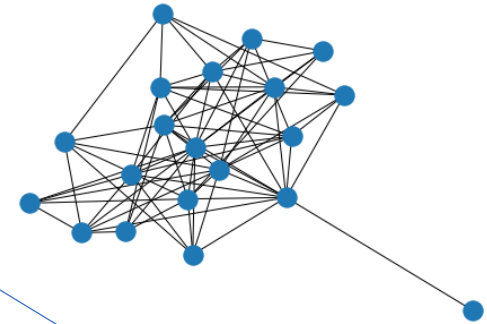
# Side note: Barabási-Albert graph

- Last report: How can there be a such a lonely node?
- Parameters:  $n$  – final num of nodes  
 $m$  – edges from new nodes

**initial\_graph** [Graph or None (default)] Initial network for Barabási–Albert algorithm. It should be a connected graph for most use cases. A copy of `initial_graph` is used. If None, starts from a star graph on  $(m+1)$  nodes.



$$m = 5$$



**barabasi\_albert\_graph** ( $n, m, seed=None, initial\_graph=None$ )  
Returns a random graph using Barabási–Albert preferential attachment

A graph of  $n$  nodes is grown by attaching new nodes each with  $m$  edges that are preferentially attached to existing nodes with high degree.

**Parameters**

- n** [int] Number of nodes
- m** [int] Number of edges to attach from a new node to existing nodes
- seed** [integer, random\_state, or None (default)] Indicator of random number generation state. See [Randomness](#).
- initial\_graph** [Graph or None (default)] Initial network for Barabási–Albert algorithm. It should be a connected graph for most use cases. A copy of `initial_graph` is used. If None, starts from a star graph on  $(m+1)$  nodes.

**Returns**

- G** [Graph]

**Raises**

- NetworkXError** If  $m$  does not satisfy  $1 \leq m < n$ , or the initial graph number of nodes  $m_0$  does not satisfy  $m \leq m_0 \leq n$ .