

The Meta-Variational Quantum Eigensolver

Alba Cervera-Lierta, Jakob S. Kottmann, Alán Aspuru-Guzik

QWorld Quantum Science Days

June 1st, 2021

Outlook







1. Tequila: a platform for rapid development of quantum algorithms
2. Variational Quantum Algorithms (VQA)
3. VQA for Quantum Machine Learning (QML)
4. The Meta-VQE: a QML algorithm for Hamiltonian simulation

Tequila

Quantum Science and Technology

PAPER

TEQUILA: a platform for rapid development of quantum algorithms

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Published 9 March 2021 • © 2021 IOP Publishing Ltd

[Quantum Science and Technology](#), [Volume 6](#), [Number 2](#)

Citation Jakob S Kottmann *et al* 2021 *Quantum Sci. Technol.* **6** 024009

The golden era of quantum languages...



The golden era of quantum languages...

...and quantum software tools...



The golden era of quantum languages...
...and quantum software tools...
...plus classical tools for the NISQ era



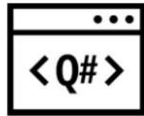


PYQUIL

Qibo



PENNY
LANE

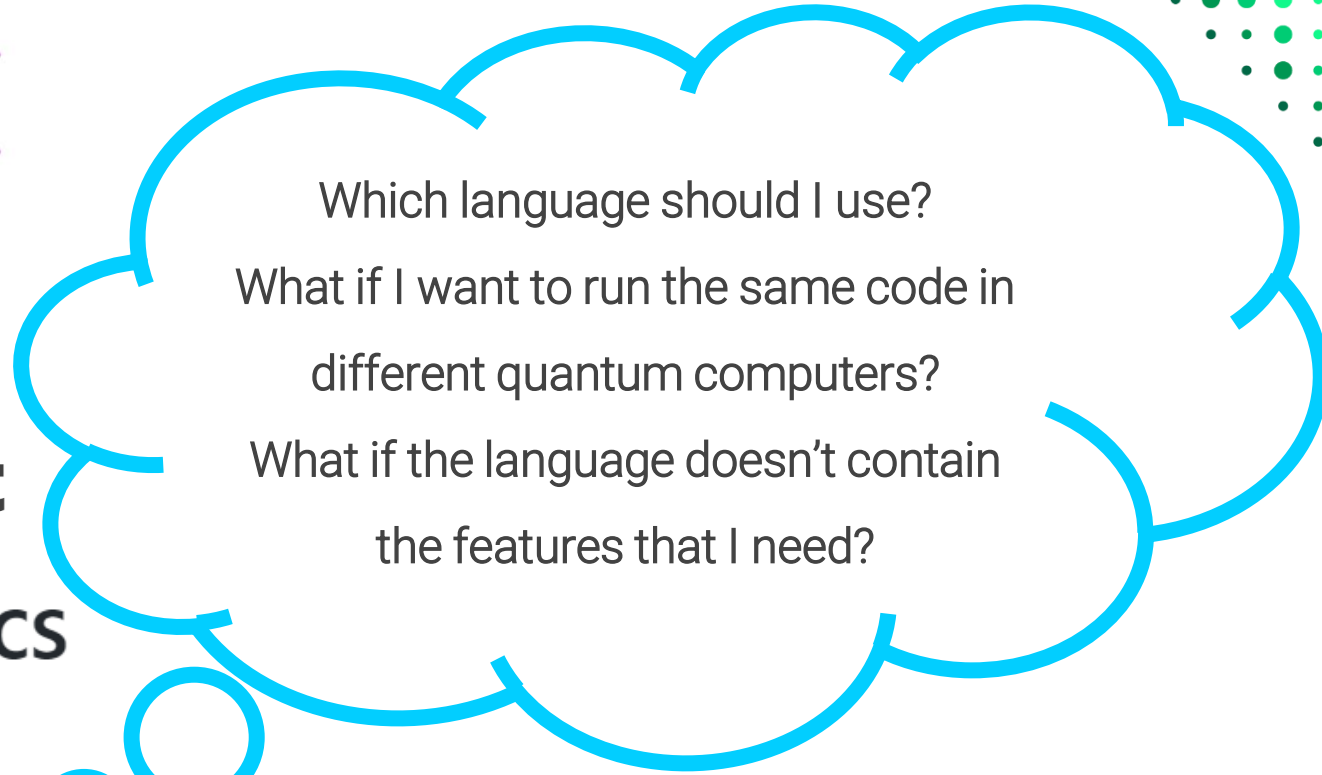
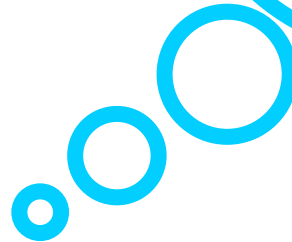


GPyOpt



Mitiq

Phoenix



Which language should I use?

What if I want to run the same code in
different quantum computers?

What if the language doesn't contain
the features that I need?



Unification, standardization, acceleration



A quantum language to simplify and accelerate implementation of new ideas for quantum algorithms.

Code

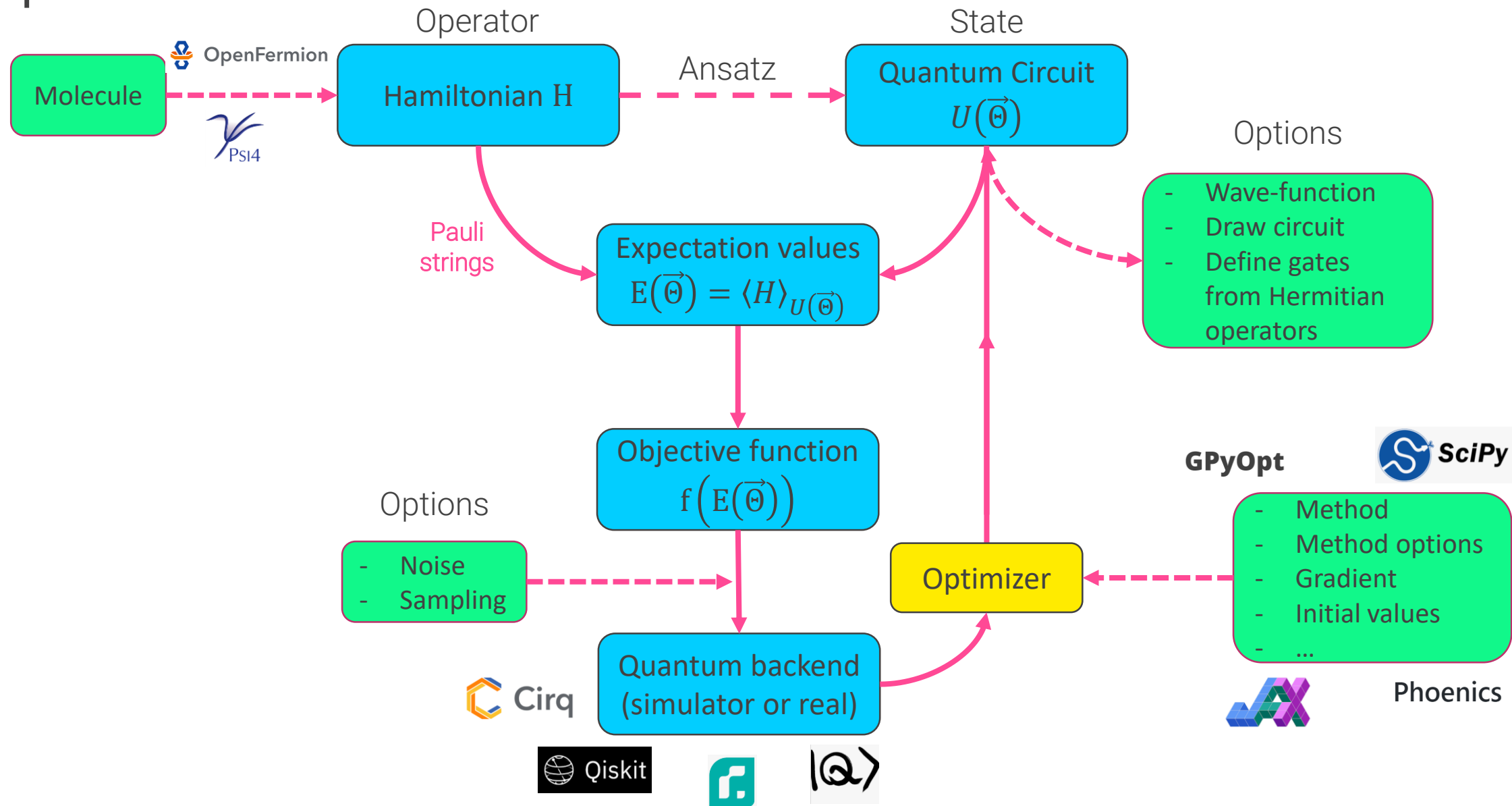
<https://github.com/aspuru-guzik-group/tequila>

Tutorials

<https://github.com/aspuru-guzik-group/tequila-tutorials>



Tequila API



Variational Quantum Algorithms

[Submitted on 21 Jan 2021]

Noisy intermediate-scale quantum (NISQ) algorithms

Kishor Bharti, Alba Cervera-Lierta, Thi Ha Kyaw, Tobias Haug, Sumner Alperin-Lea, Abhinav Anand, Matthias Degroote, Hermann Heimonen, Jakob S. Kottmann, Tim Menke, Wai-Keong Mok, Sukin Sim, Leong-Chuan Kwek, Alán Aspuru-Guzik

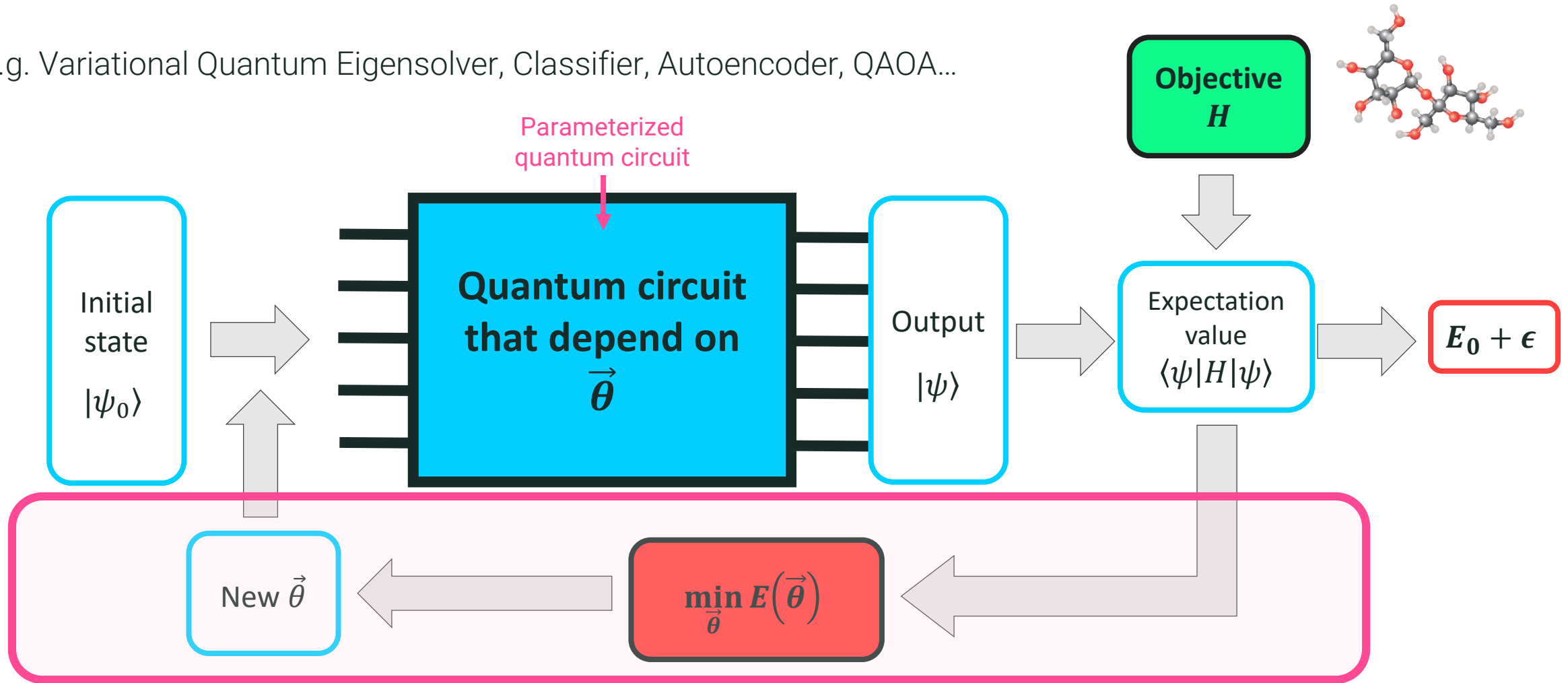
Comments: Review article, 82 pages, 7 figures, comments welcome

Subjects: **Quantum Physics (quant-ph)**; Statistical Mechanics (cond-mat.stat-mech); Artificial Intelligence (cs.AI); Machine Learning (cs.LG)

Cite as: [arXiv:2101.08448](https://arxiv.org/abs/2101.08448) [quant-ph]

Variational Quantum Algorithms

e.g. Variational Quantum Eigensolver, Classifier, Autoencoder, QAOA...



Classical optimization

Variational principle: $E = \langle\psi|H|\psi\rangle \geq E_0$

Objective function



It encodes the problem in a form of a quantum operator, e.g. a Hamiltonian

$$\langle H \rangle_{\mathcal{U}(\theta)} \equiv \langle 0 | \mathcal{U}^\dagger(\theta) H \mathcal{U}(\theta) | 0 \rangle$$

The objective is decomposed into Pauli strings which expectation value can be measured with the quantum computer.

$$H = \sum_{k=1}^M c_k \hat{P}_k \longrightarrow \langle H \rangle_{\mathcal{U}} = \sum_{k=1}^M c_k \langle \hat{P}_k \rangle_{\mathcal{U}}$$

An objective can also be the fidelity w.r.t. a particular target state that we are trying to match.

$$F(\Psi, \Psi_{\mathcal{U}(\theta)}) \equiv |\langle \Psi | \Psi_{\mathcal{U}(\theta)} \rangle|^2$$

We can use projectors or SWAP test to obtain the value of that fidelity

$$\max_{\theta} F(\Psi, \Psi_{\mathcal{U}(\theta)}) = \min_{\theta} (-\langle \hat{\Pi}_{\Psi} \rangle_{\mathcal{U}(\theta)})$$

Parameterized quantum circuits

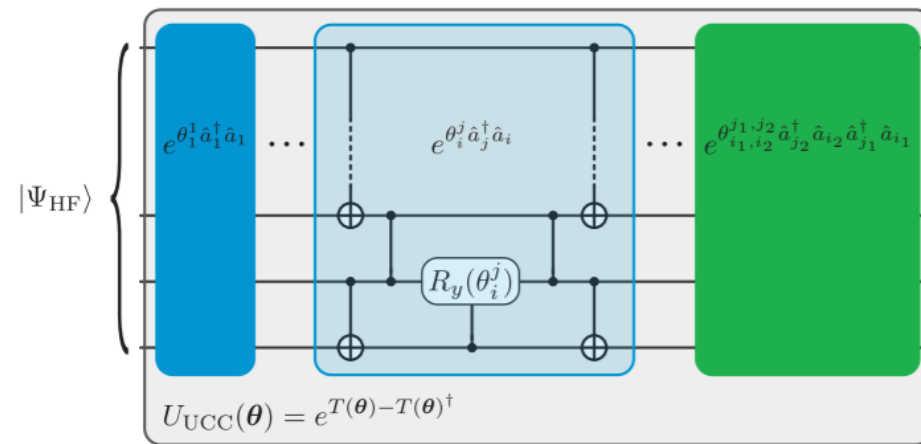


It prepares what will eventually be the approximation of the g.s. of our Objective function.

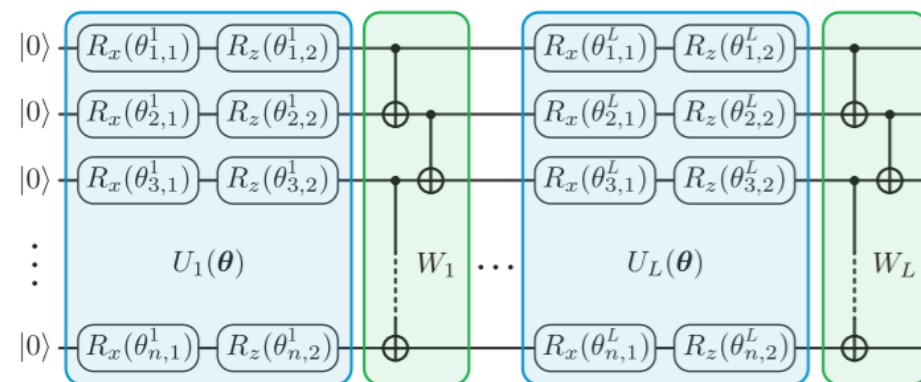
It depends on a series of parameters that have to be finetuned to minimize the objective

They can be designed from a physical point of view (e.g. UCC, QAOA,...) or from a practical point of view (using a limited set of gates and circuit topology).

a Problem-inspired ansatz



b Hardware-efficient ansatz



Classical optimization



We need to navigate the quantum circuit parameter space, e.g. by using gradient based methods

$$\theta_i^{(t+1)} = \theta_i^{(t)} - \eta \partial_i f(\boldsymbol{\theta})$$

The gradients are expectation values of the quantum circuit derivatives w.r.t. a parameter.

Example: parameter-shift rule

$$\mathcal{U}(\boldsymbol{\theta}) = V(\boldsymbol{\theta}_{-i})G(\theta_i)W(\boldsymbol{\theta}_{-i}) \quad G = e^{-i\theta_i g}$$

Eigenvalues of g are $\pm\lambda$

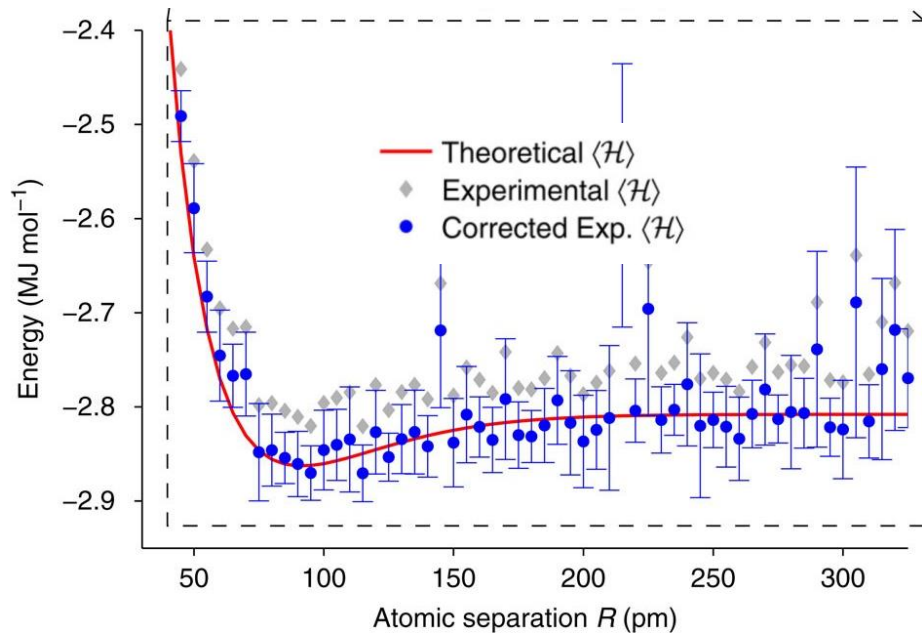
$$\partial_i \langle f(\boldsymbol{\theta}) \rangle = \lambda (\langle f(\boldsymbol{\theta}_+) \rangle - \langle f(\boldsymbol{\theta}_-) \rangle) \quad \boldsymbol{\theta}_{\pm} = \boldsymbol{\theta} \pm (\pi/4\lambda) \mathbf{e}_i$$

Gradient-free: genetic algorithms, reinforcement learning, ...



Example: the Variational Quantum Eigensolver

Bond dissociation curve of the He–H⁺ molecule.



GOAL: find $|\psi\rangle$ that minimizes $\frac{\langle \psi | \mathcal{H} | \psi \rangle}{\langle \psi | \psi \rangle}$.

Electronic structure Hamiltonian decomposed into Pauli strings

$$\langle \mathcal{H} \rangle = \sum_{i\alpha} h_{\alpha}^i \langle \sigma_{\alpha}^i \rangle + \sum_{ij\alpha\beta} h_{\alpha\beta}^{ij} \langle \sigma_{\alpha}^i \sigma_{\beta}^j \rangle + \dots$$

Quantum circuit that generates the ground state of that Hamiltonian (Unitary Couple-Cluster ansatz)

$$|\Psi(\theta)\rangle = e^{T(\theta)-T(\theta)^{\dagger}} |\Psi_{\text{HF}}\rangle$$

Unitary operation (Cluster operator) Hartree-Fock Excitations Hartree-Fock orbitals

$$T(\theta) = T_1(\theta) + T_2(\theta) + \dots$$

$$T_1(\theta) = \sum_{\substack{i \in \text{occ} \\ j \in \text{virt}}} \theta_i^j \hat{a}_j^{\dagger} \hat{a}_i$$

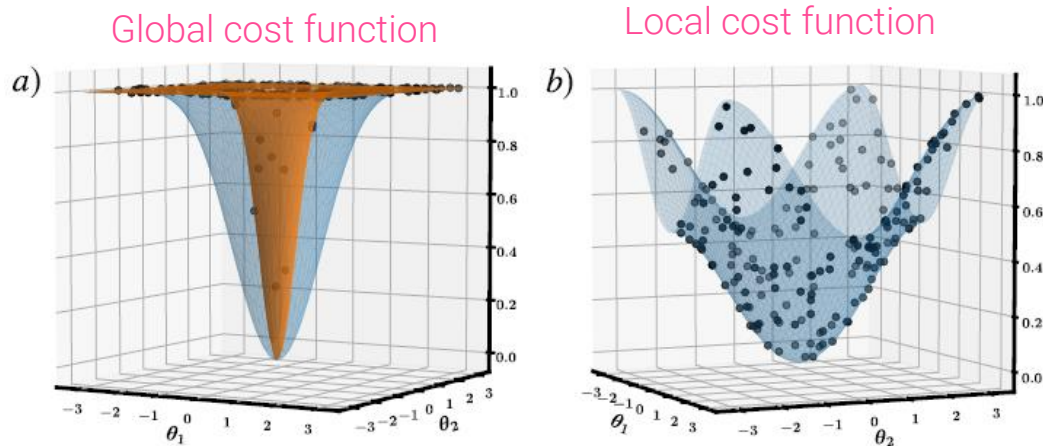
$$T_2(\theta) = \sum_{\substack{i_1, i_2 \in \text{occ} \\ j_1, j_2 \in \text{virt}}} \theta_{i_1, i_2}^{j_1, j_2} \hat{a}_{j_2}^{\dagger} \hat{a}_{i_2} \hat{a}_{j_1}^{\dagger} \hat{a}_{i_1}$$

Transform the fermionic operators to Pauli strings (e.g. Jordan Wigner) and they become the generators of the quantum gates.

The *barren-plateaux* problem

Compute the gradients with the quantum circuit and use these values to run a classical minimizer, e.g. Nelder-Mead, Adam, ...

With no prior knowledge about the solution, $\vec{\theta}$ parameters are initialized at random.



Consequence: *barren-plateaux*

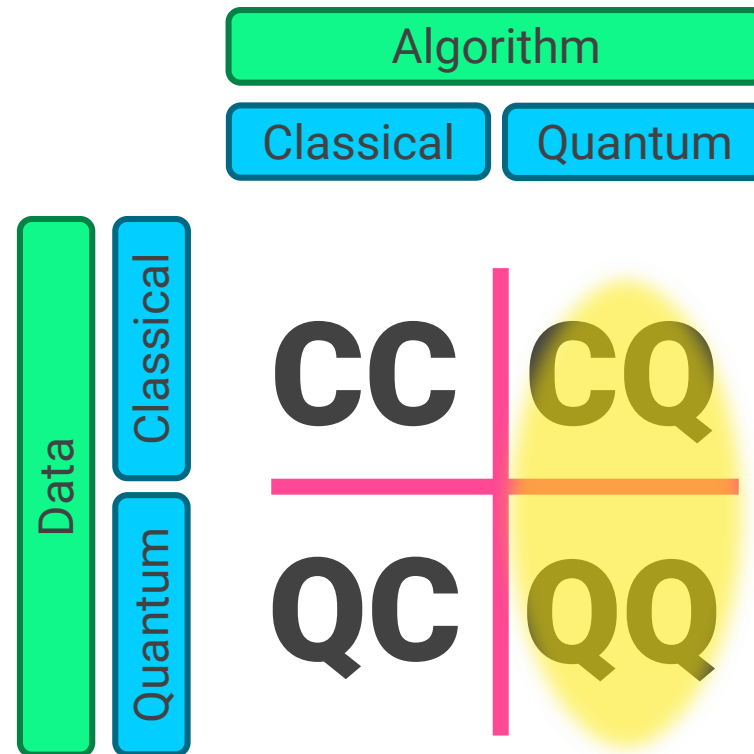
The expected value of the gradient is zero!
The expected value of the variance is also zero!

Solutions

- Use parameters close to the solution.
- Use local cost functions instead of global ones.
- Introduce correlations between parameters.

Ref.: M. Cerezo et. al. arXiv:2001.00550v2 [quant-ph]

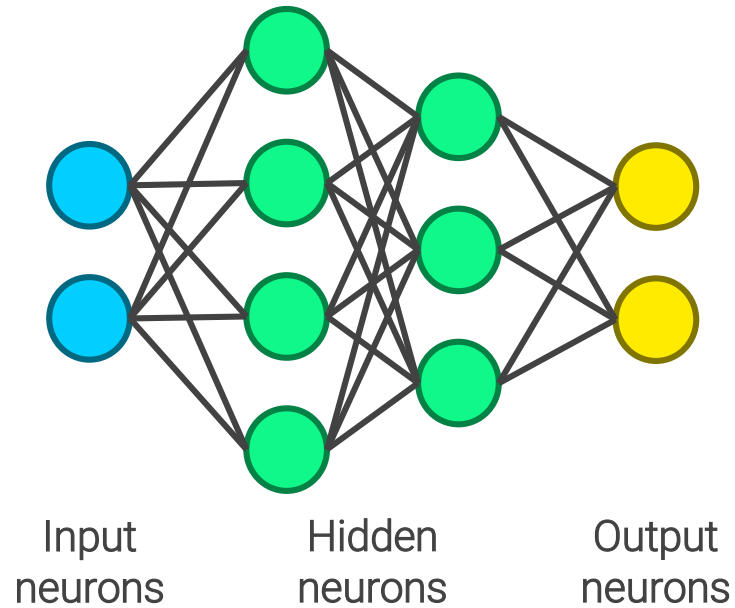
VQA for Quantum Machine Learning



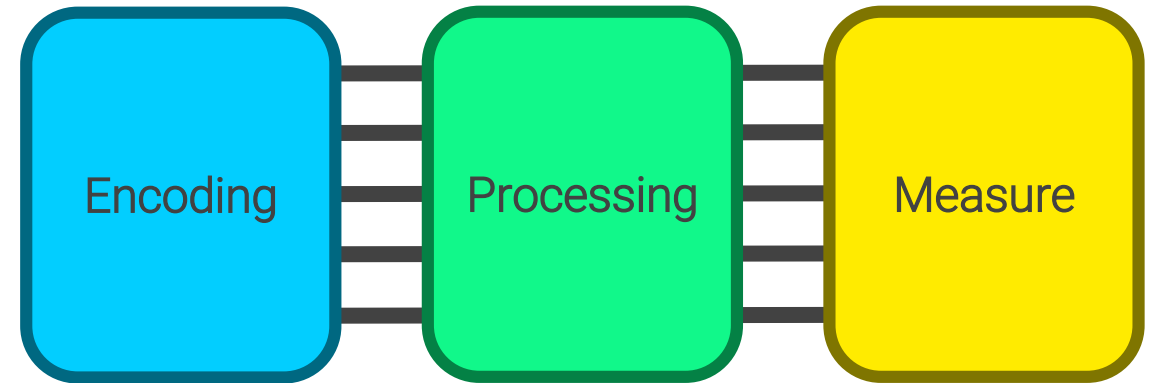
From classical to quantum NN



Classical



Quantum
(circuit centric)



K Mitarai, M Negoro, M Kitagawa, K Fujii Phys. Revs A 98 (3), 032309 (2018)

E. Farhi and H.Neven, arXiv:1802.06002 (2018)

M. Schuld and N. Killoran, Phys. Rev. Lett. 122, 040504 (2019)

M. Schuld, A. Bocharov, K. M. Svore, and N. Wiebe, Phys. Rev. A 101, 032308 (2020)



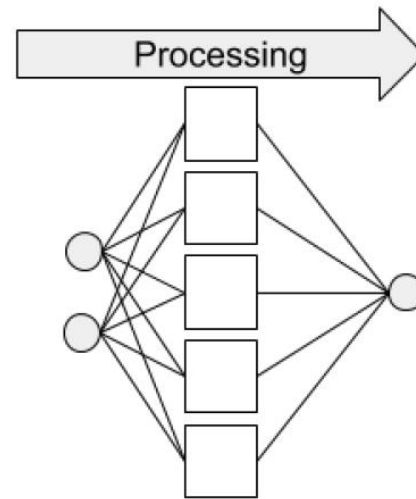
The minimal QNN

What is the most simple (but universal) NN?

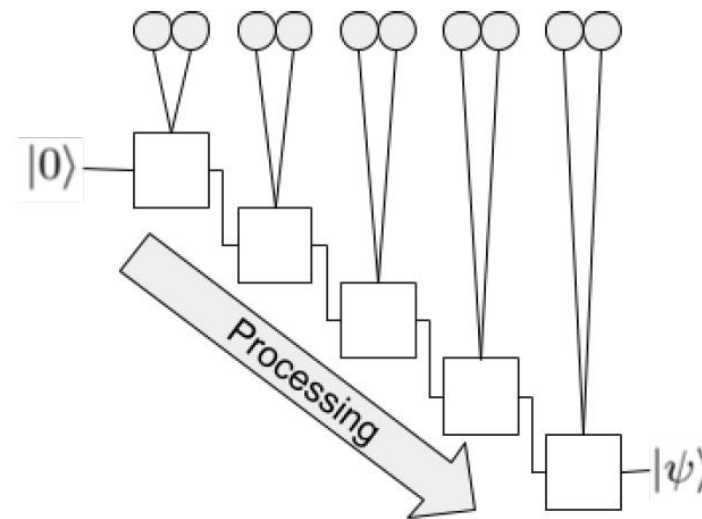
Single hidden layer NN

What is the most simple (but universal) QNN?

Single-qubit QNN



(a) Neural network



(b) Quantum classifier

Encoding the data

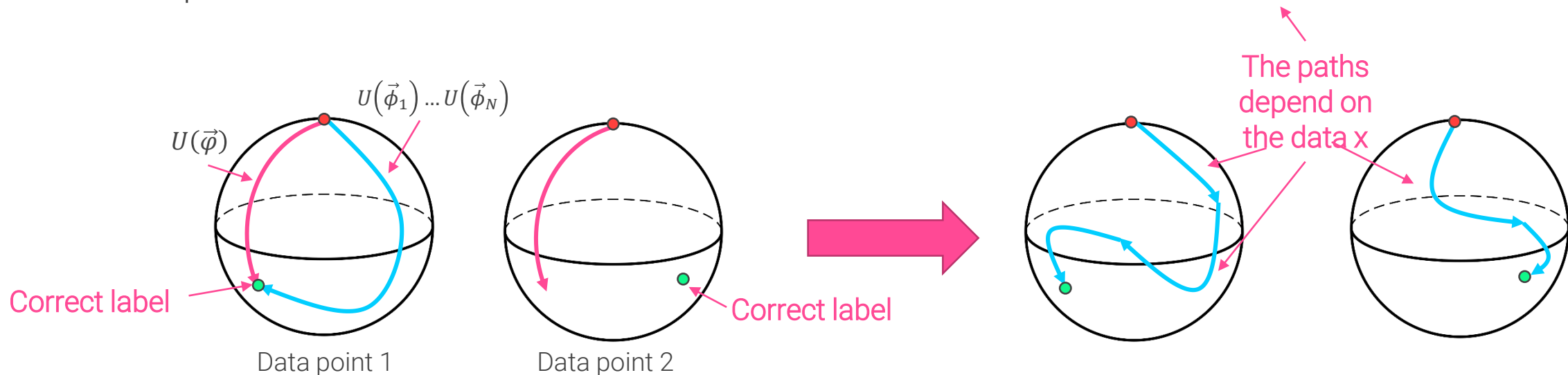
A product of unitaries can be written with a single unitary

$$U(\vec{\phi}_1) \dots U(\vec{\phi}_N) \equiv U(\vec{\phi})$$

If we add some fixed parameter dependency (the data), the operation becomes flexible and data-dependent.

Data re-uploading

$$\mathcal{U}(\vec{\phi}, \vec{x}) \equiv U(\vec{\phi}_N)U(\vec{x}) \dots U(\vec{\phi}_1)U(\vec{x})$$



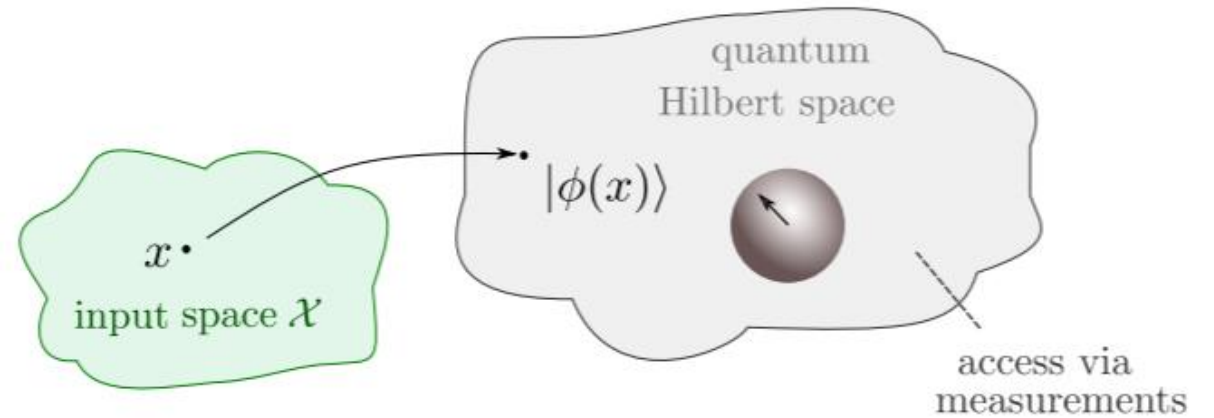
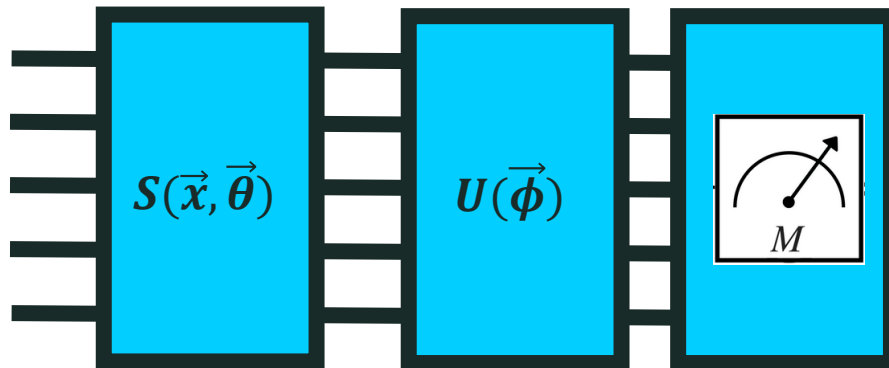
Quantum Feature Maps

$$|\psi_0\rangle \rightarrow |\psi(\vec{x}, \vec{\theta})\rangle \rightarrow |\psi(\vec{x}, \vec{\theta}, \vec{\phi})\rangle \rightarrow |gs(\vec{x})\rangle$$

Encode the data
(quantum
feature space)

Rotate to the
correct
measurement
basis

Find the parameters that
minimize the objective
(measured in the computational
basis)





The Meta-Variational Quantum Eigensolver

Open Access

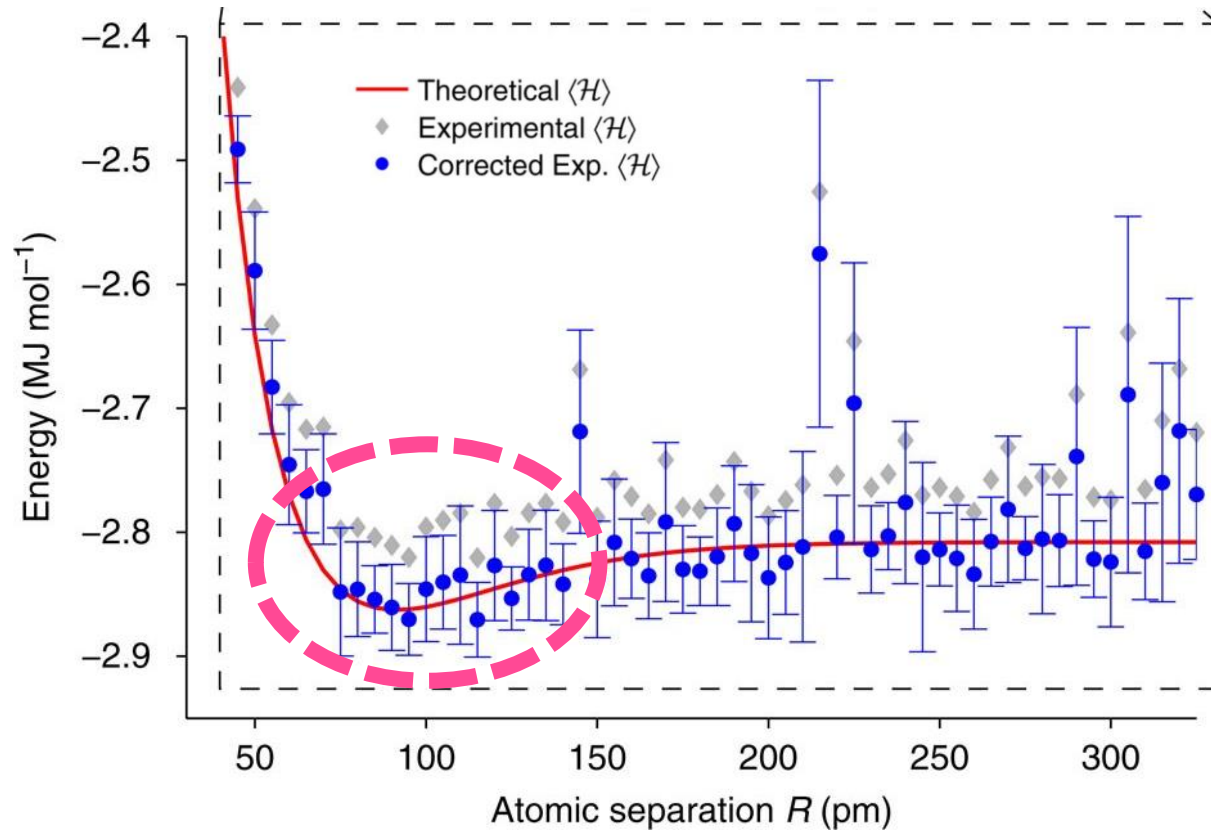
Meta-Variational Quantum Eigensolver: Learning Energy Profiles of Parameterized Hamiltonians for Quantum Simulation

Alba Cervera-Lierta, Jakob S. Kottmann, and Alán Aspuru-Guzik
PRX Quantum **2**, 020329 – Published 28 May 2021

What's the true goal of VQE?



Bond dissociation curve of the He–H⁺ molecule.



GOAL: find $|\psi\rangle$ that minimizes $\frac{\langle \psi | \mathcal{H} | \psi \rangle}{\langle \psi | \psi \rangle}$.



Find the atomic separation that minimizes the energy

$$\min \langle H(R) \rangle$$

A. Peruzzo, J. McClean, P. Shadbolt, M.-H. Yung, X.-Q. Zhou, P. J. Love, A. Aspuru-Guzik, J. L. O'Brien, Nature Comm. **5**, 4213 (2014)

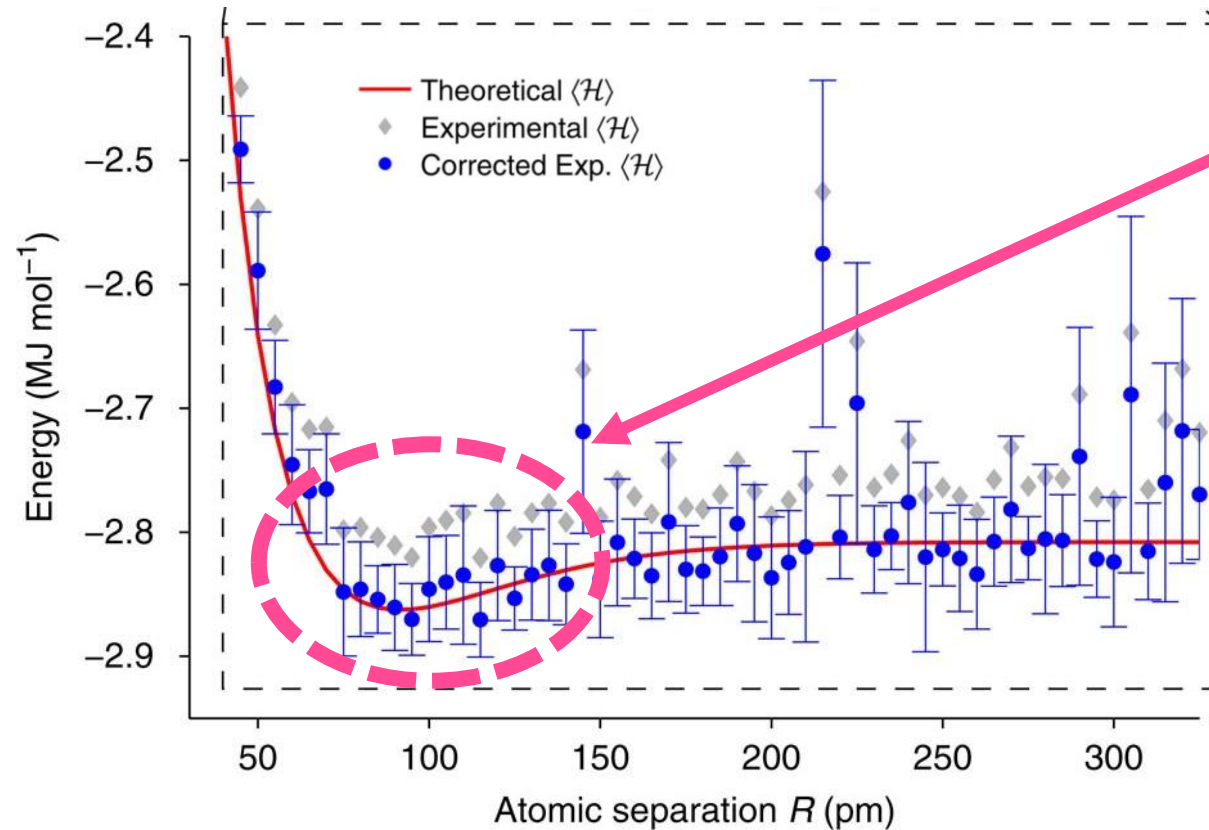
ACL, J. Kottmann, A. Aspuru-Guzik, PRX Quantum **2**, 020329 (2021)



What's the true goal of VQE?



Bond dissociation curve of the He-H⁺ molecule.



To obtain **this** you need to scan from 0 to 300.

Each blue point is a VQE, that is, you have to **prepare, run and optimize** the quantum circuit.

Can we avoid to compute the uninteresting points?

A. Peruzzo, J. McClean, P. Shadbolt, M.-H. Yung, X.-Q. Zhou, P. J. Love, A. Aspuru-Guzik, J. L. O'Brien, Nature Comm. **5**, 4213 (2014)

ACL, J. Kottmann, A. Aspuru-Guzik, PRX Quantum **2**, 020329 (2021)



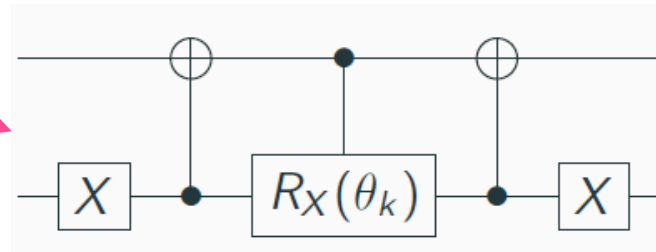
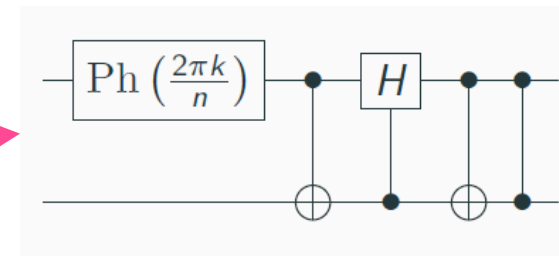
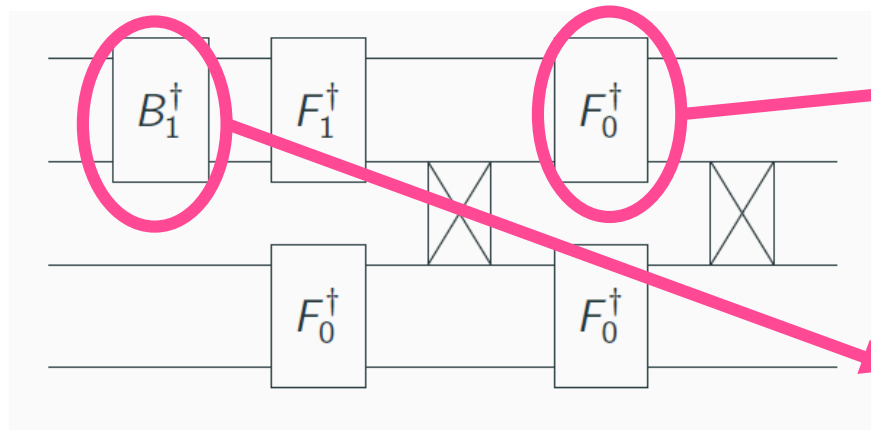
Exact simulation

If the model can be solved analytically, we can find the quantum circuit that diagonalizes the Hamiltonian (therefore, no need for any VQE)

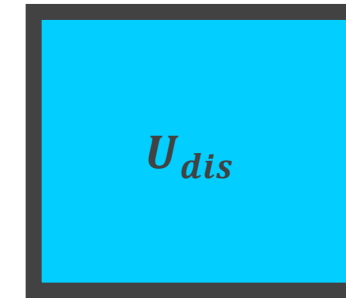
Example: Ising model

By finding the transformations that:

1. Jordan-Wigner
2. Fourier Transform
3. Bogoliubov transformation



Computational basis
state
(easy to prepare)



Eigenstate



$$U_{dis} = U_{JW}U_{FT}U_{Bog}$$

Hamiltonian parameters

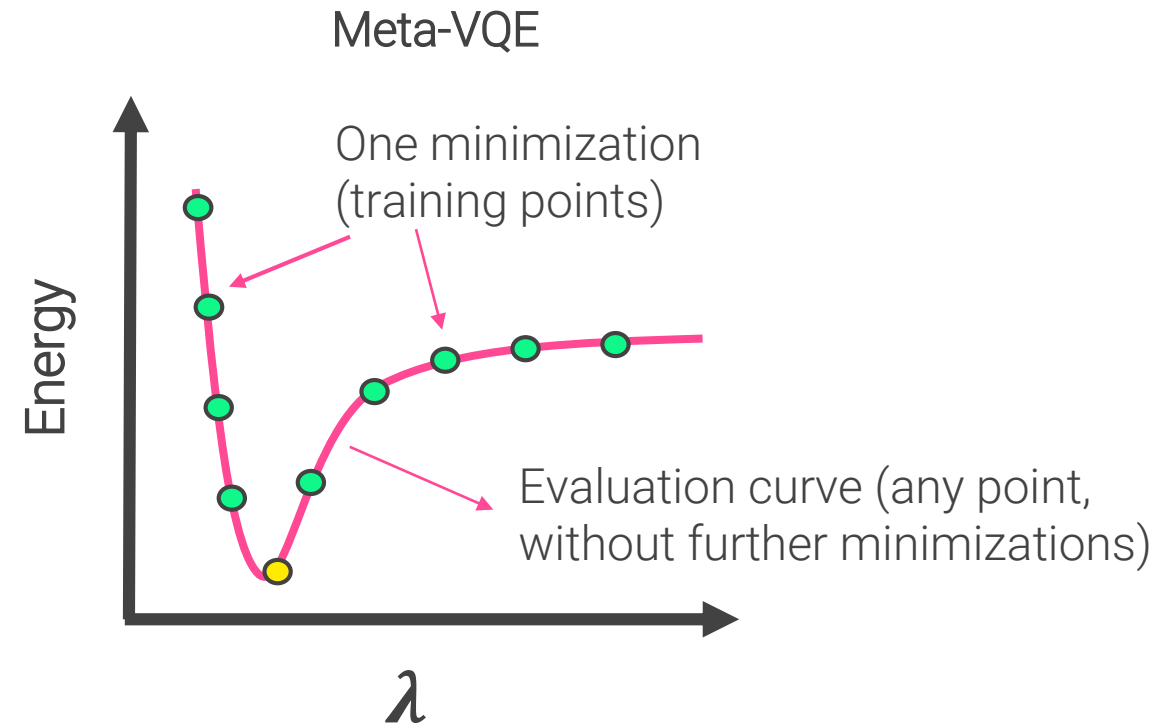
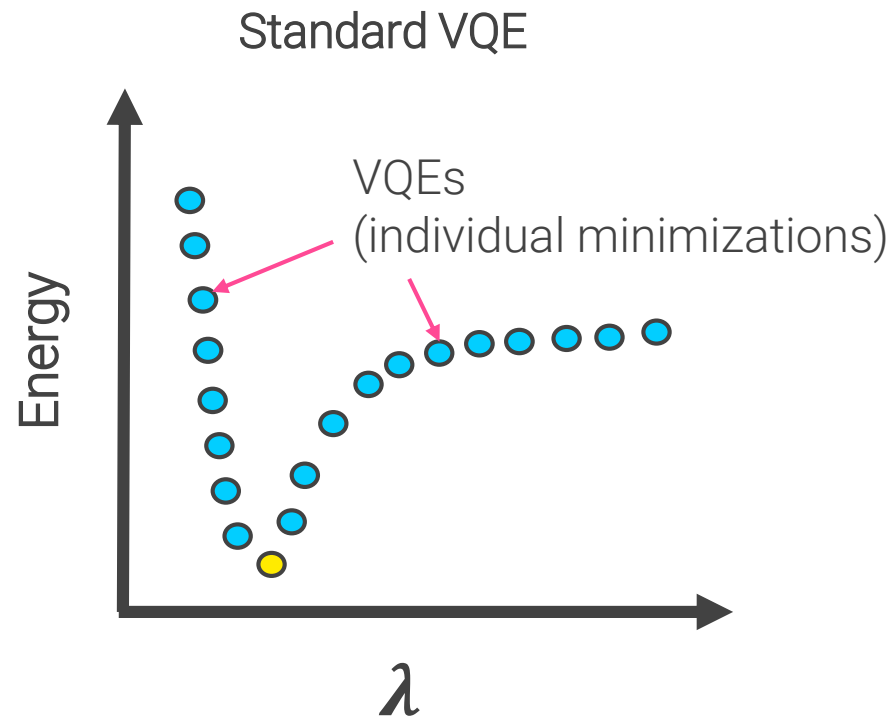
They naturally appear in the rotational gates from the Bogoliubov transformation!

$$\theta_k = 2 \arctan \left(\frac{J\gamma \sin \left(\frac{2\pi k}{n} \right)}{J \cos \left(\frac{2\pi k}{n} \right) + \lambda} \right)$$

Meta-VQE outlook

Parameterized Hamiltonian $H(\vec{\lambda})$

Goal: to find the quantum circuit that **encodes** the ground state of the Hamiltonian for any value of $\vec{\lambda}$



See also: K. Mitarai, T. Yan, K. Fujii, Phys. Rev. Applied 11, 044087 (2019)

ACL, J. Kottmann, A. Aspuru-Guzik, PRX Quantum 2, 020329 (2021)

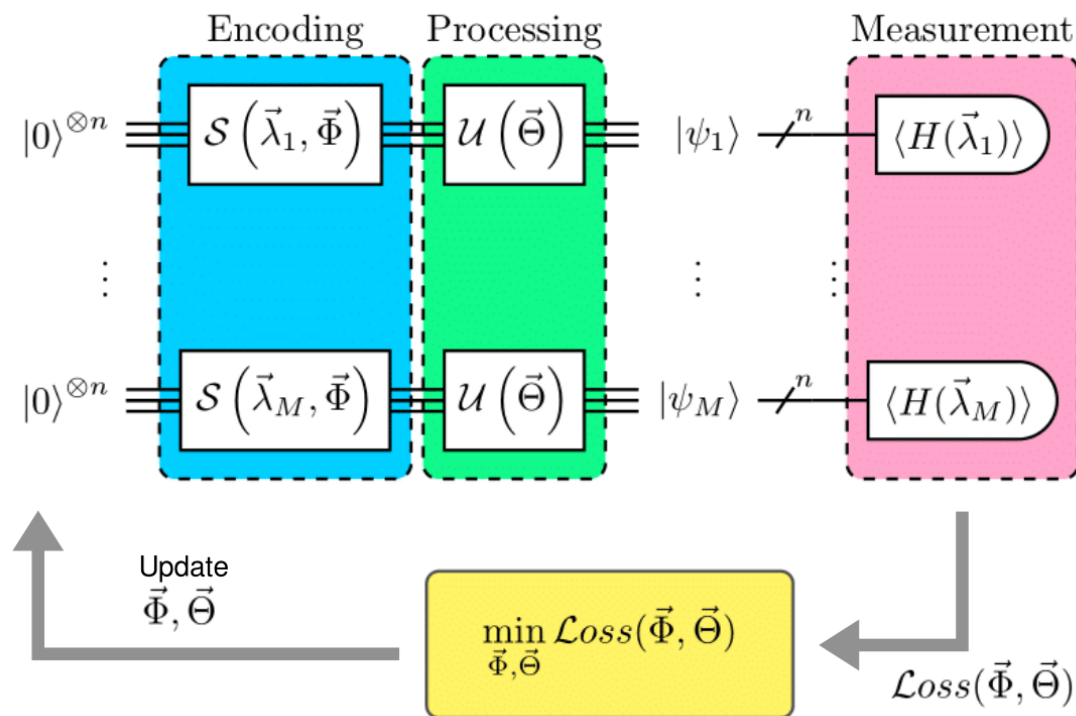
The Meta-VQE

Parameterized Hamiltonian $H(\vec{\lambda})$

Training points: $\vec{\lambda}_i$ for $i = 1, \dots, M$

Loss function with all $\langle H(\vec{\lambda}_i) \rangle$

Goal: to find the quantum circuit that **encodes** the ground state of the Hamiltonian for any value of $\vec{\lambda}$



Output: $\vec{\Phi}_{\text{opt}}$ and $\vec{\Theta}_{\text{opt}}$

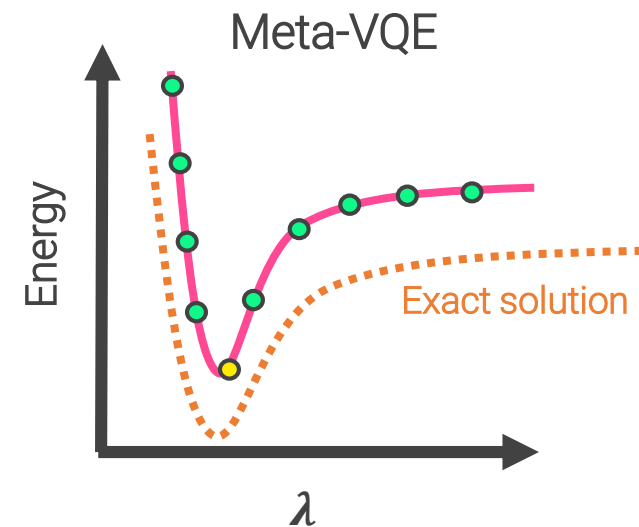
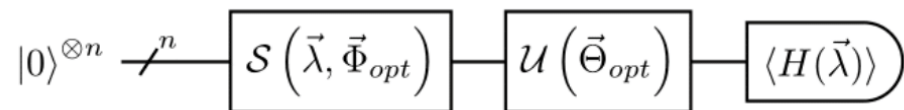
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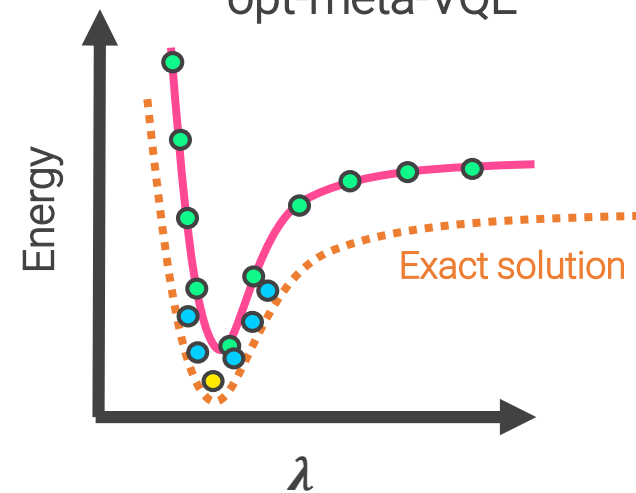
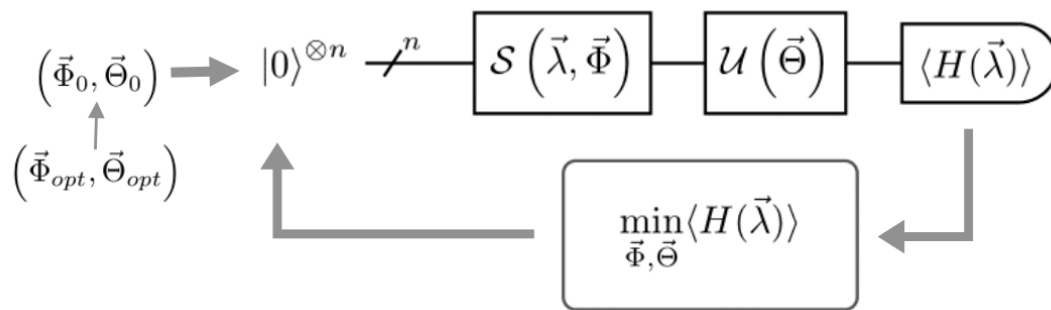
The Meta-VQE output

Output: $\vec{\Phi}_{opt}$ and $\vec{\Theta}_{opt}$

Option 1: run the circuit with test $\vec{\lambda}$ and obtain the g.s. energy profile.



Option 2: use $\vec{\Phi}_{opt}$ and $\vec{\Theta}_{opt}$ as starting point of a standard VQE optimization (opt-meta-VQE)



1D XXZ spin chain



$$H = \sum_{i=1}^n \sigma_i^x \sigma_{i+1}^x + \sigma_i^y \sigma_{i+1}^y + \Delta \sigma_i^z \sigma_{i+1}^z + \lambda \sigma_i^z$$

For $\lambda = 0$, two QPT: $\Delta = -1, \Delta = 1$

Analytical solution of the model: using the Bethe ansatz (no known quantum circuit)

Good worse-case-scenario model

- We do not know which circuit ansatz will work
- The ground state is highly entangled (that's why we need quantum computers!)
- The energy profile is not trivial: it presents a peak in the region $\Delta > -1$

1D XXZ spin chain

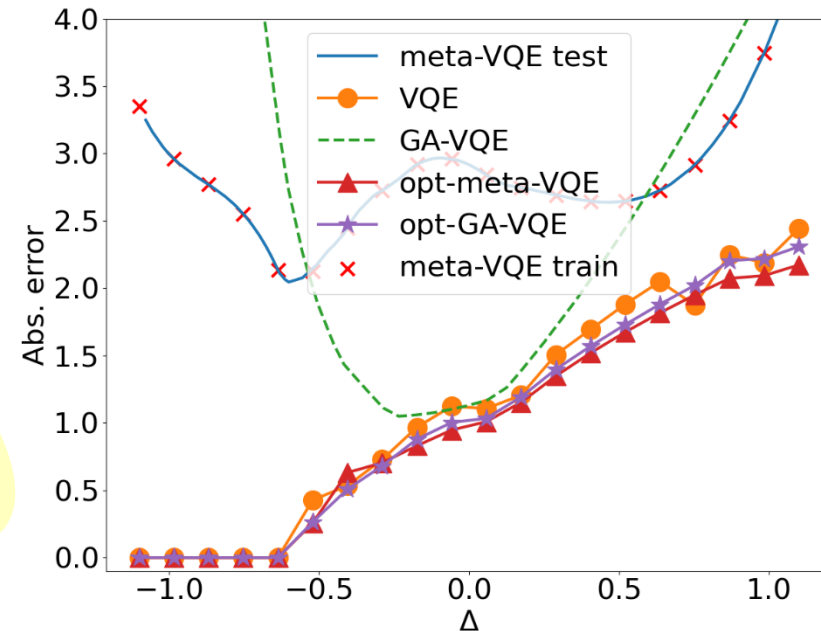
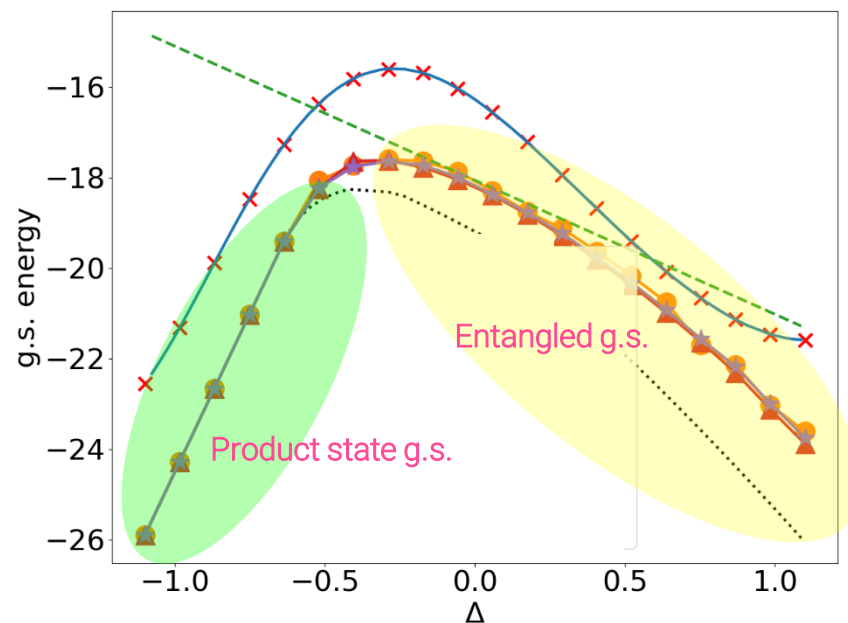
14 qubits simulation, $\lambda = 0.75$

Linear encoding: $R_z(w_1 \Delta + \phi_1) R_y(w_2 \Delta + \phi_2) \otimes$ Alternating CNOT

Processing layer: $R_z(\theta_1) R_y(\theta_2) \otimes$ Alternating CNOT

Results 2 encoding + 2 processing layers

$$H = \sum_{i=1}^n \sigma_i^x \sigma_{i+1}^x + \sigma_i^y \sigma_{i+1}^y + \Delta \sigma_i^z \sigma_{i+1}^z + \lambda \sigma_i^z$$



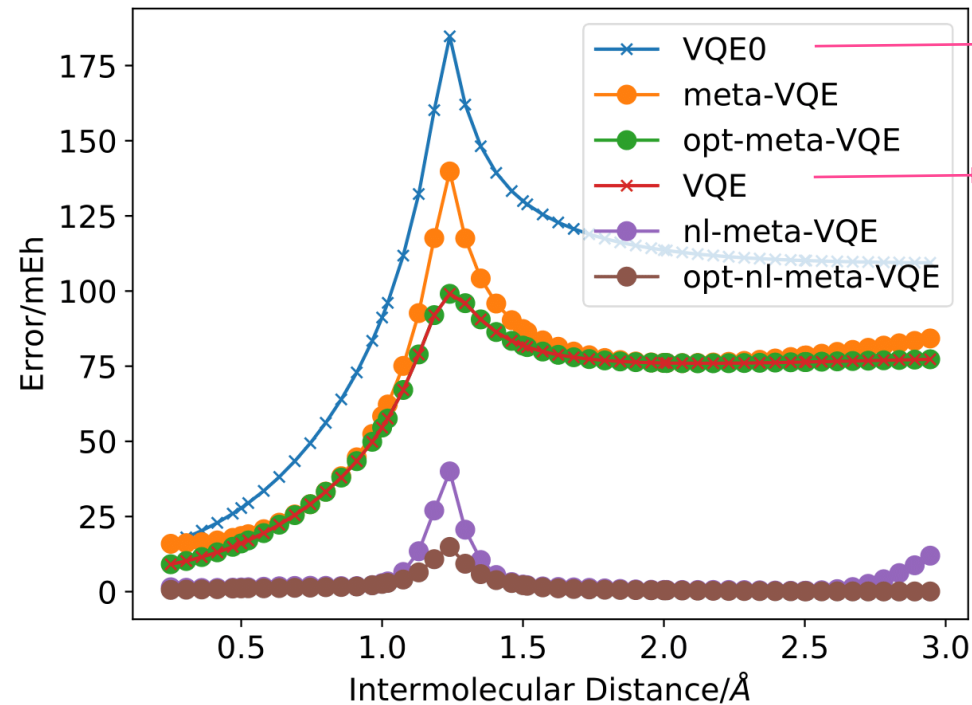
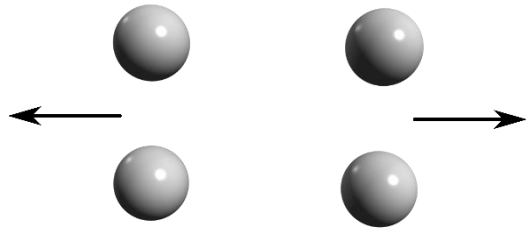
H_4 molecule

H_4 molecule in 8 spin-orbitals (STO-3G basis set)

Ansatz: k-UpCCGSD (k=2 for these results)

Linear encoding: $\theta = \alpha + d\beta$
Non-linear encoding: $\theta = \alpha e^{\beta(\gamma - d)} + \delta$ (floating Gaussians)

Hamiltonian Parameter
(intermolecular distance)



Initial state: $|0\rangle$

Initial state: $|HF\rangle$

Single transmon

Single transmon simulation using QCAD mapping

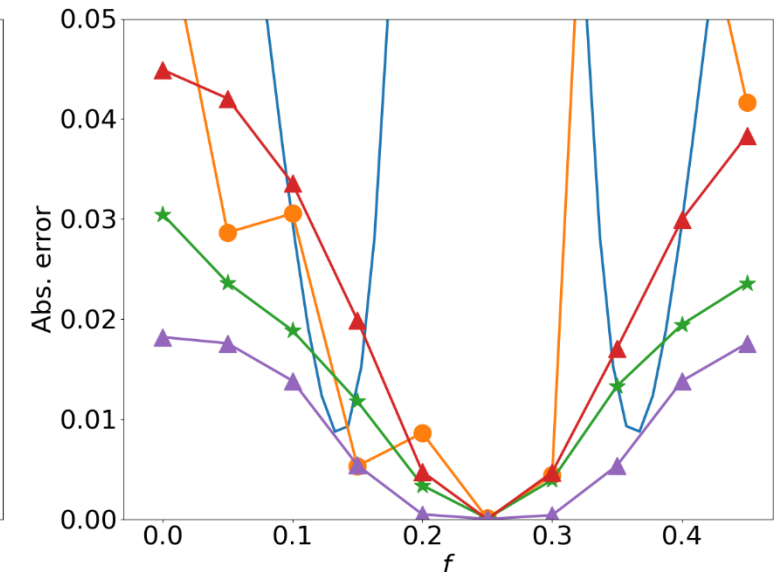
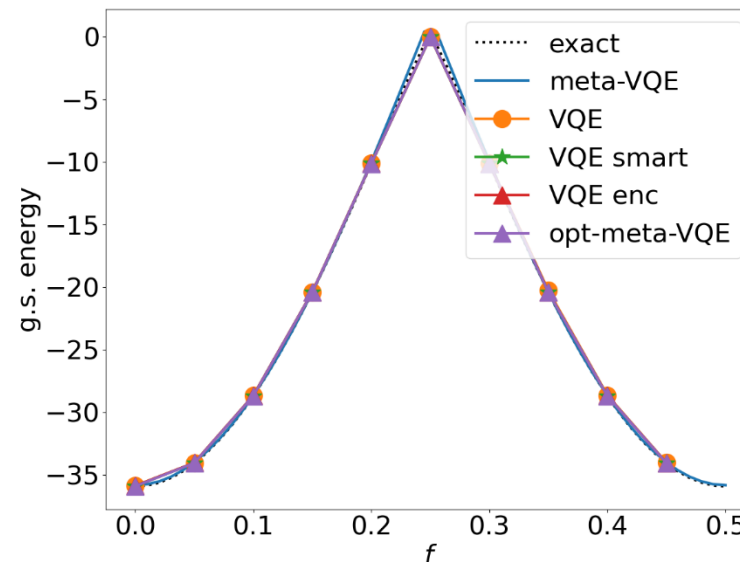
Ansatz: 1 encoding + 1 processing layers + 1 final layer of $R_x R_z$

Layer: $R_x R_z$ + all connected XX gates

Parameters of XX gates are the same in all layers (same entanglement gate)

Linear encoding: $R_x(w_1 f + \phi_1) R_z(w_2 f + \phi_2)$

Hamiltonian Parameter (flux)



Kyaw, Menke, Sim, Sawaya, Oliver,
Guerreschi, Aspuru-Guzik,
arXiv:2006.03070 (2020)

Conclusions

- Meta-VQE can be used to scan over Hamiltonian parameters to find the interesting energy regions.
 - ➡ Reduction in the total computational cost (less number of objective evaluations)
- We can use its parameter solution to run a more precise minimization (opt-meta-VQE)
 - ➡ Faster convergence, potentially avoiding barren plateaus and local minima
- The encoding strategy in VQE-type algorithms might be useful to guide the optimization towards the solution.
 - ➡ Avoiding barren plateaus (T. Volkoff, P. J. Coles, Quantum Sci. Technol. 6, 025008 (2021))

Code and demo (notebooks)

<https://github.com/aspuru-guzik-group/Meta-VQE>

Using Tequila quantum package

<https://github.com/aspuru-guzik-group/tequila>



Alán Aspuru-Guzik



Jakob Kottmann

Questions?

Special thanks to

Tequila developers

NISQ review collaboration:



Leong-Chuan Kwek



Kishor Bharti



Thi Ha Kyaw



Tobias Haug



Sumner Alperin



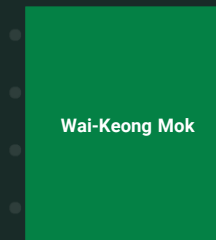
Abhinav Anand



Matthias Degroote



Hermanni Heimonen



Wai-Keong Mok



Sukin (Hannah) Sim

