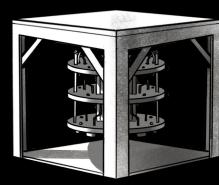


Scaling quantum computing, with the networks of **today** and **tomorrow**



Monolithic quantum computers don't scale!

Quantum computers need millions of qubits

Scaling qubit counts is **fundamentally limited**

Network stacks are required distributed computing

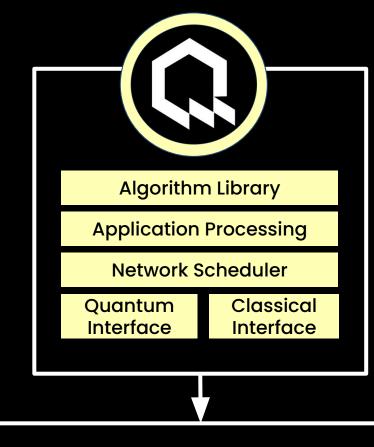


Qoro has the solution

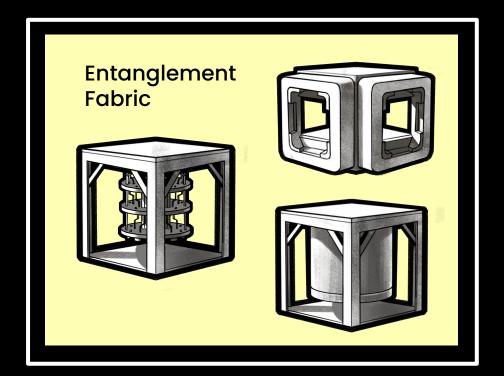
today

tomorrow

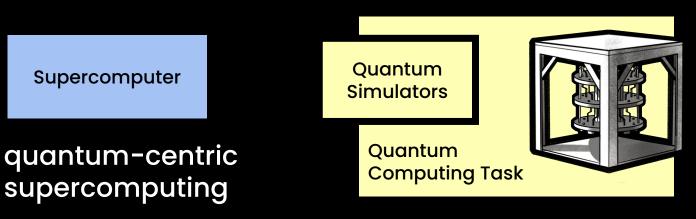
End-Users



Quantum Network Operating System



HPC



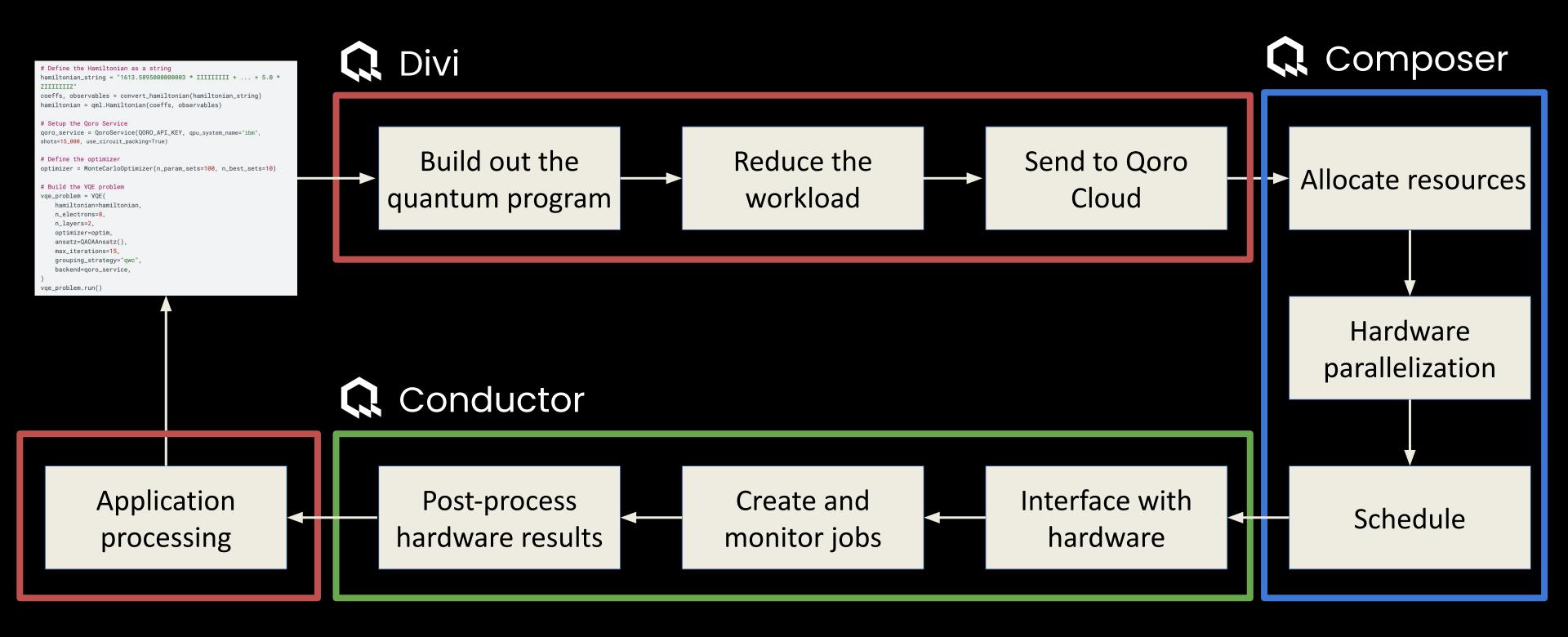


Example: Protein Folding

```
# Define the Hamiltonian as a string
hamiltonian_string = "1613.5895000000003 * IIIIIIIII + ... + 5.0 *
ZIIIIIIZ"
coeffs, observables = convert_hamiltonian(hamiltonian_string)
hamiltonian = qml.Hamiltonian(coeffs, observables)
# Setup the Qoro Service
qoro_service = QoroService(QORO_API_KEY, qpu_system_name="ibm",
shots=15_000, use_circuit_packing=True)
# Define the optimizer
optimizer = MonteCarloOptimizer(n_param_sets=100, n_best_sets=10)
# Build the VQE problem
vqe_problem = VQE(
    hamiltonian=hamiltonian,
    n_electrons=8,
    n_layers=2,
    optimizer=optim,
    ansatz=QAOAAnsatz(),
    max_iterations=15,
    grouping_strategy="qwc",
    backend=qoro_service,
vqe_problem.run()
```



What's happening?





Huge Cost Savings

Costs for running a 17 qubit protein folding problem on IBM (PAYG, \$96/min):

Conditions:

- Use Monte-Carlo optimization (gradient free)
- Using a 335 term Hamiltonian
- Using standard VQE



Huge Cost Savings

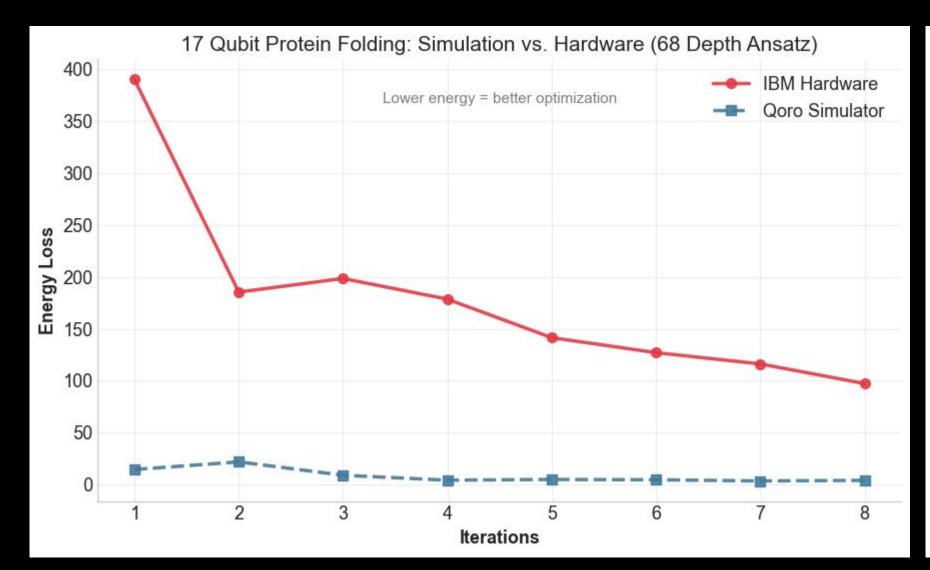
PAYG, \$96/min

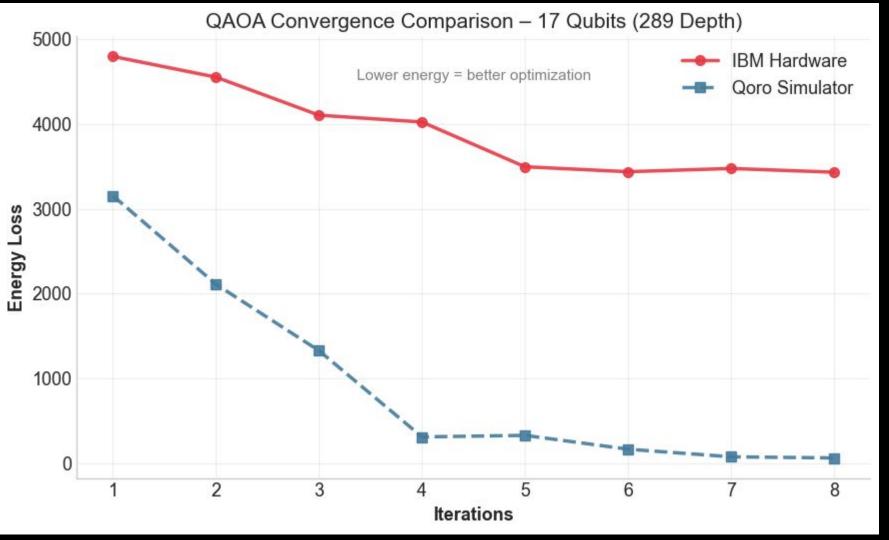
PER ITERATION!

Tool Used	Lines of code	Number of circuits	Cost for 8,000 shots	QPU Runtime	Reduction
Naive implementation	> 2,000	33,500	\$128,640	22 hours	
Divi & Grouping	20	100	\$384	4 minutes	99.7%
Composer & Packing	0	100	\$48	30 seconds	99.9%



Quality

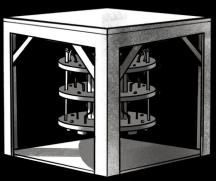








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