Kagome Lattice – Open Science 2022

A solution for the state preparation of the Kagome Lattice in a quantum computer

Our Solution

- Our solution is implemented as a python program where all the parameters of the VQE algorithm are fed as command line arguments.
- * Ansatz Options

Name	Description	Default Value
-a,ansatz_type	Ansatz type:	EfficientSU2
	ExcitationPreserving: Heuristic excitation-preserving	
	wave function.	
	• EfficientSU2: A hardware efficient SU(2) 2-local circuit.	
	PauliTwoDesign: The Pauli Two-Design ansatz.	
	TwoLocal: The two-local circuit.	
	RealAmplitudes: A heuristic trial wave function.	

Name	Description	Default Value
-ol,opt_level	Circuit optimization level (1-3)	1
-ui,uniform_interaction	Heisenberg Model uniform	Edge weight
	interaction value.	
-up,uniform_potential	Heisenberg Model uniform	0.0
	potential	
-w,weight	Lattice edge weight.	2.4
-s,shots	Number of execution shots.	2048

Optimizer Options

Name	Description	Default
		Value
-o,optimizer_type	Optimizer type:	SPSA
	SPSA: Simultaneous Perturbation Stochastic Approximation	
	SLSQP: Sequential Least Squares Programming.	
	COBYLA: Constrained Optimization By Linear Approximation.	
	UMDA: Continuous Univariate Marginal Distribution Algorithm	
	GSLS: Gaussian-smoothed Line Search.	
	GradientDescent: The gradient descent minimization routine.	
	 L_BFGS_B: Limited-memory BFGS Bound optimizer. 	
	NELDER_MEAD: Performs unconstrained optimization.	
	POWELL: The Powell algorithm with unconstrained	
	optimization	
	NFT: Nakanishi-Fujii-Todo algorithm.	
-i,max_iter	Maximum number of iterations or function evals used by the	175
	optimizer.	

Resilience and Scalability

Name	Description	Default Value
-r,resilience_type	Resilience type (1-3):	2
	• T-Rex: 1	
	• ZNE: 2	
	• PEC: 3	

Scalability

The program is designed to run in any quantum processor with any number of qubits. For example to run in Geneva (27 qubits) with an NFT optimizer and uniform potential:

\$ python3 kagome_solution.py -b ibm_geneva -t ibm_geneva -q 27 -a EfficientSU2 -w 1.0 -o NFT -up -1.0

Our Results: Solution 1 - Error 0.05%

Backend: ibmq_guadalupe

Ansatz: EfficientSU2 (reps=1,entanglement='reverse_linear')

Optimizer: **NFT**(maxiter=175)

Resilience: ZNE(2)

Shots: 2048,

Edge weight: 1.34000

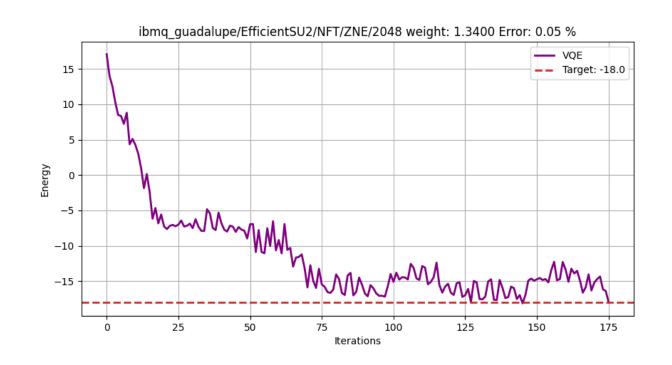
Execution time (s): 12390.63

Expected ground state energy: -18.00000000

Computed ground state energy: -17.99098307

Result eigen value: -17.99098307

Relative error: 0.05009404%



Our Results: Solution 2 - Error 0.56%

Backend: ibmq_guadalupe

Ansatz: EfficientSU2 (reps=1,entanglement='reverse_linear')

Optimizer: COBYLA (maxiter=100)

Resilience: ZNE(2)

Shots: 2048,

Edge weight: 1.80000

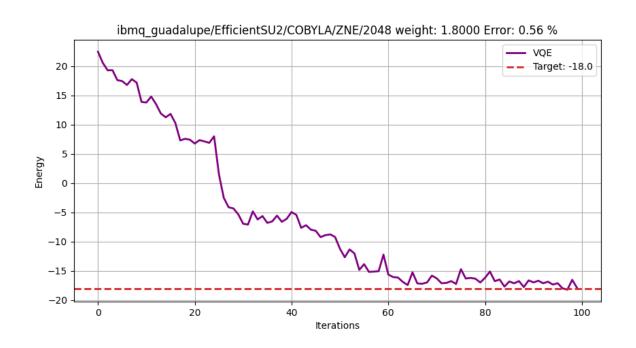
Execution time (s): 23294.63

Expected ground state energy: -18.00000000

Computed ground state energy: -17.89892578

Result eigen value: -17.89892578

Relative error: 0.5615%



Our Results: Solution 3 - Error 0.85%

Backend: ibmq_guadalupe

Ansatz: ExcitationPreserving (reps=1,entanglement='linear')

Optimizer: SPSA(maxiter=100)

Resilience: ZNE (2)

Shots: 2048

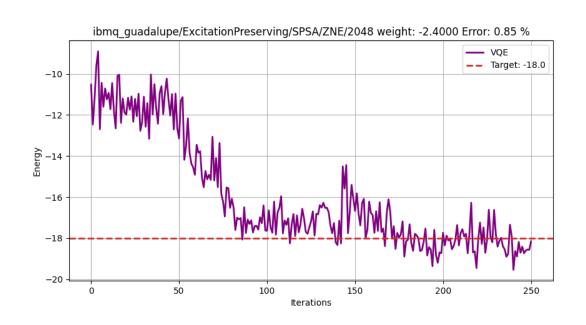
Edge weight: -2.4000

Execution time (s): 113953.17

Expected ground state energy: -18.00000000 Computed ground state energy: -18.15273438

Result eigen value: -18.15273438

Relative error: 0.848%



Other Solution - Error 7%

Backend: ibmq_guadalupe

Ansatz: EfficientSU2 (reps=1, entanglement='reverse_linear')

Optimizer: SPSA(maxiter=100)

Resilience: ZNE(2)

Shots: 2048

Edge weight: 1.9000

Execution time (s): 102658.02

Expected ground state energy: -18.00000000

Computed ground state energy: -19.38377279

Result eigen value: -19.38377279

Relative error: 7.687%

