



Novel Variational Quantum Algorithms for MaxCut

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To Kika, my loyal companion and source of endless joy.

Declaration

I declare that this document is an original work of my own authorship and that it fulfills all the requirements of the Code of Conduct and Good Practices of the Universidade de Lisboa.

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Resumo

Neste trabalho, apresentamos o *Interpolated QAOA/QEMC (iQAQE) Framework*, uma abordagem inovadora inspirada no *Quantum Approximate Optimization Algorithm (QAOA)* e no *Qubit-Efficient Max-Cut Heuristic Algorithm (QEMC)*, para desenhar múltiplos Algoritmos Quânticos Variacionais (VQAs) distintos para resolver o problema do Corte Máximo (MaxCut). Este *framework* baseia-se nos componentes centrais do QEMC enquanto integra conceitos do QAOA para aproveitar os pontos fortes de ambos os algoritmos. O *Framework* iQAQE requer menos *qubits* que o QAOA e apresenta maior resiliência à incerteza estatística associada a um pequeno número de medições (*shots*) em comparação com o QEMC.

O *framework* oferece uma gama de parâmetros ajustáveis, facilitando a criação de diversos VQAs. Introduzimos heurísticas para a seleção desses parâmetros, como o número de *qubits*, a cardinalidade das listas e os mapeamentos, e avaliamos o seu desempenho. Os nossos resultados indicam que o iQAQE frequentemente apresenta um desempenho comparável ao QEMC e pode até superar métodos clássicos de ponta, como Goemans-Williamson, em certos cenários.

Além disso, propomos duas abordagens alternativas para resolver o problema do MaxCut, derivadas da nossa investigação sobre o iQAQE. Embora esses métodos não se enquadrem diretamente no *Framework* iQAQE, oferecem *insights* valiosos e potenciais caminhos para futuras pesquisas.

Também apresentamos um pequeno modelo de aprendizagem automática, projetado para determinar o mapeamento óptimo para um grafo específico com base em propriedades estatísticas dos próprios mapeamentos. Em última análise, o *Framework* iQAQE serve como um banco de testes versátil para o desenvolvimento de novos VQAs, eventualmente abrindo caminho para resultados revolucionários no futuro.

Palavras-chave: Computação Híbrida Quântica-Clássica (HQCC), Algoritmos Quânticos Variacionais (VQAs), Problema do Corte Máximo (MaxCut), *Quantum Approximate Optimization Algorithm* (QAOA), *Qubit-Efficient MaxCut Heuristic Algorithm* (QEMC).

Abstract

In this work, we introduce the Interpolated QAOA/QEMC (iQAQE) Framework, a novel approach inspired by the Quantum Approximate Optimization Algorithm (QAOA) and the Qubit-Efficient MaxCut Heuristic Algorithm (QEMC), for designing multiple distinct Variational Quantum Algorithms (VQAs) to solve the Maximum Cut (MaxCut) problem. This framework builds on the core components of QEMC while integrating concepts from QAOA to harness the strengths of both algorithms. The iQAQE Framework requires fewer qubits than QAOA and exhibits greater resilience to statistical uncertainty associated with small shot numbers compared to QEMC.

The framework offers a range of adjustable parameters, facilitating the creation of various VQAs. We introduce heuristics for selecting these parameters, such as the number of qubits, list cardinality, and mappings, and evaluate their performance. Our findings indicate that iQAQE often performs on par with QEMC and can even surpass classical state-of-the-art algorithms like Goemans-Williamson in certain scenarios.

Additionally, we propose two alternative approaches for solving the MaxCut problem, derived from our investigations on iQAQE. While these methods do not fall directly within the iQAQE Framework, they offer valuable insights and potential avenues for future research.

We also present a small machine learning model designed to determine the optimal mapping for a specific graph based on statistical properties of the mappings themselves. Ultimately, the iQAQE Framework serves as a versatile testbed for developing new VQAs, potentially paving the way for groundbreaking results in the future.

Keywords: Hybrid Quantum-Classical Computing (HQCC), Variational Quantum Algorithms (VQAs), Maximum Cut (MaxCut) Problem, Quantum Approximate Optimization Algorithm (QAOA), Qubit-Efficient MaxCut Heuristic Algorithm (QEMC).

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Nomenclature

Greek symbols

$ \psi\rangle$	Arbitrary pure state
ρ	Density matrix or performance guarantee
$\sigma_x, \sigma_y, \sigma_z$	Pauli matrices

Roman symbols

X, Y, Z	Pauli matrices (alternative notation)
H_C or H_P	Cost or problem Hamiltonian
H_M or H_B	Mixer or bias Hamiltonian
$R_x(\theta), R_y(\theta), R_z(\theta)$	Rotation gates (of angle θ) around the x, y and z axes on the Bloch sphere
S_n	n -dimensional unit sphere
Tr	Trace [of a matrix]

Subscripts

x, y, z	Cartesian components
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Superscripts

\dagger	Hermitian adjoint
T	Transpose

List of acronyms

AR	Approximation Ratio
BSF	Best-so-far
D-CNOT	Deterministic CNOT
DV	Dependent Variable
GW	Goemans-Williamson (Algorithm)
HPC	High Performance Computing
HQCC	Hybrid Quantum-Classical Computing
iQAQE	Interpolated QAOA/QEMC (Hybrid Algorithm)
IV	Independent Variable
MaxCut	Maximum Cut (Problem)
ML	Machine Learning
ND-CNOT	Non-deterministic CNOT
NISQ	Noisy Intermediate-Scale Quantum
OLS	Ordinary Least Squares
PC	Principal Component
PCA	Principal Component Analysis
PCR	Principal Component Regression
PQC	Parameterized Quantum Circuit
QAOA	Quantum Approximate Optimization Algorithm
QEMC	Qubit-Efficient MaxCut Heuristic (Algorithm)
QKD	Quantum Key Distribution
QML	Quantum Machine Learning
SAT	Boolean Satisfiability Problem
SDP	Semidefinite Programming
SVR	Support Vector Machine Regression
TSP	Traveling Salesman Problem
VIF	Variance Inflation Factor
VQA	Variational Quantum Algorithm
VQE	Variational Quantum Eigensolver

Chapter 1

Introduction

Over the last few decades, significant progress has been achieved in the field of quantum computing. This is an entirely new computing paradigm, based on exploiting the fundamental principles of quantum mechanics to our advantage. Although the concept of a quantum computer has been around for a little over 40 years [1], only rather recently, in the present century, have we successfully engineered elementary prototypes for such cutting-edge devices. Scientists and engineers worldwide are now working towards the development of practical quantum computers, which are expected to revolutionize the way we solve complex problems in various fields, such as cryptography, optimization, drug discovery, material science, machine learning, and many more.

However, with the current generation of quantum machines, so-called Noisy Intermediate Scale Quantum (**NISQ**) devices, we are still far from achieving the promised quantum advantage. This refers to the point at which a quantum computer can outperform its classical counterpart by solving specific problems considerably faster or more efficiently. It's the "holy grail" of quantum computing and what drives the research in this field forwards.

Presently, the most promising candidates for achieving meaningful quantum advantage are anchored in what has become known as "Hybrid Quantum-Classical Computing" (**HQCC**). This approach aims to merge the strengths of both worlds: the computational power of quantum and the stability of classical computing. In this context, Variational Quantum Algorithms (**VQAs**) have been at the forefront of research, as they are particularly well-suited for near-term quantum devices, requiring fewer qubits and featuring shallower circuit depths. These algorithms are hybrid by design, using a classical optimizer to adjust the parameters of a Parameterized Quantum Circuit (**PQC**).

Amid the many prospective applications of quantum computing, notable advances have been made in recent years in solving combinatorial optimization problems. **VQAs** like the Quantum Approximate Optimization Algorithm (**QAOA**) [2] have demonstrated potential for tackling the Maximum Cut (**MaxCut**) problem, a challenging graph-based NP-hard problem in computer science. Despite its promise, however, **QAOA** has a significant drawback: it requires a large number of qubits, exceeding the capacity of current quantum devices, when scaled to larger problem instances. This constraint has prompted the development of alternative approaches such as the Qubit-Efficient **MaxCut** Heuristic Algorithm (**QEHC**)

[3], designed to address the `MaxCut` problem using fewer qubits. Meanwhile, the search for better algorithms to solve this problem continues, and the development of new hybrid quantum-classical methods is crucial to achieving quantum advantage.

1.1 Motivation

The search for efficient algorithms to solve combinatorial optimization problems is critical in numerous fields, such as logistics, finance, and telecommunications. The `MaxCut` problem, for example, has broad applications in areas like machine learning [4], statistical physics [5], circuit design [5], and data clustering [6]. Creating more efficient algorithms to solve this problem can improve the performance of these applications, driving substantial progress in their respective fields.

Moreover, there is a strong interest from the computer science community in terms of computational complexity. The `MaxCut` problem is recognized as NP-hard, and the search for efficient algorithms to solve it may offer valuable insights into the boundaries between classical and quantum computing. It might even contribute to unraveling one of the most perplexing questions in theoretical computer science: the $P = NP$ problem. Imagining a world where the $P = NP$ conjecture is proven true, albeit improbable, is intriguing. It would mean that every problem in NP could be solved in polynomial time, including `MaxCut`. This would also extend to problems like the Traveling Salesman Problem (`TSP`) and the Knapsack Problem, among others. The implications would be profound, as this would dramatically increase our ability to solve previously difficult optimization problems, with direct applications in areas like vehicle routing, job scheduling, and broader logistics. Additionally, the impact on cryptography would be as significant – if not more so – since many cryptographic techniques rely on the complexity of NP problems. For example, integer factorization is a key component of RSA (Rivest-Shamir-Adleman) encryption, a popular asymmetric encryption algorithm used extensively in public-key cryptography for secure message transmission over the internet. If $P = NP$, RSA encryption could easily be broken, leading to major security risks. Hence, the importance of studying these problems to fully understand their complexity.

The aforementioned considerations drive our efforts to develop a new algorithm for solving the `MaxCut` problem with greater efficiency and accuracy. This algorithm, the Interpolated QAOA/QEMC (`iQAQE`) Hybrid Algorithm, will be the focus of this thesis. More specifically, we will develop what we call the `iQAQE` Framework, which will serve as the foundation for developing numerous distinct `iQAQE` algorithms.

1.2 Topic Overview

In this project, we propose the Interpolated QAOA/QEMC (`iQAQE`) Framework. This will enable us to develop novel VQAs by leveraging the strengths of two existing VQAs, `QAOA` and `QEMC`, for improved performance in solving the `MaxCut` problem. `QAOA` requires a qubit for each graph vertex, making it difficult to scale. In contrast, `QEMC` uses exponentially fewer qubits by assigning one basis state to

each graph node, requiring only $\log_2(n)$ qubits (for n graph vertices). However, this compression leads to limitations in **QEMC**'s results. By interpolating both **VQAs**, we aim to create an algorithm that utilizes fewer qubits than **QAOA** and performs better than **QAOA** and **QEMC**. The new algorithm, constructed through the **iQAQE** Framework, assigns multiple basis states to each node, in contrast to **QEMC**'s single basis state approach. This design tentatively allows for a more practical implementation on present-day **NISQ** devices, thanks to its reduced qubit requirements compared to **QAOA**, fewer measurement shots than **QEMC**, and potentially greater trainability than **QAOA**.

1.3 Objectives

The primary objective of this thesis is to develop and analyze the **iQAQE** Framework. Algorithms built within this framework will be implemented and tested using classical simulations of quantum machines. The deliverables for this project include the **iQAQE** Framework's code¹, the results obtained from the simulations, and the analysis of these results. The expected outcomes are improvements in the performance of the **iQAQE** algorithm compared to **QAOA** and **QEMC**, with a focus on accuracy, efficiency, and scalability.

1.4 Thesis Outline

This thesis is structured as follows: Chapter 2 provides an overview of the background concepts related to quantum computing, variational quantum algorithms, and the **MaxCut** problem. Chapter 3 introduces the hybrid quantum-classical **iQAQE** Framework, explaining its design, advantages, and potential applications. Chapter 4 details the numerical implementation of the **iQAQE** Framework and individual algorithms **QAOA** and **QEMC**, in addition to the Goemans-Williamson (**GW**) algorithm. It also describes the benchmarking and testing methods used to evaluate the algorithms' performances. In Chapter 5, we delve into the different **iQAQE** schemes developed and the corresponding results obtained from simulations. These are compared with **QAOA**, **QEMC**, and **GW**. Chapter 6 discusses alternative approaches to solving the **MaxCut** problem, derived from our research on **iQAQE**. Finally, Chapter 7 concludes the thesis, summarizing the main findings and suggesting future research directions.

¹Reach out to the author to obtain access to the code.

Chapter 2

Theoretical background

This chapter covers the key concepts and theoretical background necessary to understand the work presented in this thesis. It begins by describing the `MaxCut` problem and discussing elements of computational complexity, reinforcing the rationale behind this research. The state-of-the-art algorithms for solving the `MaxCut` problem, both classical and hybrid quantum-classical, are then examined. After a brief introduction to quantum computing, the chapter explores hybrid quantum-classical computing and variational quantum algorithms in greater depth. Specifically, the Quantum Approximate Optimization Algorithm (`QAOA`) and the Qubit-Efficient MaxCut Heuristic (`QEMC`) are analyzed, as they play a pivotal role in this study.

2.1 The Maximum Cut Problem & Computational Complexity

The `MaxCut` problem, a fundamental problem in graph theory and combinatorial optimization, involves partitioning a graph $G = (V, E)$ into two disjoint subsets, S_1 and S_2 , such that the number of edges connecting vertices from different subsets is maximized. Formally, the objective is to find a partition (S_1, S_2) that maximizes:

$$\text{Cut}(S_1, S_2) = \sum_{(u,v) \in E} \chi(u, v), \quad (2.1)$$

where $\chi(u, v) = 1$ if u and v belong to different subsets, and $\chi(u, v) = 0$ if they belong to the same subset. This quantity is referred to as the "cut" of the partition (S_1, S_2) . Pictorially, this can be represented as cutting the edges of the graph (Figure 2.1), hence the name `MaxCut`. What we describe here is the undirected, un-weighted `MaxCut` problem. A more general formulation would involve the specific graph's adjacency matrix, W_{ij} .

Since the `MaxCut` problem is NP-hard, finding an optimal solution efficiently is a significant computational challenge, especially as the graph size grows. However, researchers have developed various approximation algorithms and heuristics to approach near-optimal solutions in a reasonable time, including both classical and quantum approaches (cf. subsections 2.1.1 and 2.1.2). These methods are

designed to tackle the intrinsic complexity of the problem, providing practical solutions that have real-world applications. As mentioned earlier, the `MaxCut` problem finds use in a variety of fields, including machine learning [4], statistical physics [5], circuit design [5], and data clustering [6].

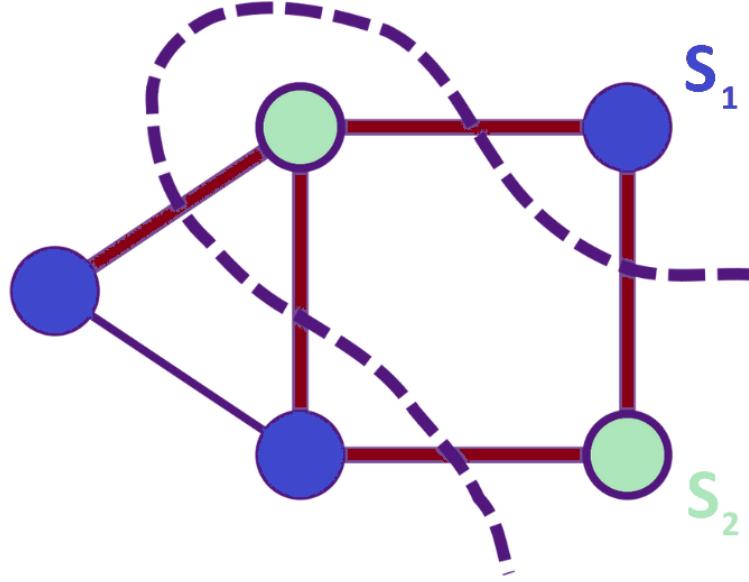


Figure 2.1: An example of a graph with a partition that maximizes the number of cut edges (red). Note that the `MaxCut` partition might not be unique.

We previously mentioned that the `MaxCut` problem is well-known to be NP-hard. (Indeed, it [decision version] even appears on Karp's original list of NP-complete problems, from 1972 [7].) Let's explore what this means in more detail. The computational complexity class of a problem is defined by the computational resources required to solve it, such as time or memory. Although there are many complexity classes (see the [Complexity Zoo](#)), the following four are key to our discussion:

1. **P (Polynomial time):** Problems that can be solved in polynomial time by a deterministic Turing machine, indicating they are computationally efficient.
2. **NP (Non-deterministic Polynomial time):** Problems that can be verified in polynomial time if given a solution, but finding the solution may not be as straightforward. Therefore, if you guess a solution, you can confirm its correctness quickly.
3. **NP-hard (Non-deterministic Polynomial time hard):** Problems that are at least as challenging as the most difficult problems in NP. If you could solve an NP-hard problem in polynomial time, you would be able to solve all problems in NP efficiently.
4. **NP-complete (Non-deterministic Polynomial time complete):** Problems that belong to NP but are also NP-hard. If you can solve one NP-complete problem efficiently, all problems in NP would become efficiently solvable.

The MaxCut problem in its **optimization form** – "Given a graph G , find the maximum cut" – is NP-hard. This implies that if you could find an efficient solution to MaxCut, you could also solve all other NP problems reliably. That would include, e.g., the boolean SATisifiability problem (**SAT**), knapsack problem, decision version of the traveling salesman problem, clique problem, and more [8]. This potential for broad application drives interest in developing high-performance algorithms for MaxCut. Alternatively, the MaxCut problem can be formulated as a **decision problem** – "Given a graph G and an integer k , determine if there is a cut of size at least k in G ." This formulation is NP-complete. NP-complete status indicates that a problem is both in NP, allowing polynomial-time verification, and at least as hard as any other problem in NP, making it central to computational theory and offering crucial insights into the boundary between efficiently solvable and intractable problems. Although our focus is on the NP-hard version of the problem, it's important to understand the significance of the NP-complete formulation.

2.1.1 Classical State-of-the-Art Algorithms for MaxCut

Over the years, a variety of algorithms have been proposed to solve the MaxCut problem, ranging from exact solutions to various approximations, including heuristics and other techniques. In this work, we focus on scaling these algorithms to handle bigger graphs, which is why exact algorithms are not discussed – they quickly become impractical for large instances. Instead, our attention is on approximation algorithms, which are more suitable for larger scales¹.

Approximation algorithms aim to find solutions that are close to optimal, often with a guaranteed Approximation Ratio (**AR**). This is a measure of the performance of an approximation algorithm, representing the worst-case deviation between the algorithm's solution and the optimal solution. Formally, this can be expressed as: (For the Goemans-Williamson (**GW**) Algorithm [9], e.g.)²

$$r_A^{GW} = \frac{\text{Cut}_{\text{GW}}}{\text{MaxCut}} \approx 0.87856,$$

where Cut_{GW} is the worst-case "cut" value obtained by the Goemans-Williamson Algorithm, and MaxCut is the MaxCut value. The **AR** is a crucial metric for evaluating the quality of an approximation algorithm, providing insight into its effectiveness. The closer the **AR** is to 1, the better the algorithm's performance. In literature, this worst-case approximation ratio might also be referred to as the *performance guarantee* ρ , often appearing alongside designations such as ρ -approximation algorithm. Using this nomenclature, the Goemans-Williamson Algorithm is a 0.87856-approximation algorithm for MaxCut (i.e., $\rho = 0.87856$).

Building on this, two commonly used **approximation algorithms** are:

- **Goemans-Williamson Algorithm** [9]: This popular approximation algorithm achieves a performance guarantee of approximately 0.87856, leveraging Semidefinite Programming (**SDP**) and a randomized rounding technique. (This will be further developed below, in subsubsection Goemans-Williamson Algorithm.)

¹In this section, "larger/bigger graphs" will often refer to graphs with many hundreds to a few thousand nodes.

²The actual value is $\alpha = \min_{0 \leq \theta \leq \pi} \frac{2}{\pi} \frac{\theta}{1 - \cos \theta} > 0.878$.

- **Spectral Methods** [10]: These techniques employ eigenvalues and eigenvectors of the adjacency matrix to identify effective cuts. While spectral methods may offer quicker computational speeds compared to other approximation strategies (e.g. GW), they may not consistently provide the optimal approximation ratios. Performance-wise, the leading spectral algorithm, Trevisan's algorithm, using Soto's refined analysis [11], features a worst-case approximation ratio of 0.614.

Heuristic and **metaheuristic algorithms** offer another approach, focusing on finding solutions more quickly, albeit without guaranteed approximation ratios. They are typically employed when computational efficiency is prioritized over achieving the best possible approximation ratio. These algorithms include:

- **Greedy Algorithms** (Described in [10]): These build solutions incrementally, following a greedy approach. Mathematically, we start with $S_1, S_2 = \emptyset$. In each step of the algorithm, we choose a vertex $v \in V$ yet to be assigned to S_1 or S_2 . Then, we add v to S_1 or S_2 by choosing the larger of the two cuts $(S_1 \cup \{v\}, S_2)$ and $(S_1, S_2 \cup \{v\})$ at this step. Repeat until all vertices are assigned to a set.
- **Simulated Annealing** [12]: Employs a probabilistic optimization approach, inspired by metallurgical annealing, to iteratively explore the solution space, gradually reducing the acceptance of worse solutions ("temperature") over time to converge towards near-optimal solutions.
- **Genetic Algorithms** [13]: These algorithms, rooted in evolutionary principles, employ selection, crossover, and mutation techniques to iteratively generate solutions for the MaxCut problem, evolving from an initial population of diverse partitions towards improved solutions.

Reiterating, these approximation and heuristic methods offer scalable solutions for the MaxCut problem, making them more suitable for larger graphs where exact methods are not feasible.

Of all the mentioned approaches, we will now elaborate on the Goemans-Williamson algorithm, which currently stands as the most effective approximation method for MaxCut, boasting the best-known performance guarantee to date (0.87856). Its detailed explanation is particularly relevant as we plan to utilize it in benchmarking our proposed algorithms later on. We opt for GW over other options for one obvious reason: Our aim is to push the boundaries of MaxCut algorithms, thus it is logical to compare against the top-performing method. (We do not use any of the heuristic/metaheuristic algorithms, as they do not provide performance guarantees, and we are interested in having a benchmark to compare against.)

2.1.1.1 Goemans-Williamson Algorithm

In this subsubsection, we shall present the Goemans-Williamson Algorithm in greater detail. This approximation algorithm, developed by Michel X. Goemans and David P. Williamson in 1995 [9], is a cornerstone in the field of combinatorial optimization, particularly for the MaxCut problem. The algorithm leverages semidefinite programming to construct a solution that is then rounded to provide a near-optimal cut. Semidefinite programming involves optimizing a linear function of a symmetric matrix while adhering to linear equality constraints and the requirement that the matrix be positive semidefinite. This programming paradigm finds utility across diverse fields such as control theory, nonlinear programming,

geometry, and combinatorial optimization, showcasing its versatility and applicability. (These applications are detailed in [9] and references therein.)

Formally, the Goemans-Williamson Algorithm can be described as follows. (We will use the original formulation, from [9], presenting a distilled version of the most important aspects of their description.) Given a graph, with vertex set $V = \{1, \dots, n\}$ and non-negative weights $w_{ij} = w_{ji}$ for each pair of vertices i and j , the weight of the maximum cut $w(S_1, S_2)$ is given by the following integer quadratic program³:

$$\begin{aligned} & \text{Maximize } \frac{1}{2} \sum_{\substack{i < j: \\ (i,j) \in E}} w_{ij}(1 - y_i y_j) \\ (Q) \quad & \text{Subject to: } y_i \in \{-1, 1\} \quad \forall i \in V. \end{aligned} \tag{2.2}$$

It is clear that the sets $S_1 = \{i | y_i = 1\}$ and $S_2 = \{i | y_i = -1\}$ correspond to a cut of weight $w(S_1, S_2) = \frac{1}{2} \sum_{(i,j) \in E} w_{ij} (1 - y_i y_j)$. This represents the same quantity as the aforementioned $\text{Cut}(S_1, S_2) = \sum_{(u,v) \in E} \chi(u, v)$ (Eq. 2.1), in the un-weighted case. Explicitly, in words, the sum in Eq. 2.2 is to be taken over all the graph's edges (i, j) , with y_i representing node- i 's associated set: $y_i = +1$, if node- $i \in S_1$, or $y_i = -1$, when node- $i \in S_2, \forall i \in V$.

Given that solving this integer quadratic program is NP-hard [9], we explore relaxations of (Q) . These are obtained by relaxing some of the constraints of (Q) , and extending the objective function to the larger space. Consequently, all solutions of (Q) are feasible for the relaxation, and the optimal value of the relaxation serves as an upper bound for (Q) 's optimal value. We interpret (Q) as constraining y_i to be a 1-dimensional unit vector. A number of possible relaxations involve allowing y_i to be a multidimensional vector v_i of unit Euclidean norm. As the linear space spanned by the vectors v_i has dimension at most n (number of graph nodes), we can assume that these vectors belong to \mathbb{R}^n . More precisely, they belong to the n -dimensional unit sphere, S_n . At this point, so as to make sure that the resulting optimization problem is, indeed, a relaxation, we must define the cost function such that it reduces to $\frac{1}{2} \sum_{i < j} w_{ij}(1 - y_i y_j)$ for vectors lying in a 1-dimensional space. One way to do this is to replace $y_i y_j$ with the inner product $v_i \cdot v_j$ in the expression for $w(S_1, S_2)$. The resulting relaxation, (P) , reads:

$$\begin{aligned} & \text{Maximize } \frac{1}{2} \sum_{\substack{i < j: \\ (i,j) \in E}} w_{ij}(1 - v_i \cdot v_j) \\ (P) \quad & \text{Subject to: } v_i \in S_n \quad \forall i \in V. \end{aligned} \tag{2.3}$$

In the context of this semidefinite programming relaxation, the GW algorithm follows these steps:

1. Solve (P) , thus obtaining an optimal set of vectors v_i ;
2. Let r be a vector uniformly distributed on the unit sphere, S_n ;
3. Set $S_1 = \{i | v_i \cdot r \geq 0\}$ and $S_2 = \{i | v_i \cdot r < 0\}$.

In other words, we choose a random hyperplane through the origin (with r as its normal), in n dimensions, and partition the vertices according to whether the vectors v_i lie "above" ($v_i \cdot r \geq 0$) or "below"

³Note that we use the general formulation here, contemplating the possibility of weighted graphs, i.e., we do not restrict ourselves to un-weighted graphs.

$(v_i \cdot r < 0)$ this hyperplane. Those above the hyperplane are assigned to S_1 , and those below to S_2 .

Furthermore, it can be shown that the **GW** algorithm has a performance guarantee of:

$$\alpha = \min_{0 \leq \theta \leq \pi} \frac{2}{\pi} \frac{\theta}{1 - \cos \theta} > 0.878. \quad (2.4)$$

A detailed proof elucidating the appearance of this 0.878 value is present in section 3 of [9].

Computationally, one defines the positive semidefinite matrix $X \in \mathbb{R}^{n \times n}$, with $X_{ii} = 1$ and $X_{ij} = v_i \cdot v_j = v_i^T v_j, \forall i, j \in V$. Then, we can solve the following semidefinite program:

$$\begin{aligned} \text{Maximize } & \frac{1}{2} \sum_{\substack{i < j \\ (i,j) \in E}} w_{ij}(1 - X_{ij}) \\ (N) \quad \text{Subject to: } & X_{ii} = 1, \forall i \in V, X \succeq 0. \end{aligned} \quad (2.5)$$

We do possess effective tools and techniques for solving semidefinite programs (e.g., interior point methods), enabling the efficient determination of a highly satisfactory X matrix. From there, we can extract the vectors v_i , by performing the square root operation on the matrix X . The resulting matrix's columns, or rows – X is symmetric – will correspond to the desired vectors v_i . This outlines the process for solving (P) as listed earlier.

2.1.2 Hybrid Quantum-Classical State-of-the-Art Algorithms for MaxCut

A number of different hybrid quantum-classical algorithms have been proposed to solve the **MaxCut** problem, leveraging the unique properties of quantum computing to potentially outperform purely classical algorithms. They are designed to exploit quantum superposition and entanglement to explore the solution space more efficiently. In this section, we will briefly mention the state-of-the-art hybrid quantum-classical algorithms for **MaxCut**, focusing on the Quantum Approximate Optimization Algorithm (**QAOA**) and the Qubit-Efficient **MaxCut** Heuristic Algorithm (**QEBC**). Due to their variational nature, these algorithms do not have a theoretical performance guarantee, but they have shown promising results in practice.

Presently, the Quantum Approximate Optimization Algorithm (**QAOA**), initially introduced in [2], is one of the top candidates for achieving meaningful quantum advantage with near-term quantum devices. Although, **MaxCut**-wise, it has a substantial limitation: it requires a number of qubits equal to the number of graph nodes. This is a significant drawback, as it severely limits the algorithm's scalability to bigger graphs. Hence, why it is believed that "*QAOA for Max-Cut requires hundreds of qubits for quantum speed-up*" [14]. After all, our classical algorithms work just fine for small graphs, which is why we do not present performance metrics for basic **QAOA**. Alternatively, different **QAOA** variations have been proposed to address the issue of high qubit requirements, such as **QAOA-in-QAOA** [15], which partitions a large graph into many smaller subgraphs, each of which is easily solved using a separate **QAOA** instance. Afterwards, these results are joined together by working through another, this time **weighted**, **MaxCut** problem. This approach in particular has proven to yield competitive or even better performance

over the best known classical algorithms, i.e. **GW**, for graphs up to 2000 nodes.

A separate problem one might face when implementing **QAOA**, that also hinders its application to larger graphs, has to do with qubit connectivity⁴, as **QAOA** might require specific connectivity patterns that are not natively available in present-day **NISQ** devices. To implement these, we are often required to utilize a number of 2-qubit **SWAP** gates, which degrade the algorithm's performance by introducing extra noise in the system, and hence errors in our results. Therefore, a parallel line of work has been to develop **QAOA** variations that have lesser connectivity requirements, in the sense that they do not depend on arbitrary qubit connectivity. For example, **Parity-QAOA** [16, 17] was designed for quantum chips with planar lattice geometry (Figure 2.2), requiring only nearest-neighbour connectivity. Although it necessitates more qubits, it entirely solves this problem.

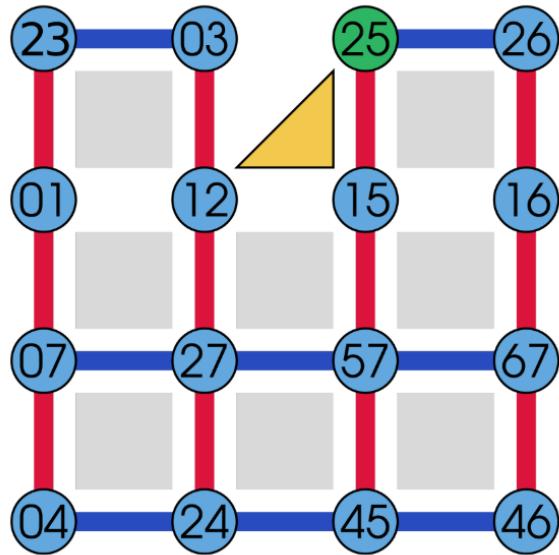


Figure 2.2: Planar lattice geometry illustration: the circular nodes represent the physical qubits, each of which can interact directly with its nearest neighbours. The lines represent terms of the driver Hamiltonian, from **Parity-QAOA** [16]. The numbers refer to edges in the original graph. Image reproduced from [16].

Furthermore, the Qubit-Efficient **MaxCut** Heuristic Algorithm (**QEMC**), introduced in [3], is a quantum-classical hybrid algorithm that aims to solve the **MaxCut** problem using fewer qubits than **QAOA**, therefore extending its applicability to larger graphs. It has shown cutting-edge performance in practice, occasionally surpassing the classical state-of-the-art, i.e. **GW**, for graphs up to 2048 nodes. The novelty of this algorithm is that it opens up the door for quantum devices to be used for large **MaxCut** instances, previously unfeasible due to qubit number limitations, which rendered **QAOA** computations for such large graphs impossible. As was already mentioned, **QEMC** allows for an exponential compression in the number of needed qubits, enabling its utilization on today's quantum hardware for considerably larger graphs than what would be possible with traditional **QAOA**. However, this exponential compression also makes it efficiently simulable classically, which effectively defeats its purpose as a quantum

⁴Qubit connectivity refers to the pattern or arrangement of connections between qubits in a quantum computing system, determining which qubits can interact directly with each other.

algorithm. For this reason, QEMC is termed a quantum-inspired classical algorithm and, by definition, will not produce any quantum advantage. Either way, we include it in this subsection, since it can always be implemented as a VQA on current-day NISQ devices.

More recently, different algorithms have been proposed in an attempt to reduce the number of qubits required for larger graphs, without falling into the trap of efficient classical simulability. (After all, we are looking for quantum advantage.) Polynomial compression-based algorithms, such as that proposed in [18], have too shown promising results. This one in specific [18] is, to the best of our knowledge, boasting the highest quality attained experimentally on sizes up to 7000 graph nodes, competitive with state-of-the-art classical solvers. Not only that, but their specific qubit-efficient encoding brings in a super-polynomial mitigation of barren plateaus⁵ as a built-in feature, which constitutes a significant advantage over other VQAs. Curiously enough, our initial idea for the iQAQE Framework is quite similar to what is explored in [18]. However, we believe our concept to be somewhat more general, by not assuming *a priori* a polynomial compression in the number of qubits.

2.2 Quantum Computing Primer

In this section, we present a brief introduction to quantum computing, focusing on its fundamental ideas and principles. We begin with an overview of the general concepts, followed by a discussion of quantum gates and circuits, crucial for understanding quantum algorithms. This groundwork will prepare us to delve into the specifics of variational quantum algorithms and their application to the MaxCut problem.

General concepts and mathematical formalism

(This description is inspired by [19].)

Quantum computing makes use of the foundational principles of quantum mechanics in an attempt to extract some sort of quantum advantage from them. Instead of using classical binary digits, quantum computers use their quantum analog, qubits. In practice, qubits can be any two-state quantum system. Said states are, unequivocally, associated with the states $|0\rangle$ and $|1\rangle$, conventionally, very much like in classical computing when one uses bits. The primary distinction lies in the fact that a qubit, being a quantum system, can exist in a superposition of both states generally denoted as: (We use Dirac notation.)

$$\mathcal{H}^2 \ni |\psi\rangle = \alpha|0\rangle + \beta|1\rangle, \quad (2.6)$$

where $\alpha, \beta \in \mathbb{C}$, with $|\alpha|^2 + |\beta|^2 = 1$. This ensures probability normalization to 1, as, according to the famous Born rule, the probability of measuring the qubit in state $|0\rangle$ is $|\langle 0|\psi\rangle|^2 = |\alpha|^2$, and in state $|1\rangle$ is $|\langle 1|\psi\rangle|^2 = |\beta|^2$.

In the same vein, one can compose (represented by the tensor product) n qubits to obtain an n -qubit

⁵In VQA analysis, barren plateaus are characterized by a vanishing expectation value of $\nabla \mathcal{L}$ over random parameter initializations and an exponential decay (in the number of qubits, n) of its variance, where \mathcal{L} refers to the loss function. These are flat regions of the parameter space, impeding optimization.

state:

$$|\psi\rangle = \bigotimes_{i=0}^{n-1} (\alpha_i |0\rangle_i + \beta_i |1\rangle_i). \quad (2.7)$$

More generally, we can express an arbitrary pure state of n qubits, $|\psi\rangle$, by a normalized linear combination of the computational basis states $|b_0 b_1 \dots b_{n-1} b_n\rangle := |b_0\rangle \otimes |b_1\rangle \otimes \dots \otimes |b_{n-1}\rangle \otimes |b_n\rangle$:

$$|\psi\rangle = \sum_{\mathbf{b} \in \{0,1\}^n} \alpha_{\mathbf{b}} |b_0 b_1 \dots b_{n-1} b_n\rangle, \quad (2.8)$$

where $\sum_{\mathbf{b} \in \{0,1\}^n} |\alpha_{\mathbf{b}}|^2 = 1$. This superposition phenomenon (mathematically characterized by the linear combination of basis states), utterly impossible in classical physics, brings unprecedented computational power under certain scenarios, as it introduces a kind of built-in parallelism in quantum computing, a highly desirable feature for any computational scheme.

In addition to this, quantum computing also exploits entanglement as a resource. Entanglement allows for intricate correlations between qubits, which have no classical counterpart, and are believed to offer computational speed-ups, although it is yet unclear exactly how. An entangled state, by definition, is one that cannot be expressed as the tensor product of individual qubit states, i.e., it cannot be represented in the form of Eq. 2.7. The most famous example of entangled states are the Bell states, representing what are known as two-qubit (or bipartite) maximally entangled states. Such Bell states are reproduced below:

$$\begin{aligned} |\Phi^+\rangle &= \frac{1}{\sqrt{2}}(|00\rangle + |11\rangle), & |\Phi^-\rangle &= \frac{1}{\sqrt{2}}(|00\rangle - |11\rangle), \\ |\Psi^+\rangle &= \frac{1}{\sqrt{2}}(|01\rangle + |10\rangle), & |\Psi^-\rangle &= \frac{1}{\sqrt{2}}(|01\rangle - |10\rangle). \end{aligned} \quad (2.9)$$

Indeed, these states cannot be expressed as a tensor product of two individual qubit states. From a practical standpoint, Bell states frequently serve as foundational elements in advanced techniques like quantum teleportation and quantum cryptography. Quantum Key Distribution (QKD) protocols, for example, frequently rely on entangled Bell pairs to distribute secure keys for encrypting and decrypting messages.

Another key topic in quantum information theory is the distinction between pure and mixed states. So far, we have only been treating pure states, which are represented by a single ket vector in some Hilbert space. Mixed states, on the other hand, are represented by density matrices, which are positive semidefinite matrices with unit trace. These matrices are used to describe a statistical ensemble of pure states. The density matrix ρ of a pure quantum state $|\psi\rangle$ is defined as:

$$\rho = |\psi\rangle\langle\psi|. \quad (2.10)$$

In the more general case of a statistical ensemble of pure states $|\psi_i\rangle$, each with probability p_i , the density matrix ρ is given by:

$$\rho = \sum_i p_i |\psi_i\rangle\langle\psi_i|. \quad (2.11)$$

It's crucial to avoid confusing this with superposition, as a mixed state comprises a statistical ensemble of pure states, while superposition pertains to a single state that is a linear combination of other states. For instance, the state $|\psi_S\rangle = \frac{1}{\sqrt{2}}(|0\rangle + |1\rangle)$ represents a superposition of $|0\rangle$ and $|1\rangle$, whereas the state $\rho_M = \frac{1}{2}(|0\rangle\langle 0| + |1\rangle\langle 1|)$ denotes a mixed state, signifying that half the states in the ensemble are in state $|0\rangle$ and the other half in state $|1\rangle$. Unlike the probabilistic mixture (ρ_M), the superposition ($|\psi_S\rangle$) can exhibit quantum interference, hence why it is important to distinguish between the two. If desired, this distinction can also be observed in the off-diagonal terms of the density matrix, referred to as coherences, which are responsible for quantum interference effects: the off-diagonal terms of ρ_M are zero, whereas those of $\rho_S = |\psi_S\rangle\langle\psi_S| = \frac{1}{4}(|0\rangle\langle 0| + |0\rangle\langle 1| + |1\rangle\langle 0| + |1\rangle\langle 1|)$ are not.

Types of qubits – physical implementations

Currently, numerous quantum computing architectures are in competition and the ultimate victor remains uncertain. Nevertheless, the most prominent architectures have qubits composed of either photons [20, 21], neutral atoms (using Rydberg states) [22, 23], trapped ions [24], semiconductors [25], or, most notably, superconducting circuits [26, 27] – employed by giants like IBM and Google in their quantum processors. Looping back to the definition of a qubit, considering a photonic quantum computer, it becomes rather straightforward and natural to map the photon's vertical and horizontal polarizations to the $|0\rangle$ and $|1\rangle$ states, respectively⁶. Once again, this mapping is analogous to how classical computing maps the current/no current states to 1 and 0.

The next step is to understand how these so-called quantum circuits act on our qubits, and how they can be used to perform quantum computations. This is where quantum gates come into play.

Quantum gates and circuits

Presently, quantum computers are being designed predominantly with a circuit-based architecture. These circuits serve as the quantum equivalent of classical logic circuits, however, instead of the typical AND, OR, and NOT gates, quantum circuits feature quantum gates. These gates apply specific transformations to qubits, based on their definitions, taking an initial state $|\psi\rangle$ to an output state $|\phi\rangle = U|\psi\rangle$, for some gate U , **designed to be unitary**, i.e., $U^\dagger U = UU^\dagger = I$, where I is the identity matrix. Unitary transformations preserve the normalization of quantum states and the inner product between states, which are essential properties in quantum mechanics. Additionally, they ensure that quantum operations are reversible, meaning that information is not lost during computation. This [unitary gates] can also be seen as a natural consequence of the Schrödinger equation, governing the time evolution of quantum systems, requiring said evolution to be described by a unitary operator.

The most frequently used quantum gates consist of rotations around each of the three Cartesian axes, represented as complex exponentials of the Pauli x , y and z matrices (σ_x , σ_y and σ_z). The Pauli matrices are defined as: (Note the alternative notation, X , Y and Z .)

$$X := \sigma_x = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}, \quad Y := \sigma_y = \begin{pmatrix} 0 & -i \\ i & 0 \end{pmatrix}, \quad Z := \sigma_z = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}. \quad (2.12)$$

⁶Note that there are alternative ways (other than polarization) to encode quantum information in photons for quantum computing.

Furthermore, when we mention rotations, these are to be seen in the Bloch sphere, which is a geometrical representation of the pure state space of a two-level quantum system, an illustration of which can be seen in Figure 2.3.

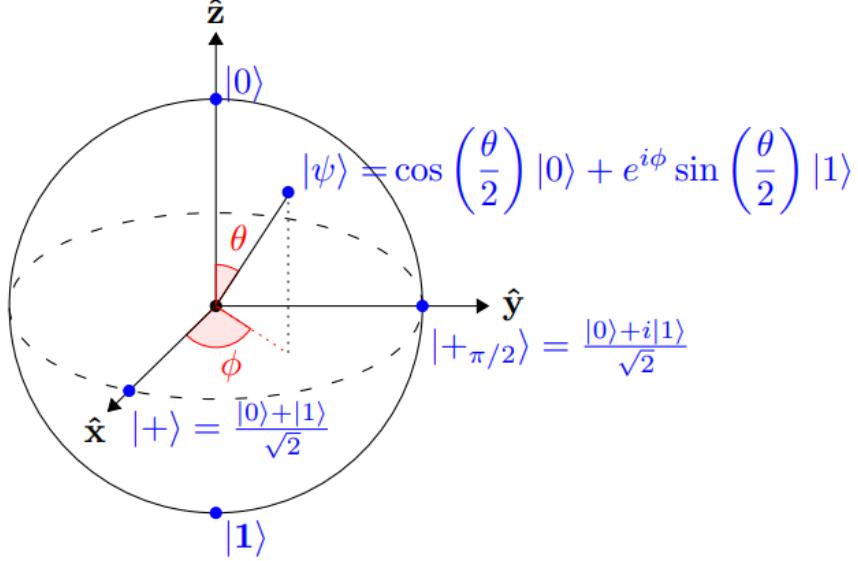


Figure 2.3: Bloch sphere representation. The sphere's north and south poles correspond to states $|0\rangle$ and $|1\rangle$, respectively. Reproduced from [28].

The aforementioned rotation gates can be written as: $R_x(\theta) = e^{-i\theta X/2}$, $R_y(\theta) = e^{-i\theta Y/2}$, and $R_z(\theta) = e^{-i\theta Z/2}$. In matrix form, this reads:

$$R_x(\theta) = \begin{pmatrix} \cos\left(\frac{\theta}{2}\right) & -i\sin\left(\frac{\theta}{2}\right) \\ -i\sin\left(\frac{\theta}{2}\right) & \cos\left(\frac{\theta}{2}\right) \end{pmatrix}, \quad R_y(\theta) = \begin{pmatrix} \cos\left(\frac{\theta}{2}\right) & -\sin\left(\frac{\theta}{2}\right) \\ \sin\left(\frac{\theta}{2}\right) & \cos\left(\frac{\theta}{2}\right) \end{pmatrix}, \quad R_z(\theta) = \begin{pmatrix} e^{-i\frac{\theta}{2}} & 0 \\ 0 & e^{i\frac{\theta}{2}} \end{pmatrix}, \quad (2.13)$$

These will frequently appear in quantum circuits, as they are the most basic quantum gates, in the sense that all other single-qubit gates can be re-constructed from them, i.e., for any unitary single-qubit gate U , one can always find a decomposition $U = e^{i\phi}R_Z(\gamma)R_X(\beta)R_Z(\alpha)$ (formally proven in [29]).

For illustrative purposes, to elucidate how single-qubit gates act on qubits, it is quite simple to verify that the σ_x matrix (alternative symbol, X) behaves exactly like a NOT gate. This correspondence is established through a 180° rotation around the x -axis. Mathematically,

$$X = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}, \quad (2.14)$$

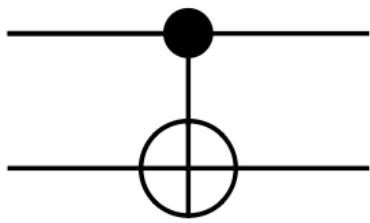
such that:

$$X |0\rangle = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix} \begin{pmatrix} 1 \\ 0 \end{pmatrix} = \begin{pmatrix} 0 \\ 1 \end{pmatrix} = |1\rangle. \quad (2.15)$$

Another essential gate found in most quantum algorithms is the Hadamard gate. It can be represented

as the following linear combination of X and Z matrices: $H = \frac{1}{\sqrt{2}}(X + Z)$, and is particularly useful for creating superpositions, as it maps the state $|0\rangle$ to $\frac{1}{\sqrt{2}}(|0\rangle + |1\rangle) := |+\rangle$, and $|1\rangle$ to $\frac{1}{\sqrt{2}}(|0\rangle - |1\rangle) := |-\rangle$.

In addition to these simple one-qubit quantum gates, there are also gates designed for multiple qubits. For instance, consider the controlled-NOT gate (CNOT), which does exactly what its name suggests: if the control qubit is in state $|1\rangle$, it applies NOT to the test qubit, whereas if it is in state $|0\rangle$ instead, nothing happens. The CNOT gate is also quite prevalent in quantum circuits, playing a vital role in creating entanglement between qubits, which is a crucial resource in quantum computing. The circuit representation and equivalent matrix form of the CNOT gate are presented below, in Figure 2.4.



(a) CNOT gate – circuit representation. The black (filled) circle identifies the control qubit. The other qubit is the test qubit.

$$\begin{bmatrix} 1 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 1 \\ 0 & 0 & 1 & 0 \end{bmatrix}$$

(b) CNOT gate – equivalent matrix form.

Figure 2.4: Two-qubit gate example – the controlled-NOT gate.

More generally, within a quantum circuit, an intricate network of interconnected qubits and gates collaborates to execute a specific algorithm. For example, the circuit depicted in Figure 2.5, below, is designed for the implementation of the renowned Grover's search algorithm [30].

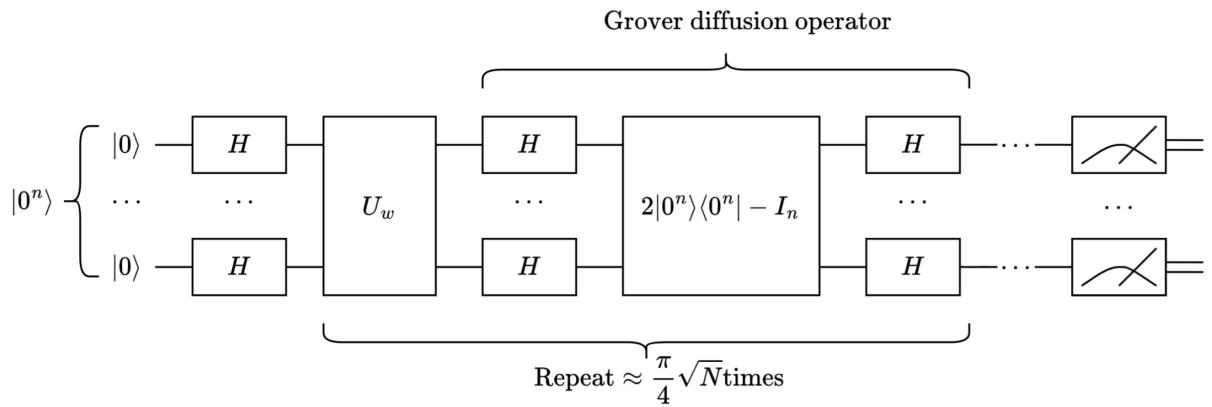


Figure 2.5: Quantum circuit utilized in the famous Grover's search algorithm [30]. Integrally reproduced from [31].

This is a purely quantum algorithm, different from what we intend to explore in this work. Nevertheless, it serves as a good example of how quantum circuits can be used to implement complex quantum algorithms.

An additional point of interest is the potential for classical simulation of quantum circuits. Quantum gates, being akin to mathematical operators, can always be expressed in matrix form. Consequently, analyzing the impact of a series of quantum gates, i.e., a quantum circuit, on a specific quantum state is

achievable by sequentially applying the relevant operators, as they appear in the circuit. At the end of this process, we arrive at a quantum state typically distinct from the initial state. This entire simulation can be executed classically. However, its scalability is severely limited, as the number of states, constituting our computational basis, representable with n qubits grows exponentially, at a rate of 2^n . As one might anticipate, this very quickly becomes unmanageable. Eventually, either memory constraints impede the storage of such extensive states, or the computations require so much time that the endeavor loses practicality. Effectively, simulating quantum computers becomes challenging beyond a few dozen qubits. Some tabletop calculations suggest that simulating 28 qubits would necessitate approximately 13 GB of memory, reaching the theoretical maximum capacity of a modern standard 16 GB personal computer. It's important to note that for each additional qubit, memory requirements double. For instance, simulating 50 qubits would demand about 54 Petabytes of memory, and for 51 qubits, approximately 108 PB, and so forth. These estimations are based on the assumption that a complex number is stored as two Python floats, each requiring 24 bytes of memory, thus a complex amplitude requires 48 bytes of memory to be stored. Then, one such amplitude is required for each of the 2^n basis states, for n qubits.

2.3 Hybrid Quantum-Classical Computing

Hybrid quantum-classical computing refers to a computational approach that combines elements of both classical and quantum computing paradigms to leverage the strengths of each. In this model, classical processors and quantum processors work in tandem to solve complex problems more efficiently than either could achieve alone. This collaborative strategy aims to harness quantum computing's unique capabilities while mitigating the challenges and limitations associated with quantum systems, such as error correction and decoherence. As of today, the synergy between classical and quantum elements holds promise for addressing complex real-world problems in areas like optimization, machine learning, and cryptography (cf. Figure 2.6).

2.3.1 Variational Quantum Algorithms

The hallmark of HQCC is what are called variational quantum algorithms (VQAs), which correspond to hybrid quantum-classical algorithms, typically realized through a parameterized quantum circuit (PQC). Additionally, as part of their implementation, the parameters are subjected to training in order to achieve the desired outcomes, by minimizing the value of a cost function. As such, VQAs can be thought of as the quantum analogue of highly successful machine-learning methods, such as neural networks. Moreover, since they outsource the circuit parameters' optimization to a classical optimizer, exterior to the quantum processor, VQAs leverage the full toolbox of classical optimization. As a whole, this approach has the added advantage of keeping the quantum circuit depth shallow, which, ultimately, helps in suppressing the overall noise level. This is crucial, since it allows for these algorithms to be implemented in the current NISQ (Noisy Intermediate Scale Quantum) era, not requiring complete fault-tolerance to produce reasonable results. A wide array of possible applications has been examined for

VQAs, essentially covering all the use cases envisioned by researchers for quantum computers (cf. Figure 2.6). Despite all this, it is important to note that VQAs are not without fault. There is still a lot of work to be done regarding their trainability, accuracy and efficiency. Either way, they are undoubtedly the most promising candidates for achieving useful quantum advantage in the near future, which is exactly why they have come under the spotlight, drawing the attention of numerous scientists over the past few years.

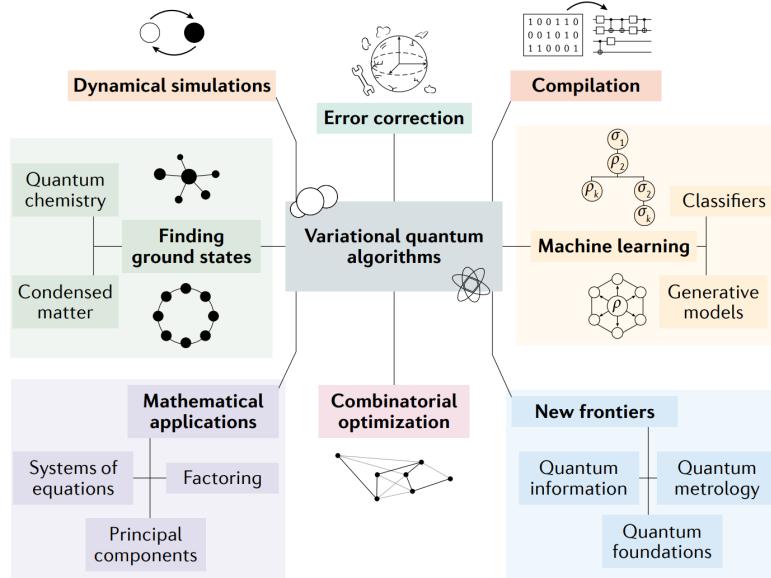


Figure 2.6: Applications of variational quantum algorithms. Sourced from [32] in its entirety.

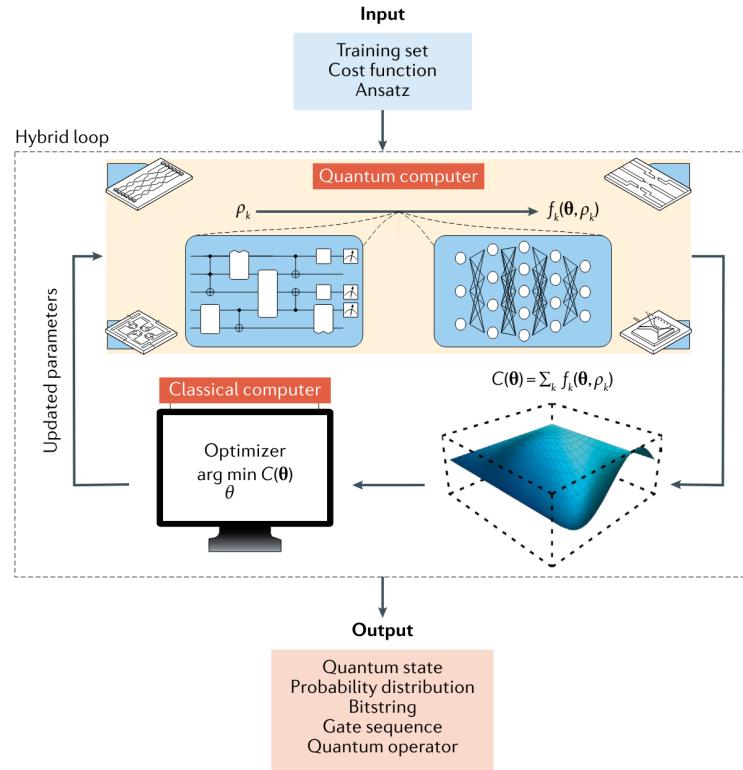


Figure 2.7: Schematic diagram of a variational quantum algorithm. Reproduced from [32].

2.3.1.1 Basic structure of a VQA

(This description is heavily inspired by [32] and mirrors their section layout.)

This section provides an overview of the essential components of VQAs. Kicking off the development of a VQA involves defining a cost (or loss) function, C , that encodes the solution to the problem. Following this, an ansatz is suggested, that is a quantum circuit that relies on a set of continuous or discrete parameters, θ , subject to optimization. In other words, we propose the functional form/structure of the parameterized quantum circuit. Afterwards, the suggested ansatz undergoes training, within a hybrid quantum-classical loop (cf. Figure 2.7), to address the following optimization task

$$\theta^* = \arg \min_{\theta} C(\theta), \quad (2.16)$$

corresponding to the minimization of the selected cost function. The distinctive feature of VQAs is their application of a quantum computer to estimate the cost function, $C(\theta)$, and its gradient, while relying on classical routines for the optimization of the parameters' values, θ . Next, we provide supplementary information for each step of the VQA framework.

Cost function

An essential aspect of VQAs involves encoding the problem into an adequate cost function. Similar to classical machine learning, this function associates values of the trainable parameters, θ , with real numbers. At a more abstract level, the cost function defines a hypersurface, often termed the cost landscape (cf. Figure 2.7), in a way such that the optimizer's task is to then navigate across this landscape and discover its global minimum, as this would correspond to the sought-after solution. In many instances, it is beneficial and feasible to express the cost in the form:

$$C(\theta) = \sum_k f_k (Tr [O_k U(\theta) \rho_k U^\dagger(\theta)]), \quad (2.17)$$

for some set of functions $\{f_k\}$, where $U(\theta)$ is a parameterized unitary, θ is composed of discrete and/or continuous parameters, ρ_k are input states from a training set and O_k are a set of observables. Note how the argument of f_k , $Tr [O_k U(\theta) \rho_k U^\dagger(\theta)]$, corresponds to the expectation value of observable O_k , in state $\rho'_k = U(\theta) \rho_k U^\dagger(\theta)$. Oftentimes, it is convenient to design the cost function to be given by the expectation value of a Hamiltonian (e.g., the cost Hamiltonian in QAOA), which is why we mention that it is beneficial for it to have this specific form. Then, the minimization of the cost corresponds, unambiguously, to the determination of the ground-state energy. Additionally, one frequently uses Pauli operators, whose expectation values are rather easy to compute in practice.

Furthermore, when designing a cost function, we want it to be as faithful as possible, in the sense that its minimum should closely correspond to the desired solution. Likewise, it is imperative that we can efficiently estimate $C(\theta)$, as this is integral to the functionality of this method. So far, we have been taking this for granted. However, one must guarantee that this is indeed verified experimentally, otherwise we are not able to proceed.

Ansätze

Another crucial element of a VQA is the ansatz. In broad terms, the ansatz's structure determines the nature of the parameters θ and, consequently, how they can be optimized to minimize the cost. The design of an ansatz often relies on the specific requirements of the task, allowing for the creation of problem-inspired ansätze. However, certain ansatz architectures are generic and problem-agnostic, making them applicable even in scenarios where relevant information is lacking. As one might expect, problem-agnostic ansätze, being more general, will also, usually, require a greater number of parameters, which unsurprisingly hinders the optimization task. However, this also grants them more flexibility. On the other hand, problem-oriented ansätze are capable of assimilating problem-specific information into the structure of quantum circuits. This custom-fit approach can result in a reduction of the parameter space, making optimization more efficient, and lead to solutions that are more meaningful and interpretable. Furthermore, adopting such an approach might (at times, not always!) enable the use of shallower quantum circuits, thereby aiding in the reduction of overall noise levels. (Greater circuit depths require more quantum gates, which results in more noise.) When it all comes down to it, one should weigh the trade-off between accuracy and generality, and carefully assess resource utilization in the choice of ansatz. (Yet, for achieving optimal results tailored to a specific problem, it's always recommended to utilize problem-inspired ansätze. The challenge lies in their design. While it's tempting to opt for a problem-agnostic ansatz and consider the task done, problem-inspired designs consistently yield superior performance.) For illustration purposes, here are a few examples of commonly considered ansatz types: hardware-efficient ansätze (aimed at reducing circuit depth), unitary coupled clustered ansätze (for quantum chemistry simulations), quantum alternating operator ansätze (used in QAOA), et cetera.

Gradients

After specifying the cost function and ansatz, the next step involves training the parameters, θ , and tackling the optimization problem described in Eq. 2.16. It is known that, for numerous optimization tasks, leveraging information from the gradient or higher-order derivatives of the cost function can enhance the speed and ensure the convergence of the optimizer [32]. A notable benefit of various VQAs is the ability to analytically compute the gradient of the cost function. This can be done using the parameter-shift rule. If the cost function has the form of Eq. 2.17, with $f_k(x) = x$, and θ_l is the l^{th} element in θ , which parameterizes a unitary $e^{i\theta_l \sigma_l}$ in the ansatz, with σ_l a Pauli operator, then the parameter-shift rule states that the equality

$$\frac{\partial C}{\partial \theta_l} = \sum_k \frac{1}{2 \sin \alpha} (Tr [O_k U(\boldsymbol{\theta}_+) \rho_k U^\dagger(\boldsymbol{\theta}_+)] - Tr [O_k U(\boldsymbol{\theta}_-) \rho_k U^\dagger(\boldsymbol{\theta}_-)]), \quad (2.18)$$

with $\boldsymbol{\theta}_\pm = \boldsymbol{\theta} \pm \alpha e_l$, holds for any real number α . In practice, one uses $\alpha = \pi/4$, since this maximizes the accuracy of the result [32]. Note that, although the expression is analytically exact, we can only ever obtain approximate results (hence the discussion about accuracy), since we are required to compute

expectation values of O_k , necessitating the quantum circuit to be sampled. This section was simply meant to shed some light on how one can experimentally compute these gradients, required for the optimization procedure.

Workflow (Hybrid Loop)

The VQA workflow is incredibly similar to a traditional machine learning pipeline. We have some input quantum state entering our ansatz, which is then processed by the quantum circuit, parameterized by θ . Afterwards, the output of this circuit is measured and the results are used to compute the cost function, $C(\theta)$. Following this, said cost is fed into a classical optimizer, which updates the parameters, θ , in order to minimize the cost⁷. This process [run quantum circuit, estimate cost, update parameters] is repeated iteratively until the optimizer converges to a minimum, at which point the optimal⁸ parameters, θ^* , are obtained. These can then be used to prepare the quantum state that solves the problem at hand, in the sense that the problem's solution can be extracted from it. This is a very high-level overview of the VQA workflow, but it should give a good idea of how the previously mentioned concepts come together to form a working algorithm. All the VQAs described in this work will use this hybrid loop (cf. Figure 2.7). Therefore, specifying the cost function and ansatz will be sufficient to characterize them.

Applications & Examples

In what follows, I will merely outline some of the most promising applications of VQAs for solving real world problems. These are (cf. Figure 2.6): (Specific algorithms for these tasks are indicated in parenthesis.)

1. **Finding ground states and excited states** (Variational Quantum Eigensolver [33], VQE, and variations thereof);
2. **Dynamical quantum simulations** - Based on iterative variational algorithms, they allow, e.g., for the simulation of open quantum systems [32];
3. **Optimization tasks** (Quantum Approximate Optimization Algorithm [2], QAOA, and Qubit-Efficient MaxCut Heuristic Algorithm [3], QEMC, both of which we shall analyse in further detail later);
4. **Mathematical applications** [32] - Linear systems, matrix-vector multiplication, non-linear equations, factoring, *et cetera*;
5. **Error correction** (Variational Quantum Error Corrector [34], QVECTOR);
6. **Machine learning and data science** - Quantum Machine Learning (QML) [35], classifiers, generative models, *et cetera*.

We are, now, finally ready to tackle the two specific variational quantum algorithms, QAOA and QEMC, which constitute the main focus of this work. Let us start with QAOA.

⁷It's a bit more complex, as we are required to compute the gradients for the classical optimization. For this, we simply use the parameter-shift rule. This requires some more shots, with different parameter values, but is entirely feasible.

⁸That is, assuming we don't get stuck in a suboptimal local minimum.

2.3.1.2 Quantum Approximate Optimization Algorithm (QAOA)

Renowned as the most prominent VQA in quantum-enhanced optimization, the QAOA [2] was originally devised to approximate solutions for combinatorial optimization problems, including constraint-satisfaction, SAT [36], and MaxCut problems [37]. In this section, we intend to explain in greater detail the inner workings of QAOA, since it is closely related to the new hybrid algorithm proposed by the HQCC project collaboration⁹. In a similar vein, the next section will delve into an equivalent level of detail, this time concentrating on the Qubit-Efficient MaxCut Heuristic Algorithm, QEMC.

Combinatorial optimization problems are formulated on binary strings, $s = (s_1, \dots, s_N)$, with the goal of minimizing a designated classical objective function, $L(s)$. In QAOA, however, the cost function is defined as the expectation value of a quantum Hamiltonian, H_C (or H_P), termed the cost (or problem) Hamiltonian, and doesn't explicitly depend on such bit-string s . This [cost] is constructed by mapping each classical variable, $s_j \in \{1, -1\}$, to a Pauli spin-1/2 operator, Z_j . Thus, the usual MaxCut objective $L(s) = \frac{1}{2} \sum_{\substack{i < j: \\ (i,j) \in E}} (1 - s_i s_j)$, denoting the cut of partition $s = (s_1, \dots, s_N)$, is transformed into the cost Hamiltonian

$$H_C = \frac{1}{2} \sum_{\substack{i < j: \\ (i,j) \in E}} (1 - Z_i Z_j), \quad (2.19)$$

from which one can extract the QAOA objective function.

Next up is the QAOA ansatz. Drawing inspiration from the quantum adiabatic algorithm, QAOA substitutes adiabatic evolution with n cycles of alternating time evolution between the cost Hamiltonian, H_C , and a suitably chosen mixer (or bias) Hamiltonian, H_M (or H_B). The role of this mixer Hamiltonian can be understood as introducing quantum fluctuations, or transitions, between different states, helping the algorithm explore the solution space more effectively, ideally preventing it from getting trapped in sub-optimal local minima. The entirety of this process forms the previously mentioned quantum alternating operator ansatz.

In practice, defining $\theta = \{\gamma, \alpha\}$, the cost function is $C(\gamma, \alpha) = \langle \psi_n(\gamma, \alpha) | H_C | \psi_n(\gamma, \alpha) \rangle$, with

$$|\psi_n(\gamma, \alpha)\rangle = e^{-i\alpha_n H_M} e^{-i\gamma_n H_C} \dots e^{-i\alpha_1 H_M} e^{-i\gamma_1 H_C} |\psi_0\rangle, \quad (2.20)$$

where $|\psi_0\rangle$ is the initial state entering the ansatz. This ansatz is illustrated in Figure 2.8, below, for $n \in \mathbb{N}$ QAOA layers.

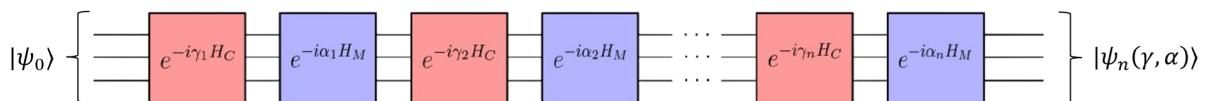


Figure 2.8: Quantum alternating operator ansatz. Adapted from [38].

Replacing adiabatic time evolution with a *Trotterized* evolution works due to the famous Trotter-Suzuki

⁹More information about this collaboration is present in the Acknowledgments section.

formulas [19], which state that (for general non-commuting H_j):

$$e^{-i\sum_{j=1}^m H_j t} = \prod_{j=1}^m e^{-iH_j t} + O(m^2 t^2). \quad (2.21)$$

If $t \ll 1$, then the error in this approximation becomes negligible. On the other hand, if t is large, Trotter-Suzuki formulas can still be used to simulate the dynamics accurately by breaking things up into a sequence of short time-steps. Let r be the number of steps taken in the time evolution, so each time-step runs for time t/r . Then, we have that

$$e^{-i\sum_{j=1}^m H_j t} = \left(\prod_{j=1}^m e^{-iH_j t/r} \right)^r + O(m^2 t^2/r), \quad (2.22)$$

which implies that if r scales as $m^2 t^2/\epsilon$, then the error can be made at most ϵ for any $\epsilon > 0$. In the QAOA case, we simply have $H = H_C + H_M$ and we treat each of the evolution time intervals as variational parameters that are optimized classically (can be seen as $t/r \equiv \gamma_l, \alpha_l$). This means we do not use this *Trotterization* approach directly, but it heavily inspires the QAOA ansatz nevertheless.

Let's now further detail the practical realization of the algorithm. Reiterating, the aim of MaxCut is to maximize the number of edges in a graph that are "cut" by a given partition of the vertices into two sets (cf. Figure 2.1). One way of implementing QAOA for MaxCut is to consider: H_C as in Eq. 2.19, where the sum is over all edges of the studied graph, connecting vertices (j, k) ; and $H_M = \sum_{j=1}^N \mathbf{X}_j$, for N qubits and graph nodes. In this scheme, the associated cost function, $\langle \psi_n(\gamma, \alpha) | H_C | \psi_n(\gamma, \alpha) \rangle$, merely rewards cases where we have edges between nodes of different sets. Furthermore, we can, now, also write¹⁰

$$U_{H_{C_l}} = e^{-i\gamma_l H_C} = \prod_{\text{Edge } (j,k)} e^{-i\gamma_l (1 - Z_j Z_k)/2} \quad (2.23)$$

$$U_{H_{M_l}} = e^{-i\alpha_l H_M} = \prod_{j=1}^n e^{-i\alpha_l \mathbf{X}_j} \quad (2.24)$$

These give us the form of the unitaries that are applied in each layer of the QAOA ansatz, representing the *Trotter*-inspired time evolution of Figure 2.8 (red and blue boxes). They correspond to the complex exponentials of the cost and mixer Hamiltonians, respectively: $e^{-i\gamma_l H_C}$ and $e^{-i\alpha_l H_M}$. Notice how $U_{H_{M_l}}$ uses Pauli- x operators, instead of the usual Pauli- z in $U_{H_{C_l}}$. In practice, the mixer Hamiltonian terms $e^{-i\alpha_l H_M}$ correspond to $R_x(2\alpha_l)$ and are easy to implement. The cost Hamiltonian terms, on the other hand, are a bit more tricky, requiring the use of 2 CNOT gates. Each of the terms $e^{-i\gamma_l (1 - Z_j Z_k)/2}$ can be transpiled into a quantum circuit, as shown in Figure 2.9, below. For each edge in the graph, we'll have one of these terms, in each of the n layers of the QAOA ansatz.

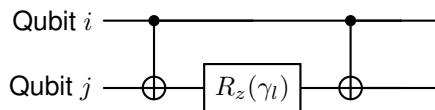


Figure 2.9: $e^{i\gamma_l Z_i Z_j / 2}$ decomposition.

This allows us to re-write Eq. 2.20 as

$$|\psi_n(\gamma, \alpha)\rangle = U_{H_{M_n}} U_{H_{C_n}} \dots U_{H_{M_1}} U_{H_{C_1}} |\psi_0\rangle \quad (2.25)$$

In QAOA, it is customary to start with a uniform superposition over the N bit-string basis states, where N represents the number of graph nodes. This is denoted as $|\psi_0\rangle = |+\rangle^{\otimes N} = \frac{1}{\sqrt{2^N}} \sum_{z \in \{0,1\}^N} |z\rangle$, and is achieved by applying a Hadamard gate to each of the qubits, at the start of the quantum circuit.

Subsequently, a classical routine is employed to optimize the values of the parameters by minimizing the negative of the cost (analogous to the "cut"), therefore enabling the extraction of the MaxCut partition. Experimentally, for each step of the classical optimization, it is necessary to compute expectation values of Pauli- z operators, which requires the quantum circuit to be sampled a certain number of times (shots). Additionally, in case it was not yet clear, note that in QAOA we use one qubit for each node of the graph, with the qubits' values indicating which set they correspond to: $|0\rangle$: "Set 0"; $|1\rangle$: "Set 1". As such, for the graph in Figure 2.1, we'd require 5 qubits. Exemplifying, if the most sampled basis state at the end of the quantum circuit is $|00011\rangle$, then nodes 1, 2 and 3 belong to "Set 0", while nodes 4 and 5 belong to "Set 1".

2.3.1.3 Variational Qubit-Efficient MaxCut Heuristic Algorithm (QEMC)

The Qubit-Efficient MaxCut Heuristic Algorithm (QEMC) [3] is somewhat similar to QAOA. After all, it was heavily inspired by it. However, it has a number of crucial differences. First, it only requires $n = \log_2(N)$ qubits, instead of N , where N is the number of nodes in the graph. Additionally, the QEMC algorithm is based on a novel probability threshold encoding scheme, a suitable cost function, and a parameterized unconstrained quantum circuit. Going in order:

Probability threshold encoding scheme

With n qubits, each of the $N = 2^n$ basis states will represent one of the graph's nodes. Following the sampling of the quantum circuit, a probability distribution is generated. Nodes with probabilities exceeding a certain threshold, $p_{th} = \frac{1}{2B}$, belong to "Set 1", while probabilities below this value indicate inclusion in "Set 0". This strongly diverges from how we encode set inclusion in QAOA.

Suitable cost function

The objective function utilized in QEMC is the following:

$$L(\{p(i)\}) = \sum_{\substack{j < k: \\ (j,k) \in E}} \left[\left(d(j, k) - \frac{1}{B} \right)^2 + \left(s(j, k) - \frac{1}{B} \right)^2 \right], \quad (2.26)$$

where $d(j, k) = |p(j) - p(k)|$ and $s(j, k) = p(j) + p(k)$ are the absolute difference and sum of the corresponding basis states' probabilities. The idea is that as both $d(j, k)$ and $s(j, k)$ tend towards $1/B$, the probability of one node approaches zero (distinctive "Set 0"), while the probability of the other node

approaches $1/B$ (distinctive "Set 1"), without specifying which is which. Ultimately, just like for QAOA, connections between nodes of different sets are favoured. Note, however, that this probability threshold encoding scheme assumes *a priori* that one of the sets ("Set 1") has B nodes. Nevertheless, this is not an issue, as we can efficiently iterate through all potential values of $B = 1, \dots, \lfloor \frac{N}{2} \rfloor$. Frequently, it is reasonable to set $B = N/2$, and we shall use this as our default starting point.

Problem-agnostic ansatz

The QEMC circuit ansatz is agnostic to specific graph instances, a departure from QAOA where the graph structure is explicitly encoded in the quantum circuit. Instead, the graph is implicitly encoded through the cost function. As a result, the QEMC quantum circuit is not bound to any particular form and only needs to be expressive enough to approximate the optimal states in the Hilbert space. Such problem-independent ansatz approach provides considerable flexibility in ansatz selection. Frequently, the circuit ansatz known as "Strongly Entangling Layers" is employed, as depicted below (Figure 2.10). As can be seen, this ansatz applies a series of parameterized single-qubit rotations interspersed with entangling gates to generate a highly entangled quantum state. We use PennyLane's `qml.StronglyEntanglingLayers` implementation for this purpose.

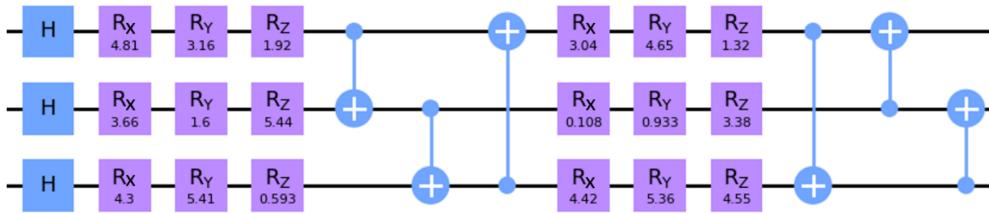


Figure 2.10: "Strongly Entangling Layers" circuit ansatz: case of $n = 3$ qubits and $p = 2$ layers. After a single layer of Hadamard gates, each subsequent layer consists of $3n$ single-qubit parameterized rotation gates and n CNOT gates (entangling gates). Reproduced from [3].

These constitute the main ingredients necessary to understand the QEMC algorithm. At this point, the same hybrid loop as before would be run, so as to optimize the "Strongly Entangling Layers" ansatz's parameters, to minimize the cost function. To read the MaxCut partition, generated by the QEMC algorithm, one would sample the circuit's output one more time, after the training, and build the basis states' probability distribution. The threshold p_{th} would then be applied to determine each nodes' set inclusion, from each of their associated basis states' probabilities.

One notable, and perhaps unfortunate, property of QEMC is that it is efficiently simulable classically. Due to the exponential compression of the number of qubits, the algorithm can be feasibly simulated on a classical computer, even for large graphs, hence defeating its purpose as a quantum algorithm. This is something the authors of the algorithm [3] realized in hindsight. For this reason, it is now termed a quantum-inspired classical algorithm.

¹⁰In reality, QAOA in this form doesn't require the use of *Trotterization*, since all the operators in the sums constituting H_C and H_M commute. Nevertheless, it is *Trotter*-inspired.

Chapter 3

iQAQE Framework

In this chapter, we present the iQAQE Framework, which interpolates between QAOA and QEMC. We begin with a comparison of the two algorithms, followed by a detailed description of iQAQE, its theoretical model, motivation, and workflow.

3.1 Comparison of QAOA and QEMC

To motivate the development of our new algorithm(s), we summarize the merits and drawbacks of both QAOA and QEMC. QEMC has several advantages over QAOA, including significantly lower qubit requirements ($n = \log_2(N)$ vs. N) and the ability to support shallower circuit depths [3]. The threshold probability encoding scheme allows QEMC to associate each graph partition with a volume of quantum states, potentially enhancing noise resilience. Nevertheless, while QEMC might be better trainable [3], it requires the entire probability distribution at each optimization step to compute the cost, increasing the number of needed shots compared to QAOA, which only requires expectation values. Additionally, it has been shown that QEMC can be efficiently simulated classically, somewhat undermining its purpose as a quantum algorithm [3].

Having thoroughly understood these two algorithms, the collaboration within the HQCC project has introduced a novel framework for designing multiple distinct VQAs. This will be the focus of our next discussion.

3.2 Interpolated QAOA/QEMC Hybrid Algorithm(s) (iQAQE)

3.2.1 Theoretical Model

Algorithms developed within the proposed framework aim to interpolate between QAOA and QEMC, combining their strengths into a more robust approach, tentatively named iQAQE (Interpolated QAOA/QEMC). In iQAQE, we depart from the QEMC approach by associating each graph node with a list (sometimes also called a sub-list) of basis states, in contrast to QEMC's consideration of a single

basis state for each node (cf. Figure 3.1). Each of these lists comprises somewhere between $[1, 2^{n-1}]$ basis states, where n represents the number of qubits. The number of basis states per list is called the list's cardinality, denoted as c . This design allows for potential overlap among states from different lists/nodes. What we formerly referred to as "mapping" involves distributing basis states among these lists. The term "mapping" can also signify a specific allocation of basis states among the lists. Additionally, it is important to note that the encoding of these states will utilize a qubit range expected to fall between the **QEMC** and **QAOA** requirements, specifically in $[\log_2(N), N]$, for an N -node graph. With all that said, the main goal of my thesis is to ascertain the optimal mapping from basis states to lists, essentially determining which basis states should be included in specific lists/nodes. Naturally, this undertaking also involves the identification of a suitable cost function and ansatz to ensure the algorithm's effective operation. For most of our work, we consider **iQAQE** to use the same ansatz and cost function as **QEMC**, adjusted for the appropriate number of qubits and with some modifications that we'll discuss below.

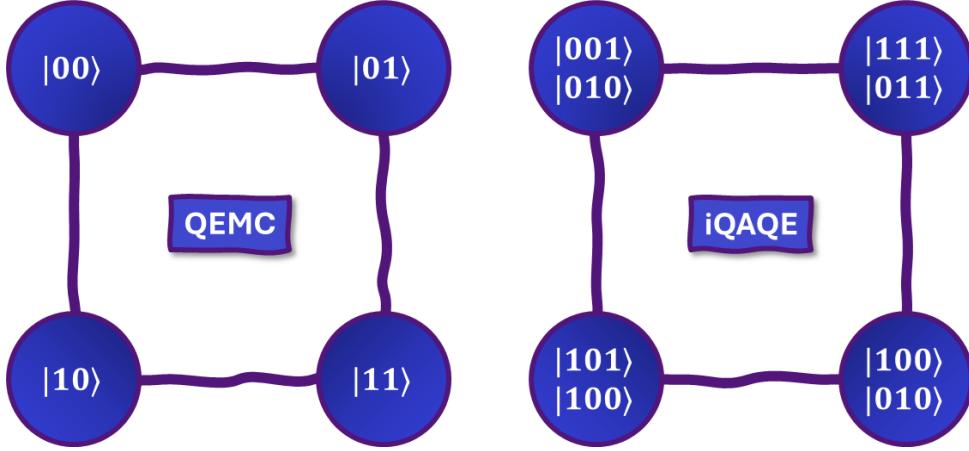


Figure 3.1: Example schematic representation of the **iQAQE** encoding scheme. Comparison between the **QEMC** and **iQAQE** encoding schemes: example for a trivial 4-node graph. In **QEMC**, each node is associated with a single basis state, while in **iQAQE**, each node is associated with a list of basis states.

3.2.2 Motivation & Outlook

At this point, it is important to mention that the development of this algorithm is entirely exploratory, although we do believe it should allow for a more practical algorithm implementation-wise. We can postulate that such a hybrid approach might have some potential advantages. Namely, it should allow for less shots than in **QEMC** and fewer qubit requirements than in **QAOA**, in addition to, arguably, being better trainable. Another interesting characteristic of such an algorithm is its tunable "quantum-ness". As we interpolate between **QAOA** (hybrid quantum-classical) and **QEMC** (classical, quantum-inspired), the ability to selectively adjust the degree of both could prove to be advantageous. As part of a more long-term perspective (not in this thesis), I may explore the possibility of generalizing this algorithm to tackle additional combinatorial problems, such as Max- k -Cut. Throughout my thesis, I will rigorously test the algorithm using classical simulations of quantum machines. The aim is to identify the most effective implementation strategies, taking into account both cost and ansatz considerations, as well as

the distribution of basis states.

3.2.3 Algorithm Description & Workflow

A sketch of the workflow is provided below (Table 3.1). Regarding the hybrid loop, I only detail the changes relative to **QEMC**. Following this, I will elaborate on each step.

Table 3.1: Core steps for algorithms built on the iQAQE Framework.

Step	Description
1. Select number of qubits	$n_{\text{qubits}} = n \in [\log_2(N), N]$, where N refers to the graph's number of nodes.
2. Select list cardinality	$\text{list_cardinality} = c \in [1, 2^{n-1}]$, where n represents the chosen number of qubits. Note that c is taken to be the same for all the lists, although this could change in future work.
3. Assignment/mapping	Map c basis states to each graph node.
4. Calculate nodes' probabilities (within the hybrid loop)	To obtain each node's probability, sum the probabilities of the basis states in the associated list and normalize the results, so they add up to 1.
5. Cost function and ansatz (within the hybrid loop)	Keep the same cost function and ansatz as in the QEMC algorithm, just adapt the ansatz to the correct number of qubits.

1. Select the number of qubits

In selecting the number of qubits, we allow for any intermediate value between the **QEMC** ($\log_2(N)$) and **QAOA** (N) limits. This should give us more flexibility regarding the degree of "quantum-ness" that we want our algorithm to exhibit. This is to be understood in the context of **QEMC** being entirely classical, whereas **QAOA** is hybrid quantum-classical. As such, by choosing a number of qubits in between these two limits, we expect our algorithm to be hybrid in terms of its quantum nature. In addition to this, the number of qubits, alongside the ansatz, will also influence the expressivity of the model.

Another key consideration, closely tied to our next workflow step, is the chosen encoding method. This refers to how we determine each node's color and map basis states to respective lists. Different encoding strategies, such as using expectation values of Pauli strings (e.g., as seen in [18]), inherently influence the number of required qubits. For instance, the Pauli string encoding in [18] automatically enforces a polynomial compression on the number of qubits. Currently, we've employed **QEMC**'s probability threshold encoding scheme, which provides flexibility in the number of qubits and deriving node probabilities from basis states. Further discussion on this will be provided in point 4 below.

2. Select list cardinality

If it hasn't been clear already, we assign a list of basis states to each graph node. The cardinality of each list, denoted as c , falls within the interval $[1, 2^{n-1}]$, where n denotes the number of qubits used. This choice represents a balance between the **QEMC** (1) and **QAOA**¹ (2^{n-1}) extremes. Determining the

¹This 2^{n-1} value is determined by what we refer to as the **QAOA** mapping for iQAQE, which we will discuss later.

optimal value for c is non-trivial, as it depends on various interconnected factors including the number of qubits, ansatz, cost function, and color encoding. To explore its impact on algorithm performance, we conduct grid searches on this parameter.

3. Assignments/mapping

The process of assigning/mapping basis states to lists adds significant complexity. *A priori*, there's no clear-cut method for this mapping. The simplest approach involves randomly assigning c basis states to each node, ensuring all states are utilized at least once to avoid discarding information. However, exploring more intricate mappings may yield potential benefits. For instance, in the QAOA mapping, a single qubit among N is fixed to 1 for each node, allowing for 2^{N-1} permutations of the remaining qubits. In other words, each list, comprising 2^{N-1} elements, corresponds to a distinct graph node, with each node having a different qubit fixed to 1. As can be noted, there's a myriad of distinct mappings one can build. In Chapter 5, we will thoroughly examine several possible mappings, each presenting its unique advantages and disadvantages. The challenge lies in identifying the most effective mapping for a given problem instance. We aim to address this in our work.

4. Calculate nodes' probabilities

Above (Table 3.1), we mentioned a straightforward method of computing node probabilities by summing the probabilities of associated basis states and normalizing the result. This approach serves as our foundational strategy throughout this work. However, there may be hidden advantages in devising more intricate methods. For instance, one could explore calculating node probabilities using the arithmetic or geometric averages of basis states' probabilities, followed by normalization. Yet, designing such schemes requires caution, as mismatches between the ansatz and encoding could result in constant probabilities for nodes, impeding variational circuit optimization². Furthermore, the normalization process itself may eliminate the probabilities' dependence on variational parameters under certain circumstances. While we explore alternative methods for calculating probabilities, our primary emphasis in this work remains on the basic approach described earlier, given its apparent robustness.

5. Cost function and ansatz

We employ the QEMC cost function, utilizing the probability threshold encoding scheme mentioned earlier. It's important to note that this scheme assumes, by default, that one of the sets ("Set 1") comprises B nodes³, which may not align with the MaxCut partition. In some cases, we may need to iterate through potential values of B , ranging from 1 to $\lfloor \frac{N}{2} \rfloor$ – although, in practice, careful selection of B should prevent this scenario. Often, setting $B = N/2$ proves to be a reasonable starting point, and we adopt this as our default.

²This occurred once during testing when we used the QAOA mapping and ansatz alongside the QEMC cost function and probability threshold encoding scheme.

³This B value is user-defined.

Given our approach to mapping basis states to node lists, it's logical for us to employ the **QEMC** cost function rather than **QAOA**'s. The latter relies on a problem Hamiltonian defined on N qubits, which may not match our choice of n . Typically, $n < N$ as we aim to reduce the number of qubits.

Similarly, we opt for the Strongly Entangling Layers ansatz from the **QEMC** scheme for its versatility regarding qubit count. This ansatz, agnostic to specific problems, offers flexibility in our algorithm's implementation compared to **QAOA**'s problem-inspired ansatz. However, there are plans to explore more tailored ansatz designs correlated with individual mappings, aiming to enhance algorithm performance by leveraging graph-specific information. Note that this may necessitate an adjusted objective function developed with these considerations in mind. While we experiment with problem-inspired ansätze in this work, our primary focus remains on the Strongly Entangling Layers ansatz.

Chapter 4

Implementation

In this chapter, we outline the numerical implementation of the models discussed in Chapters 2 and 3, focusing on pseudo-code representations. Furthermore, we highlight the Python libraries employed and provide specific code details, accompanied by a brief discussion on the benchmarks developed to evaluate the algorithms' performances.

4.1 Individual Algorithms

For numerical implementation of the previously described algorithms, we employ PennyLane [39], a versatile Python library designed for differentiable programming of quantum computers. PennyLane facilitates the execution of variational quantum circuits and their simultaneous training, akin to training a classical neural network, within the same Python environment, which is highly convenient. It offers numerous features, including automatic differentiation, crucial for optimizing variational quantum circuits.

In our implementation with PennyLane, we utilize one of two devices from its extensive selection. Firstly, we employ the `default.qubit` device, a straightforward state-vector qubit simulator implemented in Python, compatible with Autograd, JAX, TensorFlow, and Torch backends. This device is well-suited for optimizations involving a small to moderate number of qubits and parameters, offering exact expectation values. Secondly, we utilize the `lightning.qubit` device, a fast state-vector qubit simulator with a C++ backend, recommended for scenarios involving moderate numbers of qubits and parameters or when utilizing stochastic expectation values. Our simulations utilize both devices: the former excels in analytical simulations of small quantum circuits, which predominates our work, while the latter is preferable for simulations of larger quantum circuits with a finite number of shots, especially when many shots are required. Although analytical simulations are favored due to their lower computational resource requirements, they lack the output sampling aspect of a true quantum computer simulation. Nonetheless, they suffice for our goal of optimizing variational quantum circuit parameters.

Moreover, we heavily rely on the NetworkX [40] Python library, a valuable tool for creating, manipulating, and analyzing complex networks. NetworkX aids in generating the graphs utilized in the MaxCut problem, offering useful visualization tools for graph representation and partitioning visualization, essen-

tial for interpreting algorithm results.

Additionally, we utilize the CVXPY [41] Python library, specifically designed for convex optimization tasks. CVXPY facilitates solving the semidefinite programming (SDP) relaxation of the MaxCut problem within the GW algorithm framework, which inherently constitutes a convex optimization problem.

Regrettably, we didn't have the chance to test these on actual quantum hardware, which would have been invaluable for evaluating the performance of these algorithms in a real-world setting. Furthermore, the lack of access to a High-Performance Computing (HPC) cluster significantly limited our capacity to perform large-scale simulations and grid-searches necessary for hyperparameter tuning. These limitations occasionally posed a bottleneck in our research, particularly concerning our inability to evaluate the algorithms on larger graphs, which would have been crucial for comprehensive testing. All simulations presented herein were conducted using my personal computer – a system equipped with an 8-core AMD Ryzen 7 5800H CPU, integrated Radeon graphics¹, and 16 GB of RAM.

Lastly, all simulations were performed using the Adam classical optimizer [42], with default parameters except for the learning rate, which was treated as a hyperparameter and fine-tuned accordingly. The Adam optimizer is a popular optimization algorithm for variational quantum algorithms, utilizing stochastic gradient descent with adaptive learning rates and momentum.

4.1.1 Quantum Approximate Optimization Algorithm (QAOA)

Below, we present the pseudo-code representation of the Quantum Approximate Optimization Algorithm (QAOA) implementation, as described in subsubsection 2.3.1.2.

Algorithm 1 Quantum Approximate Optimization Algorithm (QAOA)

Require: graph, n_layers, parameters, Optional MaxCut value

Ensure: n_qubits = graph.n_nodes

1. Define U_C and U_M
2. Create variational quantum circuit using `@qml.qnode` with `n_layers` layers of U_C and U_M

Note: Start with `qml.Hadamard` on all qubits

3. Define objective function to minimize: $-\langle H_C \rangle$, obtained through `-1 * qml.expval(H_C)`
4. Initialize Adam optimizer with default parameters
5. Start timer & begin training:

while True **do**

- 5.1 Update parameters using Adam
- 5.2 Store cut value, cost function, and approximation ratio (if MaxCut is provided), for each iteration

if max_iter **or** abs_tol **or** rel_tol reached **then**

break

end if

end while

6. Stop timer: record training time

7. Sample circuit to get most frequent bitstring

8. Compute partition & corresponding cut value
-

¹No GPU-assisted computations were conducted.

All of the developed code has been compiled in a GitHub repository (Private), available at https://github.com/kaiuki2000/HQCC_Code_Implementations/tree/main. For access to the code, please contact the author.

4.1.2 Qubit-Efficient MaxCut Heuristic Algorithm (QEMC)

Here, we provide the pseudo-code representation of the implementation of the Qubit-Efficient Max-Cut Heuristic Algorithm (QEMC), as detailed in subsubsection 2.3.1.3.

Algorithm 2 Qubit-Efficient MaxCut Heuristic Algorithm (QEMC)

```

Require: graph, n_layers, parameters, B, Optional MaxCut value           ▷ B is a parameter of the cost function
Ensure: n_qubits = math.ceil(np.log2(graph.n_nodes))

1. Define QEMC ansatz layer (Strongly Entangling Layers ansatz)
2. Create variational quantum circuit using @qml.qnode with n_layers layers of QEMC ansatz

Note: Start with qml.Hadamard on all qubits
3. Define objective function using probability threshold encoding scheme
4. Initialize Adam optimizer with default parameters
5. Start timer & begin training:

while True do
    5.1 Update parameters using Adam
    5.2 Store cut value, cost function, and approximation ratio (if MaxCut is provided), for each iteration
    if max_iter or abs_tol or rel_tol reached then
        break
    end if
end while
6. Stop timer: record training time
7. Sample circuit to get output probability distribution
8. Compute partition using probability threshold encoding scheme & corresponding cut value

```

4.1.3 Interpolated QAOA/QEMC Hybrid Algorithm (iQAQE)

This subsection presents the pseudo-code for a general iQAQE algorithm (Algorithm 3), as discussed in Chapter 3 (iQAQE Framework). It is important to note that point 3 of the algorithm involves significant complexity. The method for mapping the states is flexible and can be chosen in various ways. We will explore several possible mappings in Chapter 5.

Additionally, note the similarities between iQAQE and QEMC in terms of pseudo-code, as iQAQE is fundamentally built on the same basis.

Algorithm 3 Interpolated QAOA/QEMC Algorithm (iQAQE)

Require: graph, n_layers, parameters, B, Optional MaxCut value ▷ B is a parameter of the cost function

1. Select number of qubits: n_qubits = $n \in [\log_2(N), N]$, with $N = \text{graph.n_nodes}$
2. Select list cardinality: list_cardinality = $c \in [1, 2^{n-1}]$
3. Assignment/mapping: Map c basis states to each graph node
4. Define iQAQE ansatz layer (adapted Strongly Entangling Layers ansatz)
5. Create variational quantum circuit using @qml.qnode with n_layers layers of iQAQE ansatz

Note: Start with qml.Hadamard on all qubits

6. Define objective function using probability threshold encoding scheme
7. Initialize Adam optimizer with default parameters
8. Start timer & begin training:

while True **do**

- 8.1 Update parameters using Adam

Note: Calculate nodes' probabilities by summing the probabilities of the basis states in each list and normalizing the result

- 8.2 Store cut value, cost function, and approximation ratio (if MaxCut is provided), for each iteration

if max_iter or abs_tol or rel_tol reached **then**

- break**
- end if**

end while

9. Stop timer: record training time
10. Sample circuit to get output probability distribution
11. Compute partition using probability threshold encoding scheme & corresponding cut value

4.1.4 Goemans-Williamson Algorithm

Here, we present the numerical implementation of the Goemans-Williamson algorithm, in the form of pseudo-code.

Algorithm 4 Goemans-Williamson Algorithm for MaxCut

Input: graph, Optional MaxCut value

Output: Partition, Cut_Value, Approximation_Ratio (if MaxCut is given), Running_Time

1. Initialize: n (number of nodes), edges (graph edges)
2. Define SDP variable X and constraints $X \gg 0$ (X positive semidefinite) and $X[i, i] == 1$ for $i \in \text{range}(n)$
3. Set up the objective function:

```
objective = sum(0.5 * (1 - X[i, j]) for (i, j) in edges)
```
4. Start the timer and solve the SDP problem with the defined objective function and constraints, using CVXPY
5. Record the running time
6. Retrieve X and extract the vectors v from the columns of sqrtm(X) (square root of matrix X).
7. Perform random-hyperplane "selection" utilizing the vectors v to derive the partition.
8. Compute the cut value and, if applicable, the approximation ratio
9. Return the partition, cut value, approximation ratio (if applicable), and running time

4.2 Benchmarking and Testing Methods

Now that we have introduced all the algorithms and their numerical implementations, we must discuss the benchmarking and testing methods used to evaluate their performance. The primary metrics for evaluation are the cost plot (cost vs. iterations) and the approximation ratio plot² (approximation ratio vs. iterations), both produced during training. These metrics help compare the algorithms' convergence rates and final cut values.

Given that we initialize our algorithms with random parameters, we devised statistically meaningful metrics to compare their performances. A single run isn't sufficient to draw conclusions about an algorithm's performance, as initial parameters can significantly impact the results. Thus, the primary statistical metric used is the average cut value (or approximation ratio). This metric averages the cut values (or approximation ratios) derived at each iteration during optimization. It enables us to evaluate the algorithm's average performance for a single run after training. Essentially, it provides the likelihood of obtaining the correct result after training by executing the circuit once.

Another metric commonly used is the average best-so-far (BSF) cut value. This metric computes the average of the best-so-far cut values obtained at each iteration during optimization. For a given training curve, with cut values vs. iterations, the BSF cut value is obtained by storing the highest cut value achieved so far. If a dip occurs in the graph, it is adjusted to match the previous graph value, ensuring that the final result is monotonically increasing. The code for this transformation is shown below:

```
1 def BSF_Transform(vec):
2     """
3         Transforms a vector into a best-so-far vector.
4     """
5     bsf_vec = vec.copy()
6     for i in range(1, len(bsf_vec)):
7         if bsf_vec[i] < bsf_vec[i-1]:
8             bsf_vec[i] = bsf_vec[i-1]
9     return bsf_vec
```

Listing 4.1: Auxiliary function: best-so-far transformation

This transformation ensures that the BSF cut value curve is monotone increasing, as it always takes the best value so far during training.

Additionally, we consider the median best-so-far cut value (cf. section 5.6), which helps eliminate outliers often present in our simulations. By using the median, we can mitigate the impact of particularly lucky or unlucky initializations and better reflect the true performance of the algorithm. Both the average BSF and the median BSF curves are usually computed using either 3, 5 or 10 training curves³, to ensure statistical robustness. Each curve is generated with different random initial parameters, obtained through `2 * numpy.pi * numpy.random.rand`, producing values in the range $[0, 2\pi]$.

Initially, we used the best-so-far of the average, but we found it less representative of the algorithms' best performances. We now prefer to first apply the BSF transform to obtain the best possible per-

²We employ the AR when the MaxCut is known. Otherwise, we simply compute the cut plot.

³Unless explicitly stated otherwise, we always employ 10 runs for computing averages/medians.

formance from each curve and then compute the average or median. This approach ensures that we account for the best performance of each curve, rather than just the best-so-far average performance. However, we initially used the best-so-far of the average for most of our work and only later switched to the new metrics. While this doesn't alter the interpretation of the results⁴, it does affect the metric values, providing a more accurate quantification of the algorithms' performances.

Additionally, we conduct grid searches on the models' hyperparameters (such as Adam's learning rate, `n_layers`, `list_cardinality`, etc.) to optimize their performance and understand how they behave under different conditions. Nevertheless, due to the unavailable access to an HPC cluster, these grid searches are often limited in scope.

To compare the algorithms' performances, we also benchmark them against the Goemans-Williamson algorithm. This comparison helps us evaluate how well our algorithms perform relative to state-of-the-art solutions. Often, our plots will include comparisons with QAOA⁵, QEMC, and the GW algorithm for a comprehensive analysis of the algorithms' performances.

⁴In general, the interpretation remains consistent, but comparing both metrics offers additional insights, hence our decision to retain the best-so-far average rather than simply using the average best-so-far.

⁵If not specified otherwise, we default to using 3 QAOA layers.

Chapter 5

iQAQE Schemes and Results

In this chapter, we present the various iQAQE schemes we have developed and the results of their numerical simulations. These schemes explore different combinations of qubit numbers, list cardinalities, basis state mappings, and circuit layers. Each scheme features a unique mapping, resulting in distinct outcomes and properties. We compare these results and discuss the potential of each scheme. This chapter is essentially a collection of various ideas, all built upon the iQAQE Framework described in Chapter 3. The majority of these schemes are essentially heuristic methods designed to aid in selecting the number of qubits and mapping.

At times, our work diverged from iQAQE, leading us to explore other ideas for MaxCut algorithms. Although these ideas originated from our work on iQAQE, they are distinct enough to warrant their own chapter. These will be presented in the next chapter (Chapter 6).

From here onwards, unless explicitly stated otherwise, all VQAs are being tested on the following 8-node graph, whose optimal cut is 10 (Figure 5.1). This graph is one of the benchmarks used in [3], which is why we employ it here. Additionally, its small size allows for reasonably quick simulations on my personal computer, which is a significant advantage.

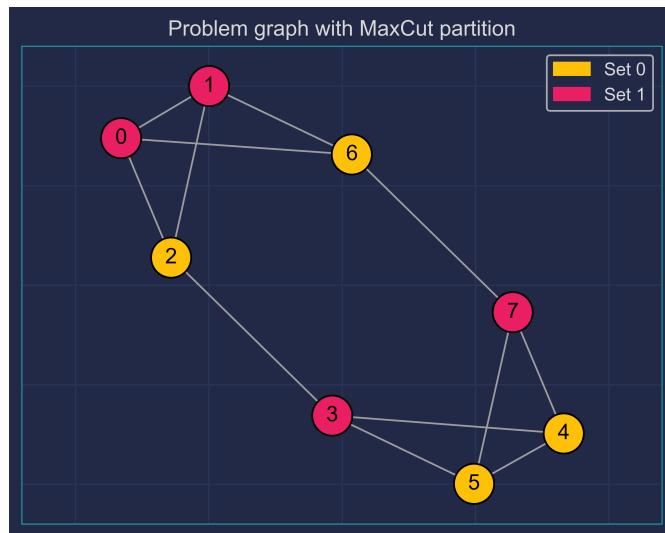


Figure 5.1: Considered 8-node graph instance: the MaxCut partition is illustrated. The optimal cut (10) was found through brute-force (exhaustive search).

5.1 Random iQAQE

First and foremost, we aim to understand how iQAQE compares with both QAOA and QEMC. In terms of iQAQE's implementation for this comparison, the number of qubits (n) and the list cardinalities (c) were selected without careful consideration. They were randomly chosen using `np.random.randint`, resulting in $n = 4$ and $c = 4$. Currently, the allocation of basis states to lists is also done randomly, without ensuring that all basis states are utilized. The goal of this initial simulation is to take a preliminary look at iQAQE's performance and understand how much the results vary between finite shot number simulations and analytical ones. (This also served as a comparison test between the two previously mentioned PennyLane devices. For this small graph, using the `default.qubit` device consistently proved to be more effective. Therefore, we will use this exclusively from now on.) The plots below display the best-so-far average approximation ratio, derived from 10 iQAQE runs, each with different random initial parameters. The results for all three algorithms are generated using a configuration of 4 ansatz layers and an Adam learning rate of 0.99.

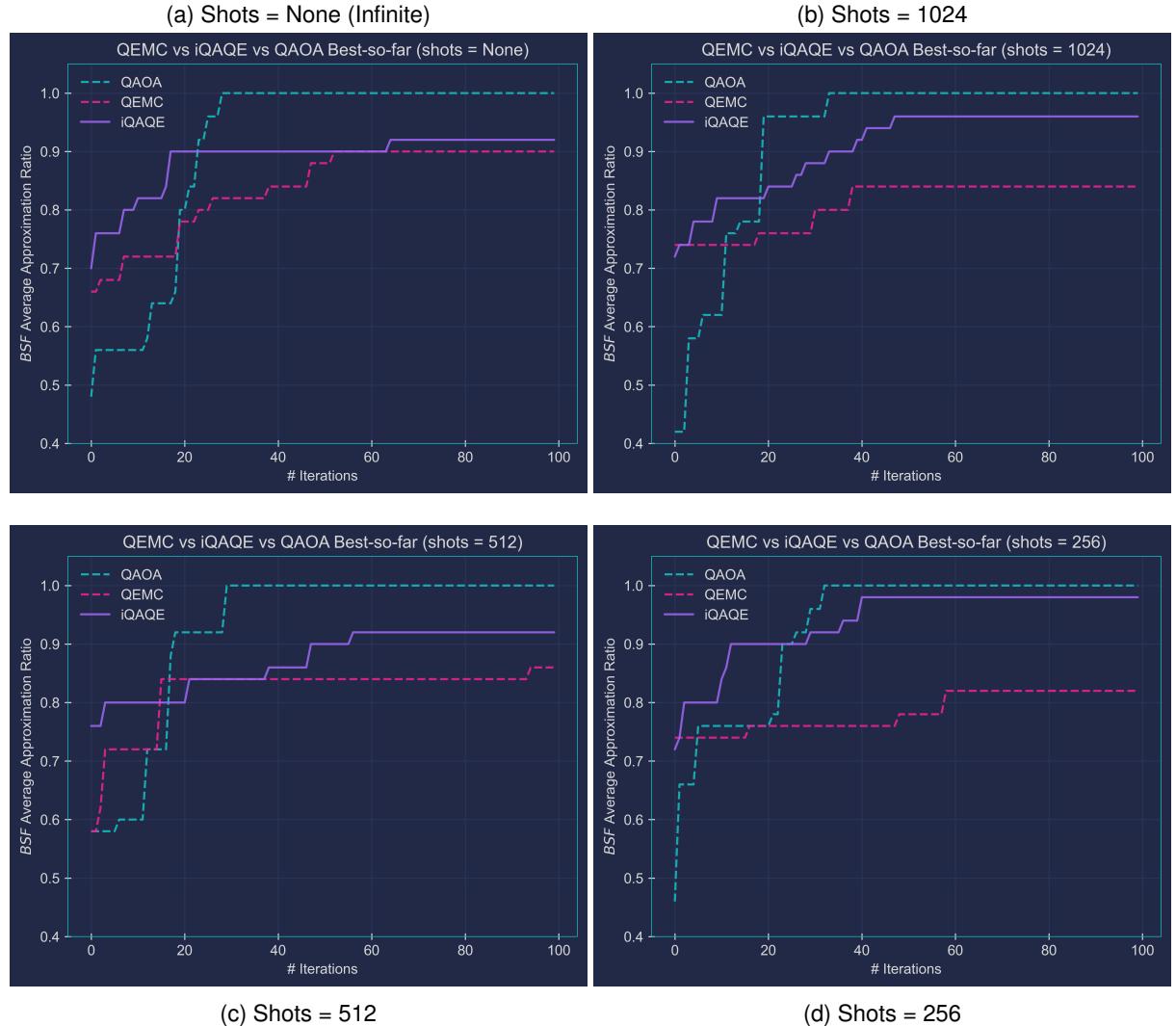


Figure 5.2: Comparison between the 3 VQAs' performances (QAOA, QEMC and iQAQE), using an infinite and finite number of shots. The best-so-far average is plotted.

These plots generally align with our theoretical expectations. As QEMC relies heavily on a large number of shots, transitioning to a finite shot number naturally impacts the results. Surprisingly, however, iQAQE seems to exhibit slightly greater resilience to this effect, despite sharing QEMC’s cost function and ansatz. This resilience might be attributed to the presence of multiple basis states associated with each graph node, potentially reducing the need for exhaustive sampling. If some basis states are harder to sample, the probabilities of the remaining states might adapt to compensate, as governed by the QEMC objective function. It is worth noting that the performance of iQAQE appears to vary with the number of shots used. This variability can be attributed to different random initial parameters for each number of shots and fluctuations from circuit sampling, which introduce more variance in the results. (This is also true for QAOA and QEMC.) Nonetheless, QAOA consistently outperforms both approaches, highlighting its superiority. For the same number of layers, QAOA’s problem-inspired ansatz and cost function consistently yield better results. To determine if iQAQE can surpass QAOA’s performance, extensive testing across various combinations of parameters such as n , c , number of layers, and Adam’s learning rate is necessary. This will motivate some of the grid searches we conduct later in this study. However, given the vast number of potential combinations, testing all of them is impractical. Therefore, some of the more targeted mappings explored in this work aim to constrain one or more of these variables, thereby reducing the total number of combinations for testing. Following this approach, conducting grid searches over the parameter values becomes more manageable and practical to determine the optimal settings. The subsequent sections of this chapter will detail these mapping strategies.

5.2 Polynomial Compression-type Encodings

At this point, driven by our earlier discussion on the nature of random-based algorithms, we begin exploring more meaningful mappings. Here, for each node/list, we opt to fix one or more (different) qubits to 1 (or 0) while allowing the remainder to vary. The number of qubits we fix (k) depends on the desired order of compression. This approach not only enables the utilization of fewer qubits (if $k > 1$) but also establishes the total number of qubits to be used, simplifying subsequent grid searches over parameters. By fixing k qubits, we achieve polynomial compression of order- k (in the number of qubits). The total number of nodes that can be encoded in this manner is determined by $\binom{n}{k} = \frac{n!}{k!(n-k)!}$, where n denotes the number of qubits. Essentially, this represents the number of ways we can select k qubits from the available n and set them to 1 (or 0), each combination encoding one graph node. Consequently, we need to ensure that $\binom{n}{k} \geq N$, where N represents the number of graph nodes. This criterion guides our selection of n and k . In practice, to determine n for any k , we select the smallest n that satisfies $\binom{n}{k} \geq N$. This yields $N = \mathcal{O}(n^k)$. Subsequently, we decide on the number of basis states to associate with each graph node, chosen from a total of 2^{n-k} (with $n - k$ free qubits). While we could potentially utilize all 2^{n-k} basis states for each node, we allow for the flexibility to choose $c \leq 2^{n-k}$. The value of c requires optimization. With the mapping established, the algorithm proceeds as before, employing QEMC’s cost function and ansatz. Notably, we cannot use QAOA’s ansatz since $n \neq N$. Due to polynomial compression, $n < N$ always, unless $k = 1$ (in which case there is no compression).

5.2.1 Basic Polynomial Compression-type iQAQE

This is the simplest type of polynomial compression-based iQAQE scheme. It involves fixing k qubits to 1 instead of 0 (e.g., $\{|11xxx\rangle\}$, for $k = 2$; this refers to a single node). The number of fixed qubits (k) is determined by the desired order of compression, as previously described. Below (Figures 5.3 and 5.4) are the results for $k = 2$ and $k = 3$, representing quadratic and cubic compression levels, respectively. In the figures, the list cardinality is set to 4 (using random mapping) for both the $k = 2$ curves (iQAQE_k2) and the $k = 3$ curves (iQAQE_k3). Additionally, the number of layers is set to 4 for $k = 2$ and 5 for $k = 3$, with an Adam learning rate of 0.98. The final curves are derived from 10 runs of the algorithms, each with different random initial parameters. The results are compared with those of QAOA and QEMC.

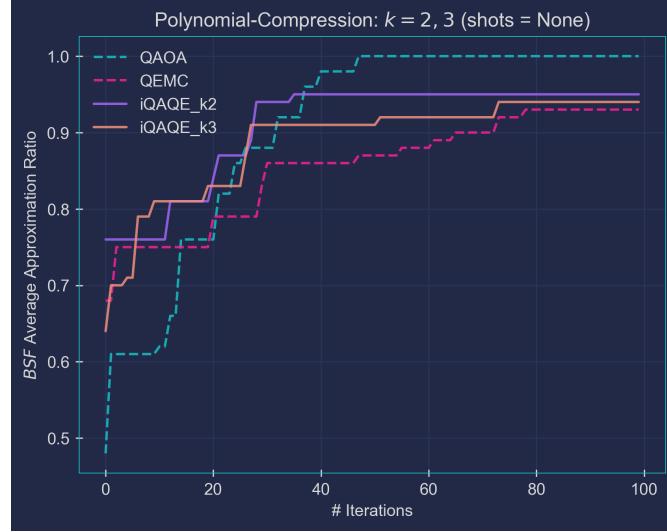


Figure 5.3: BSF Average Approximation Ratio vs. iteration number for the tested VQAs: QAOA, QEMC, and iQAQE, using the aforementioned polynomial compression scheme. The number of layers considered were 4 for $k = 2$ and 5 for $k = 3$.

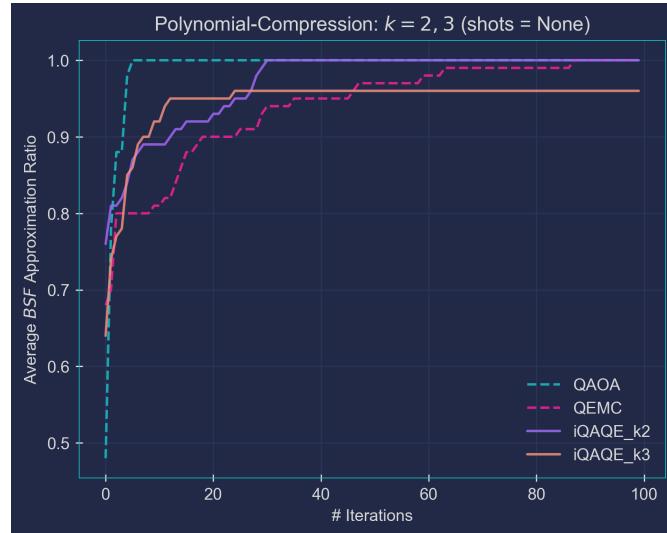


Figure 5.4: Average BSF Approximation Ratio vs. iteration number for the tested VQAs: QAOA, QEMC, and iQAQE, using the aforementioned polynomial compression scheme. The number of layers considered were 4 for $k = 2$ and 5 for $k = 3$.

Notice the difference between the two figures. In the first figure, we average the curves first and then apply the `BSF` transformation. In the second figure, we reverse the process: first applying the `BSF` transformation to each curve and then averaging the transformed curves. The results are quite different, as illustrated. This suggests that both `QEMC` and `iQAQE` are typically more sensitive to fluctuations in their performance curves, hinting at a more complex optimization landscape. This is an important consideration when interpreting the forthcoming results. Often, we will choose to discard the `BSF` average in favor of the average `BSF` (cf. section 5.6). Nonetheless, both approaches are valuable for evaluating the algorithms' properties. Additionally, in some instances, we will simply use the average. Each case will be clearly specified.

Anyways, the primary aim of this study was to assess whether different types of mappings, like the one explored here, offer advantages compared to the purely random case. The results in Figures 5.3 and 5.4 indicate that this may be the case. At the very least, they suggest that there is potential in exploring these ideas further. For example, `iQAQE_k2` achieves a perfect average `BSF` approximation ratio (Figure 5.4), although it requires more iterations [to get there] than `QAOA`. Nevertheless, our method surpasses `QEMC` for $k = 2$. The results for $k = 3$ are less encouraging. However, both mappings use fewer qubits ($n = 5$) than `QAOA` ($n = 8$), which is beneficial for implementation on practical (`NISQ`) quantum computers. The next step is to conduct grid searches over the parameters for this specific graph instance to determine if we can consistently achieve better performance than `QAOA`. This might provide further insights into the underlying principles of the algorithm's performance.

The results of the grid searches conducted for $k = 2$ and $k = 3$ are present in Figures 5.5 and 5.6. Various values were explored for the parameters: `step_size_list = [0.35, 0.4, 0.5, 0.6, 0.7, 0.8, 0.9, 0.95, 0.98, 0.99]`, `n_layers` ranging from 1 to 8, and `c` varying from 2 to $(2^n - k - 1)$, for $k = 2, 3$, where `n_layers` and `c` are integers. Notably, all simulations for the grid searches were executed with an infinite number of shots (`shots = None`). Additionally, this time, we utilized the average curve (from 10 instances) instead of the average best-so-far curve. This choice was made to evaluate the overall performance of the algorithm rather than focusing solely on the best-performing instances. Each run continued until convergence (`abs_tol1 <= 1e-5`) or `max_iter = 1000` was reached, followed by recording the final partition's cut. The objective was to identify any discernible patterns in the grid search to enhance the scheme's performance. However, no apparent pattern emerged, highlighting a recurring challenge in this work: the abundance of hyperparameters makes it challenging to optimize results through straightforward means. The best conclusion we can draw is to avoid high layer counts with high learning rates. Later, we will explore employing a more intricate machine learning scheme to address this challenge. For now, we proceed with a different approach.

Recovering the `QAOA` limit ($k = 1$)

Next, we tested whether we could recover the `QAOA` limit using the `iQAQE` Framework with what we call the `QAOA` mapping. This mapping involves using the same number of qubits as there are nodes (as in standard `QAOA`) and assigning 2^{n-1} basis states to each graph node. For each graph node, a different qubit is fixed to 1, while the remaining $n - 1$ qubits are free to vary, resulting in the aforementioned 2^{n-1}

¹Difference in costs between consecutive iterations.

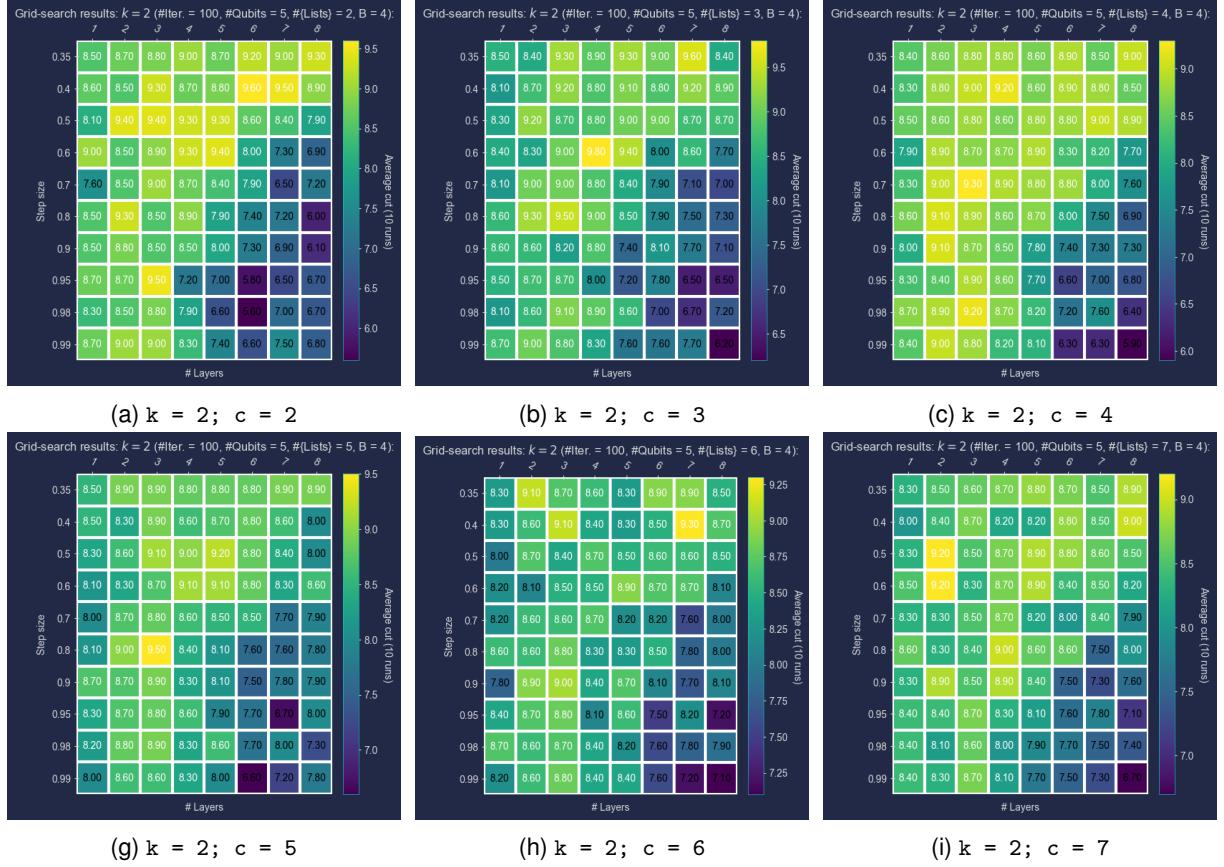


Figure 5.5: Basic Polynomial Compression-type iQAQE – grid search, using $k = 2$. All the simulations were done using an infinite number of shots (shots = None), and each value consists of the average of 10 iQAQE instances, using different random initial parameters.

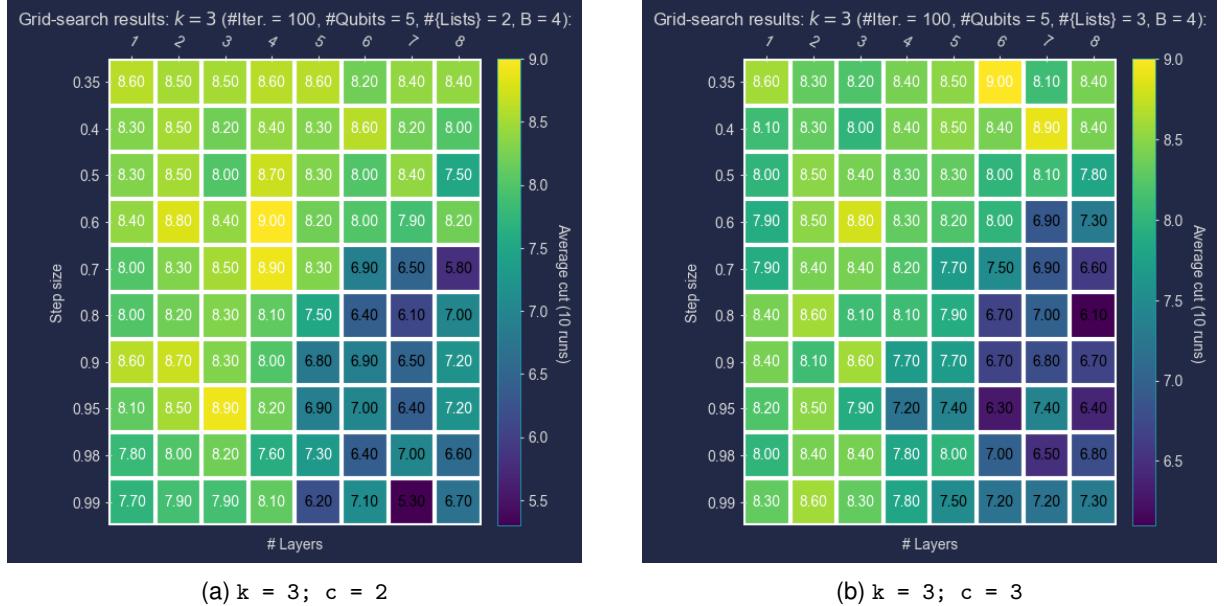


Figure 5.6: Basic Polynomial Compression-type iQAQE – grid search, using $k = 3$. All the simulations were done using an infinite number of shots (shots = None), and each value consists of the average of 10 iQAQE instances, using different random initial parameters.

basis states per node. The rest of the algorithm remains unchanged, employing `QEMC`'s ansatz and cost function. The nodes' probabilities are derived from the normalized sum of the probabilities of their associated basis states. This approach was motivated by our desire to better understand `iQAQE`'s properties and behavior. Previously, we observed (Figure 5.2) that `iQAQE`'s performance seemed to fall between that of `QAOA` and `QEMC`, which is intuitive, as it is designed as an interpolation between the two². To further test this behavior, we used this `QAOA` mapping to see how closely we could approach `QAOA`'s performance levels, which have been the best so far. The following results were obtained, focusing only on analytical simulations (i.e., infinite shots) – we employ the average approximation ratio curve obtained from 10 runs, as we did in the previous grid search:

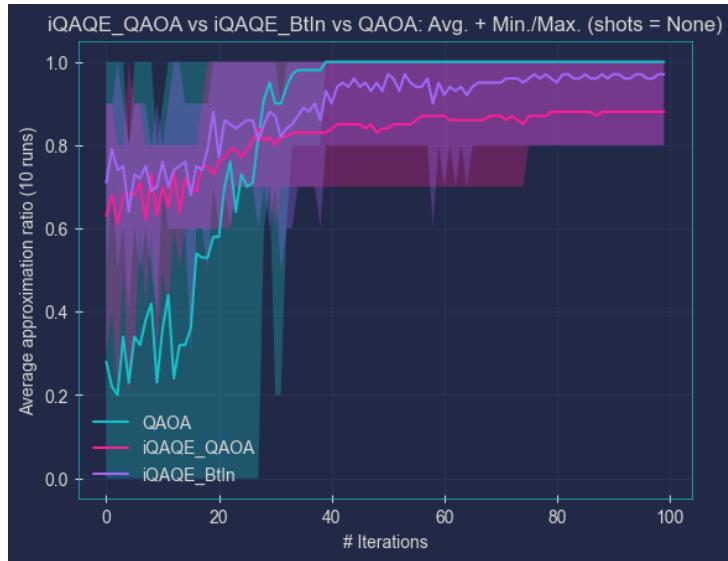


Figure 5.7: Average approximation ratio (\pm Min./Max.) vs. iteration number, for the three tested VQAs: `QAOA`, `iQAQE_QAOA` (as described above) and `iQAQE_Btn`. The latter stands for `iQAQE`, with better initialization. This corresponds to the same mapping as the one in Figure 5.2.

This realization led us to understand that we cannot expect to replicate `QAOA`'s outstanding performance on this graph simply by using a `QAOA`-type mapping. After all, we are still relying on `QEMC`'s cost function and ansatz, so this outcome is not entirely surprising. Moreover, to add to the challenge, we found that `iQAQE_QAOA`'s performance can even be surpassed by another random, non-specific mapping, highlighting how far we are from achieving `QAOA`-like results.

In an effort to bridge this gap, we attempted to integrate `QAOA`'s ansatz into the scheme, alongside the `QAOA`-type mapping. However, after extensive testing, we realized this approach is not feasible. Here's why: using both elements together with the node probability encoding (as a normalized sum of their associated basis states' probabilities) causes the `QEMC` cost function to become independent of the ansatz's parameters³. This means we cannot proceed with training, rendering the scheme ineffective. Thus, implementing this approach in its current form is not possible. For it to be successful, other aspects would need to change, such as the cost function or the method of deriving node probabilities from basis states' probabilities, which is the main issue here.

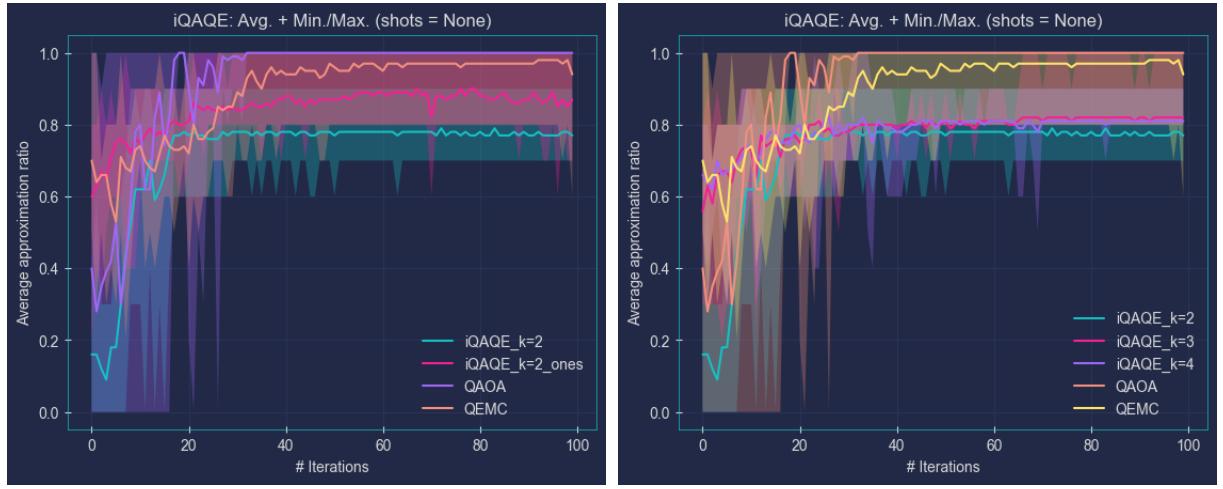
²We'll see later that this isn't entirely true.

³To verify this, we performed analytical calculations for a simple toy model: a graph with two nodes connected by a single edge. The results generalize to larger graphs.

5.2.2 Correlation-based iQAQE

This scheme closely resembles what was previously discussed in subsection 5.2.1. The key distinction lies in the inclusion of the possibility for both 0's and 1's to be fixed (e.g., $\{|11xxx\rangle, |00xxx\rangle\}$, for $k = 2$; recall, this encodes a single node). Consequently, the scheme aims to identify correlations among the different qubits, wherein "correlations" denote identical colors. In other words, the constructed lists regard the k fixed qubits as positively correlated, meaning they are either all set to 1 or all set to 0. This approach was inspired by the notion that by identifying correlations between nodes, it becomes feasible to color the entire graph as long as one node is initially colored. Although seeking correlations between qubits, as done here, differs from seeking correlations between nodes, we anticipated that it might yield promising results to some extent. In this scenario, the color of each node is determined by the probability that its k associated qubits are all positively correlated. As such, c is now fixed at 2^{n-k} for n qubits. Unlike the previous case (Basic Polynomial Compression-type iQAQE), we do not permit $c < 2^{n-k}$.

To delve deeper into the properties and characteristics of this scheme, we conducted a thorough formal analysis, exploring its limits by considering scenarios where k ranges from 1 to n qubits, covering the entire spectrum of possibilities. Our primary objective was to ascertain if we could formulate this correlation-based iQAQE scheme in a manner that truly interpolates between QAOA and QEMC. We aimed to determine if we could retrieve both algorithms by adjusting the parameters accordingly. However, with this correlation-based scheme, we not only cannot recover QAOA, but we also fail to retrieve QEMC. Thus, this does not represent an interpolation between the two algorithms. Likewise, earlier, we loosely referred to the iQAQE Framework as an interpolation. However, upon closer scrutiny, this characterization seems inaccurate. While it is possible to recover the QEMC limit⁴, the same cannot be said for QAOA, as we have observed before. Additionally, we present the results of numerical simulations applied to the standard 8-node graph using this correlation-based iQAQE approach (Figure 5.8).



(a) Comparison of correlation-based iQAQE for $k = 2$ with results from QAOA, QEMC, and the previous Basic Polynomial Compression-type iQAQE⁵ ($iQAQE_k=2_ones$).

(b) Comparison of correlation-based iQAQE results for $k = 2, 3$, and 4 with outcomes from QAOA and QEMC.

Figure 5.8: Correlation-based polynomial-type compression scheme ($k = 2, 3$ and 4): numerical simulations for the usual 8-node graph.

⁴Simply utilize the appropriate number of qubits and select one unique basis state for each node.



Figure 5.9: Correlation-based polynomial-type compression scheme ($k = 2, 3, 4, 5, 6$ and 7): numerical simulations for the usual 8-node graph.

Note that we once again compute the average approximation ratio from 10 runs, each with different random initial parameters, and generate all curves with `n_layers = 3` and `step_size = 0.95`. Additionally, we plot the average, with the shaded region representing the maximum and minimum obtained approximation ratios (out of the 10 runs). Performance-wise, this scheme does not fare well. It is significantly outperformed by QAOA, QEMC, and the basic polynomial compression-type iQAQE. We also compared the results for all possible values of k (Figure 5.9). It turns out that the specific value of k is not very meaningful, as performance is quite similar across all values.

The underperformance is likely due to our doubling the number of basis states in each sub-list. This encoding results in each pair of sub-lists having a 50% overlap, similar to the Basic Polynomial Compression-type iQAQE, specifically when $c = 2^{n-k}$ for n qubits. However, with twice⁶ the number of basis states this time, the number of overlapping basis states is significantly higher. This increase means that more states are shared by multiple nodes. When the overlap is too large, it becomes more challenging to adjust the color of one node without affecting the others. We believe this to be the main reason for the underperformance. We will be mindful of this in the design of future schemes.

⁵This refers to the scenario where only 1's are fixed, not 0's (Basic Polynomial Compression-type iQAQE). To maintain a fair comparison with the correlation-based scheme, we use $c = 2^{n-k}$ for n qubits in this method.

⁶Either 0's or 1's can be fixed.

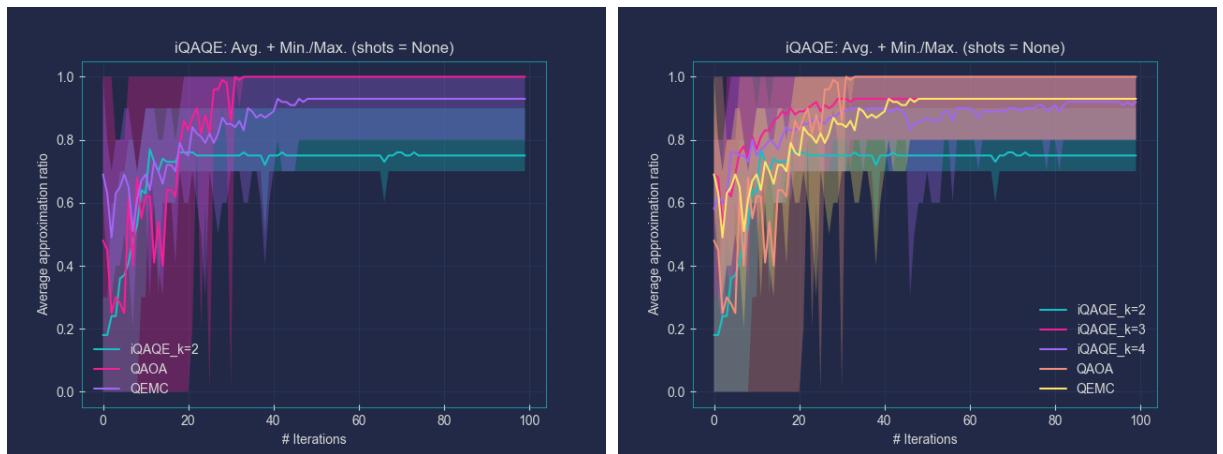
5.2.3 Fixed-Parity iQAQE

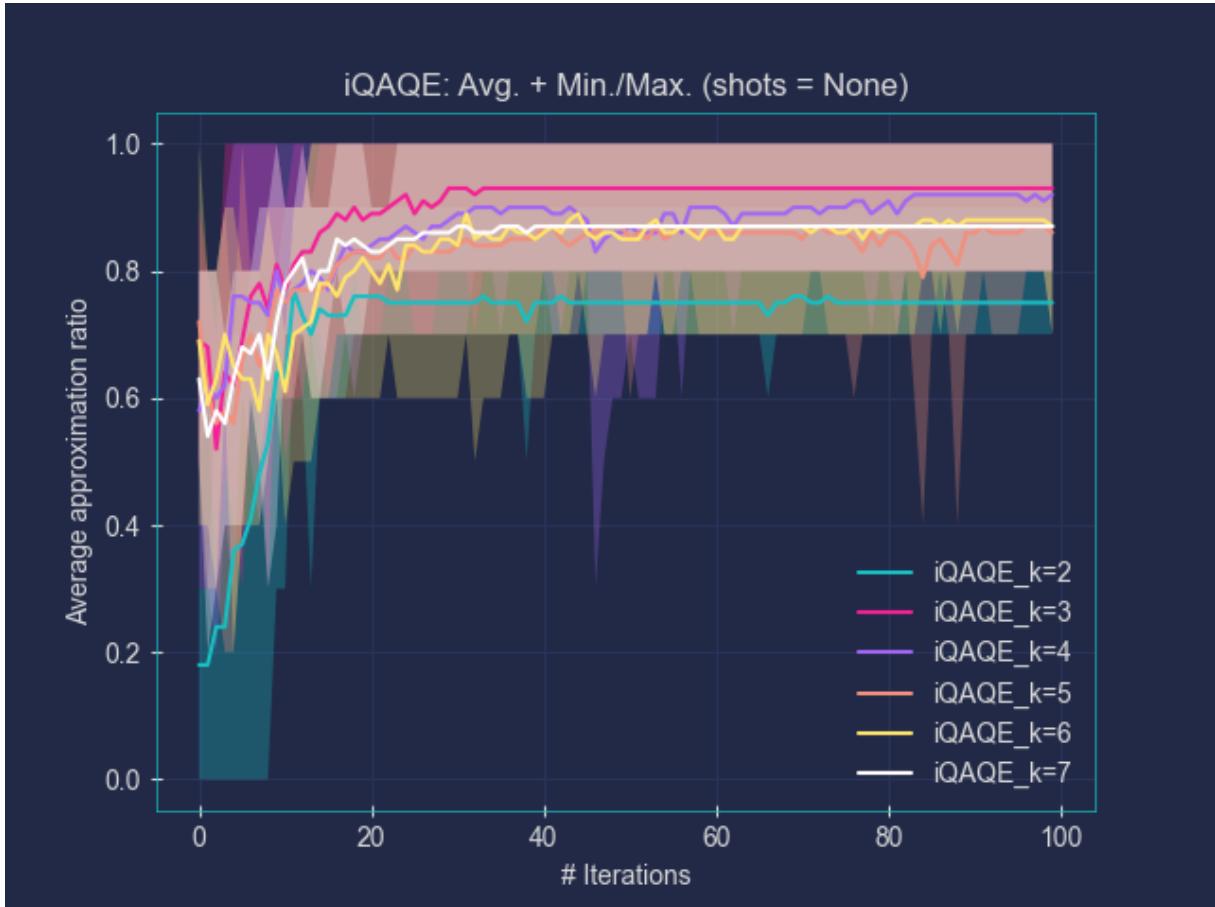
In this section, we present another heuristic method for mapping the basis states to the nodes. Although it is similar to the previous one, this time we **fixed the parity of the selected k qubits to be even**. Parity is determined by the number of 1's: if the count is even, the parity is even; otherwise, it is odd. For instance, for $k = 3$ (with $n_qubits = 5$), the lists would take the form: (Keep in mind that we are using an 8-node graph. Additionally, x denotes a free qubit: either 0 or 1. The cardinality of each list is given by $c = \sum_{i=0}^{\lfloor \frac{k}{2} \rfloor} \binom{k}{2i} \cdot 2^{(n-k)}$ for n qubits.)

1. $\{|000xx\rangle, |011xx\rangle, |101xx\rangle, |110xx\rangle\};$
2. $\{|00xx0\rangle, |01xx1\rangle, |10xx1\rangle, |11xx0\rangle\};$
3. $\{|0xx00\rangle, |0xx11\rangle, |1xx01\rangle, |1xx10\rangle\};$
4. $\{|xx000\rangle, |xx011\rangle, |xx101\rangle, |xx110\rangle\};$
5. $\{|x0x00\rangle, |x0x11\rangle, |x1x01\rangle, |x1x10\rangle\};$
6. $\{|x00x0\rangle, |x01x1\rangle, |x10x1\rangle, |x11x0\rangle\};$
7. $\{|x000x\rangle, |x011x\rangle, |x101x\rangle, |x110x\rangle\};$
8. $\{|0x0x0\rangle, |0x1x1\rangle, |1x0x1\rangle, |1x1x0\rangle\}.$

Once again, numerical simulations were performed, and the obtained results are presented below (Figure 5.10). Although the performance is still not quite on par with QAOA, it is significantly better than in the previously considered scenario. Note that the case where $k = 2$ corresponds to the previous correlation-based scheme. For $k = 2$, requiring the pair to be even is equivalent to requiring them to be the same (either both 0 or both 1), which explains the similarity in results. This similarity is apparent in the less favorable outcomes shown for $k = 2$ in Figure 5.10, which qualitatively match those in Figures 5.8 and 5.9. However, when we allow for $k > 2$, the performance improves significantly, likely due to the reduced overlap between the basis states associated with different nodes.

A keen observer might notice that this scheme's results also allow for generating the MaxCut partition, as shown by the shaded pink region in Figure 5.10b for $k = 3$. These shaded regions depict the maximum and minimum cuts achieved out of the 10 runs. This reveals that, despite the greater variability and lower average performance than QAOA, the algorithm can achieve the MaxCut partition in some runs. In practice, this is crucial. When running the algorithm 10 times, our main interest is in the best outcome. This highlights the potential of this heuristic scheme.

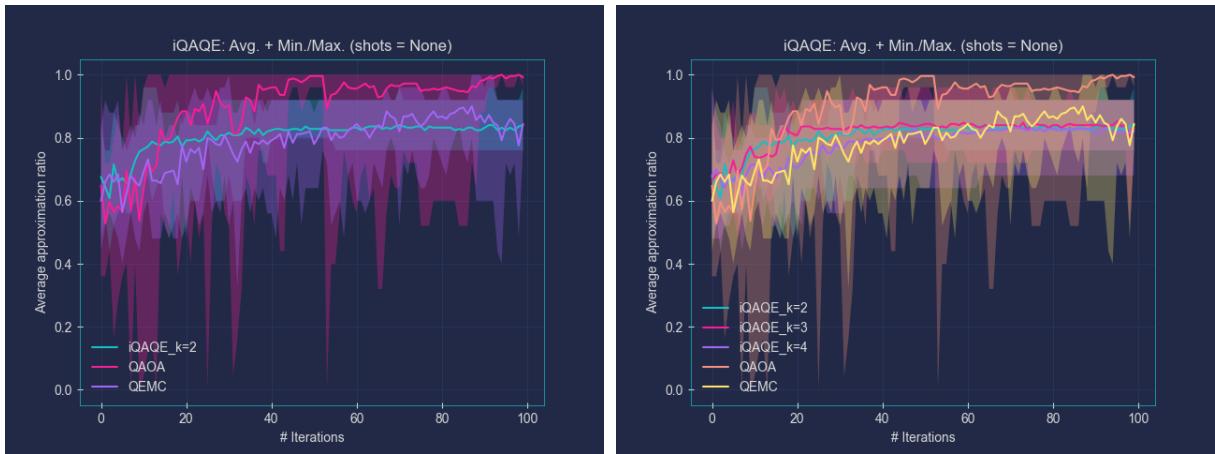




(c) Fixed-parity iQAQE for all values of $k = 2, 3, 4, 5, 6$, and 7 (8-node graph).

Figure 5.10: Fixed-parity polynomial-type compression scheme: numerical simulations for the usual 8-node graph. Please note that we utilize `n_layers = 3`, `step_size = 0.95` and $B = 4$.

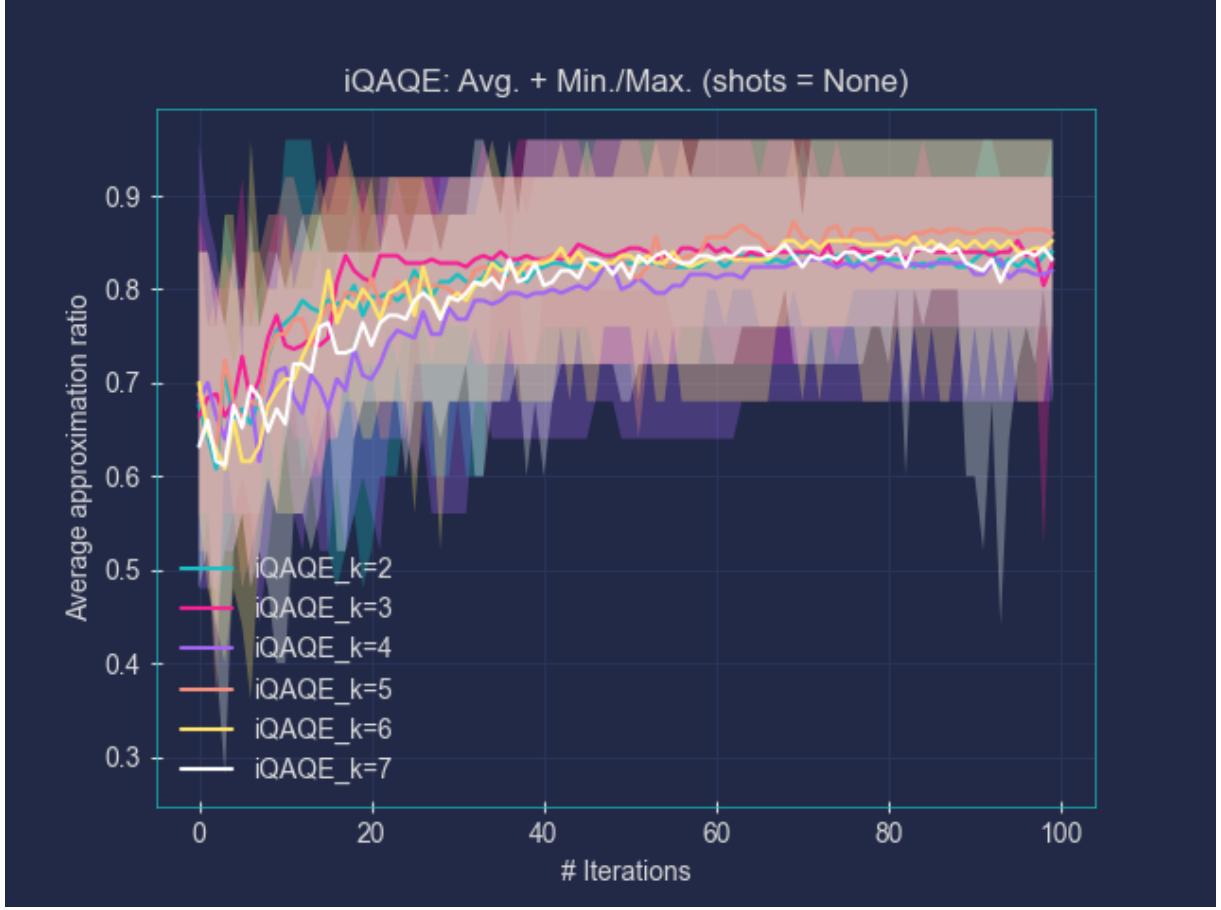
Furthermore, numerical simulations were carried out on a random 16-node graph⁷. Considering the satisfactory performance observed with the previous 8-node graph, we were interested in assessing the algorithm's performance on a larger scale. The results are now presented (Figure 5.11).



(a) Fixed-parity iQAQE for $k = 2$ compared with the results from QAOA and QEMC (16-node graph).

(b) Fixed-parity iQAQE for $k = 2, 3$, and 4 compared with the results from QAOA and QEMC (16-node graph).

⁷This graph was generated using NetworkX's `nx.gnm_random_graph(n, m)` with $n = 16$ nodes and $m = 32$ edges, using `np.random.seed` set to 44.



(c) Fixed-parity iQAQE for all values of $k = 2, 3, 4, 5, 6$, and 7 (16-node graph).

Figure 5.11: Fixed-parity polynomial-type compression scheme: numerical simulations for a 16-node graph. Please note that we utilize `n_layers = 3`, `step_size = 0.95` and $B = 8$.

The previously identified trends are also apparent here. However, the performance now seems insufficient to achieve the `MaxCut` partition (determined by brute-force), as illustrated by the shaded areas in Figure 5.11c: the maximum AR never quite reaches 1. Moreover, for this graph, varying the value of k does not significantly impact performance. Hence, it is advisable to choose the value of k that requires the fewest qubits. In this instance, $k = 3$ is optimal, needing only `n_qubits = 6`, since $\binom{6}{3} = 20 \geq 16$.

5.3 Extended-QEMC

Amid the various schemes we've attempted, I've devised an extension to the standard QEMC scheme. This new approach enhances QEMC by incorporating m additional qubits beyond the usual $\lceil \log_2(N) \rceil$ qubits required for N graph nodes.

5.3.1 Unmodified Extended-QEMC

Rather than associating a single basis state with each graph node, this scheme assigns 2^m basis states per node. Importantly, similar to QEMC, the sets of basis states associated with different nodes

do not overlap. This non-overlapping property is crucial, as we suspect that significant overlap between nodes' lists might contribute to the subpar performance observed in the correlation-based and fixed-parity (16-node graph) iQAQEs.

This extended scheme can be implemented within the iQAQE formalism by using appropriate sublists of basis states. The assignment of states to each list is currently arbitrary. For simplicity, we use a straightforward partition: the first 2^m basis states go to the first list, the next 2^m to the second list, and so on (referred to as Unmodified Extended-QEMC). However, there may be a more optimal method for distributing these states that we have not explored here.

Additionally, it is important to note that Extended-QEMC requires more qubits than regular QEMC. This increase in qubits prevents the scheme from being easily classically simulable and allows for more basis states to be associated with each graph node. Consequently, this should provide more variables to manipulate, potentially improving our chances of finding a good solution.

Now, I present the results of applying this scheme to the usual 8-node graph (Figure 5.12).

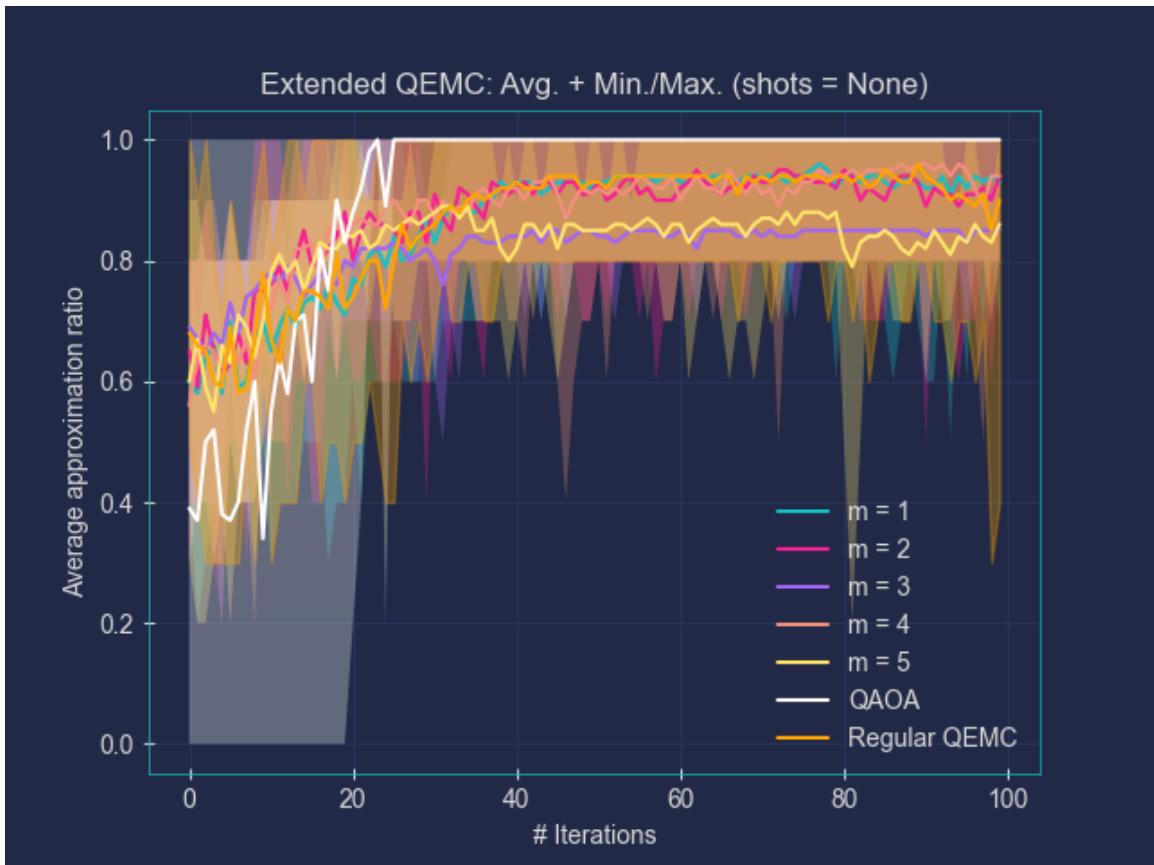


Figure 5.12: Implementation of the Unmodified Extended-QEMC scheme for the usual 8-node graph: comparison with QAOA and regular QEMC for various values of m . Note that m is not extended beyond 5, as this would exceed the number of qubits used in QAOA. Additionally, we utilize `n_layers` = 3, `step_size` = 0.95 and B = 4.

As shown in this figure, the performance of the Unmodified Extended-QEMC scheme is qualitatively very similar to that of regular QEMC. For some values of m , we achieve slightly better results, but these improvements are not particularly significant. After conducting several tests, I found no clear pattern to predict which values of m would yield the best results, supporting the earlier conclusion. Therefore, this

presents another heuristic for mapping the basis states, with performance comparable to the standard QEMC scheme. While this is a satisfactory outcome, it is not particularly remarkable.

5.3.2 Cardinality = 1 Extended-QEMC

Motivated by the results in Figure 5.17a⁸, I decided to make a slight change to the Unmodified Extended-QEMC scheme: instead of assigning 2^m basis states to each node, we assign only one basis state (randomly) to each node (Cardinality = 1 Extended-QEMC). Although this adjustment occasionally yielded marginally better results than regular QEMC, we do not consider these improvements significant (see Figure 5.13 for reference).

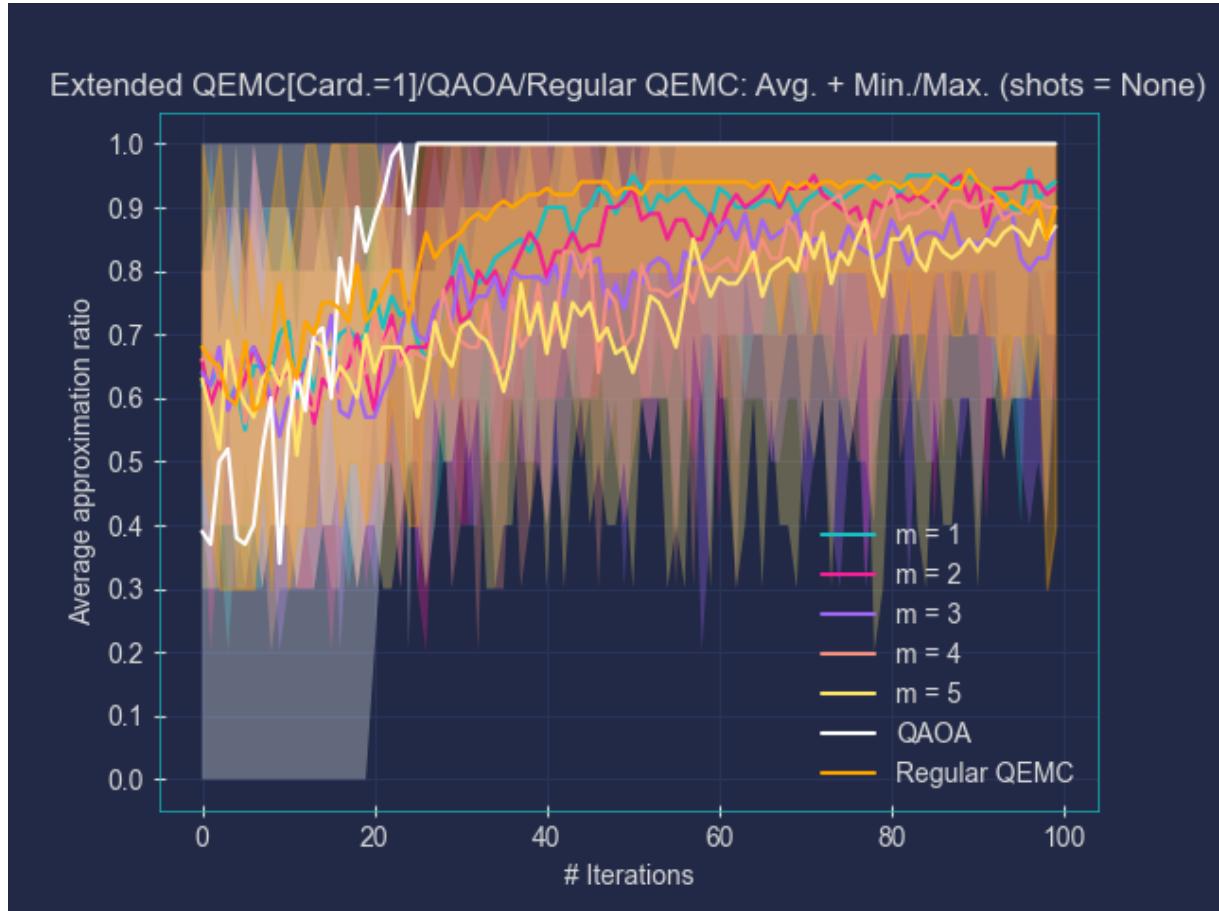


Figure 5.13: Implementation of the Cardinality = 1 Extended-QEMC scheme for the standard 8-node graph: comparison with QAOA and regular QEMC for various values of m . Please note that we utilize `n_layers = 3`, `step_size = 0.95` and `B = 4`.

5.4 Alternative ansätze

After some reflection, I began to question whether the performance of QEMC and iQAQE could be hindered by the use of a problem-agnostic ansatz. It's generally advantageous to incorporate problem-specific information into the ansatz, as exemplified by problem-inspired ansätze. Additionally, I pondered whether excessive entanglement or correlations among the system's basis states might be affecting the

⁸The trajectory of this study was far from linear. The results presented here are not in the chronological order of our research, hence this reference to later work, which was actually conducted before devising the Extended-QEMC scheme.

schemes' performances. Such phenomena could make it challenging for the model to adjust the amplitude of a specific basis state without significantly impacting others. To address this potential issue, the concept of implementing non-deterministic CNOT gates (ND-CNOTs) occurred to me⁹. This approach would allow the model to dynamically adjust the degree of entanglement. The most rudimentary way to implement this is by employing parameterized R_x gates before each CNOT's control qubit, which is precisely what we did. The R_x gate allows the control qubit to rotate in and out of the $|1\rangle$ state, thus adjusting the degree of entanglement introduced by the CNOT gate. The results for the standard 8-node graph are presented below (Figure 5.14), for both QEMC and Cardinality = 1 Extended-QEMC.

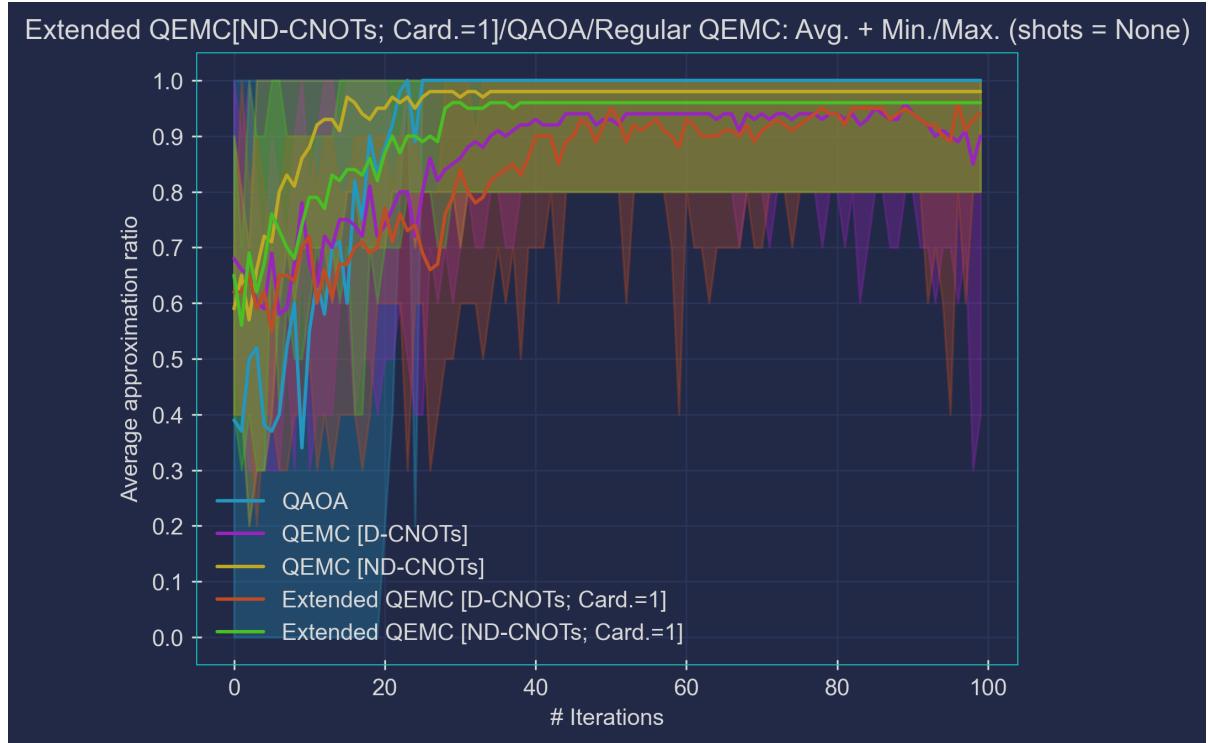


Figure 5.14: Comparison of QEMC and Cardinality = 1 Extended-QEMC implementations utilizing both D-CNOTs and ND-CNOTs for the usual 8-node graph, alongside QAOA. We utilize $B = 4$.

Several noteworthy observations can be made here. Firstly, it's apparent that ND-CNOT-based schemes exhibit significantly more stable performance with fewer fluctuations compared to deterministic CNOT-based (D-CNOT¹⁰) schemes. This discrepancy arises from differences in their implementations: D-CNOT circuits employ `n_layers = 3` and `step_size = 0.95` (Adam's learning rate), whereas ND-CNOT circuits use `n_layers = 2` and `step_size = 0.5`. The latter, `step_size = 0.5`, accounts for the smoother curves observed in Figure 5.14, as expected, and could also clarify the observed faster convergences. Secondly, the performance appears noticeably enhanced when utilizing ND-CNOTs, particularly evident in the case of regular QEMC: a distinct improvement is evident when comparing the yellow and purple lines. Despite introducing added complexity to the optimization landscape, this approach appears to yield significant performance gains, at least for this small 8-node graph. Thus, we've proposed a novel scheme that builds upon QEMC, showing great potential to outperform it.

⁹We recognize that the term "non-deterministic CNOT gates" is employed in linear optical quantum computing (KLM protocol), yet we use it differently in this context.

¹⁰This refers to the standard CNOTs.

Moreover, this analysis serves to validate the notion that deviating from purely problem-agnostic ansätze is generally beneficial. This insight will inform our approach in developing future schemes in subsequent research endeavors. Encouraged by the promising outcomes from this study, we opted to expand the testing of this scheme to larger graphs to assess its scalability. This paves the way for comparisons with the Goemans-Williamson algorithm.

5.5 Goemans-Williamson and Bigger Graphs

At one point, I began to question whether it was appropriate to directly compare our results to QAOA, which stands out as the best-performing VQA among those tested. After thorough deliberation, we concluded that, indeed, it makes sense to do so for smaller graphs. However, as graphs increase in size, the availability of quantum machines with a necessary number of qubits to support such QAOAs becomes limited. This circumstance has led to the emergence of schemes like QEMC [3] and others [18], which aim to address the limitations imposed by the number of qubits. At this stage, it becomes quite pertinent to draw comparisons with the Goemans-Williamson algorithm.

In light of this, I proceeded to evaluate the performance of our proposed schemes on larger graphs. Two such graphs were considered:

1. 32-node Erdős–Rényi graph, constructed using NetworkX's `nx.erdos_renyi_graph(n=32, p=0.2, seed=0, directed=False)`, where each edge is included in the graph with a probability of $p = 0.2$;
2. 100-node Erdős–Rényi graph, defined similarly to the previous graph, utilizing `nx.erdos_renyi_graph(n=100, p=0.2, seed=0, directed=False)`.

The results obtained from these evaluations are presented below¹¹ (Figures 5.15 and 5.16).

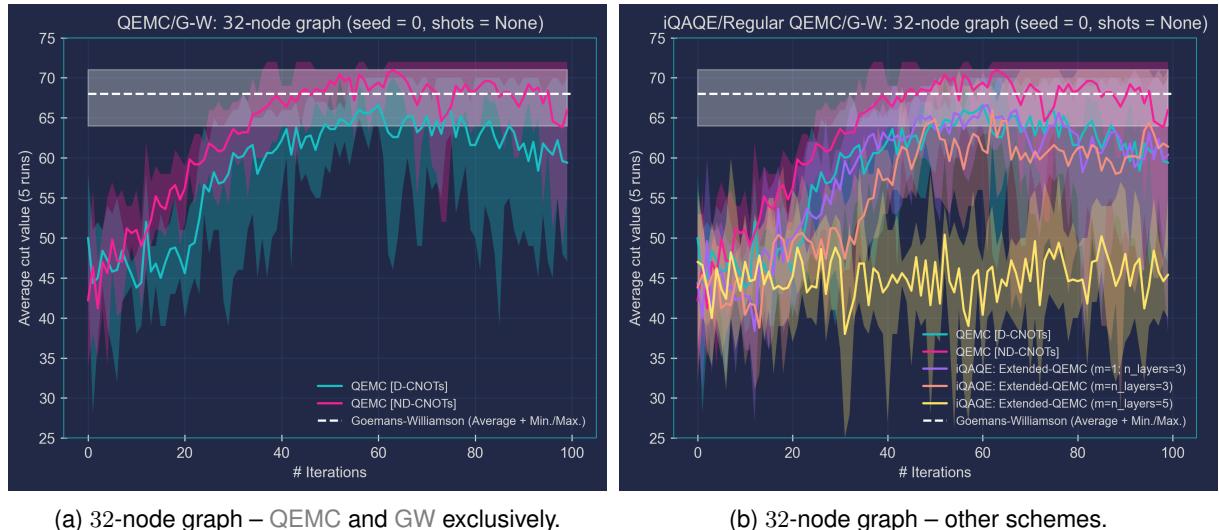
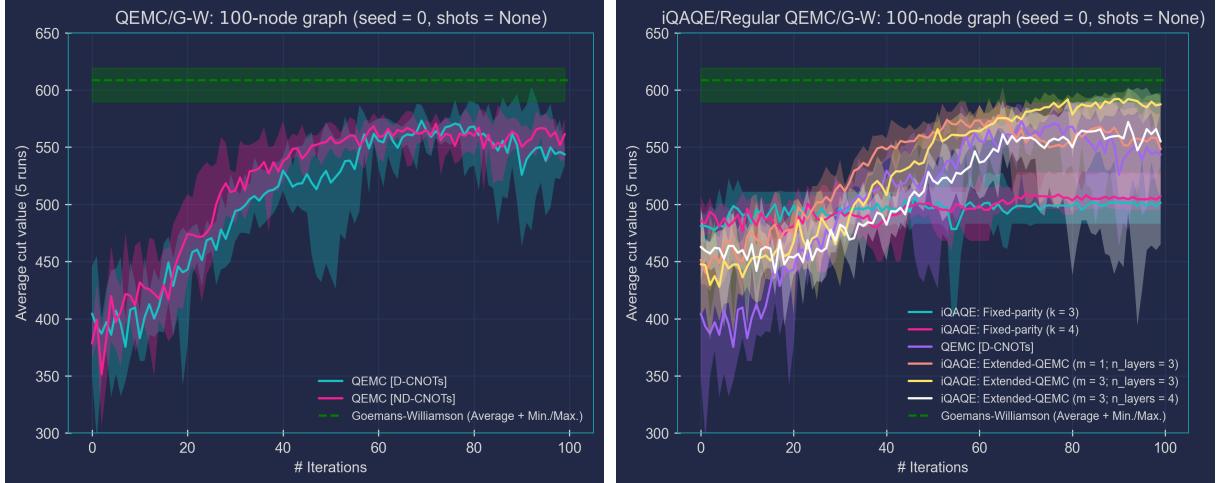


Figure 5.15: Comparison of performance (average cut values) among various VQA schemes for the specified 32-node graph. `n_layers = 3` was employed for both QEMCs, with `step_size = 0.95` for D-CNOT-QEMC and `step_size = 0.5` for ND-CNOT-QEMC. For all Extended-QEMC schemes, `step_size = 0.95` was utilized, with the number of layers specified in the legend. We utilize $B = 16$.

¹¹Please note that we have included only the most promising schemes.



(a) 100-node graph – QEMC and GW exclusively.

(b) 100-node graph – other schemes.

Figure 5.16: Comparison of performance (average cut values) among various VQA schemes for the specified 100-node graph. This time, all considered VQAs utilized `n_layers` = 3 and `step_size` = 0.95, except for Extended-QEMC with m = 3, which used `n_layers` = 4. We utilize B = 50.

Please note that we averaged the results over 5 runs instead of the typical 10. This adjustment was made to expedite the process while still achieving a reasonable averaging of statistical fluctuations. Additionally, in Figure 5.16b, I included the fixed-parity iQAQE scheme to assess its performance on a larger graph.

Now, concerning the results, several interesting observations emerge. For the 32-node graph, Extended-QEMC's performance varies considerably with different values of m and `n_layers`, as previously observed, making it inconsistent and challenging to optimize for other graphs. Once again, we observe the benefits of using ND-CNOTs in QEMC compared to the usual D-CNOTs. One might also observe that the average performance of the ND-CNOT-based QEMC is highly competitive with GW, occasionally even outperforming it. Furthermore, the maximum performance achieved by the ND-CNOT-based scheme surpasses that of GW, as evident from the comparison of the shaded pink and white regions (Figure 5.15a). This marks the first instance in this study where one of our heuristics surpasses classical state-of-the-art algorithms, representing a significant milestone and underscoring the potential for further exploration of the proposed iQAQE Framework.

Turning to the 100-node graph, the situation differs somewhat. This time, the Extended-QEMC scheme stands out, delivering remarkable performance competitive with GW. Additionally, we found that the previously mentioned fixed-parity iQAQE performs exceptionally poorly, demonstrating its inability to generalize to larger graphs. Moreover, there is no longer a significant performance gap between the ND-CNOT-based and regular QEMC schemes, suggesting that the former doesn't scale well to larger graphs. We suspect this might be due to the heightened complexity of the optimization landscape, making it more difficult for the model to locate an optimal solution. This increased complexity arises as a result of the larger number of qubits, which necessitates the optimization of a greater number of ND-CNOT gates. Regardless, for this 100-node graph, reaching the average performance level of the Goemans-Williamson algorithm remains elusive, indicating ample room for improvement. However, once again, promising results are evident in the maximum cuts (shaded regions).

5.6 Best-so-far correction

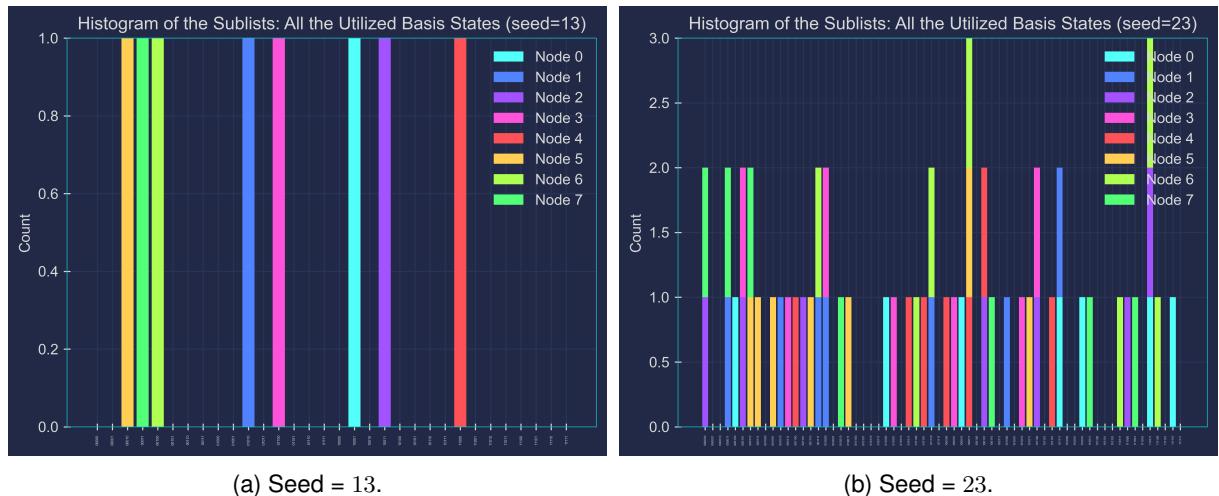
As previously noted, there exists a distinction between implementing the `BSF` transformation before or after the averaging process. We assert that the former (prior to averaging) provides the most accurate representation of the algorithm's optimal performance. Therefore, in alignment with the average `AR` plots previously showcased, we opted to rerun several earlier schemes with the `BSF` transformation applied before averaging/computing the median. Given the substantial number of resulting graphs, they will be included in Appendix A. The plots shown there feature the median best-so-far cut and are intended to complement the previous analyses.

5.7 Randomized iQAQE benchmarking

In this section, our aim is to adopt a more systematic approach to identify ways to enhance the mapping of basis states. While heuristics have served us thus far, we sought to determine whether there are additional patterns that we may have overlooked by relying solely on these heuristics.

Initial proposition

Continuing from the earlier discussion, we ran numerous iQAQE instances, each with random qubit number, list cardinalities, and assignments¹². The underlying idea was to identify optimal combinations of these by running multiple instances of iQAQE. Subsequently, we aimed to discern meaningful patterns that could inform the development of a strategy for more refined basis state selection. However, our analysis did not yield a clear indication of how the partitioning should proceed. As demonstrated below, there existed a plethora of vastly different combinations of qubit number, list cardinality, and partitioning that performed similarly well. Consequently, it proved humanly impossible to extract any coherent insights from this dataset. The results supporting this claim are presented in Figure 5.17 for the usual 8-node graph.



¹²We always employed `n_layers` = 3, `step_size` = 0.99, and `B` = 4. This time, the final selected cut is derived from the `BSF` average curve (not the average `BSF`), constructed from 10 repetitions. Although not optimal, this does not alter our interpretation of the results.

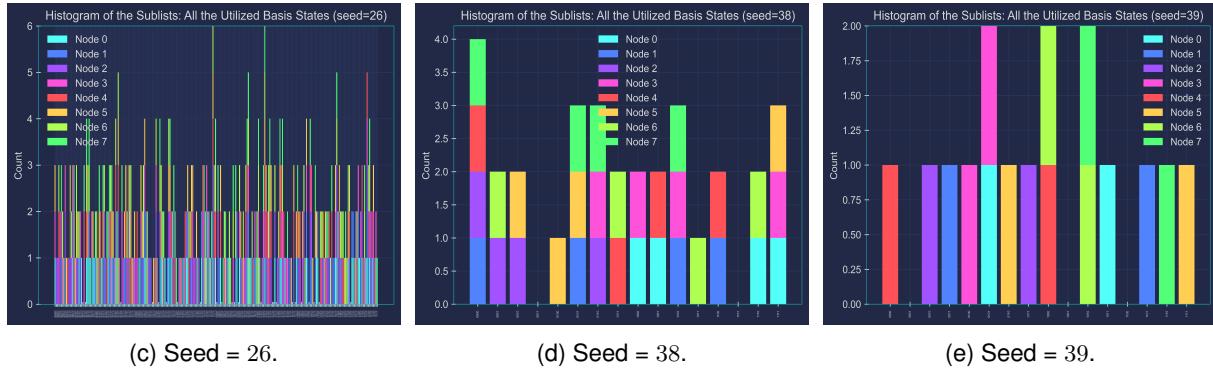


Figure 5.17: Basis states' distributions over the 8 nodes of the usual graph for many different `np.random.seed` seeds. Only those for which the best-so-far average approximation ratio after 100 iQAQE iterations was found to be ≥ 0.975 were selected. (These were the best performing seeds out of the 50 tested for $\text{seed} \in [0, 49] \cap \mathbb{Z}$.)

As can be seen, not much can be directly inferred from this. When confronted with seemingly unstructured data, a natural progression is to employ machine learning schemes to discern patterns. Indeed, this aligns with the forte of such models. Various options were considered, spanning from neural networks to regression and clustering models. Initially, we opted to implement a basic regression model for its simplicity and ease of interpretation. Establishing such a model necessitates identifying a set of statistical properties of the input data (lists of basis states) that can serve as inputs for the regression model. This discussion follows in the subsequent section.

Machine learning-based approach

Following the prior discussion regarding the iQAQE formalism (refer to Figure 5.17), our objective was to construct a small machine learning (ML) model aimed at comprehending the patterns in basis states' partitioning that contribute to the algorithm's performance variations. Naturally, such a model necessitates defining input (independent variables - IVs) and output (dependent variable - DV). As previously mentioned, our IVs will encompass statistically significant properties extracted from each specific sub-list. Alternatively, including the basis states' mapping directly as the model's input presents challenges concerning memory-efficient encoding. Traditional one-hot encoding methods are inadequate due to the vast number of potential basis states available (which scales exponentially with the number of qubits). Thus, we opted for the following statistical properties: (These make up our IVs.)

- **Type 1 variables** - these are input variables that are defined for each of the sub-lists of basis states. They are:
 - **Number of basis states** in each sub-list, the so-called cardinality of the sub-list;
 - **Average Hamming weight** of the basis states in each sub-list;
 - **Standard deviation/variance of the Hamming weights** in each sub-list;
 - **Average pair-wise Hamming distance** between the basis states in each sub-list;
 - **Standard deviation/variance of the pair-wise Hamming distances** between the basis states in each sub-list;

- **Parity score** of each sub-list: this is obtained by: #(Even basis states) – #(Odd basis states), for each sub-list.
- **Type 2 variables** - these are input variables defined for the entire set of selected basis states across all the sub-lists. They are:
 - **Average Hamming weight** of the entire set of basis states;
 - **Standard deviation/variance of the Hamming weights** of the entire set of basis states;
 - **Average pair-wise Hamming distance** between the basis states in the entire set;
 - **Standard deviation/variance of the pair-wise Hamming distances** between the basis states in the entire set;
 - **Parity score** of the entire set of basis states. (Given by the sum of the parity scores of the sub-lists);

– **Global basis-states' entropy**: this is given by the Shannon entropy of the basis states' distribution in the entire set. It is calculated as: $-\frac{1}{\log_2(N)} \sum_{i=1}^N p_i \log_2(p_i)$, where p_i is the probability of the i^{th} basis state in the set. Note that it is normalized by the maximum entropy value, which is $\log_2(N)$, where N is the total number of basis states in the set. (This entropy's maximum is achieved when we have a uniform distribution of the basis states.) Also, this is a measure of the randomness of the basis states' distribution in the whole set.

The dataset was constructed in a similar manner to that of Figure 5.17, but expanded to encompass 1000 seeds. For each seed, we computed the **median** BSF cut value after training (**DV**) and the associated independent variables (**IVs**) related to the partitioning of basis states. (It's worth noting that we employed `n_layers = 3, step_size = 0.99, B = 4, max_iter = 50, and repeat = 5`, i.e., 5 repetitions for each seed.)

With the **IVs** and **DV** finally chosen, and the data set fully constructed, we now need to define our model. We decided to start simple, with a basic multiple linear regression model. Here, the **DV** is computed as a linear combination of the **IVs**, as below:

$$y = \beta_0 + \beta_1 x_1 + \beta_2 x_2 + \dots + \beta_n x_n + \epsilon, \quad (5.1)$$

where x_i are the input variables (**IVs**), y is the dependent variable (**DV**), β_i are the coefficients, and ϵ is the error term. The coefficients will be estimated using the Ordinary Least Squares (**OLS**) method. The model's goodness-of-fit will be evaluated using the R^2 value. Additionally, it's worth noting that the independent variables were scaled using the `StandardScaler` from the `sklearn.preprocessing` module, a common practice in machine learning. This process standardizes the features by removing their mean and scaling them to unit variance, ensuring they have equal weight in the model. This standardization provides a stable foundation for both multiple linear regression and Principal Component Analysis (**PCA**).

After implementing the model, we obtained an R^2 of 0.1023, which is far from the ideal value of 1, and a 5-fold cross-validation loss¹³ of 0.0061 with a standard deviation of 0.0004. To improve the model, I analyzed the independent variables (IVs) by examining their correlation matrix and Variance Inflation Factor (VIF). The correlation matrix revealed that some IVs were highly correlated, which was expected in hindsight. High correlation among IVs is problematic because it makes it difficult to isolate the effects of each IV on the dependent variable (DV), as altering one IV affects others.

To quantify this issue, we calculated the VIF for each IV. The VIF measures the increase in the variance of parameter estimates if an additional variable is added to the linear regression, indicating the level of multicollinearity among the IVs. Put simply, it evaluates the extent to which the variance of a predicted regression coefficient increases when the predictors exhibit correlation. If the VIF is greater than 5 or 10 (depending on the user's tolerance), it suggests high multicollinearity, and the variable should probably be removed from the model. Our analysis showed that many input variables were perfectly multicollinear, meaning some were essentially redundant. I experimented with removing these redundant variables one by one to improve model performance. While this approach did lead to some improvement, it was never significant. I tried various combinations, such as removing only Type 2 variables, only Type 1 variables, or a mix of both, but none resulted in substantial performance enhancements. The predictive power of the model was also unsatisfactory. Although it could estimate the order of magnitude correctly, it was often off by more than 0.1 in approximation ratio (AR), which is unacceptable. If we could develop a model that accurately predicts the (AR) based on certain IVs, we could reverse-engineer the sub-lists that lead to the best-performing median BSF approximation ratios. This would be a significant step towards our ultimate goal of identifying effective mapping schemes for the basis states.

In any case, this poor performance could indicate two things: either the model is inadequate (i.e., the underlying relationship is not actually linear), or the chosen IVs are not the most suitable (or, we simply do not have enough data-points; we were limited by the available hardware, so we cannot do much more regarding this). We suspect the latter might be true, although we're unsure which other variables to use. This circles back to the initial challenge of encoding the sub-lists' information in a memory-efficient and representative manner. Regarding the potential non-linearity of the model, we've considered using non-linear models, such as Support Vector Machine Regression (SVR), which exploits the kernel trick to fit non-linear data. However, we have set this idea aside for now¹⁴, as our current goal is to identify which IVs most influence the DV's value.

One alternative approach that came to mind is to perform Principal Component Analysis (PCA) followed by regression (Principal Component Regression – PCR). Although PCA also assumes a linear relationship, it could help us achieve a better model. By computing the Principal Components (PCs), we can indirectly understand which IVs have the greatest influence in explaining the variance of the input data. These influential IVs would be the ones we should focus on when choosing our basis states' partitioning. By examining the loadings of the PCs, we can determine which IVs constitute each PC.

¹³The metric employed is the mean squared error.

¹⁴This would be an interesting direction for future work.

Another advantage of PCA is that the obtained PCs are orthogonal (uncorrelated), thus eliminating the multicollinearity problem among the input variables. We can then attempt linear fitting (PCR) on the PCs against the DV. The main drawback is that we lose some interpretability of the input variables, as accurately interpreting the PCs can be challenging due to their nature as linear combinations of the original IVs.

Anyways, PCR using the 5 most significant principal components also yielded subpar results. An R^2 of 0.0287 was reported, with a 5-fold cross-validation loss of 0.0059 and a standard deviation of 0.0004. As one could already predict from the R^2 value, this model's predictive powers were also unsatisfactory. Nevertheless, we can still try to extract some information from the principal components to understand which independent variables are most influential in determining the dependent variable. Notably, the first 5 principal components (out of 47) account for 65.3% of the variance in the data:

```
Explained variance ratio:  
[0.46558461 0.07637042 0.04557191 0.03879498 0.02634421]
```

Listing 5.1: Explained variance ratio for the first 5 PCs (8-node graph).

Now, if we examine the loadings of the first and most significant principal component, we notice they are quite evenly distributed across the independent variables. This makes interpretation challenging, as we cannot easily identify one IV as being more important than the others. Here are the loadings of the first principal component for the 8-node graph: (Identically uniform distributions were also verified for the other PCs.)

```
PCA_1 = [-1.75323157e-01 -9.26451207e-02 -8.98377538e-02 -8.77605361e-02 -9.18316965e-02  
-8.58000439e-02 -1.03221966e-01 -9.24948614e-02 -1.11659776e-01 -1.75791034e-01  
-1.73978061e-01 -1.70204027e-01 -1.70707746e-01 -1.73770114e-01 -1.72400120e-01  
-1.73386482e-01 -1.71136657e-01 -1.88767128e-01 -1.82462202e-01 -1.81233077e-01  
-1.81277591e-01 -1.83509589e-01 -1.80911686e-01 -1.83449672e-01 -1.85536765e-01  
-1.87937749e-01 -1.85009049e-01 -1.87993729e-01 -1.86774034e-01 -1.87581036e-01  
-1.85531361e-01 -1.86124394e-01 -1.86528612e-01 -2.09742228e-03 -3.66652955e-03  
9.18121507e-04 -9.31248048e-03 -4.63173882e-03 -1.18673071e-03 1.43514268e-02  
-1.14790537e-02 -1.48687943e-01 -1.33654722e-01 -1.54547996e-01 -1.55191025e-01  
-5.69361380e-03 -1.50033651e-01]
```

Listing 5.2: Loadings of the first PC, for the 8-node graph.

Table 5.1: Cross-validation loss (mean square error) and R^2 values for the two fitted models, Multiple Linear Regression and PCR, in the context of the usual 8-node graph.

Model (8-node graph)	5-fold cross-validation loss		R^2
	Average (\bar{x})	Std. dev. (σ)	
Multiple Linear Regression	0.0061	0.0004	0.1023
PCR (5 Principal Components)	0.0059	0.0004	0.0287

As can be seen, there are no IVs whose impact can be neglected, making the interpretation of this principal component nearly impossible, which doesn't help our case much. Ideally, we would discover that a small number of independent variables have a significantly larger impact than the others. This

would enable us to concentrate on those variables when selecting our basis states' partitioning. Unfortunately, this is not the situation here. For the sake of completeness and organization, I compile the obtained results in Table 5.1.

Furthermore, after revisiting the regression analyses' coefficients, we noticed something curious: when not performing PCA before fitting¹⁵, we observed significantly higher coefficients for four specific input variables (on the order of magnitude 10^9 or 10^{10} , compared to $[10^{-4}, 10^{-2}]$ for all other variables). These IVs are (for the 8-node graph dataset):

1. Average Hamming weight (Type 1);
2. Parity Score (Type 1);
3. Average Hamming weight (Type 2);
4. Parity Score (Type 2).

Initially, it appeared that these variables contributed the most to the dependent variable's value (median BSF AR after training). However, we no longer believe that to be the case. Instead, we attribute this observation to high multicollinearity among the independent variables (IVs), which renders the model unstable.

In light of the above, we also tested Ridge regression (L_2 regularization) as an alternative to PCR. Ridge regression is useful when the input data has obvious multicollinearity, as it adds a penalty¹⁶ to the regression coefficients, reducing their variance and stabilizing the model. Using `alpha = 1.0` in `sklearn.linear_model.Ridge`, we obtained an R^2 value of 0.1023 and an average cross-validation loss of 0.0061, with a standard deviation of 0.0004. These results were identical to those from multiple linear regression (for the 8-node graph). However, with Ridge regression, no IVs had disproportionately large coefficients. They all ranged between 10^{-4} and 10^{-2} . This regularization added stability to the model by preventing coefficient divergence. Moreover, its predictive power appeared to be slightly better than that of regular linear regression for our current dataset, often resulting in errors smaller than 0.1 in approximation ratio (AR).

This analysis reinforced our belief that the previous 10^{10} -valued coefficients were due to model instabilities from high multicollinearity. Thus, Ridge regression offers a better model with more meaningful and interpretable coefficients. However, the confirmed uniform spread of coefficients makes it difficult to pinpoint a few IVs as the most impactful, which we had hoped to identify.

We have explored several techniques to find patterns or strategies for better mapping. However, the task is more complex than initially anticipated. Capturing all intricate patterns might require a more complex model, such as a neural network¹⁷ or another non-linear approach. We have decided to leave this direction for future work. Another concern is that iQAQE's performance may depend heavily on

¹⁵If PCA is applied before fitting (PCR), the linear regression coefficients remain stable within the range $[10^{-4}, 10^{-3}]$ in absolute value and do not diverge.

¹⁶The L_2 penalty term is appended to the end of the RSS (Residual Sum of Squares) function, represented as follows: $RSS_{L_2} = \sum_{i=1}^n (y_i - \hat{y}_i)^2 + \alpha \sum_j j = \mathbf{1}^T \mathbf{B}_j^2$, where \mathbf{B}_j denotes the OLS coefficients. The symbol α represents a tunable hyperparameter.

¹⁷We need to exercise caution when implementing a neural network, as it's quite easy to compromise the interpretability of the model.

the specific graph instance, complicating the generalization of a model for predicting optimal mapping strategies. Despite this, we believe it is still worth pursuing.

Chapter 6

Alternative Schemes

In this chapter, we will present alternative schemes that, although not entirely based on the iQAQE Framework, emerged from our research on these topics. These are different algorithms proposed to solve the MaxCut problem.

6.1 Parity-like QAOA

The next idea we will explore involves "reducing" QAOA's cost Hamiltonian to use fewer qubits than usual. To achieve this, we will utilize parity encodings, which allow us to compress the representation of N nodes into $n < N$ qubits (with $N = \mathcal{O}(n^k)$, for any k^1 , as long as $\binom{n}{k} \geq N$). In this scheme, QAOA's cost Hamiltonian, the usual

$$H_C = \frac{1}{2} \sum_{i < j} (1 - Z_i Z_j), \quad (6.1)$$

gets "unfolded" into:

$$H_C = \frac{1}{2} \sum_{i \leq k < l; i < j \leq l} (1 - Z_i Z_j Z_k Z_l), \quad (6.2)$$

for example, when performing quadratic compression with $k = 2$. (The sums are taken over the graph's edges, as is customary.) Each node is now associated with a tuple of $k = 2$ qubits, and the node's set/color is determined by the parity of those k qubits (even number of 1's: Set 0; odd number of 1's: Set 1). This criterion is used to obtain the final partition, after training, from the most frequently sampled bit-string. The terms with 4 Pauli- z operators (when $k = 2$) can be understood as a product of two terms with 2 Pauli- z operators. Each of these terms represents the set/color of the corresponding nodes for each edge in the graph. It is easy to verify that this product gives -1 when the individual nodes belong to different sets, as expected. Additionally, note that any value of k can be used, so long as we can encode all necessary nodes in that many qubits (Number of possible encodings = $\binom{n}{k} \implies N = \mathcal{O}(n^k)$). Then, n is taken as the smallest value that satisfies $\binom{n}{k} \geq N$, similar to the previous polynomial compression-based algorithms. Furthermore, this approach can be generalized to other values of k by "unfolding" each Z in the $Z_i Z_j$ terms of Eq. 6.1 into products of k Pauli- z 's, according to the specified encoding. A

¹This k is equivalent to the previously mentioned k in polynomial compression-based schemes.

minimal example is provided below. This was kindly produced by Bence Bakó. This example aims to give more intuition on how this method works in practice.

Suppose we are solving **MaxCut** on a 6-node graph using this Parity-like **QAOA** with $k = 2$, resulting in a cost Hamiltonian similar to the one described in Eq. 6.2. This is achieved by compressing the usual 6-qubit representation (**QAOA**) to use only 4 qubits ($\binom{4}{2} = 6 \geq 6$) with the following mapping:

- Node 0 → Qubits (0, 1),
- Node 1 → Qubits (0, 2),
- Node 2 → Qubits (0, 3),
- Node 3 → Qubits (1, 2),
- Node 4 → Qubits (1, 3),
- Node 5 → Qubits (2, 3).

The standard **QAOA** ansatz, adapted to this scheme, can be readily implemented. The corresponding 4-qubit operators $e^{-i\alpha Z_i Z_j Z_k Z_l / 2}$ can be implemented as illustrated in Figure 6.1. Subsequently, the expectation value of H_C serves as the cost function.

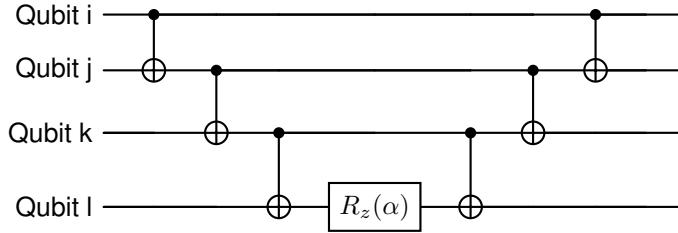


Figure 6.1: $e^{-i\alpha Z_i Z_j Z_k Z_l / 2}$ decomposition.

With this approach, we effectively reduced the number of qubits from 6 to 4, but at the cost of a deeper circuit. The mapping is flexible and does not have to be to an equal number of qubits: one can mix mappings to 1, 2, 3, or any number of qubits. The encoding aspect of this idea is essentially the same as in [18], restricted to the Z basis. This can also be seen as a type of parity encoding, since we multiply Pauli- z matrices, which have eigenvalues of ± 1 . This also shares some similarity with [16].

Now, I present the outcomes of the numerical simulation of Parity-like **QAOA** for the **MaxCut** problem on the standard 8-node graph. These results are depicted in Figure 6.2. It became apparent that, for $k = 2$, the specified encoding appeared incapable of generating the **MaxCut** partition (cf. Figure 6.2: The shaded blue region extends only up to 0.9). This sparked discussions regarding the generality of such an encoding. If the **MaxCut** partition cannot be constructed using this encoding, then achieving it through our ansatz becomes unattainable. This constitutes a significant flaw. Further insights arose from discussions with Prof. Zoltán Zimborás and others, emphasizing the presence of "cycles" in our encoding. These cycles impede our ability to encode all desired states, effectively tying certain nodes to the parity of others and automatically excluding certain partitions from our operational domain. Initially, the goal was to completely avoid such cycles. However, it has become clear that completely eliminating them is impossible. These cycles emerge as a result of confining the problem to a reduced subspace.

For instance, in a 6-node graph encoded in 4 qubits, there are 2 cycles: 0, 1, 3 and 3, 4, 5 (from the previous example encoding). The occurrence of these cycles is not incidental. Reducing from 6 qubits to

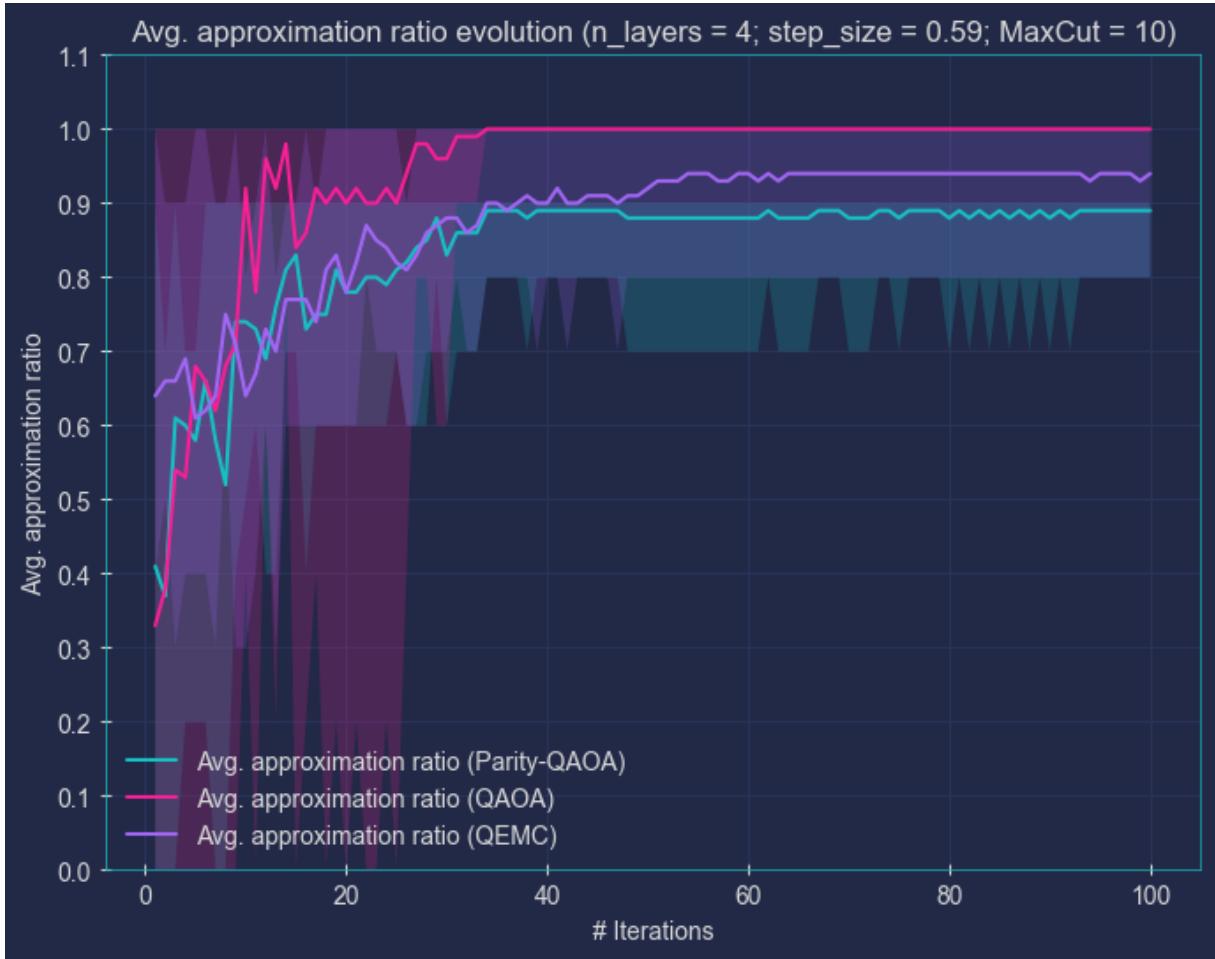


Figure 6.2: Results from the numerical simulation of Parity-like QAOA ($k = 2$) applied to the MaxCut problem on the standard 8-node graph.

4 imposes two constraints, represented by these cycles. Hence, it is theoretically impossible to encode all potential 6-node partitions in the given 4 qubits. Unless the encoding ensures the inclusion of the MaxCut partition within this reduced space, the scheme appears destined to fail in theory. Achieving such a feat would be remarkable, though, as it would effectively diminish the problem size, thereby enhancing the search process. However, I am uncertain about its feasibility. Mastering this encoding would likely necessitate specific graph instance knowledge. Given the variability in graph instances, devising a universal rule seems improbable.

Additionally, it was observed that the symmetry property, previously confirmed for standard QAOA², was only upheld when k was odd. In cases where k was even, inverting a bit-string produced the exact same partition, indicating that one possible MaxCut partition was overlooked by the algorithm.

Furthermore, while implementing the algorithm, it was found that despite the encoding leading to deeper ansätze, the outcomes were not as dire as initially anticipated. This was due to the simplification³ of terms such as $e^{Z_0 Z_1 Z_0 Z_2}$ to $e^{Z_1 Z_2}$, resulting in simpler circuits compared to those for 4 Pauli- z 's (which could be more than 4 depending on the value of k). It was also confirmed that the ordering of

²Wherein original and inverted bit-strings, e.g., 1001 and 0110, represent identical partitions (though with different labels) and are thus trained to achieve identical probabilities.

³Remember: $Z_0 Z_0 = I$.

terms in the complex exponential was irrelevant, allowing for flexibility in reordering terms to streamline implementation.

Inspired by [18], an effort was made to address the encoding issue stemming from the previously identified "cycles". These cycles were interpreted as manifestations of frustration effects, akin to those observed in antiferromagnetic spin chains. To mitigate these effects, a regularization term was introduced into the cost Hamiltonian. This term coerces the expectation values, necessary for computing the cost function, to approach zero, thereby reducing the impact of frustration. This approach entails modifying the read-out process to ensure coherence with the interpretation. Specifically, each node's color is determined by the sign of its associated expectation value, derived from the encoding. It could be said that experimental adjustments were implemented while preserving the same fundamental approach. Consequently, the cost function was revised to incorporate the regularization term: (Also, note that, in general, $\langle Z_j \rangle \langle Z_k \rangle \neq \langle Z_j Z_k \rangle$. This constitutes another deviation from the standard QAOA cost.)

$$\sum_{j < k} -\frac{1}{2} (1 - \langle Z_j \rangle \langle Z_k \rangle) + \mathcal{L}_{\text{reg}}. \quad (6.3)$$

Post-encoding, this is equivalent to:

$$\sum_{i \leq k < l; i < j \leq l} -\frac{1}{2} (1 - \langle Z_i Z_j \rangle \langle Z_k Z_l \rangle) + \mathcal{L}_{\text{reg}}, \quad (6.4)$$

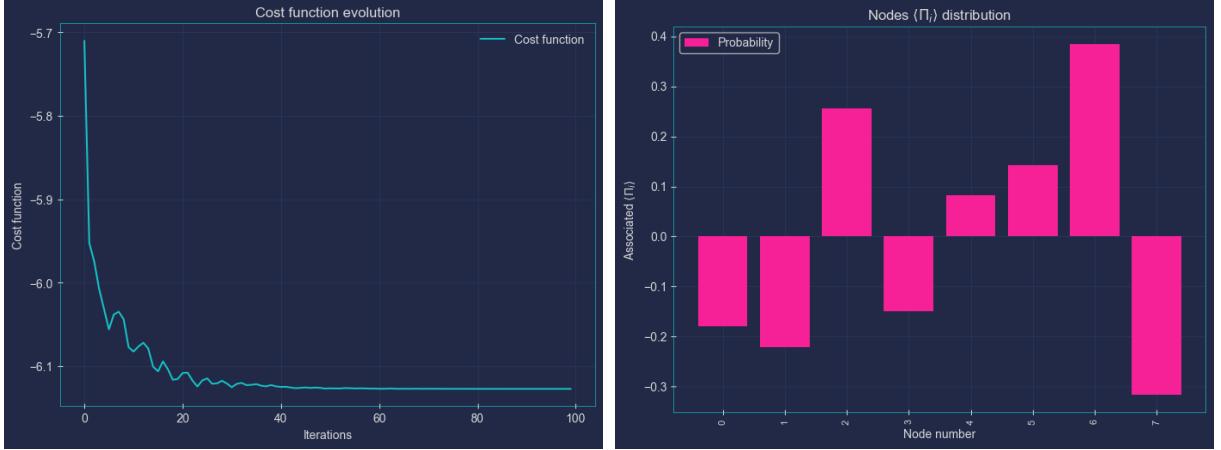
where:

$$\mathcal{L}_{\text{reg}} = \beta \nu \left[\frac{1}{N} \sum_{i \in [m]} \langle \Pi_i \rangle^2 \right]^2. \quad (6.5)$$

Here, ν represents an estimate of the MaxCut value, N denotes the total number of vertices in the graph, and β serves as a hyperparameter, adjusted on a case-by-case basis. Furthermore, $\langle \Pi_i \rangle$ represents the expectation value of the i -th correlator, where m denotes the total number of correlators, one for each graph node. These correlators correspond to the Pauli- z operators associated with each node. For instance, $Z_i Z_j$ for $k = 2$ denotes the correlator pertaining to the node whose color is determined by the parity of qubits i and j .

This approach successfully circumvented the limitations posed by the cycles, enabling the generation of partitions previously hindered by these constraints. Experimental validation on an 8-node graph demonstrated the efficacy of the method, yielding the MaxCut partition under certain parameter combinations⁴ (Figure 6.3). Notably, while these results appear promising, they largely replicate the findings of [18], thus falling short of originality. The key disparities lie in the choice of ansatz (QAOA vs. "brickwork" circuit ansatz) and the formulation of the cost function (hyperbolic tangents are utilized for non-linearity in [18]). Nonetheless, this endeavor contributed to a deeper comprehension of the referenced work and confirmed its reproducibility and effectiveness.

⁴For this demonstration, we employed 3 layers, set Adam's learning rate to 0.19, and conducted 100 iterations.



(a) Plot of the cost function values against the number of iterations during training.
(b) Plot displaying the expectation values of each node's associated correlator.

Figure 6.3: Training and results plots for the Parity-like Quantum Approximate Optimization Algorithm (QAOA), employing enforced regularization to mitigate frustration effects introduced by the aforementioned "cycles". The partition on the right (11010001) corresponds to the `MaxCut` partition of the usual 8-node graph ("cut" = 10).

6.2 Batch-based Oracle Coloring

Another idea we're interested in exploring is the following. Inspired by the previous concept that knowing the correlation between nodes is, trivially, enough to deduce the graph's partition into two colors, we can design a new scheme which makes use of this to compute the `MaxCut` partition. In short, this scheme could be implemented with a trainable Oracle, which would receive $k < n$ qubits, and return the colors of k nodes. Afterwards, we would run $\frac{n}{k}$ batches of this, to obtain all the n nodes' colors. Algorithmically (Algorithm 5), one could write:

Algorithm 5 Batch-based Oracle coloring

Fix a certain value $k < n$ (dimension of batches).

for $i = 1$ to $\frac{n}{k}$ **do**

Partition: $\{c_1^i, c_2^i, \dots, c_k^i\} \leftarrow O(|i\rangle)$

$i = i + 1$

end for

Subroutine: Utilize the sub-partitions $\{c_1^i, c_2^i, \dots, c_k^i\}$ to reconstruct the full graph's partition.

Although this scheme doesn't explicitly exploit correlations between nodes, a simple variation can be easily formulated in which the Oracle provides the correlation between the inputted nodes.

Returning to Algorithm 5, it's important to ensure that the input state going into the Oracle varies for different values of i , allowing the Oracle to distinguish between the batches. However, several nuances need to be addressed in this approach. One crucial consideration is how to train such an Oracle effectively. Training it separately on the sub-graph instances of the original graph may overlook edges connecting different sub-graphs. Thus, a clever approach is needed to incorporate this connectivity during training.

Another possibility is training multiple Oracles, each dedicated to a specific sub-graph instance. This idea emerged from discussions about what should be provided as input to the scheme. Should

it be simply $|i\rangle$ or something like $|i \bmod k\rangle$? The former approach might not utilize all 2^k available "entrance" basis states effectively, limiting the encoding of meaningful information to only the first $\frac{n}{k}$ states. Additionally, a constraint arises from the requirement that 2^k must be greater than or equal to $\frac{n}{k}$ for the scheme to be valid, which may not hold true for all combinations of n and k . This is relevant because we have k qubits entering and leaving the Oracle. If $2^k < \frac{n}{k}$, there aren't enough basis states to encode the total number of required batches, which would be problematic.

These complexities must be carefully considered in the design of the algorithm. However, we have decided to set aside this scheme for now. We found several papers, such as [43] and [15], that address this exact problem, providing numerous solutions with varying degrees of effectiveness. Notably, [15] (Zeqiao Zhou et al.), titled "*QAOA-in-QAOA: solving large-scale MaxCut problems on small quantum machines*", particularly captured our attention. It presents insightful approaches that closely align with our original intention of solving MaxCut by addressing sub-graphs individually.

Chapter 7

Conclusions

We've introduced a new framework, inspired by QAOA and QEMC, for seamlessly designing multiple distinct VQAs. This framework builds on QEMC's core components while integrating concepts from QAOA to leverage the strengths of both algorithms for improved performance. Tentatively named the iQAQE Framework, it opens the door to numerous unexplored and unknown VQAs, offering a variety of parameters to experiment with. It can be seen as a testbed for developing new VQAs. While this may not be revolutionary in itself, it has the potential to pave the way for groundbreaking results in the future.

7.1 Main Findings

The primary achievement of this work is the development of the iQAQE Framework itself. We believe this new framework has the potential to compete with current state-of-the-art algorithms, as we've demonstrated. It offers reduced qubit requirements compared to QAOA and greater resilience to statistical uncertainty associated with small shot numbers, compared to QEMC, i.e., the results don't significantly deteriorate when we use finite shot numbers.

We also introduced various heuristics for selecting the number of qubits, the cardinality of lists, and mappings. These heuristics have generally yielded results that are competitive with QEMC. In certain scenarios, such as ND-CNOT QEMC, we have surpassed the classical state-of-the-art (GW) for a reasonably large graph (32 nodes). This promising result highlights the potential of the iQAQE Framework.

Additionally, we proposed a scheme for determining the optimal mapping for a specific graph based on statistical properties of the mappings themselves, using a small machine learning model. Although the results were not entirely satisfactory, we believe this approach can inspire future research, building on this idea and our initial work.

7.2 Future Work

There are several promising directions for future research building on the current work. Key areas of exploration include improvements to the machine learning (ML) model, the development of new heuris-

tics, and testing on larger and more complex graphs. Let us take a closer look at these.

Enhancements to the Machine Learning Model: Exploring the use of neural networks could potentially yield better results. Investigating additional input variables and incorporating more data points will likely improve the model's accuracy. Another avenue worth exploring is the use of Support Vector Regression (SVR) as an alternative approach, for non-linear model fitting.

Development of New Heuristics and Exploration of Different Schemes: Developing new and creative heuristics is crucial. For example, testing different cardinalities for different nodes may reveal an optimal configuration that has not yet been identified. This exploration could uncover a "sweet spot" that enhances the performance of the algorithm. Moreover, in the context of QEMC's encoding, experimenting with various methods to derive node probabilities from the basis states' probabilities could be interesting. Exploring alternatives to the current "sum and normalize" method may yield better results. Additionally, trying different (problem-inspired) ansätze and cost functions could lead to further improvements. For instance, developing a mapping-inspired ansatz that depends on the mapping itself may offer novel insights and enhance performance. Additionally, expanding on the alternative schemes (Chapter 6) mentioned in this work could provide valuable additional avenues of research. Finally, another potential research direction could involve extending the iQAQE Framework to address other combinatorial optimization problems, such as Max- k -Cut.

Testing on Larger Graphs and Scalability Analysis: Expanding the testing of the framework to larger graphs and a greater number of instances will enhance the validation of its effectiveness. Conducting more extensive and rigorous testing will provide a better understanding of the framework's capabilities and limitations. Understanding how the framework scales with much larger graphs, potentially containing thousands of nodes, is a crucial area of investigation. This could reveal significant performance advantages over QEMC, but my current computational resources are insufficient to test this effectively at present.

Real-World Quantum Computer Implementation: Running the algorithms on an actual quantum computer is a critical step. This would enable us to assess the framework's performance in real-world scenarios.

Further Testing and Refinement: Lastly, further testing of some of the proposed heuristics is necessary. This includes experimenting with different shot numbers for some of our heuristics to determine their impact on performance. Moreover, there is potential for these heuristics to be refined and improved, leading to even better performance of the framework.

Concluding, by addressing these areas, future research can build on the foundation laid by this work, potentially leading to significant advancements in the field of Variational Quantum Algorithms (VQAs).

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Appendix A

Best-so-far correction

Here are the revised plots illustrating the median best-so-far metric for the numerous heuristics tested throughout this study.

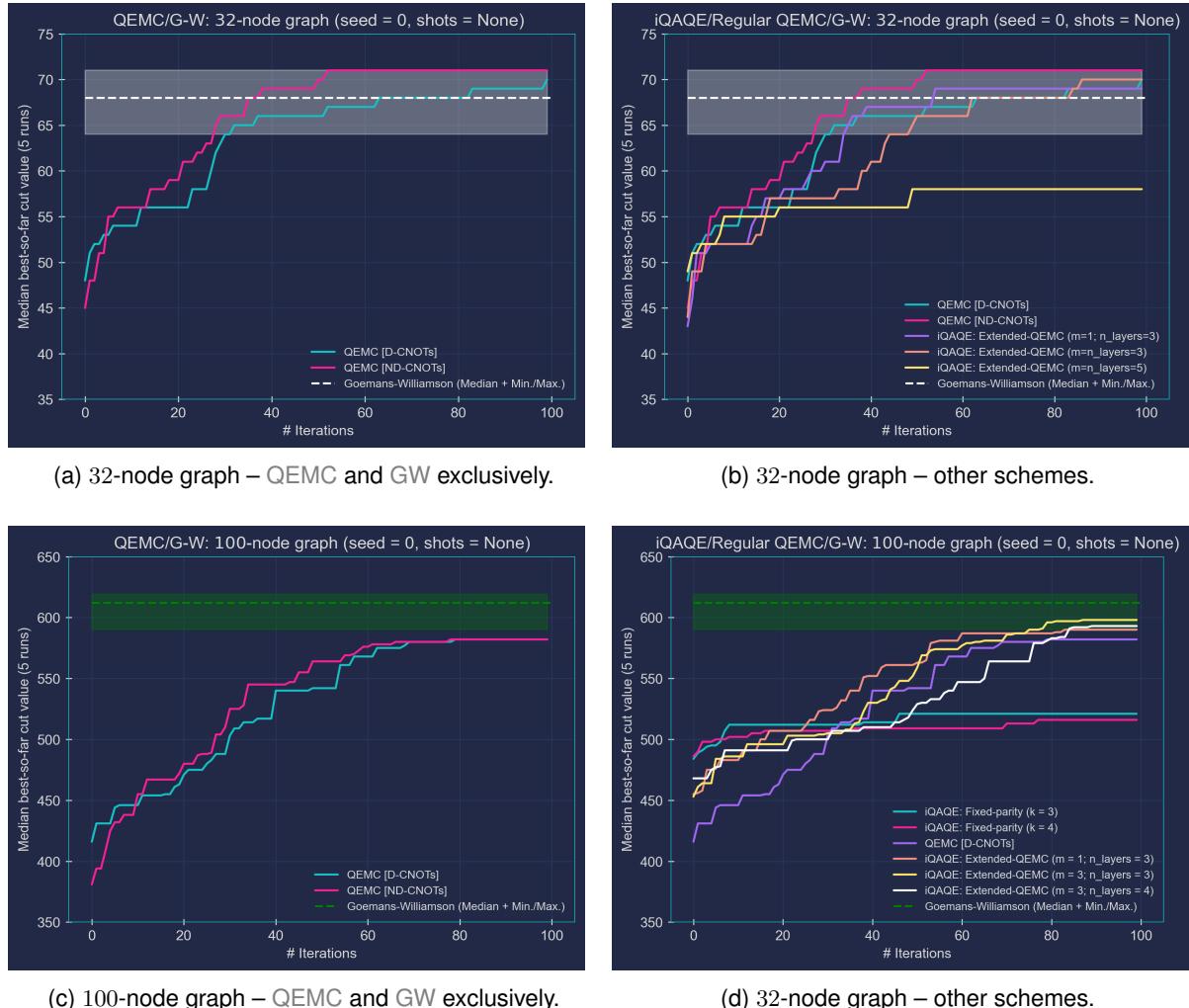


Figure A.1: Revised results using the median best-so-far metric for the 32 and 100-node graphs.

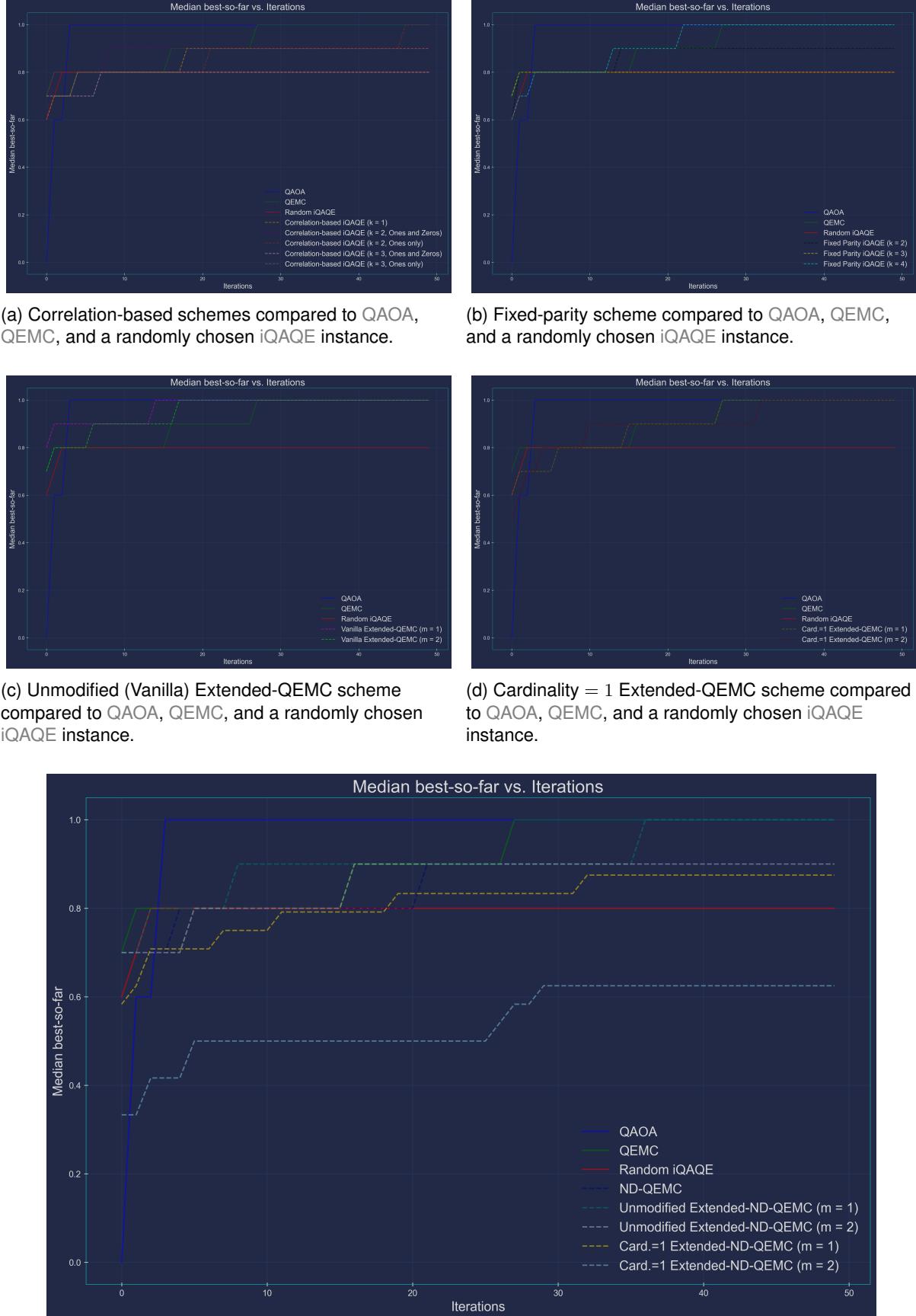


Figure A.2: Revised results using the median best-so-far AR metric for the 8-node graph.