Contents

1	Introduction	1
2	A semidefinite relaxation for the Max cut problem	3
3	A rank-two relaxation	9
4	A heuristic algorithm for MCP	18
5	Solving BQP using MCP	23
6	Experimentation	24
	6.1 Experimental Setup	24
	6.2 Instances with known optimal value	25
	6.3 Instances without known optimal value	26
	6.4 Results	26

1 Introduction

Throughout this text we consider the maximum cut problem, which is shown to be NP-hard by [GJS74]. To tackle this problem approximation algorithms and heuristic algorithms have been proposed. We will develop the theory for the heuristic presented by Burer, Monteiro and Zhang in [BMZ02], which we will call Burer heuristic.

The text divides into two conceptual parts: The first part can be considered the theoretical part, in which we try to develop the theory behind the Burer heuristic. The second part is where we present the experimental results.

In section 2, we give an overview of an approximation algorithm proposed by Goemans and Williamson. This will be the basis for the Burer heuristic. That includes introducing a semidefinite programming problem relaxation for the maximum cut problem and showing how to derive a cut from a set of points on the unit sphere. We will discuss the performance guarantee in detail. However, the proof of polynomial runtime will not be treated here, we refer to [KV18].

In section 3, we lay the theoretical foundation for the Burer heuristic (described in section 4). In particular we study the special case of the SDP relaxation introduced in section 2 for dimension two in polar coordinates. This leads to an unconstrained nonconvex optimisation problem, which has advantages and disadvantages. On the upside by using this relaxation the number of variables is not increased, i.e. the number of variables remains the number of vertices of the graph. This implies good scalability to large instances. However as the function is nonconvex, we can not expect to solve this relaxation optimally. Therefore we face a trade-off between computational runtime and a theoretical guarantee, see [BMZ02, p. 506]. We will conduct experiments on Mallach instance library to investigate this trade-off. This yields data demonstrating the effectiveness.

In section 4 we give a detailed exposition of the Burer heuristic. The heuristic combines the rank-two relaxation and Goemans-Williamson-type cuts. Given a set of points on the unit circle, we will provide a complete description of how to generate a best possible Goemans-Williamson-type cut in a deterministic manner. Furthermore we will describe an algorithm to improve a cut value by means of trial and error, where the termination criterion is given

by the number of consecutive unsuccessful attempts to improve the cut value. Lastly we will end the section with algorithm of the Burer heuristic. This is the algorithm on which the experimental data is based upon.

In section 5 we show that instances of the binary quadratic programming problem can be transformed to instances of the maximum cut problem.

In the second part we will go into computational results. We implemented the heuristic (distillation from ..) and ran it on the instance library provided by Mallach. The results described in section, gather more evidence, that the rank-two relaxation proposed by Burer et al. is highly effective. In the results part we will

2 A semidefinite relaxation for the Max cut problem

Throughout this text we consider undirected graphs G=(V,E) with $V=[1:n]:=\{1,\ldots,n\}$ and $E\subseteq\{(i,j)\in V\times V\mid i< j\}$. Note that we adopt this notation to stay coherent with [BMZ02] on which this text is based. By the condition i< j there can only be one edge between the vertices i and j, thus we could have just as well considered $E\subseteq \{\{i,j\}\mid i,j\in V:i\neq j\}$. Let the edge weights $w_{ij}=w_{ji}$ be given such that $w_{ij}=0$ if $(i,j)\notin E$. We then have $w_{ii}=0$ for all $i\in V$. We are interested in studying the weights of cuts in the graph G. Given a bipartition of (S,\overline{S}) of V, a cut is the set $\{e\in E\mid |S\cap e|=1 \text{ and } |\overline{S}\cap e|=1\}$. Clearly, a cut is not uniquely defined by (S,\overline{S}) , as (\overline{S},S) generates the same cut. Furthermore a cut is uniquely defined by a set X, as the bipartition is uniquely given by $(X,V\setminus X)$. The corresponding weight is obtained by summing the weights of the edges in the cut. The task of finding a cut of maximum weight is called the maximum cut problem, abbreviated by MCP, see [KV18]:

Maximum Weight Cut Problem

Instance: An undirected weighted graph G.

Task: Find a cut in G with maximum total weight.

It is well known that the MCP is NP-hard, see [GJS74]. Therefore there has been a lot of work on approximation algorithms and heuristics tackling the MCP. A groundbreaking contribution is the approximation algorithm described by Goemans and Williamson in [GW95]. In fact the Burer heuristic is based on this approximation algorithm. Therefore we will spend the rest of this section on developing the theory to understand the algorithm by Goemans Williamson. We will focus on the main ideas and not go into, for example proving the polynomial runtime.

Following the description given by [Vaz03, p. 268 ff], we can motivate the formulation of the maximum cut problem as a binary quadratic program:

We give a for our purposes sufficient definition of binary quadratic program.

Definition 2.1 (Binary quadratic program) Let B be either $\{0,1\}$ or $\{-1,1\}$, and $Q \in \mathbb{R}^{n \times n}$. A (unrestricted) binary quadratic program is of the form:

minimize
$$x^T Q x$$

subject to $x_i \in B \quad \forall i \in [1:n]$ (1)

To every vertex i we assign a indicator variable x_i which is constrained to be in $\{-1,1\}$. This allows us to define the cut-defining partition (S,\overline{S}) by $S:=\{i\mid x_i=1\}$ and $\overline{S}=\{i\mid x_i=-1\}$. For any edge $\{i,j\}$ in the cut we have $i\in S$ and $j\in \overline{S}$ or vice versa. Thus we have $x_ix_j=-1$ for every edge $\{i,j\}$ in the cut. On the other hand, for every edge $\{i,j\}$ that is not in the cut we have $i,j\in S$ or $i,j\in \overline{S}$, implying $x_ix_j=1$.

$$\frac{1}{2}(1 - x_i x_j) = \begin{cases} 1 & \{i, j\} \text{ in the cut defined by } S \\ 0 & \text{otherwise} \end{cases}$$

This explains, why we can write the Maximum Cut Problem, abbreviated with MCP, as the following quadratic program.

maximize
$$\frac{1}{2} \sum_{1 \le i < j \le n} w_{ij} (1 - x_i x_j)$$

subject to
$$x_i \in \{-1, 1\} \quad \forall i \in [1 : n]$$
 (2)

Alternatively, we can solve the following binary quadratic program as Lemma 2.2. This will prove useful as we progress in the section.

minimize
$$\sum_{1 \le i < j \le n} w_{ij} x_i x_j$$

subject to $|x_i| = 1 \quad \forall i \in [1:n]$ (3)

For the sake of convenience we will introduce some notation to denote sums, see [Aig07, p. 41]. Say we want to sum the even numbers between 0 and 100, we can write $\sum_{k=0}^{100} k [k \text{ even}]$. The expression in brackets, which is to be multiplied, means that

$$[k \text{ has property E}] = \begin{cases} 1 & \text{if } k \text{ satisfies property E} \\ 0 & \text{otherwise} \end{cases}$$

This expresses the same sum as

$$\sum_{\substack{k=0\\k \text{ even}}}^{100} k \quad \text{or} \quad \sum_{k=1}^{50} 2k$$

Lemma 2.2 The solutions of 2 and 3 coincide.

Proof. We first observe that the constraints of 2 and 3 coincide so it suffices to show that the target function of the BQP is minimised if and only if the target function of the MCP is maximised. The following computation uses that $x_i x_j \in \{-1, 1\}$ for all $i, j \in [1:n]$:

$$\begin{split} & \sum_{1 \leq i < j \leq n} w_{ij} x_i x_j = \sum_{1 \leq i < j \leq n} w_{ij} x_i x_j \left[x_i x_j = 1 \right] + \sum_{1 \leq i < j \leq n} w_{ij} x_i x_j \left[x_i x_j = -1 \right] \\ & = \sum_{1 \leq i < j \leq n} w_{ij} \left[x_i x_j = 1 \right] - \sum_{1 \leq i < j \leq n} w_{ij} \left[x_i x_j = -1 \right] = \sum_{1 \leq i < j \leq n} w_{ij} - 2 \sum_{1 \leq i < j \leq n} w_{ij} \left[x_i x_j = -1 \right] \end{split}$$

Using $\sum_{1 \leq i < j \leq n} w_{ij}$ is constant and $\sum_{1 \leq i < j \leq n} w_{ij} (1 - x_i x_j) = 2 \sum_{1 \leq i < j \leq n} w_{ij} [x_i x_j = -1]$. We can thus conclude, that maximising $\frac{1}{2} \sum_{1 \leq i < j \leq n} w_{ij} (1 - x_i x_j)$ (and thus $\sum_{1 \leq i < j \leq n} w_{ij} (1 - x_i x_j)$) minimises $\sum_{1 \leq i < j \leq n} w_{ij} x_i x_j$ and vice versa.

Let us introduce some notation allowing for more concise formulation. Consider $W = (w_{ij})_{1 \leq i,j \leq n}, X = (x_{ij})_{1 \leq i,j \leq n} \in \mathbb{R}^{n \times n}$. We denote the sum of their entrywise product by $W \bullet X = \sum_{i=1}^{n} \sum_{j=1}^{n} w_{ij} x_{ij}$ and vector of diagonal elements by diag $(X) = (x_{11}, \ldots, x_{nn})^T \in \mathbb{R}^n$. Further we denote the vector of all ones by e and denote a symmetric positive semidefinite matrix X by $X \succeq 0$.

The following Lemma shows a useful reformulation of 3, which will form the basis of our construction of the Goemans Williamson algorithm.

Lemma 2.3 The BQP 3 can be rewritten into the following matrix optimisation program

minimize
$$\frac{1}{2}W \bullet X$$

subject to $\operatorname{diag}(X) = e$,
 $\operatorname{rank}(X) = 1$,
 $X \succeq 0$ (4)

Proof. First of all, we are going to show that a matrix X satisfies the conditions of 4 if and only if there exists a $x \in \{-1,1\}^n$ such that $X = xx^T$:

As rank(X) = 1 implies that the columns of the matrix are scalar multiples of each other, there exist $x, y \in \mathbb{R}^n \setminus 0$ such that $X = xy^T$.

Claim: The vectors x and y are necessarily linearly dependent.

Assume otherwise. Then the Cauchy-Schwarz inequality (CS) is strict, i.e. $|\langle x,y\rangle| < \|x\| \|y\|$. Furthermore the point $\frac{x+y}{2}$ is linearly independent from x and y. By symmetry, we only show this for y: Given x,y linearly independent. Assume there exists $\lambda \in \mathbb{R}$ such that $\frac{x+y}{2} = \lambda y \Leftrightarrow x = (2\lambda - 1)y$ contradicting x,y linearly independent. Thus $|\langle x,\frac{x+y}{2}\rangle| < \|x\| \|\frac{x+y}{2}\|$ and $|\langle y,\frac{x+y}{2}\rangle| < \|y\| \|\frac{x+y}{2}\|$. Now we can show the claim using the following preparatory calculations. We consider the vector $v := y - \frac{\langle y,\frac{x+y}{2}\rangle}{\|\frac{x+y}{2}\|^2} \frac{x+y}{2}$ and calculate:

$$\langle y, v \rangle = \|y\|^2 - \frac{\langle y, \frac{x+y}{2} \rangle}{\|\frac{x+y}{2}\|^2} \left\langle y, \frac{x+y}{2} \right\rangle \stackrel{\text{CS}}{>} \|y\|^2 - \frac{1}{\|\frac{x+y}{2}\|^2} \|y\|^2 \left\| \frac{x+y}{2} \right\|^2 = 0$$

Furthermore we have:

$$\frac{1}{2} \left(\langle x, v \rangle + \langle y, v \rangle \right) = \left\langle v, \frac{x+y}{2} \right\rangle = \left\langle y, \frac{x+y}{2} \right\rangle - \frac{\left\langle y, \frac{x+y}{2} \right\rangle}{\left\| \frac{x+y}{2} \right\|^2} \left\langle \frac{x+y}{2}, \frac{x+y}{2} \right\rangle \\
= \left\langle y, \frac{x+y}{2} \right\rangle - \left\langle y, \frac{x+y}{2} \right\rangle = 0$$

Thus we have $\langle x, v \rangle + \langle y, v \rangle = 0$ and using the first calculation we get $\langle x, v \rangle < 0$ and $\langle y, v \rangle > 0$. The claim follows by observing that $v^T X v = v^T x y^T v < 0$ contradicting X positive semidefinite.

As x and y are linearly dependent, there exist an $\lambda \in \mathbb{R}$ such that $X = \lambda x x^T$ and $X_{ii} = \lambda x_i^2$ for all $i \in [1:n]$. Using the assumption $\operatorname{diag}(X) = e$ yields $1 = \lambda x_i^2$ for all $i \in [1:n]$. This implies $\lambda \neq 0$ and we can rewrite $x_i^2 = \frac{1}{\lambda}$ for all $i \in [1:n]$. In other words we have $x_i \in \left\{-\frac{1}{\sqrt{\lambda}}, \frac{1}{\sqrt{\lambda}}\right\}$ for all $i \in [1:n]$ and setting $\widetilde{x} := \sqrt{\lambda}x$ yields $|\widetilde{x}_i| = 1$ for all $i \in [1:n]$ and $X = \widetilde{x}\widetilde{x}^T$.

For the other direction: Let $x \in \{-1,1\}^n$ and set $X = xx^T$. We clearly have rank(X) = 1, also $x_i \in \{-1,1\}$ for all $i \in [1:n]$ implies $x_ix_i = 1$ for all $i \in [1:n]$, which shows $\operatorname{diag}(X) = e$. That X is positive semidefinite follows from:

$$y^T X y = y^T x x^T y = |(x, y)|^2 \ge 0 \quad \forall y \in \mathbb{R}^n$$

Finally, we conclude the proof by showing that the objective functions coincide. Let $X = xx^T$, with $x \in \{-1, 1\}$:

$$\frac{1}{2}W \bullet X = \frac{1}{2} \sum_{i,j=1}^{n} w_{ij} x_i x_j \stackrel{w_{ij}=0}{=} \frac{1}{2} \sum_{1 \le i \le j \le n} w_{ij} x_i x_j$$

Remark In the previous proof, we showed that for linearly independent $x,y \in \mathbb{R}^n$ the matrix xy^T is not positive semidefinite, by an explicit construction. We will explain the intuition behind the construction: As x and y are different, by the geometric Hahn-Banach Theorem, there exists a hyperplane separating the points. However the points being different is not to sufficient to show that there exists a separating hyperplane containing the origin. This is where the linear independence is necessary. As the construction shows, there exists a hyperplane, generated by v, that contains the origin. Containing the origin is critical as this ensures that $x^Tv, y^Tv \neq 0$ and that x^Tv and y^Tv have different signs. Only then we can deduce $v^Txy^Tv < 0$.

In order to describe the algorithm by Goemans and Williamson [GW95] we need to define what a semidefinite program is.

Definition 2.4 (Semidefinite program [Vaz03, p.258]) Let $C, D_1, \ldots, D_k \in \mathbb{R}^n$ symmetric and $d_1, \ldots, d_k \in \mathbb{R}$. The general semidefinite programming problem, abbreviated to SDP, is given by

maximize
$$C \bullet X$$

subject to $D_i \bullet X = d_i \quad \forall i \in [1:k],$
 $X \succ 0$ (5)

The following Theorem guarantees that SDP can be approximately solved in polynomial time. Showing this is a key point in showing that the algorithm presented by Goemans and Williamson has polynomial runtime. As we will not use SDP for the heuristic refer to [Vaz03, p. 258 ff] or [KV18, Theorem 16.10] for a proof.

Theorem 2.5 ([Vaz03, p. 259]) For any $\varepsilon > 0$ semidefinite programs can be solved up within an additive error of ε , in time polynomial in n and $\log(1/\varepsilon)$

By setting $C = -\frac{1}{2}W$ and $D_i = e_i e_i^T$, $d_i = 1$ for all $i \in [1:n]$ we see that we can relax the matrix optimisation program 4 into a SDP by simply dropping the rank constraint:

minimize
$$\frac{1}{2}W \bullet X$$

subject to $\operatorname{diag}(X) = e$, (6)
 $X \succ 0$

The first big idea is to relax BQP 3 in this ingenious way. We replace the unit scalars x_i by unit vectors $v_i \in \mathbb{R}^d$ (where the dimension d > 1 can be chosen freely). Furthermore the product $x_i x_j$ can be interpreted as the one dimensional scalar product, we therefore replace it by $v_i^T v_j$, which is the standard scalar product in \mathbb{R}^d . Noting $v_i \in \mathbb{R}^d$ the relaxation writes as follows:

minimize
$$\sum_{1 \le i < j \le n} w_{ij} v_i^T v_j$$

subject to $\|v_i\|_d = 1 \quad \forall i \in [1:n]$ (7)

Remark 1. We are going to show that 7 is a relaxation of 3:

Let $x \in \{-1, 1\}^n$ be a solution to 3. Let $e_1 \in \mathbb{R}^d$ be the first standard basis vector and set $v_i = x_i e_1 \in \mathbb{R}^d$ for all $i \in [1:n]$. Then the v_i 's form a feasible solution as for every $i \in [1:n]$ we have $||v_i|| = |x_i| ||e_1||_d = 1$. Furthermore values of the objective functions coincide by $v_i^T v_j = x_i x_j e_1^T e_1 = x_i^T x_j$ for all $i, j \in [1:n]$.

2. Observe that for a feasible (and optimal) solution v_1, \ldots, v_n each point v_i is located on the unit sphere \mathbb{S}^{d-1} representing the vertex i in G.

Aiming to apply Lemma 2.3, we can change the variables $x_i x_j$ to $v_i^T v_j$. Therefore X_{ij} is given by $v_i^T v_j$ instead of $x_i^T x_j$.

Lemma 2.6 ([Vaz03, p. 259f]) We can rewrite 7 into the following SDP

minimize
$$\frac{1}{2}W \bullet X$$

subject to $\operatorname{diag}(X) = e$, (8)
 $X \succ 0$

Proof. Throughout this proof we denote the i-th column of a matrix L by L_i .

Let X be a feasible solution of 7. As X is positive semidefinite there exists a $L \in \mathbb{R}^{n \times n}$ such that $X = L^T L$. Thus we have $X_{ij} = L_i^T L_j$ for all $i, j \in [1:n]$. For $i \in [1:n]$ we set $v_i := L_i$. As X is a feasible solution to 8 we have $1 = X_{ii} = L_i^T L_i = ||v_i||_n^2$, which shows $||v_i||_n = 1$ for all $i \in [1:n]$. Thus the v_i 's constitute a feasible solution to 7 of the same objective function value:

$$\frac{1}{2} \sum_{i,j=1}^{n} w_{ij} v_i^T v_j = \frac{1}{2} W \bullet L^T L = \frac{1}{2} W \bullet X.$$

Given a feasible solution v_1,\ldots,v_n to 7 we define the matrix $L\in\mathbb{R}^{n\times n}$ by $L_i:=v_i$. Now set $X=L^TL$, we then have $X_{ij}=v_i^Tv_j$. As the v_i 's are a feasible solution we have $X_{ii}=\|v_i\|_n^2=1$ for $i\in[1:n]$. Furthermore X is positive semidefinite since $y^TXy=\|Ly\|_n^2\geq 0$ for all $y\in\mathbb{R}^n$, which shows that X is a feasible solution of 8. By construction we have that the two feasible solutions have the same value:

$$\frac{1}{2}W \bullet X = \frac{1}{2} \sum_{i,j=1}^{n} w_{ij} X_{ij} = \frac{1}{2} \sum_{i,j=1}^{n} w_{ij} v_i^T v_j$$

Remark As this is essential, we reiterate how to retrieve a solution of 7 from a solution X of 8: As X is a positive semidefinite matrix, using the Cholesky decomposition we compute a matrix $L \in \mathbb{R}^{n \times n}$ such that $X = L^T L$. The columns of L are a solution to 7.

The algorithm presented by Goemans and Williamson can be stated as follows [KV18, p.424]: We emphasise that nonnegative edge weights are required. This requirement is essential in the proof of Lemma 2.10.

As we will not solve semidefinite programs in the heuristic presented later in the text, we will for simplicity's sake, assume that we can find an optimal solution to 8. That means that we have an optimal solution to 7. This inaccuracy can be absorbed into the approximation factor; as stated in [Vaz03, p. 260]. A detailed account of how to handle nearly optimal solutions is given in [KV18]. Until the end of the section we will closely follow the description provided in [Vaz03, p. 260 ff]. In order to approximate the MCP we need to relate this optimum solution to 7 to a cut. As the matrix whose columns are an optimum solution to 7 is in general not of rank 1 we are faced with a rounding problem. The question boils down to the following: How can we derive a good cut from an optimum solution to 7?

Algorithm 1 Goemans Williamson algorithm

Input: A graph with nonnegative edge weights

Output: A cut given by the set S

- 1: Find an approximately optimal solution X to 7
- 2: Find vectors $y_1, ..., y_n \in S^d$ such that $y_i^T y_j = x_{ij}$ for all $i, j \in [1:n]$ (using Cholesky decomposition)
- 3: Choose a random point $r \in \mathbb{S}^d$
- 4: Set $S := \{i \in [1:n] \mid r^T y_i \ge 0\}$
- 5: **return** $(S, V \setminus S)$

The second big idea by Goemans and Williamson, answers the question. As the heuristic presented later on in the text is based on this idea we will describe it in detail.

Let v_1, \ldots, v_n be an optimal solution and denote the optimum value by OPT. We have $v_i \in \mathbb{S}^d$ for all $i \in [1:n]$ and for all $i, j \in [1:n]$ we have $\cos(\theta_{ij}) = v_i^T v_j$. Here θ_{ij} denotes the angle between v_i and v_j . The contribution of v_i and v_j to the value of OPT is, as we simply read off from the definition of the objective function:

$$\frac{1}{2}w_{ij}\left(1-\cos(\theta_{ij})\right)$$

Therefore the contribution of v_i and v_j is the greater, the closer θ_{ij} is to π (, while considering the same w_{ij}). Geometrically this means, that diametrically opposed points have the highest contribution. This means that roughly speaking we want to separate diametrically opposed points. The cuts introduced in the next Definition achieve just that.

Definition 2.7 (Goemans-Williamson cut) A Goemans-Williamson cut for a direction $r \in \mathbb{S}^n$ is the cut given by (S, \overline{S}) , where we set $S := \{i \in [1:n] \mid v_i^T r \geq 0\}$ and $\overline{S} := \{i \in [1:n] \mid v_i^T r < 0\}$.

As the algorithm is based on rounding the values according to a uniformly chosen Goemans-Williamson cut we are naturally interested in knowing what the odds of separating two points are.

Lemma 2.8 Let $n \geq 2$ and $v_i, v_j \in \mathbb{S}^n$ different. We divide the unit sphere into two hemispheres by a hyperplane through the origin. The probability that v_i and v_j are separated is given by $P[v_i \text{ and } v_j \text{ are separated}] = \frac{\theta_{ij}}{\pi}$

Proof. Consider the plane that contains $0, v_i$ and v_j , given by span (v_i, v_j) . A hyperplane (through the origin) separates the two points if and only if its projection onto the plane intersects the arc connecting v_i, v_j on the sphere. As the hyperplane is chosen uniformly and the angle between v_i and v_j is given by θ_{ij} the probability, that the hyperplane separates the two points is $\frac{\theta_{ij}}{\pi}$.

Lemma 2.9 The expression $\alpha := \min_{0 < \theta \le \pi} \frac{2}{\pi} \frac{\theta}{1 - \cos(\theta)}$ is well defined and $\alpha > 0.87856$ holds.

Proof. The enumerator has a simple zero and the denominator a zero of order two. As the terms are nonnegative we have $\lim_{\theta \to 0} = \frac{2}{\pi} \frac{\theta}{1 - \cos(\theta)} = \infty$. Thus we have a continuous function on an interval and the As the denominator and the numerator have simple zeroes in zero,

the function can be continuously extended in 0. Thus we have a continuous function on an interval and the minimum will be attained. We take the derivative and get

$$\frac{\partial}{\partial \theta} \frac{2}{\pi} \frac{\theta}{1 - \cos(\theta)} = \frac{2}{\pi} \frac{1 - \cos(\theta) - \theta \sin(\theta)}{(1 - \cos(\theta))^2}$$

Thus we have a critical point if and only if $1 = \cos(\theta) + \theta \sin(\theta)$ In $(0, \pi]$ there is only one solutions, namely 2.33112237041442. By the first part this has to be a local minimum with value 0.878567205784851604. By $\frac{2}{\pi} \frac{\pi}{1-\cos(\pi)} = 2$ the local minimum is a global minimum and $\alpha > 0.87856$

Lemma 2.10 The expected value of the weight of a Goemans-Williamson cut is greater or equal to $\alpha \cdot \text{OPT}$, where OPT denotes the value of an optimal solution to 7.

Proof. Using Lemma 2.9 we have for all $\theta \in (0, \pi]$:

$$\frac{\theta}{\pi} \ge \alpha \frac{1 - \cos(\theta)}{2}$$

We denote the weight corresponding to the edge (i, j) by w_{ij} and set $w_{ij} = 0$ otherwise. The expected value of a Goemans-Williamson cut is given by

 $\sum_{1 \leq i \leq j \leq n} w_{ij} P[v_i \text{ and } v_j \text{ are separated}], \text{ together with Lemma 2.8 we can compute:}$

$$\sum_{1 \leq i < j \leq n} w_{ij} P\left[v_i \text{ and } v_j \text{ are separated}\right] \geq \sum_{1 \leq i < j \leq n} w_{ij} \frac{\theta_{ij}}{\pi}$$

$$\geq \frac{\alpha}{2} \sum_{1 \leq i < j \leq n} w_{ij} \left(1 - \cos(\theta_{ij})\right)$$

$$= \alpha \cdot \text{OPT}$$

Remark In particular Theorem 2.10 implies that there exists a Goemans-Williamson cut with value $0.8785 \cdot \text{OPT}$.

Therefore we have discussed all the, for our purposes, important aspects and can state, as in [KV18, Theorem 16.12]:

Theorem 2.11 The Goemans-Williamson Max-Cut-Algorithm returns a set S for which the expected value of the associated cut is at least 0.878 times the maximum possible value.

3 A rank-two relaxation

In this section we will investigate relaxation 7 in dimension two. Instead of using Cartesian coordinates we will use polar coordinates. We will investigate the implications of these changes, following [BMZ02, Sec. 3]. We will introduce the notion of angular representation of cuts, which are certain vectors that correspond to cuts. The section will culminate in giving a classification of cuts as stationary points of the objective function of the relaxation.

Using polar coordinates we can represent any point on the unit sphere by their angle. In other words for every $v \in S^1$, there exists a $\theta \in \mathbb{R}$ such that

$$v = \begin{pmatrix} \cos(\theta) \\ \sin(\theta) \end{pmatrix}.$$

Thus we can represent n points v_1, \ldots, v_n on the unit sphere, by a vector $\theta = (\theta_1, \ldots, \theta_n) \in \mathbb{R}^n$, such that the *i*-th vector corresponds to the *i*-th coordinate:

$$v_i = \begin{pmatrix} \cos(\theta_i) \\ \sin(\theta_i) \end{pmatrix}$$
 for all $i \in [1:n]$ (9)

By using the addition Theorem the inner product simplifies to:

$$v_i^T v_j = \cos(\theta_i)\cos(\theta_j) + \sin(\theta_i)\sin(\theta_j) = \cos(\theta_i - \theta_j)$$
(10)

Next we define the following map:

$$T: \mathbb{R}^n \to \mathbb{R}^{n \times n}, \ \theta \mapsto (\theta_i - \theta_j)_{\substack{i \in [1:n] \\ j \in [1:n]}}$$

For every $\theta \in \mathbb{R}^n$ the resulting matrix $T(\theta)$ is skew-symmetric, since for all $i, j \in [1:n]$

$$T(\theta)_{ij} = \theta_i - \theta_j = -(\theta_j - \theta_i) = -T(\theta)_{ii}$$

Throughout the text, the application of a scalar function onto a matrix, is to be understood as entrywise application of the scalar function. We define the following function, which will turn out to be of central importance:

$$f: \mathbb{R}^n \to \mathbb{R}, \ \theta \mapsto \frac{1}{2}W \bullet \cos(T(\theta)) = \frac{1}{2} \sum_{i=1}^n \sum_{j=1}^n w_{ij} \cos(\theta_i - \theta_j)$$

Lemma 3.1 The function f is:

- 1. invariant with respect to simultaneous, uniform rotation on every component. That means for every $\theta \in \mathbb{R}^n$ and every $\tau \in \mathbb{R}$ we have $f(\theta) = f(\theta + \tau e)$.
- 2. 2π -periodic with respect to each variable, i.e., for all $\theta \in \mathbb{R}^n$ we have $f(\theta) = f(\theta + 2\pi e_i)$, where e_i denotes the *i*-th standard basis vector.

Proof. We begin by proving the first property. Let $\theta \in \mathbb{R}^n$ and $\tau \in \mathbb{R}$, it is sufficient to show $T(\theta + \tau e) = T(\theta)$, as this is the only term in f that depends on θ . This holds, as for any $i, j \in [1:n]$ we have $T(\theta + \tau e)_{ij} = \theta_i + \tau - (\theta_j + \tau) = \theta_i - \theta_j = T(\theta)_{ij}$. To show the second property let $\theta \in \mathbb{R}^n$ and $i \in [1:n]$ and compute using 2π -periodicity of the cosine:

$$\cos(T(\theta + 2\pi e_i))_{ij} = \cos(\theta_i + 2\pi - \theta_j) = \cos(\theta_i - \theta_j) = \cos(T(\theta))_{ij} \text{ and } \cos(T(\theta + 2\pi e_i))_{ji} = \cos(\theta_j - \theta_i - 2\pi) = \cos(\theta_j - \theta_i) = \cos(T(\theta))_{ji}$$

As all other entries of $\cos(T(\theta))$ remain unchanged, $f(\theta) = f(\theta + 2\pi e_i)$ is established.

Recalling 10 and 9 we see that minimising f is nothing but solving 7 in polar coordinates. This observation is of central importance and we denote the relaxation of the MCP:

$$\min_{\theta \in \mathbb{R}^n} f(\theta) \tag{11}$$

This point of view has advantages and disadvantages. Since f is nonconvex we have no general tool to find a global minimum or even to decide whether a local minimum is a global minimum. In general, there can be multiple local but nonglobal minima. However, formulation 11 is an unconstrained minimisation problem, with a fairly easy analytical description of f. This allows for efficient computation of local minima. Making use of its simplicity we compute its partial derivatives by hand.

Lemma 3.2 The partial derivatives of f for any $j \in [1:n]$ is given by

$$\frac{\partial f}{\partial \theta_j}(\theta) = \sum_{k=1}^n w_{kj} \sin(\theta_k - \theta_j)$$

and the gradient can be written as:

$$\nabla f(\theta) = e^T \left(W \circ \sin(T(\theta)) \right) \tag{12}$$

Here the notation \circ stands for the Hadamard product, i.e. the entrywise product of W and $\sin(T(\theta))$

Proof. We fix $j \in [1:n]$ as the index of the variable for which we compute the partial derivative. To avoid ambiguity we rename the indices of f yielding:

$$f(\theta) = \frac{1}{2} \sum_{k=1}^{n} \sum_{i=1}^{n} w_{ki} \cos(\theta_k - \theta_i)$$

From this expression we see that the term k = i = j is equal to zero and we only need to distinguish between the following two cases:

In case of k = j and $i \neq j$ we have:

$$\frac{\partial}{\partial j} w_{ji} \cos(\theta_j - \theta_i) = -w_{ji} \sin(\theta_j - \theta_i) \stackrel{w_{ij} = w_{ji}}{=} w_{ij} \sin(\theta_i - \theta_j)$$

In case of i = j and $k \neq j$ we have:

$$\frac{\partial}{\partial j} w_{kj} \cos(\theta_k - \theta_j) = w_{kj} \sin(\theta_k - \theta_j)$$

In the case of $k \neq j \neq i$ the derivative is zero as the term is constant with respect to j. By linearity we therefore conclude:

$$\frac{\partial}{\partial j} f(\theta) = \frac{1}{2} \left(\sum_{k=1}^{n} w_{kj} \sin(\theta_k - \theta_j) + \sum_{j=1}^{n} w_{ij} \sin(\theta_k - \theta_j) \right) = \sum_{k=1}^{n} w_{kj} \sin(\theta_k - \theta_j)$$

To justify the gradient, we only need to observe that $\frac{\partial}{\partial j} f(\theta)$ is the sum of the entries of the *j*-th column of the matrix $W \circ \sin(T(\theta))$. This coincides with the *j*-th entry of $(W \circ \sin(T(\theta)))^T e = (e^T (W \circ \sin(T(\theta)))^T)$.

To be able to classify stationary points we now compute the Hessian matrix.

Lemma 3.3 The Hessian matrix H of f is given by

$$H(\theta) = W \circ \cos(T(\theta)) - \operatorname{diag}\left(\left(W \circ \cos(T(\theta))\right)e\right) \tag{13}$$

and the partial derivatives of second order are explicitly give by:

$$\frac{\partial^2}{\partial \theta_i \partial \theta_j} f(\theta) = \begin{cases} w_{ij} \cos(\theta_i - \theta_j) & \text{if } i \neq j \\ -\sum_{k \neq j} w_{kj} \cos(\theta_k - \theta_j) & \text{if } i = j \end{cases} \quad \forall i, j \in [1:n]$$

Proof. We begin by computing the partial derivatives of second order: Case $i \neq j$:

By $i \neq j$ we have that $w_{kj}\sin(\theta_k - \theta_j)$ is not constant with respect to θ_i if and only if k = i. Noting that $\frac{\partial}{\partial \theta_i}w_{ij}\sin(\theta_i - \theta_j) = w_{ij}\cos(\theta_i - \theta_j)$ we get

$$\frac{\partial^2}{\partial \theta_i \partial \theta_j} f(\theta) = w_{ij} \cos(\theta_i - \theta_j)$$

Case i = j:

In this case we have $w_{jj} = 0$ and $\frac{\partial}{\partial \theta_j} \sin(\theta_k - \theta_j) = -\cos(\theta_k - \theta_j)$. By linearity we thus get

$$\frac{\partial^2}{\partial \theta_j \partial \theta_j} f(\theta) = \sum_{k \neq j} -w_{kj} \cos(\theta_k - \theta_j)$$

The result follows, since the partial derivatives of second order coincide with the entries of

$$W \circ \cos(T(\theta)) - \operatorname{diag}((W \circ \cos(T(\theta))e).$$

Next we want to investigate the relationship between the function f and the cuts of a graph.

Definition 3.4 (Angular representation of a cut) A vector $\theta \in \mathbb{R}^n$ is called an angular representation of a cut, or simply a cut, if there exist integers k_{ij} such that $\theta_i - \theta_j = k_{ij}\pi$ for all $i, j \in [1:n]$.

Let $\overline{\theta}$ be an angular representation of a cut. Using that $\cos(\overline{\theta}_i - \overline{\theta}_j)$ evaluates to 1 if k_{ij} is even and to -1 if k_{ij} is odd, there exists a binary vector $\overline{x} \in \{-1, 1\}^n$ such that

$$\cos(\overline{\theta}_i - \overline{\theta}_j) = \overline{x}_i \overline{x}_j = \pm 1 \quad \text{for all } i, j \in [1:n].$$
 (14)

This can be seen by setting $x_1 = 1$ and

$$\overline{x}_i = \begin{cases} 1 & \text{if } \overline{\theta}_i - \overline{\theta}_1 \in 2\pi \mathbb{Z} \\ -1 & \text{otherwise} \end{cases} \quad \text{for } i \in [2:n]$$

This construction is correct, as for any $i, j \in [1:n]$ we have k_{ij} is even if and only if k_{i1} and k_{j1} are both even or both odd if and only if $\overline{x}_i \overline{x}_j = 1$. The first equivalence holds by

$$k_{ij}\pi = \overline{\theta}_i - \overline{\theta}_j = \overline{\theta}_i - \overline{\theta}_1 - (\overline{\theta}_i - \overline{\theta}_1) = (k_{i1} - k_{i1})\pi$$

Then \overline{x} can be viewed as the cut corresponding to $\overline{\theta}$. Moreover, the cut value corresponding to $\overline{\theta}$ is

$$\psi(\overline{\theta}) = \frac{1}{2} \sum_{1 \le i \le j \le n} w_{ij} \left(1 - \cos(\overline{\theta}_i - \overline{\theta}_j) \right)$$
 (15)

Furthermore this equation makes sense for all $\theta \in \mathbb{R}^n$ and we can interpret its value as a relaxed cut value. In the next Proposition we give a detailed account of this correspondence.

Proposition 3.5 Modulo the uniform rotation and the periodicity for each variable, there is an one-to-one correspondence between binary cuts and angular representations of a cut, given by

$$\overline{\theta}_i = \begin{cases} 0 & \text{if } \overline{x}_i = 1\\ \pi & \text{if } \overline{x}_i = -1 \end{cases} \tag{16}$$

In the same manner we can set

$$\overline{x}_i = \begin{cases} 1 & \text{if } \overline{\theta}_i = 0\\ -1 & \text{if } \overline{\theta}_i = \pi \end{cases}$$

Proof. It is clear that assignment 16 defines a one-to-one correspondence between binary cuts and angular representations in $\{0,\pi\}^n$. On the other hand every angular representation of a cut has a representative, where every variable is either 0 or π . Let $\overline{\theta}$ be an angular representation of a cut. By the second property of f, we can assume $\overline{\theta}_i \in [0,2\pi)$ for all $i \in [1:n]$. We only need to show that there exists a $\tau \in \mathbb{R}$ such that $\overline{\theta}_i - \tau \in \{0,\pi\}$ for all $i \in [1:n]$. If $\overline{\theta}_i \in \{0,\pi\}$ for all $i \in [1:n]$ we simply pick $\tau = 0$ and are done. Otherwise there exists an $i \in [1:n]$, such that $\overline{\theta}_i \notin \{0,\pi\}$. If $\overline{\theta}_i \in (0,\pi)$ we choose $\tau := \overline{\theta}_i$, else we have $\overline{\theta}_i \in (\pi,2\pi)$ and we choose $\tau := \overline{\theta}_i - \pi$. Now let $j \in [1:n]$. With this choice we just have to show that

$$\overline{\theta}_j = \begin{cases} \tau & \text{if } \overline{\theta}_j \in (0, \pi) \\ \pi + \tau & \text{if } \overline{\theta}_j \in (\pi, 2\pi) \end{cases}$$
 (17)

holds. As $\overline{\theta}$ is an angular representation of a cut there exists a $k \in \mathbb{Z}$ such that $\overline{\theta}_i - \overline{\theta}_j = k\pi$. By rearranging the terms there exists a $\widetilde{k} \in \mathbb{Z}$ such that $\overline{\theta}_j = \overline{\theta}_i - k\pi = \widetilde{k}\pi + \tau$. This shows equation 17. This shows $(\overline{\theta} - \tau e)_i \in \{0, \pi\}$ for all $i \in [1:n]$.

Using this correspondence we can represent a cut by both $\overline{\theta}$ and \overline{x} . Given a binary representation of a cut \overline{x} (or an angular representation $\overline{\theta}$), we denote the angular representation by $\theta(\overline{x})$ (or the binary representation by $x(\overline{\theta})$ of that same cut. The next Proposition states an import property of angular representations of a cut.

Proposition 3.6 Every angular representation of a cut $\overline{\theta} \in \mathbb{R}^n$ is a stationary point of the function f.

Proof. As $\overline{\theta}$ is a cut we have $\overline{\theta}_i - \overline{\theta}_j \in \mathbb{Z}\pi$ for all $i, j \in [1:n]$. This directly implies that every entry of $\sin(T(\overline{\theta}))$ is zero. Plugging this into 12 yields $\nabla f(\overline{\theta}) = 0$. This shows that $\overline{\theta}$ is a stationary point.

Definition 3.7 (Nonnegatively summable matrix) A matrix $M \in \mathbb{R}^{n \times n}$ is called nonnegatively summable if the sum of the entries in every principal submatrix of M is nonnegative, or equivalently, if $u^T M u \geq 0$ for every binary vector $u \in \{0,1\}^n$.

Put in words a principal submatrix is a matrix that we get by crossing out rows and corresponding columns, i.e. crossing out the *i*-th row implies crossing out the *i*-th column as well and vice versa. By definition, every semidefinite matrix is nonnegatively summable. However the other implication does not hold as the following example shows.

Example Let n > 1 and denote the identity matrix as $I \in \mathbb{R}^{n \times n}$. The matrix $ee^T - I$ is nonnegatively summable but not positive semidefinite. As all the entries of $ee^T - I$ are nonnegative the matrix is clearly nonnegatively summable. To show that $ee^T - I$ is not positive semidefinite we consider the vector $u := e_1 - e_n$, where e_1 is a standard basis vector. By n > 1 we have $1 \neq n$ and we get $u^T (ee^T - I) u = u^T (-1, 0, \dots, 0, 1) = -2 < 0$.

Before we can provide a characterisation for maximum (and minimum) cuts in Lemma 3.9 we prove a little calculation rule:

Lemma 3.8 Let $W \in \mathbb{R}^{n \times n}$ be a symmetric matrix, $x \in \mathbb{R}^n$ and $\delta, \delta' \in \{0, 1\}^n$. Then we have

$$(\delta \circ x)^T W (\delta' \circ x) = \delta^T (W \circ xx^T) \delta' = \delta'^T (W \circ xx^T) \delta$$

Proof. First we observe that for any $\delta \in \{0,1\}^n$ we have

$$(\delta \circ x)_i = \begin{cases} 0 & \text{if } \delta_i = 0 \\ x_i & \text{if } \delta_i = 1 \end{cases} = \delta_i x_i$$

Using this we can compute:

$$(\delta \circ x)^T W (\delta' \circ x) = \sum_{i=1}^n (\delta \circ x)_i \sum_{j=1}^n w_{ij} (\delta' \circ x)_j = \sum_{i=1}^n \delta_i x_i \sum_{j=1}^n w_{ij} \delta'_j x_j$$
$$= \sum_{i=1}^n \delta_i \sum_{j=1}^n w_{ij} x_i x_j \delta'_j = \delta^T A \delta'$$

where $A_{ij} = w_{ij}x_ix_j$ for all $i, j \in [1:n]$. Thus we have $A = W \circ xx^T$ and the first equality holds. Having established the first equality the second follows by noting that W is symmetric:

$$(\delta \circ x)^T W (\delta' \circ x) = (\delta' \circ x)^T W^T (\delta \circ x) = \delta'^T (W \circ xx^T) \delta$$

Lemma 3.9 Let $\overline{x} \in \{-1,1\}^n$ be given and consider the matrix $M(\overline{x}) \in \mathbb{R}^{n \times n}$ defined as

$$M(\overline{x}) = W \circ (\overline{x}\overline{x}^T) - \operatorname{diag}((W \circ (\overline{x}\overline{x}^T)) e)$$
(18)

Then, \overline{x} is a maximum (respectively, minimum) cut if and only if $M(\overline{x})$ (respectively, $-M(\overline{x})$) is nonnegatively summable.

Proof. Consider the quadratic function $q: \mathbb{R}^n \to \mathbb{R}$, $x \mapsto x^T W x/2$. By Lemma 2.2 we know that $\overline{x} \in \{-1,1\}^n$ is a maximum cut if and only if \overline{x} satisfies $q(\overline{x}) \leq q(x)$ for all $x \in \{-1,1\}^n$. Consider a maximum cut $\overline{x} \in \{-1,1\}^n$ and an arbitrary $x \in \{-1,1\}^n$. As W is symmetric $\overline{x}^T W x = x^T W \overline{x}$ holds. Using this and the previous equivalence we have:

$$0 \le q(x) - q(\overline{x}) = \frac{1}{2} \left(x^T W x - \overline{x}^T W \overline{x} \right) = \frac{1}{2} \left(x^T W x + \overline{x}^T W x - x^T W \overline{x} - \overline{x}^T W \overline{x} \right)$$

$$= \frac{1}{2} (x + \overline{x})^T W (x - \overline{x}) = \left(\overline{x} + \frac{1}{2} (x - \overline{x}) \right)^T W (x - \overline{x})$$

$$= \overline{x}^T W (x - \overline{x}) + \frac{1}{2} (x - \overline{x})^T W (x - \overline{x})$$
(19)

As we have $\overline{x}, x \in \{-1, 1\}^n$ we know that either $x_i = \overline{x}_i$ or $x_i = -\overline{x}_i$ for all $i \in [1 : n]$. With this observation we see that

$$(x - \overline{x})_i = \begin{cases} 0 & \text{if } x_i = \overline{x}_i \\ -2\overline{x}_i & \text{if } x_i \neq \overline{x}_i \end{cases} \text{ for all } i \in [1:n]$$

and by defining $\delta \in \mathbb{R}^n$ as

$$\delta_i = \begin{cases} 0 & \text{if } x_i = \overline{x_i} \\ 1 & \text{if } x_i \neq \overline{x_i} \end{cases}$$

we get the identity

$$x - \overline{x} = -2\delta \circ \overline{x}$$

With this we continue computation 19:

$$0 \leq \overline{x}^T W (x - \overline{x}) + \frac{1}{2} (x - \overline{x})^T W (x - \overline{x})$$

$$= -2\overline{x}^T W (\delta \circ \overline{x}) + 2 (\delta \circ \overline{x})^T W (\delta \circ \overline{x})$$

$$= -2 (e \circ \overline{x})^T W (\delta \circ \overline{x}) + 2 (\delta \circ \overline{x})^T W (\delta \circ \overline{x})$$

$$\stackrel{3.8}{=} -2\delta^T (W \circ \overline{x}\overline{x}^T) e + 2\delta^T (W \circ \overline{x}\overline{x}^T) \delta$$

$$= -2\delta^T \operatorname{diag} ((W \circ \overline{x}\overline{x}^T)e) \delta + 2\delta^T (W \circ \overline{x}\overline{x}^T) \delta \stackrel{\text{def } M(\overline{x})}{=} 2\delta M(\overline{x}) \delta$$

In the second to last equality we used $\delta^T v = \delta^T \operatorname{diag}(v) \delta$, which follows immediately from

$$(\operatorname{diag}(v)\delta)_i = \begin{cases} 0 & \text{if } \delta_i = 0 \\ v_i & \text{if } \delta_i = 1 \end{cases} \quad \text{for all } i \in [1:n].$$

So far we have shown $0 \le \delta^T M(\overline{x}) \delta$, however for nonnegative summability we need to argue $0 \le y^T M(\overline{x}) y$ for all $y \in \{0,1\}^n$.

Claim: For any $y \in \{0,1\}^n$ there exists a $x \in \{-1,1\}^n$ such that $\delta = y$.

First recall that $\delta \in \{0,1\}^n$ depends on \overline{x} , which is fixed, and x, which is a variable. To clarify this dependence we denote δ as $\delta_{\overline{x}}(x)$. As $x \in \{-1,1\}^n$ can be chosen arbitrarily, we choose

$$x_i = \begin{cases} \overline{x_i} & \text{if } y_i = 0\\ -\overline{x_i} & \text{if } y_i = 1 \end{cases}$$

This choice directly implies $\delta_{\overline{x}}(x) = y$. The claim shows that $M(\overline{x})$ is nonnegatively summable. Now we want to show that \overline{x} is minimum cut if and only if $-M(\overline{x})$ is nonnegatively summable. We do this in the same way as above only outlining the slight changes. Let q be the quadratic function we defined above, $\overline{x} \in \{-1,1\}^n$ a minimum cut and $x \in \{-1,1\}^n$ arbitrary. Thus we have $0 \ge q(x) - q(\overline{x})$ and the same computations as above yield $0 \ge \delta^T M(\overline{x}) \delta \Leftrightarrow 0 \le \delta^T (-M(\overline{x})) \delta$. The nonnegative summability of $-M(\overline{x})$ follows in the same manner as above.

Remark 1. Let $\overline{x} \in \{-1, 1\}^n$. Then the map

$$\delta_{\overline{x}}: \{-1,1\}^n \to \{0,1\}^n, \ x \mapsto \delta_{\overline{x}}(x), \text{ where } \delta_{\overline{x}}(x)_i = \begin{cases} 0 & \text{if } x_i = \overline{x_i} \\ 1 & \text{if } x_i \neq \overline{x_i} \end{cases} \text{ for all } i \in [1:n]$$

is a bijection

In the previous proof we have shown that $\delta_{\overline{x}}$ is surjective and as $\{-1,1\}^n$, $\{0,1\}^n$ are finite sets of the same cardinality we have bijectivity.

2. The matrix $M(\overline{x})$ is symmetric for every $\overline{x} \in \{0,1\}^n$: The matrix diag $(W \circ (\overline{x}\overline{x}^T) e)$ is symmetric as it is a diagonal matrix. As W is symmetric, we have for every $i, j \in [1:n]$:

$$(W \circ \overline{x}\overline{x}^T)_{ij} = w_{ij}\overline{x}_i\overline{x}_j = w_{ji}\overline{x}_j\overline{x}_i = (W \circ \overline{x}\overline{x}^T)_{ji}$$

This shows that $(W \circ \overline{x}\overline{x}^T)$ is symmetric and, as a difference of symmetric matrices, $M(\overline{x})$ is a symmetric matrix.

Comparing the Hessian of f as given in 13 to 18, we see that the two look similar. The difference is that $\cos(T(\theta))$ in 13 is replaced by xx^T in 18. Comparing the entries we see that $\cos(\theta_i - \theta_j) = x_i x_j$ for all $i, j \in [1:n]$ is a sufficient condition for $H(\theta) = M(x)$. This observation leads to the next Theorem providing a classification of angular representations of cuts as stationary points of the function f.

Theorem 3.10 Let $\overline{\theta}$ be a angular representation of a cut, or simply a cut, and let $\overline{x} \equiv x(\overline{\theta})$ be the associated binary cut. If $\overline{\theta}$ is a local minimum (respectively, local maximum) of $f(\theta)$, then \overline{x} is a maximum (respectively, minimum) cut. Consequently, if \overline{x} is neither a maximum cut nor a minimum cut, then $\overline{\theta}$ must be a saddle point of f.

Proof. First recall that every angular representation of a cut is a critical point of f and that $\cos(\overline{\theta_i} - \overline{\theta_j}) = \overline{x_i}\overline{x_j}$ for all $i, j \in [1:n]$. Therefore $H(\overline{\theta}) = M(\overline{x})$ holds. Let $\overline{\theta}$ be a local minimum of f. Since f is smooth, $H(\overline{\theta})$ is positive semidefinite and in particular nonnegatively summable. By $H(\overline{x}) = M(\overline{x})$ we have that $M(\overline{x})$ is nonnegatively summable. Now let $\overline{\theta}$ be a local maximum of f. Then $H(\overline{\theta})$ is negative semidefinite, that means that $-H(\overline{\theta})$ is nonnegatively summable. Therefore $-H(\overline{\theta}) = -M(\overline{x})$ implies that $-M(\overline{x})$ is nonnegatively summable. Lastly let \overline{x} be neither a maximum cut nor a minimum cut. By Lemma 3.9 we know that neither $M(\overline{x})$ nor $-M(\overline{x})$ is nonnegatively summable. By $M(\overline{x}) = H(\overline{\theta})$ we have that $H(\overline{\theta})$ is neither positive semidefinite nor negative semidefinite. By the necessary condition for a local extremum we conclude that $\overline{\theta}$ is neither a local minimum nor a local maximum. As $\overline{\theta}$ is a critical point of f, it must be a saddle point.

Although all cuts are stationary points of f (by Proposition 3.6) Theorem 3.10 shows that only the maximum cuts can be local minima of f, see [BMZ02, p. 508].

Example The converse of the two implications in the above Theorem do not hold. We will give an example showing that the angular representation of a maximum cut is not necessarily a local minimum of f.

Consider the complete graph with three vertices, where every edge has weight 1. Then the weight of a maximum cut is 2 and is achieved by $\overline{x} = \begin{pmatrix} 1 & -1 & -1 \end{pmatrix}$. We then have

$$W \circ \overline{x}\overline{x}^T = \begin{pmatrix} 1 & -1 & -1 \\ -1 & 1 & 1 \\ -1 & 1 & 1 \end{pmatrix}$$
 and $(W \circ \overline{x}\overline{x}^T) e = \begin{pmatrix} -1 \\ 1 \\ 1 \end{pmatrix}$

Using 18 we get

$$M(\overline{x}) = \begin{pmatrix} 1 & -1 & -1 \\ -1 & 1 & 1 \\ -1 & 1 & 1 \end{pmatrix} - \begin{pmatrix} -1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{pmatrix} = \begin{pmatrix} 2 & -1 & -1 \\ -1 & 0 & 1 \\ -1 & 1 & 0 \end{pmatrix}$$

Straight forward computation shows nonnegative summability, namely $u^T M(\overline{x})u \geq 0$ for all $u \in \{0,1\}^n$. However, $M(\overline{x})$ is not positive semidefinite, since

$$\begin{pmatrix} 1 & 0 & 2 \end{pmatrix} M(\overline{x}) \begin{pmatrix} 1 \\ 0 \\ 2 \end{pmatrix} = \begin{pmatrix} 1 & 0 & 2 \end{pmatrix} \begin{pmatrix} 0 \\ 1 \\ -1 \end{pmatrix} = -2 < 0$$

Using $H(\theta(\overline{x})) = M(\overline{x})$, we know that $H(\theta(\overline{x}))$ is not positive semidefinite. Thus $\theta(\overline{x})$ is not a local minimum of f.

However there are special classes of instances, where for every maximum cut \overline{x} , the angular representation $\theta(\overline{x})$ is a local minimum. The following Proposition gives a class for which this holds:

Proposition 3.11 For a bipartite graph with nonnegative edge weights, the global minimum value of f is attained by a maximum cut.

Proof. We denote the bipartite graph G = (A, B, E). As all the edge weights are nonnegative, the cut (A, B), that cuts through all edges is a maximum cut. For that cut we have $\cos(\theta_i - \theta_j) = -1$ for all $\{i, j\} \in E$. For that cut f evaluates to $-\frac{1}{2}e^TWe$. This value must be a global minimum of f as all the entries of W are nonnegative.

For instances, where a maximum cut \overline{x} corresponds to a local minimum $\theta(\overline{x})$ of f, the optimality of x can be checked in polynomial time. That is because $\theta(\overline{x})$ local minimum implies $H(\theta(\overline{x}) = M(\overline{x}))$ positive semidefinite. This in turn implies $M(\overline{x})$ nonnegatively summable, which shows \overline{x} maximum cut by 3.9. A detailed proof showing that we can decide whether a symmetric matrix (- $M(\overline{x})$ is symmetric-) is positive semidefinite in polynomial time is provided in [KV18, Theorem 16.8].

Theorem 3.10 directly implies the following Corollary, which plays an important role for the heuristic we will develop in the next section.

Corollary 3.12 Let $x \in \{0,1\}^n$ be a nonmaximum cut. Then $\theta(x)$ can not be a local minimum.

Proof. The statement is the contraposition of the first statement in Theorem 3.10 \Box

Therefore, as written in [BMZ02, p.509], a good minimisation algorithm would not be attracted to stationary points which are not local minima. We construct such an algorithm in section 4.

4 A heuristic algorithm for MCP

In this section we will describe the heuristic described by [BMZ02]. Again we assume non-negative edge weights, in order to mirror the analysis done in 2.

Consider a local minimum θ of f. As θ is not necessarily an angular representation of a cut, we need to develop a method that associates a cut $x \in \{-1,1\}$ to any $\theta \in \mathbb{R}^n$. Using polar coordinates we can associate the entries of θ to points on the unit circle. Furthermore we can assume $\theta_i \in [0,2\pi)$ for all $i \in [1:n]$. Inspired by the Goemans-Williamson cuts in section 2 we can generate a cut by cutting the unit circle into two halves. For any angle $\alpha \in [0,\pi)$ we can set

$$x_i = \begin{cases} 1 & \text{if } \theta_i \in [\alpha, \alpha + \pi) \\ -1 & \text{otherwise} \end{cases}$$
 (20)

generating a cut. The weight of the obtained cut is given by

$$\gamma(x) = \frac{1}{2} \sum_{1 \le i < j \le n} w_{ij} \left(1 - x_i x_j \right)$$

Remark 1. Let $\theta \in [0, 2\pi)^n$ and $\alpha \in [0, \pi)$. Let x^{α} denote the cut generated by α and $x^{\alpha+\pi}$ the cut generated by $\alpha+\pi$. We then have $x^{\alpha+\pi}=-x^{\alpha}$. For $\alpha+\pi$ assignment 20 is to be understood as:

$$x_i = \begin{cases} 1 & \text{if } \theta_i \in [0, \alpha) \cup [\alpha + \pi, 2\pi) \\ -1 & \text{otherwise} \end{cases}$$

By observing $[0, 2\pi) = [\alpha, \alpha + \pi) \cup ([0, \alpha) \cup [\alpha + \pi, 2\pi))$ we deduce $-x^{\alpha} = x^{\alpha + \pi}$. As x^{α} and $x^{\alpha + \pi}$ are of the same weight, it suffices to consider $\alpha \in [0, \pi)$ instead of $\alpha \in [0, 2\pi)$.

- 2. The cut generated by assignment 20 can also be generated by a Goemans-Williamson cut. This is achieved by simply slightly turning the diving line in the mathematically negative sense. Slightly means more than zero but strictly less than the smallest angle between two distinct points. As we are given a finite amount of points we know that such a turn exists.
- 3. Choosing the halfopen interval $[\alpha, \alpha + \pi)$ ensures, that diametrically opposed points are separated, i.e. the corresponding vertices are on different sides of the cut.

By increasing α we have a simple way to examine all cuts obtained by assignment 20, called Goemans-Williamson-type cuts. As we will describe in algorithm 2 we have a deterministic and computationally inexpensive way of finding a best possible Goemans-Williamson-type cut. The algorithm works by rotating the separating line in the mathematically positive sense, starting with angle 0. In the update step the angle is set to the angle of the line that first intersects a new point on the circle, while rotating. If the rotated line has not encountered a new point yet it generates the same cut as the line obtained by the update step. Therefore the algorithm finds all possible Goemans-Williamson-type cuts.

We will make two assumptions without loss of generality. First of all we note that by the 2π -periodicity in each variable of f, see 3.1, we can assume $\theta_i \in [0, 2\pi)$ for all $i \in [1:n]$. Furthermore we assume $\theta_1 \leq \theta_2 \leq \cdots \leq \theta_n$.

Algorithm 2 Procedure-Cut

```
Input: \theta \in [0, 2\pi)^n satisfying \theta_1 \leq \theta_2 \leq \cdots \leq \theta_n, undirected weighted graph G
Output: A cut x^*
 1: Let x^* trivial cut, i.e. all vertices on one side
 2: Let \alpha = 0, \Gamma = -\infty, i = 1. Let j be the smallest index such that \theta_j > \pi if there is one;
    otherwise set j = n + 1. Set \theta_{n+1} = 2\pi.
 3: while \alpha < \pi do
         Generate cut x by 20 and compute \gamma(x)
         If \gamma(x) > \Gamma, then let \Gamma = \gamma(x) and x^* = x.
 5:
         if \theta_i \leq \theta_j - \pi then
 6:
             Let \alpha = \theta_i and increment i by 1.
 7:
 8:
         else
             Let \alpha = \theta_j - \pi and increment j by 1.
 9:
         end if
10:
11: end while
12: return x^*
```

Remark The assumption $\theta_1 \leq \theta_2 \leq \cdots \leq \theta_n$ is not a real constraint, we just have to keep track of the original indices, as they determine the cut. Let σ be a permutation such that $\theta_{\sigma(1)} \leq \theta_{\sigma(2)} \leq \cdots \leq \theta_{\sigma(n)}$. We can now apply algorithm 2, with the only slight change that the cut in line 2.4 needs to be generated in the following manner:

$$x_i = \begin{cases} 1 & \text{if } \theta_{\sigma(i)} \in [\alpha, \alpha + \pi) \\ -1 & \text{otherwise} \end{cases}$$

Remark Algorithm 2 finds all Goemans-Williamson-type cuts.

Recall that our rank-two relaxation has the form of the Goemans and Williamson relaxation in dimension 2 and the Goemans-Williamson-type cuts can be generated by Goemans-Williamson cuts. Hence we can analyse the performance of our algorithm in a similar manner as in section 2.

Lemma 4.1 Let $\theta \in [0, 2\pi)^n$ such that $\theta_1 \leq \theta_2 \leq \cdots \leq \theta_n$. Then the cut value generated by algorithm 2 is at least 0.878 times the relaxed cut value $\psi(\theta)$ as defined in 15, written in a formula:

$$\gamma(x^*) \ge 0.878\psi(\theta) \tag{21}$$

Proof. The entry θ_i represents the point $v_i = \left(\cos(\theta_i) \sin(\theta_i)\right)^T$ for $i \in [1:n]$. We review the probabilities that two points v_i and v_j are separated by a Goemans-Williamson-type cut, for any $i, j \in [1:n]$. In dimension 2, we say that a Goemans-Williamson cut is of angle β , if it is generated by $r = \left(\cos(\beta + \frac{\pi}{2}) \sin(\beta + \frac{\pi}{2})\right)$. A Goemans-Williamson cut of angle β induces the cut

$$x_i = \begin{cases} 1 & \theta_i \in [\beta, \beta + \pi] \\ -1 & \text{otherwise} \end{cases}$$

Clearly, if two points are separated by a Goemans-Williamson cut they are also separated by a Goemans-Williamson-type cut. Thus for an angle β the points v_i and v_j are separated by

a Goemans-Williamson-type cut but not a Goemans-Williamson cut if and only if $\beta = \theta_i - \frac{\pi}{2}$ or $\beta = \theta_j - \frac{\pi}{2}$. Since the set of separating angles only changes on a nullset the probabilities remain unchagned. Therefore we can redo the proof of Lemma 2.10 by substituting Goemans-Williamson cut for Goemans-Williamson-type cut and OPT for $\psi(\theta)$. This yields that the expected value of a Goemans-Williamson-type cut for a given θ is greater or equal to $\alpha\psi(\theta)$. Since algorithm 2 finds all Goemans-Williamson-type cuts it finds a Goemans-Williamson-type cut of value greater or equal to $\alpha\psi(\theta)$.

This is not a performance guarantee, as we can not guarantee that $\psi(\theta)$ is an upper bound on the maximum cut value. Recall, that in Goemans-Williamson algorithm, the guarantee comes from (approximately) solving the SDP 8. The next Lemma can be interpreted as a weak performance guarantee

Lemma 4.2 Let x_a^* and x_b^* be two cuts generated by algorithm 2 from θ_a and θ_b respectively. If $\gamma(x_a^*) \leq \psi(\theta)$ and $\psi(\theta_b) > \psi(\theta_a)/0.878$, then x_b^* has a higher cut value than x_a^* .

Proof. Using the assumptions we confirm

$$\gamma(x_b^*) \stackrel{21}{\ge} 0.878 \cdot \psi(\theta_b) > \psi(\theta_a) \ge \gamma(x_a^*)$$

The next paragraph, describes the idea of algorithm 3. We minimise the function f and obtain a minimiser called θ^1 . We can now start algorithm 2 and obtain a best possible cut x^1 associated with θ^1 . At this point we could return the cut x^1 . If θ^1 happens to be an angular representation of a cut we know that x^1 is a maximum cut, by Theorem 3.10. Otherwise, we can attempt to improve the cut value by spending more computational resources. By Theorem 3.10 we know that the angular representation of x^1 , denoted as $\theta(x^1)$, is a stationary point of f, most likely a saddle point. Therefore we can hope to find a deeper local minimum close to $\theta(x^1)$. As we use a gradient descend, restarting the minimisation from a stationary point is of no use. Thus we restart the minimisation of f from a slight perturbation of $\theta(x^1)$, in hopes of finding another local minimum $\theta^2 \neq \theta^1$ from which we hope to obtain a cut x^2 with a higher cut value than x^1 , i.e. $\gamma(x^2) > \gamma(x^1)$. In case $\gamma(x^2) > \gamma(x^1)$ is achieved we deem our attempt successful, and unsuccessful otherwise. We can set the termination criterion to be N consecutive unsuccessful attempts of improving the value of the cut.

In algorithm 3 we have θ^0 as input, which is the initial seed for the minimisation of f. We can try and improve our chances of finding a high quality heuristic solution by running algorithm 3 from different starting points, in other words increase the number of seeds. The number of seeds, will be given by the input M.

Assuming that we do not find a maximum cut, algorithm 3 can also be interpreted in the following way. From a seed we start minimising f. By means of Goemans-Williamson-type cuts, we look for nearby saddle points. Here it is important to note, that nearby is to be understood by the cuts that we can generate from the minimum. The so found saddle points is not necessarily the closest saddle point in the euclidean metric. The saddle point is chosen by the associated cut with the highest weight. From a slight perturbation of that point we search for a local minimum, which has a saddle point with lower f-value in its neighbourhood.

Algorithm 3 ImprovedCut

```
Input: Integer N \in \mathbb{N}, a seed \theta^0 \in \mathbb{R}^n, undirected weighted graph G
Output: A cut x^*
 1: procedure Improved Cut(input N, \theta^0)
        Given \theta^0 \in \mathbb{R}^n and integer N \geq 0, let k = 0 and \Gamma = -\infty
         while k \leq N do
 3:
             Starting from \theta^0, minimise f to get \theta
 4:
             Compute a best cut x associated with \theta by 2
 5:
             if \gamma(x) > \Gamma then
 6:
                 Let \Gamma = \gamma(x), x^* = x and k = 0
 7:
 8:
             else
                 k = k + 1
 9:
             end if
10:
             Set \theta^0 to a random perturbation of the angular representation of x.
11:
12:
        end while
        return x^*
13:
14: end procedure
```

Algorithm 4 BurerStub

```
Input: Undirected weighted graph G
Output: A cut x^*
 1: Let x^* trivial cut, i.e. all vertices on one side
 2: for i \leftarrow 0, n do
         Generate random \theta^0 \in \mathbb{R}^n
 3:
         x \leftarrow \text{IMPROVEDCUT}(N, \theta^0)
 4:
         if \gamma(x) > \gamma(x^*) then
 5:
             x^* \leftarrow x
 6:
 7:
         end if
 8: end for
 9: return x^*
```

Generally speaking the bigger M and N are the longer the algorithm will run and the better the returned cut will be. However as the termination criterion in IMPROVEDCUT relies on the number of unsuccessful attempts we can not predict the runtime. Since there are elements of randomness, namely in the choice of the seed and the random pertubation there can always be exceptions to the rule.

5 Solving BQP using MCP

Closely following [JM21, Section 2] we are going to give a cursory explanation, that binary quadratic problems of the form: Let $Q \in \mathbb{R}^{n \times n}$

minimize
$$x^T Q x$$

subject to $x_i \in \{0, 1\} \forall i \in [1:n]$ (22)

can be solved by solving the MCP on a suitable graph with appropriately chosen edge weights, see Corollary 5.2.

To this end we formulate the MCP as an integer linear programming problem:

minimize
$$\sum_{e \in E} w_e z_e$$
 subject to
$$\sum_{e \in F} z_e - \sum_{e \in C \backslash F} z_e \le |F| - 1 \quad F \subseteq C \subseteq E \text{ subject to } C \text{ cycle and } |F| \text{ odd,}$$

$$0 \le z_e \le 1 \qquad \forall e \in E,$$

$$z_e \in \mathbb{Z} \qquad \forall e \in E$$

$$(23)$$

An optimal solution to 23 gives us the following maximum cut $\{\{i,j\} \mid z_{ij} = 1\}$. Any cut of G is a feasible solution to 23 and vice versa. The cut polytope of a graph G can thus be denoted as

$$P_{\text{CUT}}^{G} := \left\{ z \in [0, 1]^{E} \mid z \text{ is a feasible solution to } 23 \right\}.$$

For the BQP 22 consider the following undirected graph G=(V,E). The vertices $V=\{v_1,\ldots,v_n\}$, where v_i corresponds to the variable x_i for every $i\in[1:n]$ and $E=\{\{v_i,v_j\}\mid i,j\in[1:n]:q_{ij}+q_{ji}\neq 0\}$ where each edge $\{v_i,v_j\}$ corresponds to the product $y_{ij}:=x_ix_j$. We admit that the Boolean quadric polytope is given by

$$P_{\text{BQP}}^{G} = \left\{ (x, y) \in \{0, 1\}^{|V|} \times \{0, 1\}^{|E|} \mid (x, y) \text{ satisfies } y_{ij} = x_i x_j \right\}.$$

Without proof we state the following Theorem proven in [De 90].

Theorem 5.1 Let $Q \in \mathbb{R}^{n \times n}$ and consider G = (V, E) with V := [1:n] and $E := \{\{i, j\} \mid i, j \in [1:n] : q_{ij} + q_{ji} \neq 0\}$. Define H = (W, F) where $W := \{v_0, \dots, v_n\}$ and $F := \{\{v_i, v_j\} \mid \{i, j\} \in E\} \cup \{\{v_0, v_i\} \mid i \in [1:n]\}$. Let $f : \mathbb{R}^F \to \mathbb{R}^{V \cup E}$ be the bijective linear map defined by

$$x_v = z_{0v}$$
 for all $v \in V$, and $y_{vw} = x_v x_w = \frac{1}{2} (z_{0v} + z_{0w} - z_{vw})$ for all $\{v, w\} \in E$.

Then $P_{\text{BQP}}^{\text{G}} = f\left(P_{\text{CUT}}^{\text{H}}\right)$.

Using this Theorem we can show a fundamental result allowing us to solve a BQP as a MCP.

Corollary 5.2 ([Mal22, p. 2]) Let $Q \in \mathbb{R}^{n \times n}$ and $E = \{\{i, j\} \mid i, j \in [1:n] \text{ such that } q_{ij} + q_{ji} \neq 0\}$. Then the corresponding BQP with the objective function x^TQx to be minimised can be solved as a MCP on H = (W, F) where $W := \{v_0, \dots, v_n\}$ and $F := \{\{v_i, v_j\} \mid \{i, j\} \in E\} \cup \{\{v_0, v_i\} \mid i \in [1:n]\}$, with the objective function

$$\max \sum_{\substack{\{v_i, v_j\} \in F \\ v_i \neq v_0 \neq v_j}} \frac{1}{2} (q_{ij} + q_{ji}) z_{ij} - \sum_{i \in W \setminus \{v_0\}} \widetilde{c}_i z_{0i},$$

where $\widetilde{c}_i := q_{ii} + \sum_{j=1}^n \frac{1}{2} (q_{ij} + q_{ji}) [\{v_i, v_j\} \in F, v_i \neq v_0 \neq v_j]$ for all $i \in W \setminus \{v_0\}$.

Proof. We have the following chain of equalities:

$$\begin{aligned} xQx^T \\ &= \sum_{i,j=1}^n q_{ij}x_ix_j + \sum_{i=1}^n q_{ii}x_ix_i \\ &\stackrel{x_ix_i=x_i}{=} \sum_{\{i,j\}\in E} \left(q_{ij} + q_{ji}\right)x_ix_j + \sum_{i=1}^n q_{ii}x_i \\ &\stackrel{=}{=} \sum_{\{i,j\}\in E} \left(q_{ij} + q_{ji}\right)\left(\frac{1}{2}\left(z_{0i} + z_{0j} - z_{ij}\right)\right) + \sum_{i\in W\setminus\{v_0\}} q_{ii}z_{0i} \\ &= -\sum_{\{i,j\}\in E} \frac{1}{2}\left(q_{ij} + q_{ji}\right)z_{ij} + \frac{1}{2}\sum_{\{i,j\}\in E} \left(q_{ij} + q_{ji}\right)\left(z_{0i} + z_{0j}\right) + \sum_{i\in W\setminus\{v_0\}} q_{ii}z_{0i} \\ &\stackrel{*}{=} -\sum_{\{i,j\}\in E} \frac{1}{2}\left(q_{ij} + q_{ji}\right)z_{ij} + \sum_{i\in W\setminus\{v_0\}} \left(q_{ii} + \frac{1}{2}\sum_{\substack{j\neq 0\\\{v_i,v_j\}\in F}} \left(q_{ij} + q_{ji}\right)\right)z_{0i} \\ &= -\sum_{\{i,j\}\in E} \frac{1}{2}\left(q_{ij} + q_{ji}\right)z_{ij} + \sum_{i\in W\setminus\{v_0\}} \tilde{c}_iz_{0i} \end{aligned}$$

In * we use that for for all $\{i, j\} \in E$ we have $i, j \neq 0$ and $i \neq j$ and move the middle term to the right term. The previous computation shows that the two objective functions are the negation of each other under the bijection f from Theorem 5.1. This shows the result. \square

6 Experimentation

6.1 Experimental Setup

The goal of our experimental study is to compare the performance of the Burer heuristic for a range of different parameters. We recall that the Burer heuristic depends on two parameter: The parameter N, which determins the number of consecutive tries for improving a cut and the parameter M, which controls the number of starting points. We are going to compare 21 solvers, which are the combinations of $M \in \{1, 5, 10, 15, 20, 30, 40, 50\}$ and $N \in \{5, 10, 15\}$. As these parameters are characteristic we have the following naming convention: The solver

with the name Mxx.Nyy is the Burer heuristic with M=xx and N=yy. The solvers are run on the instances from [Mal23]. We will give a broad overview of the library for a detailed description we refer to [Mal23]. The library comprises 431 instances, split into 8 classes. The number of vertices ranges from 21 to 3000 with a median of 101 and a mean of 394.69 vertices. There are sparse and dense graphs, (though most graphs are sparse?) For

We will distinguish between instances for which the optimal value is known and those for which the optimal value is not known. In the analysis we will consider three indicators for a given instance. The smallest and biggest returned cut value over the 21 solvers. Lastly we will investigate the standard deviation, which indicates how sensitive the heuristic is to different choices of parameters.

6.2 Instances with known optimal value

There are 332 instances with known optimum value. Out of those 146 instances are solved optimally by each of the 21 solvers. Furthermore there are 154 instances for which at least one of the solvers but not all finds the optimal value. Lastly there are 32 instances for which none of the 21 solvers finds the optimal value. For these instances we first take a look at the highest returned cut value ratio over all the solvers. Doing this wee see that there are only two instances namely gka10b 0.863636 and gka9b with 0.875912 for with the maximal ratio is less than 0.878. Other than those every other instance reaches a maximal ratio of greater than 0.99. Regarding the minimum the same phenomen occurs. For the two instances gka10b and gka9b the minimal ratios are 0.000000 and 0.277372 respectively. For the other instance the minmal ratio is greater than 0.97 exceeding 0.878. This means that for 30 out of those 32 instances every solver returns a solution satisfying the performance guarantee of the Goemans-Williamson algorithm, although there is no theoretical guarantee to do so for the Burer heuristic, see section 4. For the standard deviations we have much of the same. The standard deviation for gka9b and gka10b is with 0.331967 and 0.154788 respectively orders of magnitude larger than for the rest which have stadard deviation less or equal to 0.004201.

We will take a deeper look at the instances gka9b and gka10b, starting with gka9b: First of all we notice that the quality of the returned cut strongly varies and that it keeps oscilating even for big parameters. This is very interesting as usually the quality of the solution funnels towards high quality solutions for increasing parameters. Furthermore for 5 solvers the returned cut has the value of the trivial cut. Notably 4 out of those 5 solvers have N = 5.

We observe that both instances gka9b and gka10b arose by transformation from a BQP to MCP. The transformed graphs are dense with density 1 and the maximum cut values are within the order of magnitude of the edge weights, except for the edges connected to the sun vertex. It needs to be mentioned, that the absolute value of the edges connected to the sun vertex have a much higher weight, this stems from the transforming the BQP to a MCP. However, the instances gka*b with $* \in [1:8]$ the same characteristics hold true but their performance is much better. Another explanation might be that the Therefore the previous observations can not be a thorough explanation of the performance.

6.3 Instances without known optimal value

Although there is no optimal value we are still interested in the quality of a solution as a ratio. For each instance we therefore normalise the returned values by dividing by the maximal returned value. By construction the values will be between 0 and 1 with at least one value being 1. Performing a similar analysis as above

6.4 Results

The experiments gather further data that the Burer heuristic is highly effective in returning high quality approximations. Out of the 332 test instances with known optimum value, the instances for which the heuristic returns a cut with ratio less than 0.878 (the performance guarantee in Goemans Williamson) are so few we can name them here. They are name the instances In particular the heuristic return high quality cuts for instances with mixed sign weight, these are This is very impressive, as the performance guarantee of the Goemans Williamson algorithm only holds for graphs with nonnegative weights. Therefore there is no theoretical reason for the approximation algorithm to perform well, let alone the Burer heuristic.

We point out that the standard deviation of the returned cut values of the Burer heuristic with different parameters is very low in general.

The data indicates, that if a run with big parameters, i.e. expecting longer runtime finds a very good solution, then a high quality solution can already be found with small values.

The data clearly suggests that using the Burer heuristic to generate high quality approximate solutions is highly effective.

Looking at this from the perspective that the optimal value is not known, the following procedure might be an informed choice:

Run the Burer heuristic for very small parameters, i.e M=1 and N=5 and gradually increase them. If the approximate solution barely changes there is a good chance that it will not change for bigger parameters (within reason) either. Therefore we can accept the returned solution as an approximation. However, if there is a lot of deviation in the solutions, then the heuristic should be rerun for bigger parameters as this might lead to better solutions.

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