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Problem

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ALGORITHMS AND TESTS FOR THE COLOURFUL FEASIBILITY PROBLEM

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ABSTRACT. We study a colourful generalization of the linear programming feasibility problem, comparing the algorithms introduced by Bárány and Onn with new methods. This is a challenging problem on the borderline of tractability, its complexity is an open question. We perform benchmarking on generic and ill-conditioned problems, as well as recently introduced highly structured problems. We show that some algorithms can lead to cycling or slow convergence and we provide extensive numerical experiments which show that others perform much better than predicted by complexity arguments. We conclude that the most efficient method is a proposed multi-update algorithm.

1. INTRODUCTION

Given colourful sets S_1, \dots, S_{d+1} of points in \mathbb{R}^d and a point p in \mathbb{R}^d , the *colourful linear programming* problem is to express p as a convex combination of points x_1, \dots, x_{d+1} with $x_i \in S_i$ for each i . This problem was presented by Bárány and Onn in 1997 [BO97b], it is still not known if a polynomial-time algorithm for the problem exists. The monochrome version of this problem, expressing p as a convex combination of points in a set S , is a traditional linear programming feasibility problem.

In this paper, we study algorithms for colourful linear programming with a core condition from an experimental point of view. We learn several things. First, in our experience this problem is easy – we expend more effort to generate difficult examples than to solve them. Second, while the classical algorithms for this problem already perform quite well, we introduce modifications that achieve a substantial improvement in practical performance. Third, we construct examples where ill-conditioning leads to slow convergence for the some otherwise very effective algorithms. And finally, we remark that a simple greedy heuristic provides competitive results in practice but we find a case where it fails to solve the problem at all. Additionally we provide benchmarking that we hope will encourage research on this attractive problem.

2. DEFINITIONS AND BACKGROUND

We call a system of $d + 1$ sets of $d + 1$ points a *configuration*, and often denote it as $\mathbf{S} = \{S_1, \dots, S_{d+1}\}$. Such configurations are the simplest non-degenerate cases of colourful linear programming. We define the *core* of a configuration to be $\bigcap_{i=1}^{d+1} \text{conv}(S_i)$. In this paper we consider the *colourful feasibility* problem of expressing a given p in the interior of the core as a colourful convex combination of points in the configuration. By Bárány’s colourful Carathéodory theorem [Bár82], a solution is guaranteed to exist, and the problem is to exhibit one. This problem is described in [BO97b] as “an outstanding problem on the border line between tractable and intractable problems”.

Several close relatives of the colourful feasibility problem are known to be difficult. For example, the case where we have d colours in \mathbb{R}^d and no restriction on the size of the sets has

been shown to be strongly NP-complete through a reduction of 3-SAT. We refer to [BO97b] for more details.

In [Bár82], Bárány proposed a finite algorithm **A1** to solve colourful feasibility, and in [BO97b] Bárány and Onn analyzed the complexity of **A1** and a second algorithm **A2**. (See Section 3 for a detailed description of these and other algorithms.) Both these algorithms are essentially geometric, and the complexity guarantees depend crucially on having the point p in the *interior* of the core. In effect, the distance between p and the boundary of the core can be considered as a measure of the conditioning of the problem. Thus for a configuration \mathbf{S} we define ρ to be the radius of the largest ball around p that is contained in the core. The results for **A1** and **A2** are effectively that they are polynomial in d and $1/\rho$. While this is not polynomial in the input, it suggests that a polynomial algorithm may be possible. We remark that for configurations of $d + 1$ points in $d + 1$ colours on the unit sphere $\mathbb{S}^d \subseteq \mathbb{R}^d$, ρ will be small even if the problem has a favourable special structure, and quite small otherwise.

It is helpful to preprocess the problem by translating the point p to be the vector $\vec{0}$ in \mathbb{R}^d . If $\vec{0}$ is a point in one of the S_i 's, then the solution to the colourful feasibility problem is trivial. Otherwise, we can also scale the points of the S_i 's so that they lie on the unit sphere \mathbb{S}^d . The coordinates in any resulting convex combination can then be unscaled as a post-processing step.

We remark that restricting the sets to have size $d + 1$ is not a burden since, given a larger set, solving a monochrome linear feasibility problem allows us to efficiently find a basis of size $d + 1$ with $\vec{0}$ in its convex hull.

The colourful feasibility problem models a data mining situation where we want to select a set of points that is both diverse, in the sense that it includes representatives from predetermined classes (colours), and representative, in the sense that the selected points surround a specified point common to all the classes [Lu06]. Application of this problem to combinatorics are discussed in [BO97a].

3. SEVEN ALGORITHMS

In this paper we consider the theoretical and practical performance of seven algorithms for finding a colourful basis. The algorithms considered are the algorithms of Bárány **A1** and of Bárány and Onn **A2**, modifications of these algorithms which update multiple colours at each stage, which we will call **A3** and **A4** and a hybrid **A5** of these designed to take advantage of the strengths of both algorithms. For purposes of comparison, we also consider two simple approaches that perform well under certain circumstances: a greedy heuristic where we choose the adjacent simplex of maximum volume **A6** and a random sampling approach **A7**. All our implementations are initialized with the first points from each colour. Following are descriptions of the algorithms, see [Hua] for MATLAB implementations of each. Besides **A7**, they are implemented as pivoting algorithms with the respective pivot selection rule.

3.1. Bárány's Algorithm A1. We begin with the algorithm proposed by Bárány [Bár82], which is a pivoting algorithm. It begins with say a random colourful simplex Δ . The point x nearest to $\vec{0}$ in Δ is computed. If $x \neq \vec{0}$, then x must lie on at least one facet of Δ . Consider the colour i of the vertex of Δ that is not on this facet. Look for the point t of colour i minimizing the inner product $\langle t, x \rangle$. Then we replace the point of colour i from Δ with the point t to get a new simplex. The algorithm then repeats beginning with the new simplex.

The convergence of this algorithm relies on the fact that $\vec{0}$ is in the core of the configuration. For this reason the affine hyperplane perpendicular to the vector x cannot separate $\vec{0}$ from the points of colour i . Thus the next simplex will have a point closer to $\vec{0}$ than Δ did, and the algorithm will converge in finitely many steps. If, additionally, the core has radius at least ρ around $\vec{0}$, then there is a guarantee a given step will decrease the squared norm of the nearest point by at least a factor of $(1 - \rho^2/4)$. Using this, it is possible to show that **A1** will approach the solution in $O(1/\rho^2)$ iterations. Since an iteration can be done in polynomial time, this proves that **A1** runs in time polynomial in the input data and $1/\rho$. Consult [BO97b] for details and a proof.

We note that the complexity of a single iteration is dominated by the cost of the nearest point subroutine. This can be solved as a continuous convex quadratic optimization problem, but involves numerical issues: It can be solved to less or greater precision, either risking numerical error or increasing the running time. For the purposes of our benchmarking, we used the MATLAB built-in `quadprog()` which gave fairly good results, see Section 5.2.

3.2. Bárány and Onn’s Algorithm A2. The reliance of **A1** on nearest point calculations is a disadvantage. Partly motivated by this, Bárány and Onn proposed an alternate algorithm for the colourful feasibility problem whose calculations involve only linear algebra. This algorithm, **A2**, is described in [BO97b].

The key idea is to replace the closest point x to $\vec{0}$ on the simplex Δ by a point y on the boundary of Δ that can be computed algebraically. The initial choice of y could be one of the vertices of the initial simplex. In subsequent iterations, a colour j corresponding to a zero coefficient in y is chosen. An improving vertex v of colour j is found, and y_{new} is updated by projecting $\vec{0}$ onto the line segment between y and v and finding where the resulting vector enters the new simplex. As with **A1**, this algorithm takes $O(1/\rho^2)$ iterations, and hence is polynomial in the input data and $1/\rho$, see [BO97b].

The implementation of **A2** proposed in [BO97b] takes time $\Theta(d^4)$ for a single iteration. The bottleneck is computing y_{new} , which is the intersection of the line segment from $\vec{0}$ to a point p and the new simplex. In fact we observe that this can be done in time $O(d^3)$. First, compute the defining equations for the simplex $Ay_{\text{new}} \geq b$ by inverting the homogenized matrix of the vertices. We know the intersection point will be of the form $y_{\text{new}} = \alpha p$. We can substitute this into the above inequalities to get $\alpha(Ap) \geq b$ and simply take α to be the maximum value of $b_i/A_i p$ for $i = 1, 2, \dots, d + 1$. This is implemented in [Hua].

As noted by Maurice Queyranne, it is possible to modify **A2** to compute the nearest point on the simplex using Wolfe’s algorithm for finding the nearest point on a polytope [Wol76]. While it does not have a polynomial time guarantee, it may work well for this problem. Like **A2**, Wolfe’s algorithm uses simple linear algebra to pivot through faces; it could be adapted to use y_{new} as a warm start.

3.3. Multi-update Bárány A3. We propose the following modification of **A1**: if it happens that the nearest point x to $\vec{0}$ of the current simplex Δ lies on a low-dimensional face of Δ - i.e., on more than one facet - then we update *every* colour that is not a vertex of that face. After finding each new point, we replace x by x_{new} , the projection of $\vec{0}$ onto the line segment from x to the vertex we are adding to the simplex. The advantage of this new algorithm, which we call **A3**, is that when possible it updates several colours without recomputing a nearest point.

Since this algorithm makes at least as much progress as **A1** at each iteration, we get convergence in at most the same number of iterations. A given iteration may take longer, since it has to update multiple points. However, aside from the nearest point calculation,

all steps in an iteration of **A1** can be performed in $O(d^2)$ arithmetic operations. Hence the additional work per iteration of **A3** is $O(d^3)$, and the bottleneck remains the single nearest point calculation.

3.4. Multi-update Bárány and Onn A4. Similarly, we can adjust algorithm **A2** to pivot multiple colours when y lies on a low-dimensional face. As in **A3** we update y by setting y_{new} to the projection of $\vec{0}$ onto the line from y to the new vertex. This is faster than the computation of y from **A2** at the end of the iteration, which remains the bottleneck. We call this algorithm **A4**. It is particularly useful at the start of the algorithm since the initial point y is a vertex of Δ . This algorithm will take no more iterations than **A2**, and each iteration costs at most a constant factor more than an iteration of **A2**.

3.5. Hybrid A5. In Section 5 we describe a situation where **A2** and **A4** are slow because they repeatedly return to the same simplex, see the example in Section 6.1. A practical solution to this is to run **A4**, but use a computationally heavy step from **A3** if we detect that **A4** is returning to the same simplex. We implemented such a hybrid algorithm **A5**.

3.6. Maximum Volume A6. We also considered the performance of some greedy heuristics. The most effective of these was to pivot from Δ to an adjacent simplex of maximum volume given that the pivoting hyperplane separates Δ from $\vec{0}$. This heuristic, which we call **A6**, uses simpler linear algebra than **A2**, and by taking large simplices often gets to $\vec{0}$ in a small number of steps. We can perform an iteration of this algorithm in $O(d^4)$ time.

3.7. Random Sampling A7. Finally, we remark on a very simple guess and check algorithm where we sample simplices at random and check to see if they contain $\vec{0}$. Intuitively we would not expect such an algorithm to work well. However, as discussed in [DHST06] solutions to a given colourful feasibility problem may not be all that rare, and in some cases can be quite frequent. Since guessing and checking are relatively fast operations, it is worth considering the possibility that this naive algorithm may perform well in special cases or low dimension. We call this algorithm **A7**.

One attractive feature of **A7** is that the cost of an iteration is low – we only have to generate a random simplex and then test if it contains $\vec{0}$. The test can be done in $O(d^3)$ time by solving a linear system.

4. RANDOM, ILL-CONDITIONED AND EXTREMAL PROBLEMS

To better understand how various algorithms perform in practice, we produced a test suite of challenging colourful feasibility problems, which includes unstructured random problems, ill-conditioned problems and problems with a restricted number of solutions. In this section we describe three types of colourful feasibility problems that we consider when evaluating the practical performance of an algorithm. See [Hua] for a MATLAB implementation of each of these problem generators.

4.1. Unstructured Random Problems. The first class of problems we consider are unstructured random problems. We take $d + 1$ points in each of $d + 1$ colours on \mathbb{S}^d . The only restriction we require is that $\vec{0}$ is in the core. This is achieved by taking the last point to be a random convex combination of the antipodes on \mathbb{S}^d of the first d points. We call this generator **G1**.

4.2. Ill-conditioned Random Problems. Next, we consider ill-conditioned problems. We place d points of a given colour on the spherical cap around the point $(0, 0, \dots, 0, 1)$ and the final point of that colour in the opposite spherical cap, again as a convex combination of the antipodes. The maximum angle between a chosen vector and the final coordinate axis is a parameter, and points are concentrated towards the centre rather than uniformly distributed on the cap. Since the points all lie in a tube around the final coordinate axis, we call these *tube* generators. We implemented two tube generators: **G2** randomly places either 1 or d points of colour i on the positive side of the axis, while **G3** always places d points of colour i on the positive side of the axis.

4.3. Problems with a Restricted Number of Solutions. Finally, we consider problems where we control the number of colourful simplices containing $\vec{0}$. The paper [DHST06] provides new bounds for the number of possible solutions to a colourful feasibility problem with $\vec{0}$ in the interior of the core. It turns out that the number of simplices containing $\vec{0}$ in dimension d can be as low as quadratic in d , but not lower, see [BM07] and [ST05], or as high as $d^{d+1} + 1$ (with $\rho > 0$), which is more than one third of the total number of simplices. Constructions are given for colourful feasibility problems attaining both these values.

The probability that a simplex generated by $d+1$ points chosen randomly on \mathbb{S}^d contains $\vec{0}$ is $1/2^d$, see for example [WW01]. Thus in a uniformly generated random problem of the type generated by **G1**, we would expect about $1/2^d$ of the $(d+1)^{d+1}$ colourful simplices to contain $\vec{0}$. This is not a large fraction, but in the context of an effective pivoting algorithm such as **A1** which may pivot several neighbours to a given solution, and pivot several neighbours of the first neighbour onto it, etc., we can entertain the idea that for a random configuration most simplices are close to a solution. See Section 6.4 for further discussion.

We might expect that the difficulty of a colourful feasibility problem increases as the number of solutions, i.e. simplices containing $\vec{0}$, decreases, so we wrote three problem generators based on the constructions in [DHST06]. The first, **G4** generates perturbed versions of the configuration from [DHST06] with many solutions. These problems have $d^{d+1} + 1$ of the $(d+1)^{d+1}$ simplices containing $\vec{0}$, many more than random configurations, and we expect them to be quite easy. The second, **G5**, generates configurations where one point of each colour is close to each vertex of a regular simplex on \mathbb{S}^d . There are $d!$ solutions corresponding to picking a different colour from each vertex, this is still much less than the $(d+1)^{d+1}/2^d$ expected in a random configuration. Finally, we have **G6**, which generates perturbed versions of the configuration from [DHST06] which has only $d^2 + 1$ solutions. The generators **G4**, **G5** and **G6** randomly permute the order of the points that appear within each colour.

All these problems are ill-conditioned in the sense that points are clustered closely together. Also ρ will be quite small for **G4** and **G6**, although the construction **G5** maximizes ρ for configurations on \mathbb{S}^d , with $\rho = 1/d$.

5. BENCHMARKING AND RESULTS

In this section, we describe the results of computational experiments in which we run the colourful feasibility algorithms against our problem generators. We focus on the number of iterations that an algorithm takes to find a solution, but in Section 5.2 we also include information about the cost of iterations. The two particularly difficult, but fragile, examples of Sections 6.1 and 6.2 are not included in these results.

5.1. Iteration Counts. For each type of problem we ran tests of the algorithms in dimensions 3×2^n for $n = 0, 1, 2, 3, 4, 5, 6, 7$. Dimension 3 is our starting point since the seven algorithms degenerate to three simple and effective algorithms in dimension 2. We use the

factor 2 increase to sample higher dimensions with less frequency as we get higher. We believe this yields a reasonable sample of low, intermediate and high dimensional problems.

Note that a colourful feasibility problem instance in dimension d consists of $(d+1)^2$ points in dimension d . Thus the size of the input is cubic in d . At present it is logistically difficult to generate and store a colourful feasibility problem in dimension $d = 1000$. After dimension 100, it also becomes increasingly difficult to cope with numerical errors, especially for the algorithms that include nearest point calculations, namely **A1**, **A3** and **A5**. For this reason we do not include results for these algorithms beyond $d = 96$ for except for the relatively well-conditioned **G1** problems where we stopped at $d = 192$. As one would expect, the guess-and-check algorithm **A7** performs badly as d increases, except on problems from the **G4** generator which have an abundance of solutions. We only include results from the **A7** algorithm when they can be completed in a reasonable amount of time.

The results of our computational experiments are presented in the graphs below and the tables in Appendix C. Each graph presents results for a single random generator on a log-log scale with the average iteration count of each algorithm plotted against the dimension. Additionally, the tables contain the values of the largest iteration count observed in each type of trial; these show similar trends to the averages, although we notice that **A2** and **A4** sometimes perform substantially worse than the average, especially in the presence of ill-conditioning. The reasons for this are discussed in Section 6.2.

For each generator at $d = 3$ we sampled 100,000 problems, at $d = 6$ and $d = 12$ we sampled 10,000 problems, at $d = 24$ and $d = 48$ we sampled 1,000 problems and finally for $d \geq 96$ we sampled 100 problems. Because of the varying sample sizes, it may not be entirely fair to compare the maxima listed in Appendix C between dimensions. The results are plotted on as log-log graphs in Figures 1–6. We remark that polynomials appear asymptotically linear in log-log plots, with the slope of the asymptote being the exponent of the leading term of the polynomial and the y -intercept of the asymptote representing the lead coefficient.

In Figure 1 we see that **A1** and **A2** appear to be taking a polynomial number of iterations to solution, while **A6** and **A7** do not appear to be polynomial. Since each algorithm takes a polynomial time per iteration, the graphs of time versus dimension show similar trends.

For tube experiments **G2** and **G3**, we used an angle parameter of $\pi/6$, that is, all the vectors in the configuration made an angle of at most $\pi/6$ with the x -axis.

The tube experiments summarized in Figures 2 and 3 show the impact of ill-conditioning on all the algorithms. For **A1**, **A3**, **A5** and **A6**, convergence is slightly slower and numerical errors become more common. With these algorithms, our experiments began to crash at dimension 192. By contrast for the better conditioned problems from **G1**, the three algorithms with minimum distance calculations crashed only at dimension 384 and **A6** would in any case take too long on problems of this size. Nevertheless, these algorithms remain effective at $d = 96$.

The algorithms **A2** and **A4** are more robust in the sense that they are not as prone to crashes due to numerical errors. This is the advantage of relying entirely on straightforward linear algebra computations rather than considering nearest points or volumes. At the same time, they converge much more slowly due to problems of the type described in Section 6.2 and Appendix 6.2.

If we decrease the angle parameter which controls the width of the tube and hence the conditioning, all the algorithms become less stable numerically and experience a degradation in performance. In the cases of **A2** and **A4** they become substantially slower.

We comment that the **A7** algorithm performs about the same on **G2** problems as it did on **G1** problems. This indicates that **G2** problems typically have a similar number of solutions to **G1** problems. As one would expect, solutions to the one-sided tube problems generated

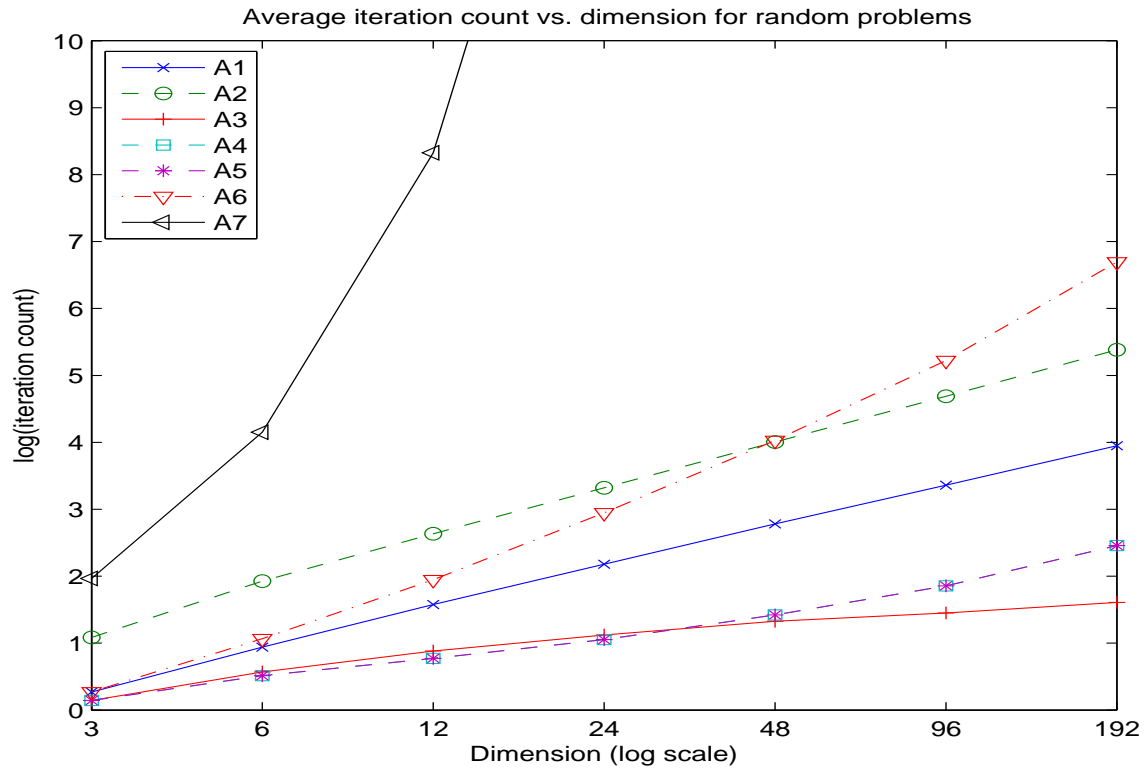


FIGURE 1. Results for **G1**.

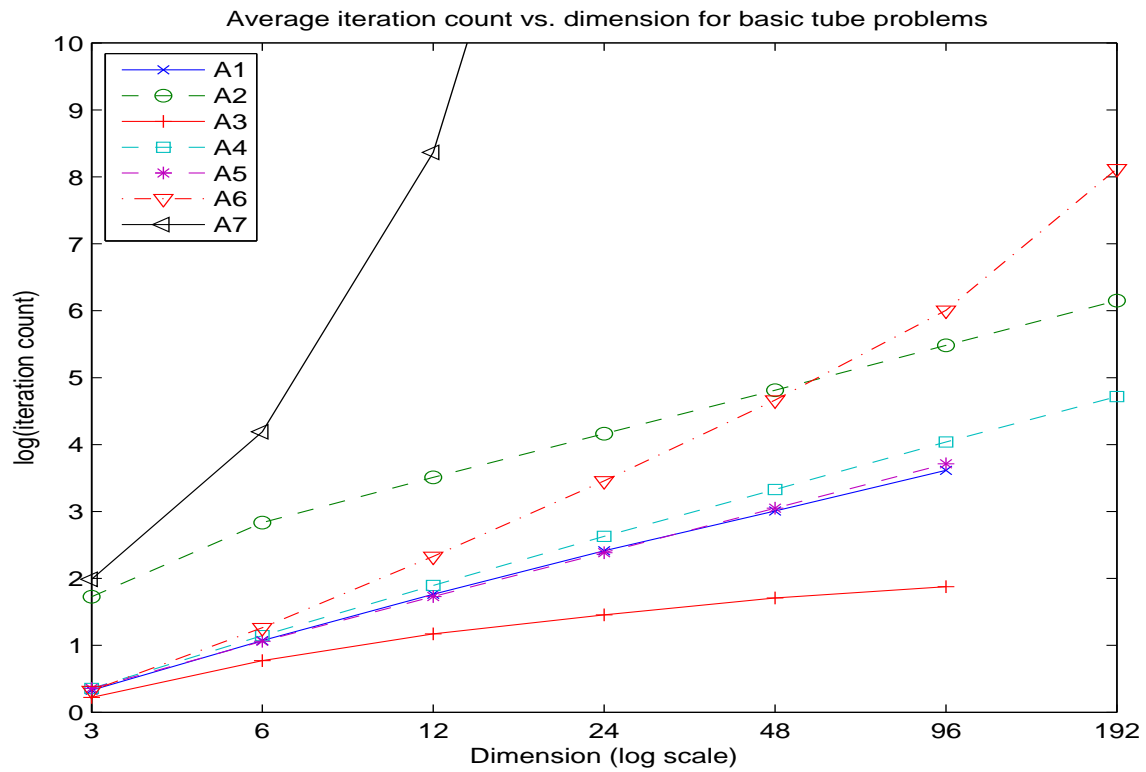
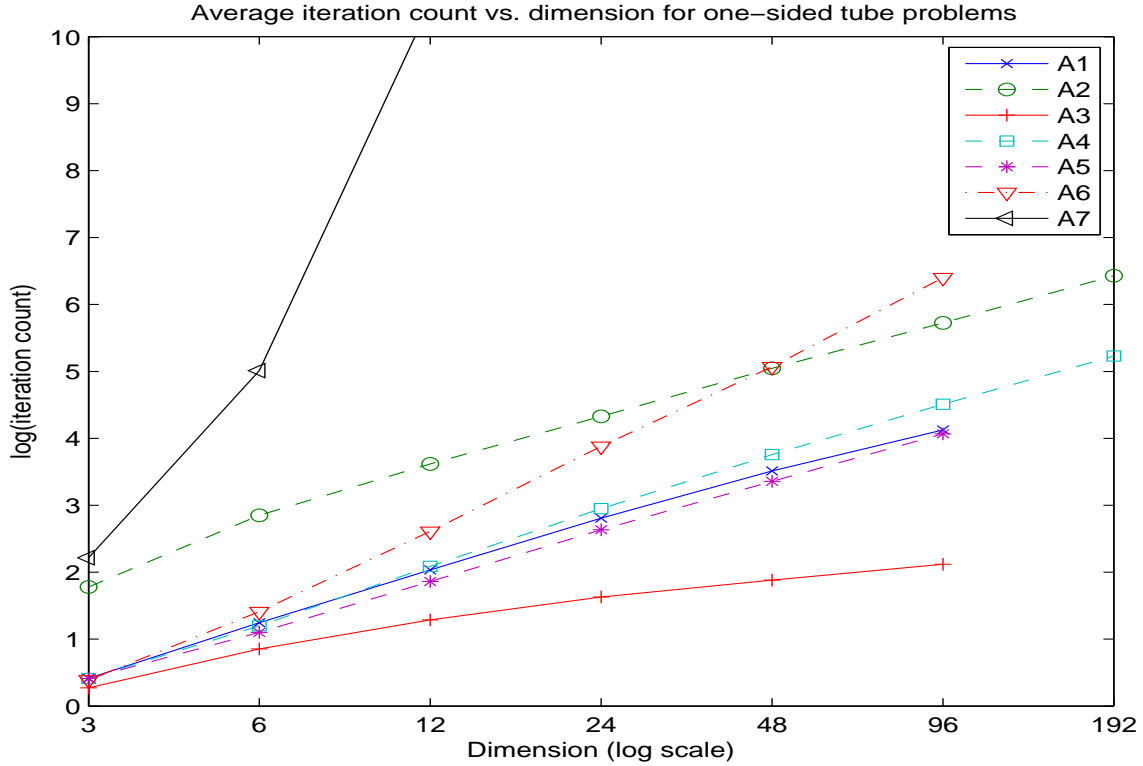


FIGURE 2. Results for **G2**.

FIGURE 3. Results for **G3**.

by **G3** are rarer than solutions to **G1** and **G2** problems since the most of the points are clustered on one side. Hence **A7** performs much worse on this type of problem.

The problems with many solutions produced by **G4** are solved very quickly by all the algorithms, as illustrated in Figure 4. In this case the random sampling algorithm **A7** offers excellent performance. With the abundance of solutions, most of the algorithms solve such problems in an expected constant number of iterations. The exception is **A2** which needs $\Theta(d)$ iterations at the start to unwind the nearest point substitute y from a vertex to an interior point on a facet. Since all the algorithms begin by checking the feasibility of the initial simplex, the **G4** problems are often solved in 0 iterations.

For the simplex structured problems of **G5**, we see all the algorithms except **A7** perform very well, despite the relative scarcity of solutions. We see that the other algorithms have exactly the proper response to this structure – they systematically take points near vertices that are not part of the current set. In the case of **A1**, a new vertex of the simplex will be added at each step to give convergence in at most d iterations, for **A2** it takes one pass through the $d + 1$ colours, and for the multi-update algorithms **A3**, **A4** and **A5** one or two passes through the colours. Algorithm **A6** also solves these problems in a reasonable number of iterations.

Finally, we see that the problems from **G6** where solutions are scarce are indeed more difficult than random problems, but that, except for the **A7** algorithm, the impact on algorithmic performance is mild. See Figure 6. Curiously, the **G6** problems are the most difficult problems for the **A1** algorithm. The multi-update algorithms **A3**, **A4** and **A5** perform extremely well.

5.2. Cost per Iteration. In Figure 7 we present the average iteration times observed for all seven algorithms on problems from the **G1** generator. The raw data for this graph is in

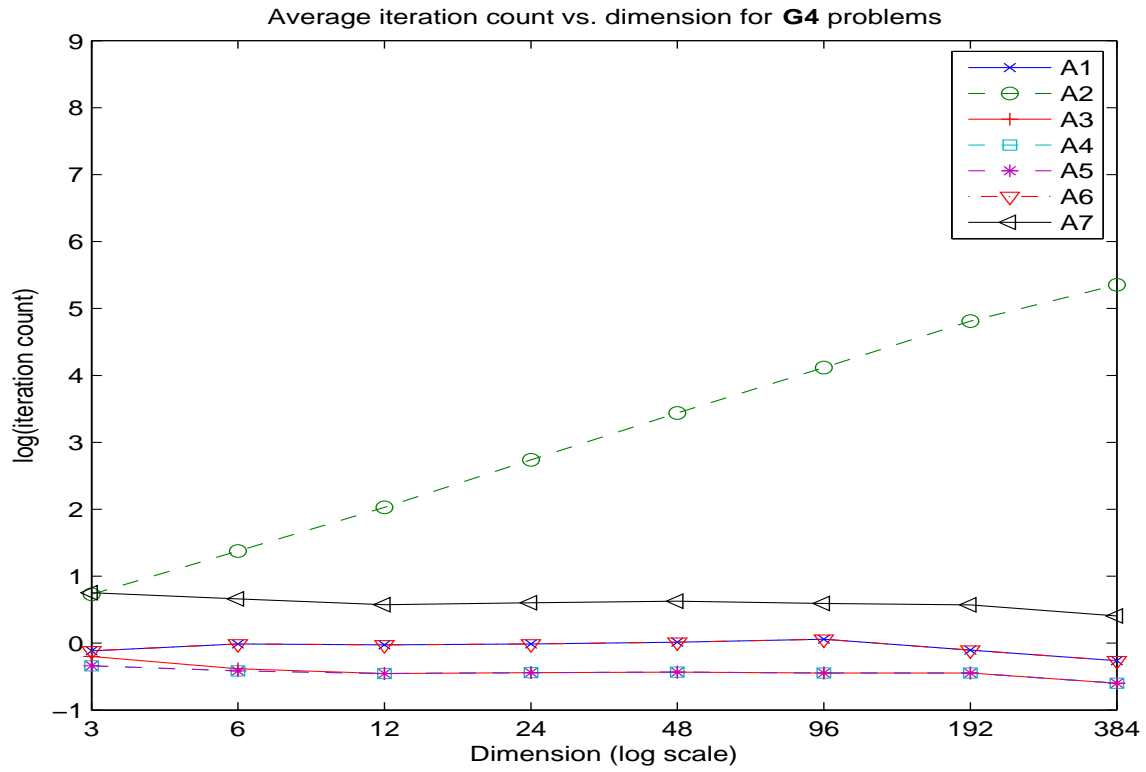


FIGURE 4. Results for **G4**.

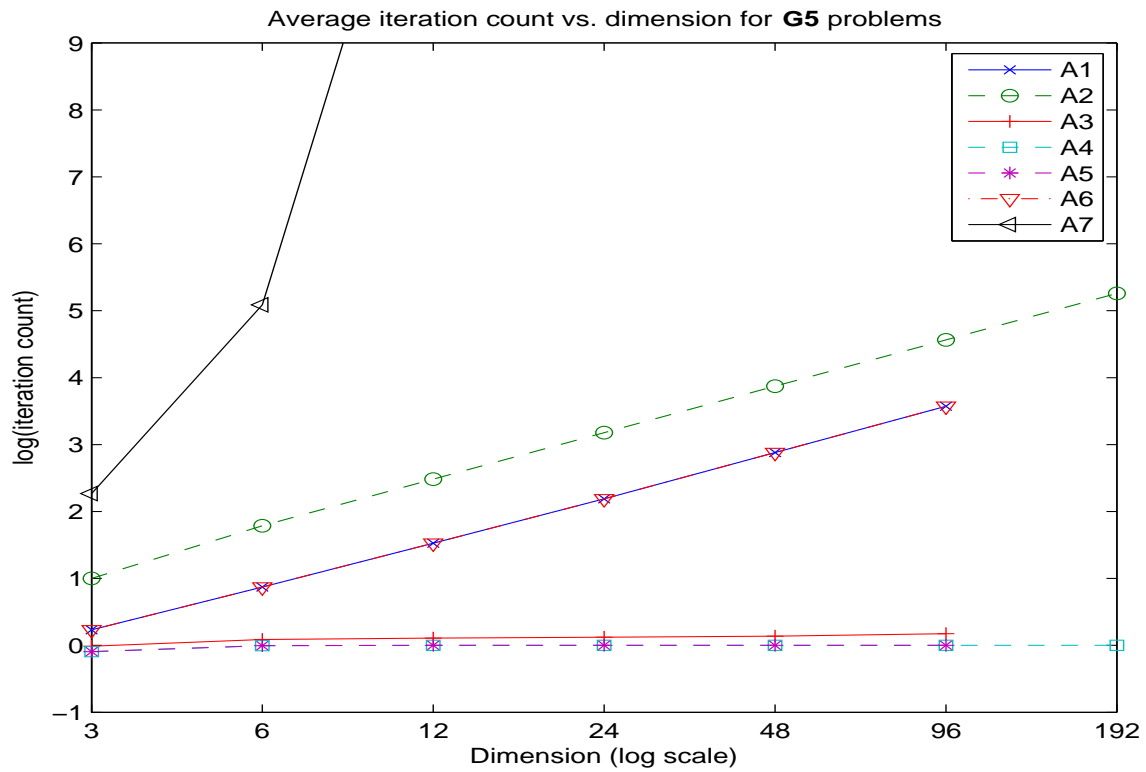
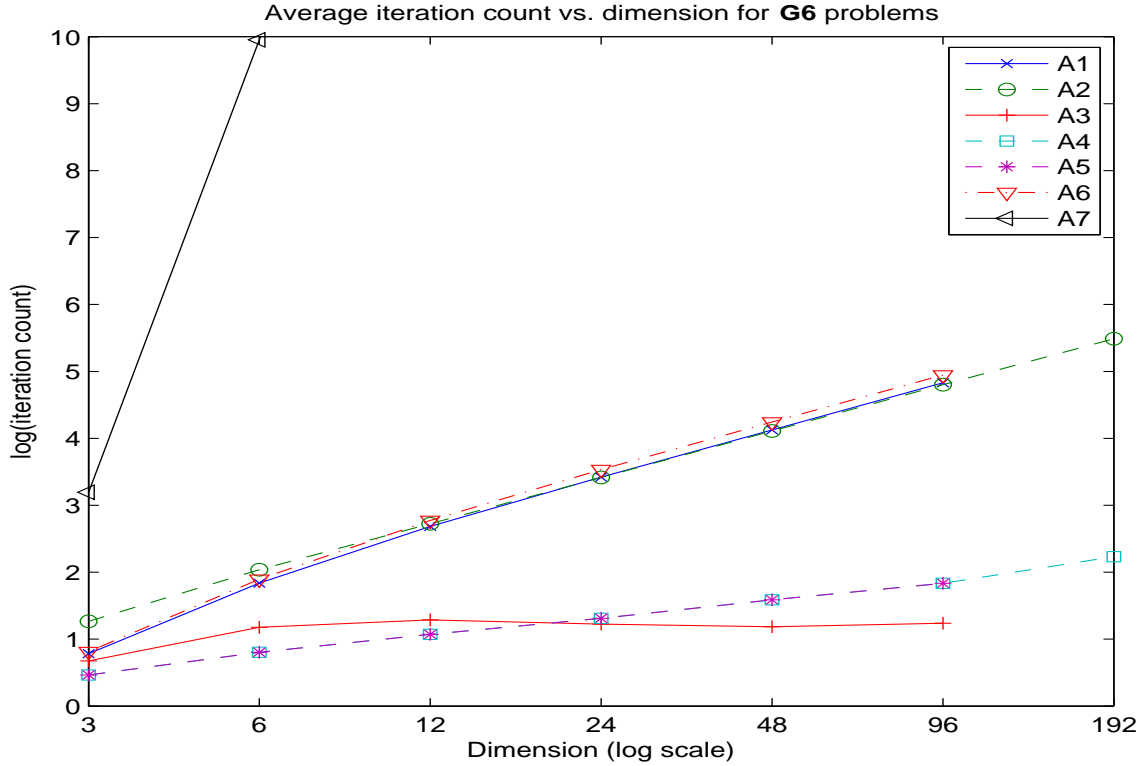


FIGURE 5. Results for **G5**.

FIGURE 6. Results for **G6**.

Appendix D. We comment that the average time to complete an iteration does not change significantly with the problem type, so we have not included the similar graphs for other generators. The data shows that in our implementation of these algorithms, the average time for an iteration is never very large. For the slowest algorithms in the highest dimensions the average iteration took less than 2 seconds.

We see some interesting trends in the graphs. In low dimensions all the iteration times are very fast and are presumably dominated by fixed startup costs. As the dimension increases, we begin to see the asymptotic behaviour. The algebraic algorithms **A2** and **A4** show the expected $\Theta(d^3)$ behaviour, which appears linear in the log-log plot. Asymptotically, the average time for an iteration of **A4** is about 10 times longer for an iteration of **A2**.

The algorithms **A1** and **A3**, which depend on a minimum distance calculation, take longer on average to complete an iteration than **A4**. The extra cost for the multiple updates in **A3** is relatively small. However, the asymptotic slope of these lines appear higher than for **A2**, which means that the nearest point calculations are causing the iterations to take time $\Omega(d^3)$. The algorithm **A6** has iteration times not much worse than **A2** in low dimension, but its asymptotics look close to $O(d^4)$ as suggested in Section 3.6. Algorithm **A7** exhibits $\Theta(d^3)$ iteration time and is asymptotically about twice as fast on average per iteration than **A2**.

Unlike the other algorithms, the average iteration time for **A5** will be substantially affected by the conditioning of the problem. Using the well-conditioned **G1** problems, **A5** usually degenerates to **A4** and has a very similar average iteration time. As the problems become more ill-conditioned, **A5** will begin to use **A3** steps as well, and the average iteration time will increase towards the average iteration time for **A3**.

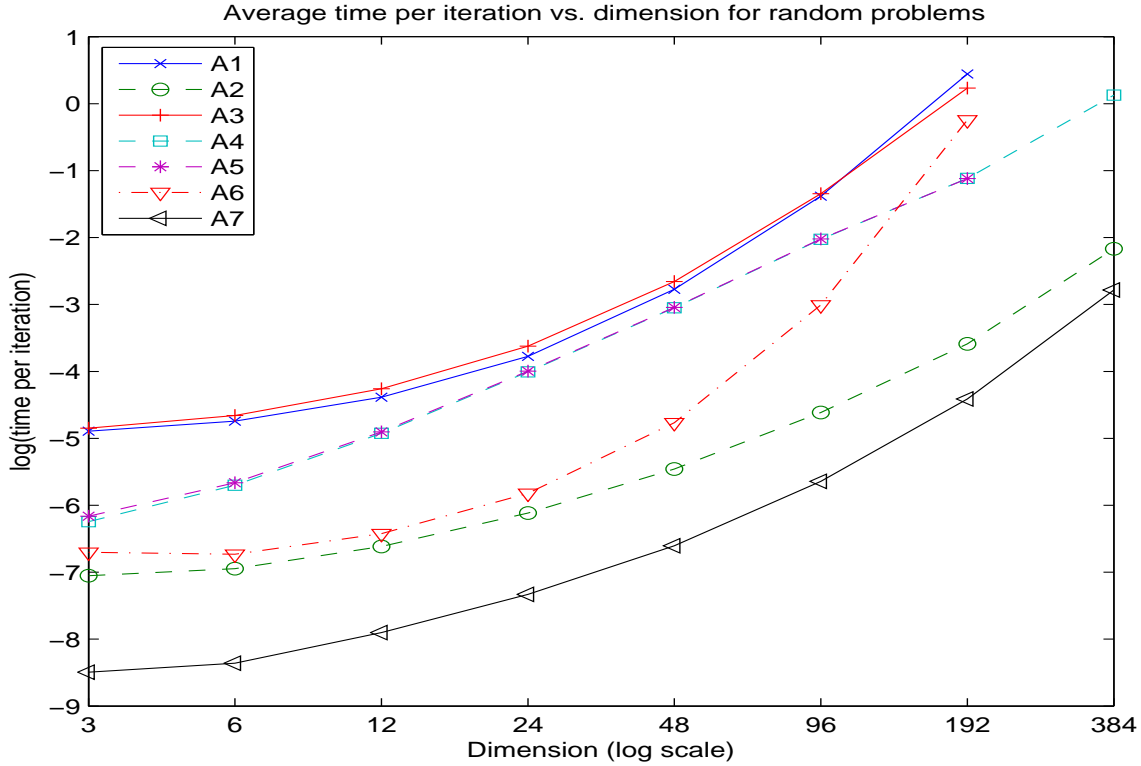


FIGURE 7. Average iteration time of the algorithms.

6. DISCUSSION AND WORST-CASE CONSTRUCTIONS

Our experiments reveal several features of colourful feasibility algorithms. After considerable searching, we found a problem instance which caused **A6** to cycle. We also found that **A2** and **A4** can converge extremely slowly in the face of ill-conditioning although **A1** and **A3** continue to perform reasonably well on the same examples. We conclude that computationally the best algorithms are the multi-update variants and remark that these tightened algorithms do yield substantial gains over the originals.

6.1. A Cycling Example for A6 in Dimension 4. In Appendix A we exhibit an example in dimension 4 for which the maximum volume heuristic cycles. Since this example shows that **A6** can cycle, it is remarkable that it happens so rarely. It did not occur in the entire test suite of Section 5. We were unable to find any examples of cycling in dimension 3 or any examples of cycling in dimension 4 with cycle length shorter than 6. Higher dimensions and longer cycle lengths do occur.

One explanation for the results is that as one might expect, **A6** is an effective heuristic in a typical situation. The distinguishing feature of the few bad examples is that the points are placed in such a way that the simplices cluster into a few groups of similar shape and volume. The heuristic of taking the maximum volume is then not very helpful in choosing promising simplices. We note that this example is solved easily by the other algorithms.

6.2. Flip-flopping During Convergence for A2: 40,847 Iterations in Dimension 3. We constructed an example of a colourful feasibility problem in dimension 3 that takes 40,847 iterations to solution using a basic implementation of **A2**. The exact points we used are contained in Appendix B. The algorithm is initialized with the simplex that uses the first point of each colour. At the fifth iteration, the algorithm reaches a situation where

the current point y lies on a facet F of colours 2, 3 and 4 very close to $\vec{0}$. Using this point the algorithm will pick the point of colour 1 that has minimum dot product with y . The second and third points of colour 1 lie almost in the directions of y and $-y$, however neither of these forms a simplex with F containing $\vec{0}$. In fact the fourth point of colour 1 does form a simplex containing $\vec{0}$ with F , but it is nearly orthogonal to y . As a result, after two iterations, **A2** returns to the same simplex. The point y will be recomputed at each step, and is slightly closer to $\vec{0}$ when the algorithm returns to the previous simplex. However, the improvement is quite small. Of course ρ is also very small, so this is consistent with the performance guarantee described in Section 3.2. The algorithm then proceeds to return to the same simplex more than 20,000 times, with an incremental improvement to y at each iteration before finally taking the fourth point of colour 1 and terminating.

As one would expect with a very ill-conditioned problem, this example is numerically fragile – the current version of our code normalizes the coordinates before starting and does not suffer the same fate. However bad behaviour is fairly typical. The tube generator for ill-conditioned problems in [Hua] produces problems whose ill-conditioning depends on a parameter defining the width of the tube. As the width decreases, we get an increasing number of cases where **A2** and **A4** take enormous numbers of iterations.

We remark that, in contrast, **A1** never returns to the same simplex, so it cannot suffer from this type of flip-flopping. Indeed in dimension 3 it could do no worse than visiting all $4^4 = 256$ simplices. At least 10 of these must contain $\vec{0}$, see [BM07], so the algorithm must terminate in at most 246 iterations. It is quite hard to see how this limit could be approached. The authors wonder if a Klee-Minty-like example, see [KM72], of worst-case behaviour for Barany’s pivoting algorithm could be constructed.

6.3. Advantages of Multiple Updates and Initialization. The multi-update algorithms **A3** and **A4** do provide substantial gains over their single update counterparts, **A1** and **A2**. In the case of **A3**, we get a large reduction in iteration count at very little cost in terms of iteration time. In our benchmarking experiments, this produced times that were competitive with **A2** and much better than **A1**. The gains for **A4** relative to **A2** are less impressive. In our benchmarking experiments, **A4** consistently averaged a 10% to 40% savings in total time to solution.

In fact, **A2** is not as well suited as **A1** to take advantage of multiple updates. The point y close to $\vec{0}$ computed by **A2** will almost always lie in the interior of a facet of Δ , meaning that **A2** will only have a single candidate colour to pivot. In contrast, in high dimension, the closest point x to $\vec{0}$ will often lie on a relatively low dimensional face of Δ , allowing multiple updates throughout the algorithm.

One difficulty for **A2** is that it begins with y at a vertex. In a normal situation, the first d steps of **A2** will each increase the dimension of the smallest face containing y by one until y lies in the interior of a facet, without necessarily yielding a much better current simplex. The multi-update **A4** does this all in the first iteration in less time than it takes **A2** to do d steps.

We have not discussed the effects of the initial simplex in this paper, but we can employ various heuristics to choose a good initial simplex. A few of these are implemented in [Hua]. We found that the most useful initialization heuristic was to run the first iteration of **A4**. This runs in $O(d^3)$ time and improves the subsequent iteration counts of the algorithms, with the obvious exception of **A7**. Even **A4** experiences a reduced iteration count, since the point y found by the initialization is not passed to the algorithm.

6.4. Theoretical Complexity of the Algorithms. In Section 3, we remarked that Bárány and Onn proved a worst-case bound for **A1** and **A2** of $O(1/\rho^2)$ iterations up to numerical considerations and we improved their iteration time for **A2** from $O(d^4)$ to $O(d^3)$. We also mentioned that we do not expect the multi-update and hybrid algorithms to improve the theoretical bounds. From the example of Section 6.1, we see that **A6** is not guaranteed to converge. The expected running time of **A7** is 1 over the probability that random simplex contains $\vec{0}$, i.e. around 2^d for random problems, and as bad as $(d+1)^{d+1}/(d^2+1)$ for the type of problems generated by **G6**.

The poor performance of **A2** on ill-conditioned problems and examples like that of Section 6.2 confirm the worst-case predictions of Bárány and Onn’s analysis. On the other hand, we did not see this type of behaviour for **A1**, and it is hard to see how it could occur.

The model proposed in Section 4.3 is that a pure pivoting algorithm such as **A1**, defines a set of rooted trees on the $(d+1)^{d+1}$ simplices. Each simplex which contains $\vec{0}$ is the root of a tree, and we draw an edge between the vertices representing simplices Δ_1 and Δ_2 if when **A1** encounters Δ_1 it pivots to Δ_2 . Then the worst performance of the algorithm in terms of the number of iterations would be the height of the highest tree. A smart algorithm will produce short trees by pivoting several simplices to a given simplex at a lower level.

Consider a situation where trees have a constant expansion factor k near the base, that is, low level vertices are connected to roughly k vertices in the level above. The number of trees is $p(d+1)^{d+1}$ where p is the probability that a simplex contains $\vec{0}$. If the trees expand up to height h , each tree will contain on the order of k^h vertices. Then we must have $k^h p(d+1)^{d+1} \leq (d+1)^{d+1}$, the total number of vertices. Rearranging, we get $h \leq -\log_k(p)$. This expression predicts the average iteration count for **A1** to grow linearly for **G1** problems, to be constant for **G4** problems and to grow at $\Theta(d \log d)$ for **G6** problems. All of these match very well with our observed results. The **G5** problems are predicted to be more difficult than they are observed to be, but that is not surprising given their simple structure.

7. SUMMARY AND FUTURE WORK

Despite the examples of Sections 6.1 and 6.2, the results presented in Section 5 show that, except for **A7** and to a lesser degree **A6**, all the algorithms did a good job of solving all the problems. We did find that the methods which include nearest point calculations were more vulnerable to numerical errors than **A2** and **A4**, since our implementations began to crash once we got past $d = 100$, especially on ill-conditioned problems. For the most part, the reduced iteration counts of the nearest point algorithms do not offset the extra time spent per iteration compared to **A2** and **A4**. In some cases of extreme ill-conditioning, such as in Section 6.2, **A2** and **A4** will take many additional iterations and be much slower compared to the nearest point algorithms. In this situation either a hybrid algorithm such as **A5**, or the basic **A1** or **A3** would work better.

We had hoped that the hybrid algorithm **A5** would offer the benefits of **A4**, namely speed and robustness in high dimensions, while stopping long periods of flip-flopping from occurring. This did happen to a degree, but in our benchmarking experiments the net time savings were negligible, while **A5** retained **A3**’s tendency to crash due to numerical errors in high dimension.

We finish by returning to the motivating question of Bárány and Onn: Is there a polynomial time algorithm for colourful feasibility? By improving the implementation of **A2**, we have improved the worst case for this algorithm from $O(d^4/\rho^2)$ to $O(d^3/\rho^2)$, however the dependence on ρ has not improved. Indeed our experiments give strong evidence that the analysis for **A2** is tight.

The situation for **A1** is less clear. We do not see the same bad behaviour with ill-conditioned problems that we found for **A2**, so it is possible that a better guarantee exists for this algorithm. In light of the model suggested in Section 6.4 it is quite difficult to see how to construct a Klee-Minty-like bad case for **A1** as discussed in Section 6.2. We view this as an appealing challenge.

8. ACKNOWLEDGMENTS

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APPENDIX A. EXAMPLE IN DIMENSION 4 WHERE **A6** CYCLES

This example consists of 5 points in each of the 5 colours in \mathbb{R}^4 . The points are presented in Table 1. They are grouped by colour, with the rows representing x, y, z and w coordinates, respectively.

Red points				
$7/52$	$1/89$	$-1/60$	$-1/28$	$4/127$
$1/176$	$-8/65$	$5/49$	$6/35$	$9/118$
$4/29$	$1/961$	$-8/191$	$1/40$	$-1/75$
$-\frac{\sqrt{4238906047}}{66352}$	$\frac{\sqrt{30434652805951}}{5559385}$	$\frac{\sqrt{11360296502737439}}{107254140}$	$-\frac{\sqrt{69789743}}{31640}$	$-7\frac{\sqrt{25600756871}}{1123950}$
Green points				
$3/85$	$-5/71$	$8/45$	$3/88$	$-1/114$
$-1/67$	$1/10$	$-38/155$	$-2/131$	$-24/185$
$1/173$	$-2/101$	$1/95$	$3/53$	$7/85$
$\frac{\sqrt{29008089867051134}}{170445655}$	$-\frac{\sqrt{5063381959}}{71710}$	$2\frac{\sqrt{159502559}}{26505}$	$5\frac{\sqrt{14863381455}}{610984}$	$\frac{\sqrt{125498719055}}{358530}$
Blue points				
$-3/77$	$4/141$	$3/22$	$16/111$	$-3/46$
$-3/20$	$-4/63$	$-3/17$	$5/29$	$3/47$
$-2/71$	$-3/173$	$-5/79$	$-1/210$	$1/33$
$-\frac{\sqrt{470161115387}}{694309}$	$-8\frac{\sqrt{122080034994545}}{88619769}$	$-\frac{\sqrt{826050579}}{29546}$	$-\frac{\sqrt{48208184671}}{225330}$	$\frac{\sqrt{5043188147}}{71346}$
Tan points				
$1/59$	$6/151$	$8/45$	$-3/29$	$11/76$
$1/29$	$-1/122$	$-7/32$	$4/43$	$-1/8$
$3/56$	$1/536$	$8/97$	$-1/14$	$9/59$
$25\frac{\sqrt{14625287}}{95816}$	$\frac{\sqrt{554855708771634695}}{745501496}$	$\frac{\sqrt{17827555751}}{139680}$	$-\frac{\sqrt{297327743}}{17458}$	$\frac{\sqrt{75612155}}{8968}$
White points				
$1/167$	$3/43$	$11/52$	$-19/65$	$-3/100$
$1/241$	$-1/244$	$-5/134$	$2/129$	$1/62$
$1/53$	$2/9$	$13/142$	$1/4386$	$-4/73$
$-5\frac{\sqrt{1201121068645021462891}}{173320847963}$	$-\frac{\sqrt{8432767415}}{94428}$	$-\frac{\sqrt{57852799351}}{247364}$	$-\frac{\sqrt{74312211919}}{285090}$	$\frac{\sqrt{50998516979}}{226300}$

TABLE 1. Coordinates of points of an example where **A6** cycles in dimension 4.

The points obtained were initially found as floating point numbers, we rounded them to rational numbers and verified the cycling with rational arithmetic.

The initial simplex is taken to be $(1,1,1,1,1)$, i.e., the first point of each colour. The algorithm proceeds to visit simplices $(1,1,4,1,1)$, $(3,1,4,1,1)$, $(3,1,4,3,1)$, $(3,1,1,3,1)$ and $(1,1,1,3,1)$ before returning to the original simplex and repeating. At steps one, three and five, there are two candidate colours for pivoting, the candidates that are not chosen for pivoting are 1, 3 and 4 respectively. In the even numbered steps there is a single candidate colour for pivoting.

We noted that 170 of the 3125 colourful simplices in this example contain zero, slightly less than the expected $3125/16 \approx 195$. The average volume of a zero containing simplex was about 0.000427, whereas the average volume of a not zero containing simplex was about 0.000201. All the simplices were quite small, the largest had volume about 0.002024. The largest two simplices did not contain zero.

The sequence of simplex volumes seen in our cycling example is: 0.0001035, 0.0001958, 0.0001175, 0.0001350, 0.0001435, 0.0000821.

APPENDIX B. EXAMPLE IN DIMENSION 3 WHERE **A2** TAKES 40,847 ITERATIONS

This example consists of 4 unnormalized points in each of the 4 colours in \mathbb{R}^3 . The points are presented in Table 2. They are grouped by colour, with the rows representing x , y and z coordinates, respectively.

Red points			
1.00000320775369	-0.01000436049274	-0.01000129525998	1.00000089660284
0.00000340785030	0.99999739350954	-1.00000497855619	0.00000051797159
0.00999859615603	0.00000371775824	0.00000030149139	-0.01999639732055
Green points			
1.00000363763560	-0.00999644886160	-0.00999943004295	1.00000335962280
-0.00000325123594	1.00000064545156	-1.00000169806216	-0.00000080450760
0.01000493174811	-0.00000024088601	0.00000009099437	-0.01999811804365
Blue points			
0.99999949817337	-0.00999587145461	-0.00999627213896	0.99999551963712
-0.00000260397964	1.00000485455718	-1.00000419710665	-0.00000024626161
0.00999854691703	0.00000123671997	-0.00000381812529	-0.01999801526314
Tan points			
0.99999980645233	0.10000000280522	-0.60000327600988	0.99999642880542
0.00000024487465	-0.98999719313413	0.79999695643245	-0.00000429109491
0.01000455311709	-0.00000405877812	0.00000372117690	-0.01000272055280

TABLE 2. Coordinates of points of an example taking 40,847 iterations of **A2** in dimension 3.

The initial simplex is taken to be $(1,1,1,1)$, i.e., the first point of each colour. It then updates to $(1,3,1,1)$, $(1,3,2,1)$, $(1,3,2,3)$, $(1,3,2,2)$ and reaches $(3,3,2,2)$ on the fifth iteration. At this point, it begins to flip between $(3,3,2,2)$ and $(2,3,2,2)$ with y initially alternating between values close to $(0.2, \pm 0.00200, 0.00285)$. The values of all these coordinates decrease very slowly as the algorithm continues. At iteration 40,847 it chooses fourth point of colour 1 instead of the third. This makes the current simplex $(4,3,2,2)$ which contains $\vec{0}$.

APPENDIX C. ITERATION COUNTS FROM OUR EXPERIMENTS

In this Appendix we present the raw data from our computational experiments. Each table presents results for a single random generator. The entries give the average number of iterations to solution for each algorithm at the given dimension. For each generator at $d = 3$ we sampled 100,000 problems, at $d = 6$ and $d = 12$ we sampled 10,000 problems, at $d = 24$ and $d = 48$ we sampled 1,000 problems and finally for $d \geq 96$ we sampled 100 problems.

	A1	A2	A3	A4	A5	A6	A7
$d = 3$	1.31	2.96	1.15	1.15	1.15	1.31	7.15
$d = 6$	2.56	6.87	1.77	1.67	1.67	2.90	63.48
$d = 12$	4.84	13.93	2.42	2.16	2.16	7.01	4133.15
$d = 24$	8.84	27.70	3.07	2.87	2.87	19.07	Large
$d = 48$	16.14	54.88	3.77	4.14	4.14	56.12	Large
$d = 96$	28.80	108.71	4.26	6.39	6.39	185.57	Large
$d = 192$	51.96	217.59	4.99	11.68	11.68	808.78	Large
$d = 384$	Unstable	425.26	Unstable	21.63	Unstable	Large	Large

TABLE 3. Average iteration counts in **G1** generator tests.

	A1	A2	A3	A4	A5	A6	A7
$d = 3$	5	136	4	4	4	5	102
$d = 6$	7	21	5	5	5	12	579
$d = 12$	10	30	6	6	6	20	47362
$d = 24$	15	37	6	8	8	43	Large
$d = 48$	22	67	6	9	9	105	Large
$d = 96$	39	120	6	10	10	269	Large
$d = 192$	63	241	7	19	19	1574	Large
$d = 384$	Unstable	472	Unstable	30	Unstable	Large	Large

TABLE 4. Maximum iteration counts found in **G1** generator tests.

	A1	A2	A3	A4	A5	A6	A7
$d = 3$	1.39	5.62	1.25	1.43	1.43	1.38	7.30
$d = 6$	2.92	17.00	2.17	3.14	2.89	3.54	66.02
$d = 12$	5.83	33.48	3.23	6.65	5.64	10.26	4296.66
$d = 24$	11.18	64.30	4.29	13.86	10.86	31.75	Large
$d = 48$	20.24	123.02	5.51	27.91	21.11	106.11	Large
$d = 96$	37.12	240.49	6.54	56.70	40.91	406.10	Large
$d = 192$	Unstable	468.52	Unstable	111.84	Unstable	3367.60	Large
$d = 384$	Unstable	909.82	Unstable	220.50	Unstable	Large	Large

TABLE 5. Average iteration counts in **G2** generator tests.

	A1	A2	A3	A4	A5	A6	A7
$d = 3$	5	4783	4	5	5	6	109
$d = 6$	8	2880	6	44	10	14	1079
$d = 12$	13	842	8	60	14	33	78418
$d = 24$	21	217	9	36	23	78	Large
$d = 48$	31	249	9	55	41	258	Large
$d = 96$	47	323	9	77	76	840	Large
$d = 192$	Unstable	561	Unstable	140	Unstable	11784	Large
$d = 384$	Unstable	1013	Unstable	260	Unstable	Large	Large

TABLE 6. Maximum iteration counts found in **G2** generator tests.

	A1	A2	A3	A4	A5	A6	A7
$d = 3$	1.51	5.93	1.31	1.51	1.51	1.48	9.16
$d = 6$	3.48	17.26	2.35	3.31	3.01	4.10	150.31
$d = 12$	7.64	37.22	3.62	8.06	6.43	13.61	Large
$d = 24$	16.59	75.73	5.11	19.11	13.92	48.51	Large
$d = 48$	33.51	155.48	6.57	42.81	28.70	159.29	Large
$d = 96$	61.97	306.64	8.32	90.98	58.44	602.07	Large
$d = 192$	Unstable	619.55	Unstable	186.86	Unstable	9607.73	Large
$d = 384$	Unstable	1221.43	Unstable	382.10	Unstable	Large	Large

TABLE 7. Average iteration counts in **G3** generator tests.

	A1	A2	A3	A4	A5	A6	A7
$d = 3$	6	2756	5	6	6	6	127
$d = 6$	9	3704	7	38	9	14	1709
$d = 12$	16	689	8	55	16	46	Large
$d = 24$	28	195	9	52	27	124	Large
$d = 48$	50	257	10	83	47	505	Large
$d = 96$	78	374	11	133	83	2023	Large
$d = 192$	Unstable	736	Unstable	226	Unstable	72317	Large
$d = 384$	Unstable	1399	Unstable	454	Unstable	Large	Large

TABLE 8. Maximum iteration counts found in **G3** generator tests.

	A1	A2	A3	A4	A5	A6	A7
$d = 3$	0.89	2.07	0.82	0.71	0.71	0.89	2.12
$d = 6$	0.99	3.96	0.68	0.66	0.66	0.99	1.94
$d = 12$	0.97	7.61	0.63	0.63	0.63	0.97	1.78
$d = 24$	0.99	15.46	0.64	0.64	0.64	0.99	1.83
$d = 48$	1.01	31.15	0.65	0.65	0.65	1.01	1.87
$d = 96$	1.06	61.44	0.64	0.64	0.64	1.06	1.81
$d = 192$	0.90	122.88	0.64	0.64	0.64	0.90	1.77
$d = 384$	0.77	211.20	0.55	0.55	0.55	0.77	1.50

TABLE 9. Average iteration counts in **G4** generator tests.

	A1	A2	A3	A4	A5	A6	A7
$d = 3$	2	5	2	3	3	2	38
$d = 6$	3	7	2	2	2	3	17
$d = 12$	6	12	1	1	1	6	30
$d = 24$	6	24	1	1	1	6	19
$d = 48$	5	48	1	1	1	5	16
$d = 96$	5	96	1	1	1	5	14
$d = 192$	3	192	1	1	1	4	15
$d = 384$	4	384	1	1	1	4	9

TABLE 10. Maximum iteration counts found in **G4** generator tests.

	A1	A2	A3	A4	A5	A6	A7
$d = 3$	1.26	2.72	0.99	0.91	0.91	1.26	9.67
$d = 6$	2.39	5.97	1.09	0.99	0.99	2.39	161.93
$d = 12$	4.61	12.00	1.12	1.00	1.00	4.61	Large
$d = 24$	8.94	24.00	1.13	1.00	1.00	8.94	Large
$d = 48$	17.82	48.00	1.15	1.00	1.00	17.82	Large
$d = 96$	35.58	96.00	1.19	1.00	1.00	35.58	Large
$d = 192$	71.15	192.00	1.47	1.00	1.00	71.15	Large

TABLE 11. Average iteration counts in **G5** generator tests.

	A1	A2	A3	A4	A5	A6	A7
$d = 3$	3	5	3	2	2	3	128
$d = 6$	5	6	3	1	1	5	1371
$d = 12$	9	12	3	1	1	9	Large
$d = 24$	14	24	2	1	1	14	Large
$d = 48$	24	48	2	1	1	24	Large
$d = 96$	41	96	2	1	1	41	Large
$d = 192$	81	192	3	1	1	81	Large

TABLE 12. Maximum iteration counts found in **G5** generator tests.

	A1	A2	A3	A4	A5	A6	A7
$d = 3$	2.19	3.54	1.96	1.59	1.59	2.26	24.39
$d = 6$	6.27	7.67	3.24	2.23	2.23	6.65	21041.05
$d = 12$	14.64	15.23	3.63	2.92	2.92	16.03	Large
$d = 24$	30.55	30.42	3.40	3.71	3.71	34.25	Large
$d = 48$	61.96	60.95	3.27	4.89	4.89	69.65	Large
$d = 96$	125.31	121.73	3.45	6.26	6.26	140.79	Large
$d = 192$	Unstable	242.06	Unstable	9.31	Unstable	Unstable	Large

TABLE 13. Average iteration counts in **G6** generator tests.

	A1	A2	A3	A4	A5	A6	A7
$d = 3$	5	7	5	4	4	6	242
$d = 6$	12	15	7	6	6	12	173941
$d = 12$	25	25	8	9	9	25	Large
$d = 24$	47	49	9	13	13	51	Large
$d = 48$	101	94	13	22	22	95	Large
$d = 96$	154	174	6	35	35	183	Large
$d = 192$	Unstable	331	Unstable	69	Unstable	Unstable	Large

TABLE 14. Maximum iteration counts found in **G6** generator tests.

APPENDIX D. AVERAGE TIME PER ITERATION

In Table 15 we give the average CPU time per iteration for our **G1** experiments. This was computed using the MATLAB `cputime` function.

	A1	A2	A3	A4	A5	A6	A7
$d = 3$	0.0075	0.0009	0.0078	0.0019	0.0021	0.0012	0.0002
$d = 6$	0.0087	0.0010	0.0095	0.0033	0.0035	0.0012	0.0002
$d = 12$	0.0124	0.0013	0.0141	0.0073	0.0074	0.0016	0.0004
$d = 24$	0.0229	0.0022	0.0267	0.0182	0.0184	0.0030	0.0007
$d = 48$	0.0625	0.0043	0.0702	0.0474	0.0477	0.0085	0.0014
$d = 96$	0.2510	0.0099	0.2608	0.1318	0.1324	0.0495	0.0035
$d = 192$	1.5592	0.0277	1.2623	0.3275	0.3268	0.7843	0.0121
$d = 384$	Unstable	0.1144	Unstable	1.1381	Unstable	Unstable	0.0619

TABLE 15. Average iteration times on **G1** generator tests.

The time per iteration is fairly constant across problem types so we do not include data from the other generators. One difference that will occur is that **A5** will have a higher average iteration time as that **A4** for ill-conditioned problems. In random problems, we rarely see slow convergence of **A4** so it is unnecessary to use the slower steps from **A3**. With ill-conditioned problems the **A3** steps become more frequent and increase the average time per iteration.

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