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A quasilinear algorithm for linear programs with large binary size and few number of vertices.

Adrien CHAN-HON-TONG

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Abstract

This paper presents the first algorithm proven quasilinear for linear programs with large binary size and few number of vertices.

In addition to this theoretical property, toy numerical experiments also highlights a possible practical interest despite complexity for generic linear programming is not established. Precisely, this algorithm is somehow simplex related but with the interesting features of always exiting a vertex (even degenerated) in strongly polynomial time thank to Chubanov algorithm for homogeneous linear feasibility.

1 Introduction

Linear programming is the very studied task of solving $\min_{x \ / \ Ax \geq b} c^T x$ for given $A \in \mathbb{Q}^{M \times N}$ a matrix, $b \in \mathbb{Q}^M$ and $c \in \mathbb{Q}^N$ some vectors.

Let L be the binary size required to write the given input (i.e. A, b, c), then, this problem can be solved in polynomial times (i.e. in a number of binary operations bounded by a L^{γ} with a fixed γ) since [9, 6, 8], with interior point method [11] being currently the state of the art.

Yet, there is no known algorithm proven to have a strong polynomial complexity i.e. able to solve the problem in a number of binary operations bounded by $L \times \log(L)^{\kappa} \times (MN)^{\gamma}$. And, [1] shows that major interior point algorithms are not strong polynomial. Today, only, some families of linear programs can be solved in strongly polynomial times:

- linear program with 0/1 matrix A [13] (by specific algorithm)
- linear program with at most two variables per inequality [7] (by specific algorithm)
- Markov chain [12] (by simplex algorithm [4])
- system having binary solution [2] (by specific algorithm)

• homogeneous linear feasibility i.e. $\exists ?x \in \mathbb{Q}^N \ / \ Ax = \mathbf{0}, \ x > \mathbf{0}$ given $A \in \mathbb{Q}^{M \times N}$ a full rank matrix [3] by Chubanov algorithm.

As no known algorithm is strongly polynomial, there is an interest to design an algorithm to solve linear programs with very large binary size using prior on the geometry of the problem (this claim is formalized in next section). Then, section 3 presents an algorithm based on Chubanov algorithm [3] for linear programming. Despite there is little hope that this algorithm is polynomial in generic case, it solves theoretically the issue of linear programs with very large binary size when number of vertices is small. This algorithm shares some feature with simplex but exits vertex in strongly polynomial time. These claims are discussed in section 4.

Notations

 \mathbb{N} , \mathbb{Q} are the sets of integer and rational numbers. \ is the ensemble subtraction. For all integers i, j, [i, j] will symbolize the **integer** range i.e. $\{i, i+1, ..., j\}$ which is empty if i > j (there will be no ambiguity with the interval in \mathbb{R} as there is no real range in this paper).

For all integers i, j, I, J, \mathbb{Q}^I is the set of I dimensional vectors on \mathbb{Q} , and, $\mathbb{Q}^{I \times J}$ is the set of matrix with I rows and J columns, with values in \mathbb{Q} , and, $._i$ designs the i component: a row for a matrix and a rational for vector or a row. \mathbb{Q}^I would be matched with $\mathbb{Q}^{I \times 1}$ i.e. vectors are seen as columns, and, row of a matrix are matched with $\mathbb{Q}^{1,J}$. For all sets $S \subset \mathbb{N}$, A_S, b_S is the submatrix or subvector obtained when keeping only components indexed by $s \in S$. T is the transposition operation i.e. $A_{j,i}^T = A_{i,j}$. $\mathbf{0}$ and $\mathbf{1}$ are the 0 and 1 vector i.e. vector contains only 0 or only 1, and \mathbf{I} is the identity matrix.

If $A \in \mathbb{Q}^{I \times J}$, the null vector space of A (i.e. the kernel) is written $Ker(A) = \{v \in \mathbb{Q}^J / Av = \mathbf{0}\}$, with the convention that Ker of empty A is all space.

Finally, it is written that complexity is O(f(size)) when there exists λ such that for all size size, worse case complexity of the algorithm on all instances of size size is less than $\lambda f(size)$. The same stands for $\tilde{O}(f(size))$ except that it means that there exists both λ and κ such that complexity is less than $\lambda f(size) \log(f(size))^{\kappa}$.

2 Limit of state of the art

There is no known algorithm to solve linear programming in strongly polynomial time, and, major family of interior point are proven unable to do so [1]. Thus, there is a practical and theoretical issue for the state of the art to deal with linear programs with very large integer. Let formalize theoretically this issue.

Let first recall the classical primal dual trick to go from optimization problem into a decision one. Let assume original goal is to solve $\max_{A_{raw}x \leq b_{raw}, x \geq \mathbf{0}} c_{raw}^T x$. It is well known that the dual problem is $\min_{A_{raw}^T y \geq c_{raw}, y \geq \mathbf{0}} b_{raw}^T y$. Now, the primal

dual is formed by combining all constraints: $A_{raw}x \leq b_{raw}$, and, $x \geq \mathbf{0}$, and, $A_{raw}^Ty \geq c_{raw}$, and $c_{raw}x = b_{raw}y$, and finally, $y \geq \mathbf{0}$.

So, the problem $\max_{A_{raw}x < b_{raw}, x > \mathbf{0}} c_{raw}x$ can be folded into $A_{big}x_{big} \ge b_{big}$ with

$$A_{big} = \begin{pmatrix} -A_{raw} & 0 \\ I & 0 \\ 0 & A_{raw}^T \\ 0 & I \\ c_{raw} & -b_{raw} \\ -c_{raw} & b_{raw} \end{pmatrix} \text{ and } b_{big} = \begin{pmatrix} -b_{raw} \\ 0 \\ c_{raw} \\ 0 \\ 0 \\ 0 \end{pmatrix}.$$

Then, this problem $A_{big}x \geq b_{big}$ can be transformed by adding a variable z into $A_{big}x+z \geq b_{big}$, $z \geq 0$. This way, there is an equivalence between feasibility $\exists ?x \ / \ A_{big}x \geq b_{big}$ and minimization of z with constraints $A_{big}x+z \geq b_{big}$, $z \geq 0$. Now, let introduce the large integer linear programming problem:

Definition 2.1 (large integer linear programming). The large integer linear programming problem is to decide if: $\exists x \in \mathbb{Q}^N \ / \ Ax \ge b, c^T x \ge 0, c^T x = 0$ when $NM \le \log(L)$, given $A \in \mathbb{Q}^{M \times N}$, $b \in \mathbb{Q}^M$ and $c \in \mathbb{Q}^N$ with $Ac > \mathbf{0}$.

For all instances for which $NM \geq \log(L)$, the output is trivially true. So, all algorithms could be derived to solve these instances trivially. So, the interesting instances are those with $NM \leq \log(L)$. Any strongly polynomial algorithm will have an quasilinear complexity i.e. $\widetilde{O}(L) = O(L \times \log^{\gamma}(L))$ with a constant γ , but, any non strongly polynomial algorithm will be super linear i.e. with a L^{γ} complexity with $\gamma > 1$.

The interesting point is that no algorithm are known quasilinear to solve large integer linear programming and major state of the art algorithm are known not to be (counter examples from [1] have unbounded L, eventually, $NM \leq \log(L)$).

Now, one could consider simplex with classical pivoting rule [4] as a candidate for large integer linear programming as simplex complexity depend on the number of pivots which could be bounded depending on N and M and not L. However, it is not quasilinear because it could need an exponential (in N, M) number of integer operations to terminate. Precisely, let introduce the concept of vertex:

Definition 2.2 (vertex). Given $A \in \mathbb{Q}^{M \times N}$, $b \in \mathbb{Q}^M$ and $c \in \mathbb{Q}^N$ with Ac > 0 and $A_M = c$ and $b_M = 0$, a set of index $D \subset [1, M]$ is called a vertex if and only if $\exists x \in \mathbb{Q}^N$ such that $A_D x = b_D$ and $A_{[1,M] \setminus D} x > b_{[1,M] \setminus D}$, and, there is no $v \in \mathbb{Q}^N$ such that $A_D v = 0$, $c^T v < 0$.

In non singular case, a vertex is just a point x such that $A_Dx = b_D$ with A_D square and non singular, plus, $A_{[1,M]\setminus D}x > b_{[1,M]\setminus D}$ (so, obviously, there is no $v \in \mathbb{Q}^N$ such that $A_Dv = 0$, $c^Tv < 0$). But, in degenerated case, $A_Dx = b_D$ may define a sub vectorial space included in a iso-c hyperplane. Let also stress that this definition requires $A_{[1,M]\setminus D}x > b_{[1,M]\setminus D}$, this way, a point x can not belong to more than one vertex. Also, the number of vertices is at most bounded by 2^M .

Let write $\Gamma(A,b,c)$ the number of vertices given A,b,c.

Now, the idea is that vertex based algorithm are efficient if number of vertices is small.

Definition 2.3 (few vertices large integer linear programming). The few vertices large integer linear programming problem is to decide if: $\exists x \in \mathbb{Q}^N \ / \ Ax \geq b, c^T x = 0$ when $NM \leq \log(L)$ and $\Gamma(A, b, c) \leq \log(L)$, given $A \in \mathbb{Q}^{M \times N}$, $b \in \mathbb{Q}^M$ and $c \in \mathbb{Q}^N$ with Ac > 0, $A_M = c$, $b_M = 0$.

Obviously, this problem is unrealistic and not interesting in practice. But from theoretical point of view, this problem is well defined.

Now, simplex algorithm (for example) may become interesting for this problem compared to interior point because during the exploration either there is less than $\log(L)$ vertices and the simplex will output a solution after having explored these only $\log(L)$ vertices, or, there is more vertices but the algorithm could stop anyway after the $\log(L)$ first explored vertices. So complexity will be $\widetilde{O}(\log(L) \times \Upsilon)$ with Υ being the complexity of the algorithm to go from a vertex to an other. Yet, the problem with the simplex (with all known pivoting rules) is that it can require exponential time to exit a vertex, so, Υ is not just bounded by $\widetilde{O}(L \times (NM)^{\gamma})$.

However, an algorithm similar to the simplex, but, with the feature of exiting a vertex with strongly polynomial complexity will be especially interesting from the few vertices large integer linear programming (yes, obviously, this problem is designed in this purpose).

Currently, the question about the existence of simplex algorithm able to exit degenerated vertex in strongly polynomial times is not new e.g. [10]. However, it seems [10] only prove that exiting a vertex is equivalent to solve two vertices linear program, but, no more. Indeed, obviously, solving linear program allows to solve the exiting a vertex problem with the same complexity. But, knowing to exit any vertex is not sufficient to solve linear program as there can be an exponential number of vertices. Currently, [10] designs for each linear program a two vertices equivalent linear program. Of course, this two vertices equivalent linear program can be solve just by exiting 1 vertex. Yet, the construction of the two vertices equivalent linear program use the solution of the original linear program! So [10] makes no wrong statements, but, overclaims the equivalency.

The main contribution of this paper is to present an algorithm, built from [3], able to exit any vertex in strongly polynomial time.

3 The slide and jump algorithm

3.1 Key idea

The starting point of this algorithm is Chubanov algorithm [3] which solves homogeneous linear feasibility $(\exists ?x \in \mathbb{Q}^N \ / \ Ax = \mathbf{0}, \ x > \mathbf{0}$ when $A \in \mathbb{Q}^{M \times N}$ has full rank) and/or it dual $\exists ?x \in \mathbb{Q}^N \ / \ Ax > \mathbf{0}$ in strongly polynomial time.

Then, it is easy to get a small improvement from a not optimal admissible solution for linear program by solving a dual homogeneous linear feasibility

problem. Indeed, if x verifying $Ax \ge b$ has not minimal cx under assumption that $Ac > \mathbf{0}$, then it is possible (see proof in subsection 3.3) to find v such that $\begin{pmatrix} A_D \\ -c^T \end{pmatrix} v > \mathbf{0}$ with D the current saturated constraints (i.e. D such that $A_Dx = b_D$).

Now, it could take infinite time to reach the optimal solution by computing such v and updating $x = x + \varepsilon v$. Yet, if x is not on a vertex, it is always possible to decrease $c^T x$ just by moving in the kernel of current saturated constraints. So, the algorithm does sliding moves to reach a vertex from a non vertex point, and, then, a jumping move to go from a vertex to an interior point. The good feature is that *sliding* moves allow to explore completely a particular combinatorial situation (D) while *jumping* moves allow to exit such structure. The termination will be guarantee by the impossibly to see a vertex D twice, and, the correctness, by capacity to jump to next D until optimal solution is reached.

3.2 Pseudo code

Pseudo code for linear program is presented in algorithm 1.

Importantly, the algorithm assumes the input linear program is $c \in \mathbb{Q}^N$, $b \in \mathbb{Q}^M$, $A \in \mathbb{Q}^{M \times N}$ with $Ac > \mathbf{0}$, $Ax \ge b \Rightarrow c^T x \ge 0$, and $Ax_{start} \ge b$ (e.g. $x_{start} = \lambda c$). Yet, these assumptions do not restrict generality because all linear programs can be pushed in this shape using classical primal dual trick.

```
Algorithm 1 Slide and jump algorithm
```

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Require: c \in \mathbb{Q}^N, b \in \mathbb{Q}^M, A \in \mathbb{Q}^{M \times N} with Ac > \mathbf{0}, Ax \ge b \Rightarrow c^T x \ge 0, and
Ax_{start} \geq b (e.g. x_{start} = \lambda c) this does not restrict generality
Ensure: return x: Ax \ge b, cx is minimal
  1: while True do
            D = \{ m \in [1, M] / A_m x = b_m \}
  2:
            compute u the orthogonal projection of -c on Ker(A_D)
  3:
            if u = 0 then
  4:
                 call subsolver \exists ?v \ / \ \begin{pmatrix} A_D \\ -c^T \end{pmatrix} v \ge \mathbf{1}
if v not exists then return x
chose \delta with 0 < \delta < h = \min_{m \in [1,M]/A_m v < 0} \frac{A_m x - b_m}{-A_m v} (e.g. \delta = \frac{h}{4})
  5:
  6:
  7:
                  x = x + \delta v
  8:
  9:
                  g = \min_{m \in [1,M] \ / \ A_m u < 0} \ \frac{A_m x - b_m}{-A_m u}
10:
11:
```

3.3 Correctness and termination

This subsection presents a proof that this slide and jump algorithm is well defined, terminates, and produces an exact optimal solution of linear program.

Lemma 3.1. Algorithm is well defined

Proof. Problematic steps are step 3 (non standard operation) and step 10 (set should not be empty). All other steps are standard operations.

step 10:

By assumption cx is bounded by 0 i.e. $\forall x \in \mathbb{Q}^N$, $Ax \geq b \Rightarrow cx \geq 0$. If there was ω such that $c\omega < 0$ and $A\omega \geq \mathbf{0}$, then, one could produce an unbounded admissible point $x + \lambda \omega$ as $A(x + \lambda \omega) \geq Ax \geq b$ and $c(x + \lambda \omega) \xrightarrow[\lambda \to \infty]{} -\infty$. So, if $c\omega < 0$ then $\exists m \in [1, M] / A_m\omega < 0$.

step 3:

Step 3 is the projection on a vectorial space i.e. it consists to solve $\underset{A_Du=0}{\arg\min}(c+u)^T(c+u)$. This step can be done by Gram Schmidt algorithm. This procedure can not fail (returns -c on empty input), and, always returns a vector with a strictly positive scalar product with -c or $\mathbf{0}$.

Lemma 3.2. Algorithm keeps the current point in the admissible space

Proof. Steps 8 and 11 move the current point.

Step 8:

Seeing the test in step 6, $A_D v > \mathbf{0}$ so for all $m \in D$, $A_m(x + \lambda v) > b_m$. Now, for all $m \notin D$, $A_m x > b_m$, so $\exists \delta > 0$ such that $A(x + \delta v) > b$ (the offered δ works but anyway it could be less).

Step 11:

First, $\forall m / A_m u \geq 0$, $A_m(x + \lambda u) \geq b$.

Then, seeing the definition of u from step 3, if $m \in D$, $A_m u = 0$ and if $m \notin D$, $A_m x > b_m$. So $\forall m \in [1, M]$, $A_m u < 0 \Rightarrow m \notin D \Rightarrow A_m x > b_m$, and so g > 0. Now, for $m / A_m u < 0$, g is a minimum, so $\frac{A_m x - b_m}{-A_m u} \geq g$. When multiplying by a negative: $\frac{A_m x - b_m}{-A_m u} A_m u \leq g A_m u$. And, so $\forall m \in [1, M]$, $/A_m u < 0$: $A_m x + g A_m u - b_m \geq A_m x + \frac{A_m x - b_m}{-A_m u} A_m u - b_m = 0$.

Lemma 3.3. algorithm outputs optimal exact solution

Proof. The proof consists to assume the algorithm returns non optimal admissible point x, while optimal solution was x^* ($cx > cx^*$). Then, it is possible to built \hat{x} (by adding εc) which are in the interior of the admissible space and still better than x for c. But, it means that $\hat{x}-x$ is both better for the objective and for saturated constraints. So, the subsolver should have returned something, and so, there is a contradiction.

Precisely, let consider $\phi = x^* - x + \frac{c^T(x-x^*)}{2\sqrt{c^Tc}}c$, then, $A_D\phi = A_D(x^* - x + \frac{c^T(x-x^*)}{2}c) = A_D(x^* + \frac{c^T(x-x^*)}{2}c) \ge \frac{c^T(x-x^*)}{2}A_Dc > \mathbf{0}$ and $c(\phi - x) = -\frac{c^T(x-x^*)}{2} < 0$.

Lemma 3.4. algorithm can not loop more then M+1 times without calling the subsolver (equivalently reaching a vertex)

Proof. g is exactly build such that the k corresponding to the minimum enters in D. Let consider $k \notin D$ such that $\frac{A_k x - b_k}{-A_k u} = g$. So, $A_k(x + gv) - b_k = A_k x + \frac{A_k x - b_k}{-A_k v} A_k v - b_k = 0$.

So, D strictly increases (for inclusion, as set) each time the algorithm reaches step 11, but, D is bounded by [1, M], so, test 4 can not return true more than M consecutive times.

Lemma 3.5. All moves strictly decrease the objective function.

Proof. Steps 8 and 11 move the current point.

step 11: By construction $c^T u < 0$, and, g > 0. So, cost decreases when moving along u.

step 8: $c^T v < 0$ so moving a little along v decreases the cost.

Lemma 3.6. A value for set D observed in step 5 can not be observed again in step 2.

Proof. This lemma is proven by contradiction: if D is seen again, one can prove that exploration of D should have been continued (i.e. the test step 3 should have been false) instead of calling the subsolver on step 5.

Let consider $x_2 - x_1$ with x_1 corresponding to the observation of D in step 5 and x_2 to any ulterior observation. As all moves strictly decreases $c^T x$, it means $c^T(x_2-x_1) < 0$, but, by definition of D, $A_D x_2 = A_D x_1 = b_D$ so $A_D(x_2-x_1) = \mathbf{0}$. So, projection of -c on $Ker(A_D)$ is not null (at least it could be $x_2 - x_1$), so, algorithm should not have passed the test step 3.

Lemma 3.7 (Slide and jump). The slide and jump algorithm solves linear programs because iteration between two vertices is strongly polynomial, and, vertices are never reached twice.

Proof. From all previous lemmas, observing that D is bounded (so looping without reaching a vertex is impossible), and that, number of subsets from [1, M] is finite an never see twice (i.e number vertices is bounded and vertex are never explored twice), algorithm terminates. And, independently, algorithm is well defined, works in admissible space and can not output something else than an optimal solution. So both correction and termination are proven.

4 Quasilinear complexity of few vertices large integer linear programming

The main claim of this paper is the following:

Theorem 4.1. The slide and jump algorithm solves in quasilinear time the few vertices large integer linear programming. *Proof.* Complexity of the slide and jump algorithm to decide a few vertices large integer linear programming instance is bounded by $\log(L) \times \Upsilon$ binary operation where Υ is the binary complexity to exit a vertex, but, here, $\Upsilon = \widetilde{O}(L) \times (NM)^{\kappa}$ as exiting a vertex consists basically calling Chubanov algorithm (strongly polynomial), yet, as $NM \leq \log(L)$ on non trivial instance, the total complexity is $\widetilde{O}(L)$ binary operations.

In addition, this algorithm is the first with this property because current interior point algorithms are super linear (in L) independently from the number of vertices, and, simplex requires exponential (in NM) number of operations to exit a vertex.

Obviously, the few vertices large integer linear programming, could be solved directly by any algorithm able to enumerate vertices. Yet, enumerating of all sets D is easy, but, enumerating vertices is harder as a set D can be unsatisfiable, or, incomplete (i.e. face but not vertex). This way, enumerating sets is exponential (in NM) because it does not stop after the $\log(L)$ first. Currently, it may be possible to use geometric prior to design ad-hoc algorithm. But it is not that trivial because any algorithm linear program instance could be meet (just if there is large number of vertices it is possible to decide directly instead of solving it).

5 Discussion

5.1 Complexity

Despite that the main claim of this paper is somehow interesting from theoretical point of view, it may have almost none practical implication as the few vertex large integer linear programming is a very unrealistic problem.

So an other question is about the complexity of the slide and jump algorithm for standard linear programming. However, depending on the implementation of inner subsolver (to solve $A_D v \geq \mathbf{0}, c^T v < 0$), slide and jump can behave more or less like a simplex for a new pivoting rule. The behaviour would be different from simplex when $A_D v > \mathbf{0}$, as, it is even possible that two consecutive vertices have no constraint in common (but it depends not just on how x exit the vertex, but also, how x collects constraints to reach a vertex). But, the behaviour would be very close when the produced vector v verifies $A_{D\setminus\{d\}}v = \mathbf{0}$ i.e. moves follow an edge to directly go from a vertex to an adjacent vertex exactly like a simplex.

In other words, depending on the implementation, slide and jump could be just a particular simplex for a new pivoting rule with the important feature that exiting a vertex is strongly polynomial. So, the optimistic point of view is to stress that complexity of the simplex with no standard pivoting rules is still opened (polynomial Hirsch conjecture). But, the realistic point of view is that there is little hope that slide and jump algorithm is strongly polynomial for standard linear program.

A Harder point to advance on this question is that link between consecutive vertices in slide and jump does not depend only on the exiting condition of jumping moves, but, also on how constraint are collected during sliding moves.

5.2Non singular case

To advance on this complexity question, one can consider a non singular vertex: A_D (written A here) is square and non singular. The goal is to find $Av \geq$ $\mathbf{0}, c^T v < 0 \text{ (with } Ac > \mathbf{0}).$

As the vertex is non singular, one can consider A^{-1} and $\phi_1, ..., \phi_N$ the vector such that $(A\phi_n)_i = \begin{cases} 1 & \text{if } i = n \\ 0 & \text{otherwise} \end{cases}$ (i.e. the transposed of the row of A^{-1}). Let $\Omega = \{n \in [1, N] / c^T \phi_n >= 0\}$ and $\Upsilon = \{n \in [1, N] / c^T \phi_n < 0\}$.

 Υ is important because it will never again be included in any vertices. Indeed, there is no ψ such that $A_{\Upsilon}\psi = \mathbf{0}$, $A_{\Omega}\psi \geq \mathbf{0}$, and, $c^T\psi < 0$. If it was, by removing the corresponding vector from Ω , one could form a vector such that $A\omega = 0$ and $c^T\omega < 0$ (this is impossible both because it is a vertex and because A is non singular).

The basic idea of the simplex (with greedy descent pivot rule) is to move

following ϕ_k with $c^T\phi_k = \min_{n \in \Upsilon} c^T\phi_n$.

Now, if one is interested to push slide and jump in the simplex framework, it seems that the underlying pivoting rule should be to move along $\sum_{n \in \Upsilon} \phi_n$ (corresponding to solve $v = \min_{\mathbf{0} \le A\nu \le \mathbf{1}} c^T \nu$ in singular case - but it is not relevant to solve such linear problem just to exit a vertex...). Typically, if Υ is just a pair i, j, greedy descent pivot rule will select either i either j while it seems interesting to move on the bisector of both. These way, when hitting a constraint k, k will give a direction which may lead to hit i xor j. But the good point is that if i is the one meet after k, then, at least the vertex j, k has been avoided.

Now, some related pivoting rule have already be studied without closing the Hirsch conjecture. In addition, again, this is not a complete pivoting rule because the resulting vertex also depend on the sliding moves.

5.3 Implementation

As complexity is not established, some toy numerical experiments are offered. Good feature of the offered implementation is to tackle exact arithmetic to produce exact solution. Except this point, these experiments are quite limited due to external and internal limitations.

Main external limitation is that algorithm depends on the efficiency of the Chubanov implementation, and, on the average number of encountered vertices - but - no public implementation of Chubanov algorithm are available today. to bypass the miss of public implementation of Chubanov algorithm, offered source code implements subsolver using projection on -c on kernel of A_S with S a random subset of D in the spirit of a Kaczmarz-Motzkin algorithm [5]

(This is equivalent as adding ϵc provides v such that $A_D v > \mathbf{0}$ - currently the algorithm does not require functionally that $A_D v > \mathbf{0}$ and $c^T v < 0$, but, only that $A_D v \geq \mathbf{0}$ and $c^T v < 0$ - or to prove there is not). Of course, using Chubanov algorithm should be better: the current implementation does not even guarantee termination. This would be changed as soon as an implementation of Chubanov will be made public. Let stress that, in the offered code, either problem are feasible, or, the primal dual of the primal dual is computed. This way, optimal solution always exists and always has 0 cost. This allows a much simpler check of termination (just checking is $Ax \geq b$ and $c^T x = 0$) that using Chubanov algorithm to wait a certificate of impossibility. This is, in fact, required as the current simulation of Chubanov algorithm only admits feasible instance. Yet, as computing the primal dual of the primal dual is always possible this does not restrict generality or validity of the offered experiments.

Main internal limitations are that the algorithm has been implemented in python without parallelism (which is a critical issue for pratical performances), and, that no public benchmark like http://www.netlib.org/lp/data/ have been evaluated. Indeed, evaluation only relies on two types of problem are considered: Klee Minty cubes and random instances. And, tested number of dimension and number of variables are quite low (around 100) compared to state of the art test. So, obviously, there is no big challenge in solving these instances. Clearly some additional work is required to produce an interesting numerical contribution, the main contribution of the paper being theoretical.

Offered experiments are not sufficient to answer whether this algorithm is or not relevant in practice. But, they highlights that the algorithm is not purely theoretical: slide and jump algorithm solves the basic but not completely trivial tests:

- the Klee Minty cube (up to dimension 50) by exploring very low number of vertices
- random instances by exploring low number of vertices
- using exact arithmetic to produce exact solution

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Some source code

```
def opposedvector(v):
    return [-e for e in v]
\begin{array}{lll} \texttt{def saturated constraints} \, (A,b,x) \colon \\ & \texttt{return [m for m in range(len(A)) if produits calaire} \, (A[m]\,,x) = = b\,[m]\,] \end{array}
return pu
\begin{array}{ll} def & resteprojection\,(u\,,BOG\,)\colon\\ & return & combinaisonlineaire\,(u\,,Fraction\,(\,-1\,)\,,projection\,(u\,,BOG\,)\,) \end{array}
def gramschimdBOG(H,BOG):
   BOG = BOG.copy()
   while True :
        H = [resteprojection(h,BOG) for h in H]
        H = [h for h in H if produitscalaire(h,h) != Fraction()]
                          H! = []:
BOG. append (H. pop ())
                  else:
#ASSUME exist v: Av>0, av<0
#return v such that Av>=0, av<0
## COULD BE IMPLEMENTED AS CHUBANOV ALGORITHM ##
#BUT CURRENTLY TERMINATION IS NOT GUARANTEE ##
def no_guarantee_termination_subsolver(A,a):
         counter=0
         counter=0
while True:
   index = [i for i in range(len(A))]
   random.shuffle(index)
   subsetI = index[0:random.randint(0,len(index))]
                  \begin{array}{l} {\rm subset} \, = \, [\, A \, [\, i \, ] \  \, {\rm for} \  \, i \  \, {\rm in} \  \, {\rm subsetI} \, ] \\ {\rm v} \, , BOG \, = \, {\rm projection\_on\_ker} \, (\, {\rm subset} \, , a \, , [\, ] \, ) \end{array}
                  AV = [produitscalaire(A[i],v) for i in range(len(A))]
FORBID = [A[i] for i in range(len(A)) if AV[i] < Fraction()]
while FORBID! = [1:
    subset += FORBID
    v,BOG = projection_on_ker(subset,c,BOG)
    AV = [produitscalaire(A[i],v) for i in range(len(A))]
    FORBID = [A[i] for i in range(len(A)) if AV[i] < Fraction()]
                  counter+=1
value = produitscalaire(v,a)
if value<Fraction():
    print("nb trial in false Chubanov",counter)
    return v</pre>
#ASSUME Ax>=b, xc>0, Ac>0, Ay>=b with cy=0
#let d in D <=> A_dx=b_d
#return v such that A_Dv>=0 and cv<0
def jump(A,b,c,x):
    D = [i for i in range(len(A)) if produitscalaire(A[i],x)==b[i]]
    block = [A[i] for i in D]
    return no-guarantee_termination_subsolver(block,c)</pre>
S = [m \text{ for } m \text{ in range(len(A))} \text{ if } AV[m] < Fraction()]  if S = = []:
                S==[]:
print("maximalmoves: S==[]")
quit()
         \begin{array}{lll} R \, = \, \left[ -AXb\left[m\right]/AV\left[m\right] & \text{for m in S} \right] \\ 1 \, = \, \min\left(R\right) \end{array}
```

```
if l==Fraction():
    print("maximalmoves: l==0")
                                     quit()
#ASSUME Ax>=b #ASSUME Ax>=b => cx>=0 #returns y such that cy<=cx, Ay>=b and projection_on_ker(A_D,c)=0 # with d in D <=> A_Dy=b_D def slide(A,b,c,x):
    print("entering slide",datetime.datetime.now())
BOC = []
    while True:
        print("$", end="", flush=True)
        D = saturatedconstraints(A,b,x)
                                   \begin{array}{lll} v,BOG = & projection\_on\_ker\left(\left[A[m] & for \ m \ in \ D\right],c,BOG\right) \\ if & produits calaire\left(v,v\right) == Fraction\left(\right): \\ & print\left("exiting slide",datetime.datetime.now()\right) \\ & return \ x \\ l & = & maximal moves\left(A,b,x,v\right) \\ x & = & combinais on lineaire\left(x,l,v\right) \end{array}
 while True:
    print("sliding move")
    x = slide(A,b,c,x)
    D = saturatedconstraints(A,b,x)
    print(D)
                                    if produitscalaire(x,c)==Fraction():
    print("found optimal")
    return x, counter
                                    \begin{split} & \text{print} \left(\text{"chubanov jump"}\right) \\ & v = \text{jump}\left(A, b, c, x\right) \\ & 1 = \text{maximalmoves}\left(A, b, x, v\right) / 4 \\ & x = \text{combinaisonlineaire}\left(x, 1, v\right) \end{split}
                                     counter+=1
  def normalize(rawA, rawb, rawxoptimal):
#input rawA rawx >= rawb
#return A,b,c such that min{cx / Ax>=b} if equivalent
#+ A is normalized, c is normalized, cx is 0 bounded, Ac = 3/4 vector(1)
M = len (rawA)
N = len (rawA[0])
A = [[Fraction() for n in range(N+3)] for m in range(M+4)]
b = [Fraction()] * (M+4)
c = [Fraction()] * (N+3)
c[-1] = Fraction(1)
                  normRawA = [Fraction()]*M
normRawAtrick = [Fraction()]*M
for m in range(M):
    normRawA[m] = produitscalaire(rawA[m],rawA[m])
    normRawAtrick[m] = normRawA[m]/Fraction(2)+Fraction(1)
                    \begin{array}{lll} & for \ m \ in \ range (M): \\ & for \ n \ in \ range (N): \\ & A[m] \ [n] = rawA \ [m] \ [n]*Fraction (4,5)/normRawAtrick \ [m] \\ & A[m] \ [-3] = normRawA \ [m]*Fraction (4,5*2)/normRawAtrick \ [m] \\ & A[m] \ [-2] = Fraction (4,5*2)/normRawAtrick \ [m] \\ & A[m] \ [-1] = Fraction (3,5) \\ & b \ [m] = rawb \ [m]*Fraction (4,5)/normRawAtrick \ [m] \\ \end{array} 
                   \begin{array}{lll} A[-4][-3] &= & \operatorname{Fraction}\left(4,5\right) \\ A[-4][-1] &= & \operatorname{Fraction}\left(3,5\right) \\ A[-3][-3] &= & -\operatorname{Fraction}\left(4,5\right) \\ A[-3][-1] &= & \operatorname{Fraction}\left(4,5\right) \\ A[-2][-2] &= & \operatorname{Fraction}\left(4,5\right) \\ A[-1][-2] &= & -\operatorname{Fraction}\left(4,5\right) \\ A[-1][-1] &= & \operatorname{Fraction}\left(4,5\right) \\ A[-1][-1] &= & \operatorname{Fraction}\left(3,5\right) \\ \end{array} 
                    xoptimal = [Fraction()]*(N+3)
for n in range(N):
    xoptimal[n] = rawxoptimal[n]
```

```
return A, b, c, xoptimal
def primaldual(rawA,rawb,rawc):
#primal: max {rawc rawx / rawA rawx<= rawb, rawx>=0}
#dual: min {rawb rawy / transpose(rawA) rawy>= rawc, rawy>=0}
#primal dual: {rawx / rawA rawx<=rawb, rawx>=0,
# transpose(rawA) rawy>=rawc, rawy>=0, rawc rawx=rawb rawy}
#unfolded into A x >= b
M = len(rawA)
N = len(rawA)
N = len(rawA[0])
A = [[Fraction() for n in range(N+M)] for m in range(M+N+N+M+2)]
b = [Fraction()]*(M+N+N+M+2)
                           offsetY = N
offset = 0
for m in range(M):
    for n in range(N):
        A[m+offset][n] = -rawA[m][n]
    b[m+offset] = -rawb[m]
                            offset += M
for n in range(N):
   A[n+offset][n] = Fraction(1)
                            offset += N
for n in range(N):
   for m in range(M):
       A[n+offset][m+offsetY] = rawA[m][n]
   b[n+offset] = rawc[n]
                            \begin{array}{lll} offset \ += N \\ for \ m \ in \ range (M): \\ A [m\!+\!offset] \left[m\!+\!offsetY\right] \ = \ Fraction (1) \end{array}
                            \label{eq:continuous_section} \begin{split} & \text{for n in range}\left(N\right)\colon\\ & & A[-2][n] = \text{rawc}\left[n\right]\\ & \text{for m in range}\left(M\right)\colon\\ & A[-2][m+\text{offset}Y\,] = -\text{rawb}\left[m\right] \end{split}
                            \begin{array}{lll} \text{for n in range(N):} & & & & & \\ & & & & & \\ & & & & & \\ & & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & \\ & & \\ & & \\ & & \\ & & \\ & & \\ & & \\ & & \\ & & \\ & & \\ & & \\ & & \\ & & \\ & & \\ & & \\ & & \\ & & \\ & & \\ & & \\ & & \\ & & \\ & & \\ & \\ & & \\ & \\ & & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ & \\ &
                             return A, b
  \begin{array}{ll} def & cubeproblem Primal (N): \\ & twopower = [Fraction ()]*N \\ & twopower [0] = Fraction (2) \\ & for & n & in & range (1,N): \\ & & twopower [n] = & Fraction (2)*twopower [n-1] \end{array} 
                           \begin{array}{ll} b = & [\, Fraction\,(\,)\,] * N \\ b\,[\,0\,] = & Fraction\,(\,5\,) \\ for & n \ in \ range\,(\,1\,,N\,) : \\ b\,[\,n\,] = & Fraction\,(\,5\,) \ * \ b\,[\,n\,-1\,] \end{array}
                           c = twopower[::-1]
                          \begin{split} A &= \left[\left[ \operatorname{Fraction}\left(\right) \text{ for } n \text{ in } \operatorname{range}\left(N\right) \right] \text{ for } m \text{ in } \operatorname{range}\left(N\right) \right] \\ \text{for } n \text{ in } \operatorname{range}\left(N\right) \colon \\ \text{for } k \text{ in } \operatorname{range}\left(n\right) \colon \\ A[n][k] &= \operatorname{twopower}[n-k] \\ A[n][n] &= \operatorname{Fraction}\left(1\right) \end{split} 
                            return A.b.c
  \begin{array}{ll} def & cubeproblem \, (N) \colon \\ & Araw \, , braw \, , craw \, = \, cubeproblem \, Primal \, (N) \\ & A,b \, = \, primal dual \, (Araw \, , braw \, , craw \, ) \end{array}
                            \begin{array}{ll} \text{xoptimal} &= [\, \text{Fraction} \, (\,) \,] * (\, 2 * N) \\ \text{xoptimal} \, [\, -1] &= \text{Fraction} \, (\, 2\,) \\ \text{xoptimal} \, [\, N-1] &= \text{braw} \, [\, -1\,] \end{array}
                            return A, b, xoptimal
  return [Fraction(random.randint(-100,-1)) for n in range(N)]
  \begin{array}{lll} def & randomMatrix\left(M,N\right): \\ & return & \left[ \; randomVector\left(N\right) \; \; for \; m \; \; in \; \; range\left(M\right) \right] \end{array}
  \begin{array}{ll} \text{def random problem} \, (N, M) \colon \\ & \text{xoptimal} \, = \, \text{random} \, Vector \, (N) \end{array}
```