

CENTER FOR MACHINE PERCEPTION



# Minimal Problem Solver Generator

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Acknowledge grants here. Use centering if the text is too short.

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# Czech Technical University in Prague Faculty of Electrical Engineering

#### **Department of Cybernetics**

## **BACHELOR PROJECT ASSIGNMENT**

Student: Pavel Trutman

**Study programme:** Cybernetics and Robotics

Specialisation: Robotics

**Title of Bachelor Project:** Minimal Problem Solver Generator

#### **Guidelines:**

- 1. Review the state of the art in solving the polynomial systems using linear algebra [1, 2] and the automatic generator of the polynomial solvers [3, 4].
- 2. Implement the improvement [4] of [3] into the existing automatic generator of solvers.
- 3. Implement a variation of algorithm [1], review its behavior, and suggest how to take over some of its elements to the automatic generator and implement it.
- 4. Demonstrate the functionality of the new solver generator and compare it with the original solver.

#### Bibliography/Sources:

- [1] Faugere, J.-C. (June 1999): "A new efficient algorithm for computing Gröbner bases (F4)". Journal of Pure and Applied Algebra (Elsevier Science) 139 (1): 61–88. doi:10.1016/S0022-4049(99)00005-5. ISSN 0022-4049.
- [2] Faugere, J.-C. (July 2002): "A new efficient algorithm for computing Gröbner bases without reduction to zero (F5)". Proceedings of the 2002 international symposium on Symbolic and algebraic computation (ISSAC) (ACM Press): 75–83.
- [3] Kukelova, Z.: Algebraic Methods in Computer Vision. PhD Thesis. CTU in Prague 2013. http://cmp.felk.cvut.cz/~kukelova/webthesis/docs/Kukelova-phd-2013.pdf
- [4] Kukelova, Z.; Bujnak, M.; Heller, J.; Pajdla, T.: Singly-Bordered Block-Diagonal Form for Minimal Problem Solvers. ACCV 2014.

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**Valid until:** the end of the summer semester of academic year 2015/2016

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#### České vysoké učení technické v Praze Fakulta elektrotechnická

#### Katedra kybernetiky

# ZADÁNÍ BAKALÁŘSKÉ PRÁCE

Student: Pavel Trutman

**Studijní program:** Kybernetika a robotika (bakalářský)

Obor: Robotika

Název tématu: Generátor řešení minimálních problémů

#### Pokyny pro vypracování:

- 1. Prozkoumejte současné metody řešení polynomiálních rovnic [1, 2] a automatický generátor postupu řešení [3, 4].
- 2. Implementujte vylepšení [4] generátoru [3] do existujícího generátoru postupů řešení.
- 3. Implementujte variaci algoritmu [1], experimentujte s jeho chováním a navrhněte, co převzít z [1] do implementace generátoru a implementujte to.
- 4. Demonstrujte funkčnost nového generátoru postupu řešení a porovnejte ji se starým generátorem.

#### Seznam odborné literatury:

- [1] Faugere, J.-C. (June 1999): "A new efficient algorithm for computing Gröbner bases (F4)". Journal of Pure and Applied Algebra (Elsevier Science) 139 (1): 61–88. doi:10.1016/S0022-4049(99)00005-5. ISSN 0022-4049.
- [2] Faugere, J.-C. (July 2002): "A new efficient algorithm for computing Gröbner bases without reduction to zero (F5)". Proceedings of the 2002 international symposium on Symbolic and algebraic computation (ISSAC) (ACM Press): 75–83.
- [3] Kukelova, Z.: Algebraic Methods in Computer Vision. PhD Thesis. CTU in Prague 2013. http://cmp.felk.cvut.cz/~kukelova/webthesis/docs/Kukelova-phd-2013.pdf
- [4] Kukelova, Z.; Bujnak, M.; Heller, J.; Pajdla, T.: Singly-Bordered Block-Diagonal Form for Minimal Problem Solvers. ACCV 2014.

Vedoucí bakalářské práce: Ing. Tomáš Pajdla, Ph.D.

Platnost zadání: do konce letního semestru 2015/2016

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Author's declaration
I declare that I have work out the presented thesis independently and that I have listed all information sources used in accordance with the Methodical Guidelines about Maintaining Ethical Principles for Writing Academic Theses.
Prohlášení autora práce
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V Praze dne

## **Abstract**

Many problems in computer vision lead to polynomial systems solving. Therefore, we need an easy way how to generate an efficient solver for each problem. On this purpose the automatic generator has been presented. In this thesis we improve the automatic generator so we will be able to generate more efficient and numerically stable solvers.

To improve the automatic generator we review and implement some methods used in the state of the art Gröbner basis solvers. Especially, we focus on the  $F_4$  Algorithm by Jean-Charles Faugère. Solvers, generated by the automatic generator, can be speed up when efficient methods are used to work with sparse matrices. We describe and implement method which is based on matrix partitioning. This method significantly speeds up the Gauss-Jordan elimination of sparse matrices.

We demonstrate the enhancements of the automatic generator on several important minimal problems. We show that the solvers generated by the new automatic generator are faster and numerically more stable than the solvers generated by the old version of the automatic generator.

**Keywords:** computer vision, robotics, minimal problems, polynomials equations, Gröbner basis

## **Abstrakt**

Mnoho problémů v počítačovém vidění vede na řešení polynomiálních rovnic. Proto potřebujeme jednoduchý způsob, jak generovat efektivní postupy řešení každého z problémů. Z tohoto důvodu byl představen automatický generátor. V této práci vylepšíme automatický generátor, takže budeme schopni generovat ještě rychlejší a numericky stabilnější postupy řešení polynomiálních systémů.

Abychom mohli vylepšit automatický generátor, prozkoumáme a následně implementujeme některé metody používané v současných nástrojích na řešení soustav polynomiálních rovnic pomocí Gröbnerových bází. Zaměříme se zejména na algoritmus  $F_4$  představený Jean-Charlesem Faugèrem. Postupy řešení problémů, vygenerované pomocí automatického generátoru, mohou být ještě dále zrychleny, pokud použijeme efektivní metody na práci s řídkými maticemi. Popíšeme a implementujeme metodu, která je založená na rozkladu matic. Tato metoda výrazně urychluje Gauss-Jordanovu eliminaci řídkých matic.

Vylepšení automatického generátoru předvedeme na několika významných minimálních problémech. Ukážeme, že postupy řešení problémů vygenerované novým automatickým generátorem jsou rychlejší a numericky stabilnější než postupy vygenerované původní verzí automatického generátoru.

**Klíčová slova:** počítačové vidění, robotika, minimální problémy, polynomiální rovnice, Gröbnerovy báze

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# List of Algorithms

1.	Simple Buchberger Algorithm
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# List of Symbols and Abbreviations

C(f), C(F)	Set of all coefficients of the polynomial $f$ or of all
	polynomials from the set $F$ .
$\deg(f)$	The total degree of $f$ .
$\overline{f}^F$	Remainder of the polynomial $f$ on division by $F$ .
$\operatorname{gcd}$	Greatest common multiple.
lcm	Least common divisor.
LC(f), LC(F)	Leading coefficient(s) of the polynomial $f$ or of all
	polynomials from the set $F$ .
LM(f), LM(F)	Leading monomial(s) of the polynomial $f$ or of all
	polynomials from the set $F$ .
LT(f), LT(F)	Leading term(s) of the polynomial $f$ or of all poly-
	nomials from the set $F$ .
M(f), M(F)	Set of all monomials of the polynomial $f$ or of all
	polynomials from the set $F$ .
$S(f_1, f_2)$	S-polynomial of polynomials $f_1$ and $f_2$ .
T(f), T(F)	Set of all terms of the polynomial $f$ or of all polyno-
	mials from the set $F$ .
$x \mid y$	x divides $y$ .
$\lfloor x \rfloor$	$= \max\{m \in \mathbb{Z} \mid m \leq x\}; \text{ floor function.}$
$\mathbb{Z}_p$	Prime field of characteristic $p$ .

### 1. Introduction

#### 1.1. Motivation

Many problems in computer vision can be formulated using systems of algebraic equations. Examples are the minimal problems [17] which arise when computing geometrical models from image data. The polynomial systems arising from this problems are often not trivial and they consists of many polynomial equations of higher degree in many unknowns and therefore, general algorithms for solving polynomial systems are not efficient for them. So special solvers for each problem have been developed to solve these system efficiently and numerically robustly.

Minimal problems have a wide range of applications, for example in 3D reconstruction, recognition, robotics and augmented reality. In these applications, the solvers of minimal problems are only small parts of large computation systems which are supposed be fast or even to work in real-time applications. Moreover, these systems need to compute the solutions of the minimal problems repeatedly for large number of parameters. Therefore, very efficient solvers are required in computer vision.

Many solvers for minimal problems have been designed ad hoc for concrete problems and therefore, they can not be used or easily modified to solve different or even similar problems. So the automatic generator [12] has been proposed. This tool generates Gröbner basis solvers automatically which enables us to generate an efficient solver for each problem we want to solve.

There are several ways, how the solver using Gröbner basis methods can be generated. The implementation presented in [12] generates polynomials that are required for solving the system systematically. But other methods can be used. In this thesis, we review the state of the art methods for solving polynomials systems and suggest which methods can be taken over to improve the automatic generator and how to implement them.

The automatic generator deals with sparse matrices in most cases. Therefore, we may consider to implement some methods which enable us to work with sparse matrices in a efficient way to save computation time and memory. In this thesis, we focus on how to improve the Gauss-Jordan elimination of sparse matrices. We use the recent work [11] which presents the significant speed up of Gauss-Jordan elimination of sparse matrices. The speed up is caused by transforming matrices into the singly-bordered block-diagonal forms by the paritioning tool PaToH [3]. This method is based on the fact that more eliminations of smaller matrices are faster than one elimination of a big matrix.

#### 1.2. Thesis structure

In this thesis, we firstly review the state of the art methods of computing Gröbner basis of polynomial systems. We start with describing simple, but easily understandable, algorithms and continue with more difficult, but also more efficient, algorithms. It is crucial to us to better understand these algorithms because we will use some techniques from them to improve the automatic generator lately in this thesis.

#### 1. Introduction

Secondly, we briefly describe the automatic generator. Then, we suggest some improvements of the automatic generator to generate efficient and numerically stable solvers. Some techniques implementened in the automatic generator may be efficient for some minimal problem, but may be inefficient for another. Therefore, we present a benchmark which enables us to choose the best methods to generate an efficient solver in the end.

Thirdly, we run some experiments to show how the implemented improvements have enhanced the automatic generator. We compare the solvers generated with the new automatic generator with the solvers generated by the old implementation.

At last, we conclude by rewieving the contributions of this thesis.

#### 1.3. Notation used

We have decided to use the notation from the [4] in the whole thesis. We just remind that polynomial is a sum of terms and term is a product of coefficient and monomial. Be aware that in some literature [1, 6, 7] the meanings of words term and monomial are switched.

# 2. Polynomial system solving

Firstly we review the state of the art algorithms for computing Gröbner basis. Better understanding of these algorithms helps us to integrate them into polynomial solving algorithms based on Gröbner basis computation more efficiently.

## 2.1. Buchberger Algorithm

Buchberger Algorithm [2], which was invented by Bruno Buchberger, was the first algorithm for computing Gröbner basis. The algorithm is described in details in [1, 4].

#### 2.1.1. First implementation

The first and easy, but very inefficient implementation of the Buchberger Algorithm, Algorithm 1, is based on the observation that we can extend a set F of polynomials to a Gröbner basis only by adding all non-zero remainders  $\overline{S(f_i, f_j)}^F$  of all pairs from F into F until there is no non-zero remainder generated.

The main disadvantage of this simple algorithm is that so constructed Gröbner basis are often bigger than necessary. This implementation of the algorithm is also very inefficient because many of the S-polynomials that are constructed from the critical pairs are reduced to zero so after spending effort on computing them, there is nothing to add to the Gröbner basis G. How to decide which pairs need not be generated is described next.

#### Algorithm 1 Simple Buchberger Algorithm

#### Input:

F a finite set of polynomials

#### **Output:**

G a finite set of polynomials

```
1: G \leftarrow F
 2: B \leftarrow \{\{g_1, g_2\} \mid g_1, g_2 \in G, g_1 \neq g_2\}
 3: while B \neq \emptyset do
           select \{g_1, g_2\} from B
           B \leftarrow B \setminus \{\{g_1, g_2\}\}\
 5:
           h \leftarrow S(g_1, g_2) \\ h_0 \leftarrow \overline{h}^G
 6:
 7:
           if h_0 \neq 0 then
 8:
                 B \leftarrow B \cup \{\{g, h_0\} \mid g \in G\}
 9:
                 G \leftarrow G \cup \{h_0\}
10:
           end if
11:
12: end while
13: return G
```

#### 2.1.2. Improved Buchberger Algorithm

The combinatorial complexity of the simple implementation of the Buchberger Algorithm can be reduced by testing out certain S-polynomials which need not be considered. To know which pairs can be deleted without treatment, we use the first and the second Buchberger's criterion [1]. Sometimes, we can even delete certain polynomials from the set G completely, knowing that every critical pair they will generate will reduce to zero and hence these polynomials themselves will be superfluous in the output set. In the next few paragraphs we will describe the implementation of the Improved Buchberger Algorithm, and of the function Update, which deletes the superfluous polynomials from G according to Gebauer and Möller [8].

The Improved Buchberger Algorithm, Algorithm 2, has the same structure as the Simple Algorithm. The function Update is used at the beginning of the Improved Buchberger Algorithm to initialize the set B of critical pairs and the Gröbner basis G from the input set F of polynomials and at every moment when a new non-zero polynomial  $h_0 = \overline{h}^G$  of an S-polynomial h has been found and the sets B and G are about to be updated.

#### Algorithm 2 Improved Buchberger Algorithm

```
Input:

F a finite set of polynomials

Output:

G a finite set of polynomials

1: G \leftarrow \emptyset
2: B \leftarrow \emptyset
3: while F \neq \emptyset do
4: select f from F
5: F \leftarrow F \setminus \{f\}
```

```
(G,B) \leftarrow Update(G,B,f)
 6:
 7: end while
     while B \neq \emptyset do
          select \{g_1, g_2\} from B
 9:
          B \leftarrow B \setminus \{\{g_1, g_2\}\}\
10:
          h \leftarrow S(g_1, g_2)
11:
          h_0 \leftarrow \overline{h}^G
12:
          if h_0 \neq 0 then
13:
               (G,B) \leftarrow Update(G,B,h_0)
14:
```

end if

16: end while 17: return G

Now, let us look at the function Update, Algorithm 3. First, it makes pairs from the new polynomial h and all polynomials from the set  $G_{old}$  and puts them into the set C. The first while loop (lines 3-9) iterates over all pairs in the set C. In each iteration it select a pair  $\{h, g_1\}$  from the set C and removes it from the set. Then, it looks for another pair  $\{h, g_2\}$  from the set C or the set D. If there is no pair  $\{h, g_2\}$  such that  $(h, g_2, g_1)$  is a Buchberger triple, then the pair  $\{h, g_1\}$  is put into the set D. The triple

15:

 $(h, g_2, g_1)$  of polynomials  $h, g_1$  and  $g_2$  is a Buchberger triple if the equivalent conditions

$$LM(g_2) \mid lcm(LM(h), LM(g_1))$$
 (2.1)

$$\operatorname{lcm}(\operatorname{LM}(h), \operatorname{LM}(g_2)) \mid \operatorname{lcm}(\operatorname{LM}(h), \operatorname{LM}(g_1)) \tag{2.2}$$

$$\operatorname{lcm}(\operatorname{LM}(g_2), \operatorname{LM}(g_1)) \mid \operatorname{lcm}(\operatorname{LM}(h), \operatorname{LM}(g_1)) \tag{2.3}$$

are satisfied. From the second Buchberger's criterion, we know that if a Buchberger triple  $(h, g_2, g_1)$  shows up in the Buchberger Algorithm and the pairs  $\{g_1, g_2\}$  and  $\{h, g_2\}$  are amongs the critical pairs, then the pair  $\{h, g_1\}$  need not be generated. That means that such a pair is not moved from the set C to the set D but it is only removed from the set C. Moreover, this while loop keeps all pairs  $\{h, g_1\}$  where LM(h) and  $LM(g_1)$  are disjoint, i.e. LM(h) and  $LM(g_1)$  have no variable in common, even if the pairs could be removed. The reason of this is that if two or more pairs in C have the same lcm of their leading monomials, then there is a choice which one should be deleted. So we keep the pair where the leading monomials are disjoint. Pairs with disjoint leading monomials are removed in the second while loop, so we eventually remove them all.

The second while loop (lines 11-17) eliminates all pairs with disjoint leading monomials. We can remove such pairs thanks to the first Buchberger's criterion. All remaining pairs are stored in the set E.

The third while loop (lines 19-25) eliminates pairs  $\{g_1,g_2\}$  where  $(g_1,h,g_2)$  is a Buchberger triple from the set  $B_{old}$ . Then the updated set of the old pairs and the new pairs are united into the set  $B_{new}$ .

Finally, the last while loop (lines 28-34) removes all polynomials g whose leading monomial is a multiple of the leading monomial of h from the set  $G_{old}$ . We can eliminate such polynomials for two reasons. Firstly,  $LM(h) \mid LM(g)$  implies  $LM(h) \mid lcm(LM(g), LM(f))$  for arbitrary polynomial f. We can see that (g, h, f) is a Buchberger triple for any f which in future appears in the set G. Moreover, polynomial g will not be missed at the end because in the Gröbner basis G, polynomials with leading monomials that are multiples of leading monomials of another polynomial from G are superfluous, i.e. they will be eliminated in the reduced Gröbner basis.

In the end of the function, the polynomial h is added into the Gröbner basis  $G_{new}$ . The output of the function Update is the Gröbner basis  $G_{new}$  and the set  $B_{new}$  of critical pairs.

# 2.2. $F_4$ Algorithm

The  $F_4$  Algorithm [6] by Jean-Charles Faugère is an improved version of the Buchberger's Algorithm. The  $F_4$  replaces the classical polynomial reduction found in the Buchberger's Algorithm by a simultaneous reduction of several polynomials. This reduction mechanism is achieved by a symbolic precomputation followed by Gaussian elimination implemented using sparse linear algebra methods.  $F_4$  speeds up the reduction step by exchanging multiple polynomial divisions for row-reduction of a single matrix.

#### **2.2.1.** Improved Algorithm $F_4$

The main function of the  $F_4$  Algorithm, Algorithm 4, consists of two parts. The goal of the first part is to initialize the whole algorithm.

### Algorithm 3 Update

```
Input:
      G_{old} a finite set of polynomials
      B_{old} a finite set of pairs of polynomials
      h a polynomial such that h \neq 0
Output:
      G_{new} a finite set of polynomials
      B_{new} a finite set of pairs of polynomials
  1: C \leftarrow \{\{h,g\} \mid g \in G_{old}\}
 2: D \leftarrow \emptyset
 3: while C \neq \emptyset do
           select \{h, g_1\} from C
 4:
 5:
           C \leftarrow C \setminus \{\{h, g_1\}\}\}
           if LM(h) and LM(g_1) are disjoint or
 6:
                     (\operatorname{lcm}(\operatorname{LM}(h), \operatorname{LM}(g_2)) \nmid \operatorname{lcm}(\operatorname{LM}(h), \operatorname{LM}(g_1)) \text{ for all } \{h, g_2\} \in C \text{ and }
                     \operatorname{lcm}(\operatorname{LM}(h), \operatorname{LM}(g_2)) \nmid \operatorname{lcm}(\operatorname{LM}(h), \operatorname{LM}(g_1)) for all \{h, g_2\} \in D) then
                 D \leftarrow D \cup \{\{h, g_1\}\}\
 7:
           end if
 8:
 9: end while
10: E \leftarrow \emptyset
11: while D \neq \emptyset do
           select \{h, g\} from D
12:
           D \leftarrow D \setminus \{\{h,g\}\}
13:
14:
           if LM(h) and LM(g) are not disjoint then
                 E \leftarrow E \cup \{\{h,g\}\}
15:
           end if
16:
17: end while
18: B_{new} \leftarrow \emptyset
      while B_{old} \neq \emptyset do
           select \{g_1, g_2\} from B_{old}
20:
21:
            B_{old} \leftarrow B_{old} \setminus \{\{g_1, g_2\}\}\
22:
           if LM(h) \nmid lcm(LM(g_1), LM(g_2)) or
                     \operatorname{lcm}(\operatorname{LM}(g_1), \operatorname{LM}(h)) = \operatorname{lcm}(\operatorname{LM}(g_1), \operatorname{LM}(g_2)) or
                     \operatorname{lcm}(\operatorname{LM}(h), \operatorname{LM}(g_2)) = \operatorname{lcm}(\operatorname{LM}(g_1), \operatorname{LM}(g_2)) then
23:
                 B_{new} \leftarrow B_{new} \cup \{\{g_1, g_2\}\}\
           end if
24:
25: end while
26: B_{new} \leftarrow B_{new} \cup E
27: G_{new} \leftarrow \emptyset
28: while G_{old} \neq \emptyset do
           select g from G_{old}
29:
           G_{old} \leftarrow G_{old} \setminus \{g\}
30:
           if LM(h) \nmid LM(g) then
31:
                 G_{new} \leftarrow G_{new} \cup \{g\}
32:
33:
           end if
34: end while
35: G_{new} \leftarrow G_{new} \cup \{h\}
36: return (G_{new}, B_{new})
```

First, it generates the set P of critical pairs and initializes the Gröbner basis G. This is done by taking each polynomial from the input set F and calling the function Update on it, which updates the set P of pairs and the set G of basic polynomials.

The second part of the algorithm generates new polynomials and adds them into the set G. In each iteration, it selects some pairs from P using the function Sel. Many selection strategies are possible and it is still an open question how to best select the pairs [6]. Some selection strategies are described in the section 2.2.6 on page 13. Then, it splits each selected pair  $\{f_1, f_2\}$  into two tuples. The first tuple contains the first polynomial  $f_1$  of the pair and the monomial  $m_1$  such that  $LM(m_1f_1) = lcm(LM(f_1), LM(f_2))$ . The second tuple is constructed in the same way from the second polynomial  $f_2$  of the pair. All tuples from all selected pairs are put into the set L, i.e. duplicates are removed.

Next, function Reduction is called on the set L. It stores its result in the set  $\tilde{F}^+$ . In the end of the algorithm it iterates through all new polynomials in the set  $\tilde{F}^+$  and calls the function Update on each of them. This generates new pairs into the set P of critical pairs and extends the Gröbner basis G.

This algorithm terminates when the set P of pairs is empty. Then the set G is a Gröbner basis and it is the output of the algorithm.

#### **Algorithm 4** Improved Algorithm $F_4$

```
Input:
```

```
F a finite set of polynomials Sel a function List(Pairs) \to List(Pairs) such that Sel(l) \neq \emptyset if l \neq \emptyset
```

#### Output:

G a finite set of polynomials

```
1: G \leftarrow \emptyset
 2: P \leftarrow \emptyset
 3: d \leftarrow 0
 4: while F \neq \emptyset do
           select f form F
           F \leftarrow F \setminus \{f\}
 6:
 7:
           (G,P) \leftarrow Update(G,P,f)
 8: end while
 9: while P \neq \emptyset do
           d \leftarrow d + 1
10:
           P_d \leftarrow Sel(P)
11:
           P \leftarrow P \backslash P_d
12:
           L_d \leftarrow Left(P_d) \cup Right(P_d)
13:
           (\tilde{F}_d^+, F_d) \leftarrow Reduction(L_d, G, (F_i)_{i=1, \dots, (d-1)})
14:
           for h \in \tilde{F}_d^+ do
15:
                (G, P) \leftarrow Update(G, P, h)
16:
17:
           end for
18: end while
19: \mathbf{return} \ G
```

#### 2.2.2. Function Update

In the  $F_4$  Algorithm, the standard implementation of the Buchberger's Criteria such as the Gebauer and Möller installation [8] is used. Details about the function *Update* can be found in the section 2.1.2. The pseudocode of the function is shown in Algorithm 3.

#### 2.2.3. Function Reduction

Function Reduction, Algorithm 5, performs polynomial division using methods of linear algebra.

The input of the function Reduction is a set L containing tuples of monomial and polynomial. These tuples were constructed in the main function of the  $F_4$  Algorithm from all selected pairs.

First, the function Reduction calls the function Symbolic Preprocessing on the set L. This returns a set F of polynomials to be reduced. To use linear algebra methods to perform polynomial division, the polynomials have to be represented by a matrix. Each column of the matrix corresponds to a monomial. Columns have to be ordered with respect to the monomial ordering used so that the most right column corresponds to "1". Each row of the matrix corresponds to a polynomial from the set F. The matrix is constructed as follows. On the (i,j) position in the matrix, we put the coefficient of the term corresponding to j-th monomial from the i-th polynomial from the set F.

We next reduce the matrix to a row echelon form using, for example, Gauss-Jordan elimination. Note that this matrix is typically sparse so we can use sparse linear algebra methods to save computation time and memory. After elimination, we construct resulting polynomials by multiplying the reduced matrix by a vector of monomials from the right.

In the end, the function returns the set  $\tilde{F}^+$  of reduced polynomials such that their leading monomials are not leading monomials of any polynomial from the set F of polynomials before reduction.

#### **Algorithm 5** Reduction

#### Input:

L a finite set of tuples of monomial and polynomial

G a finite set of polynomials

 $\mathcal{F} = (F_i)_{i=1,\dots,(d-1)}$ , where  $F_i$  is finite set of polynomials

#### Output:

 $\tilde{F}^+$  a finite set of polynomials

F a finite set of polynomials

```
1: F \leftarrow Symbolic\ Preprocessing(L, G, \mathcal{F})
```

2:  $\tilde{F} \leftarrow \text{Reduction to a Row Echelon Form of } F$ 3:  $\tilde{F}^+ \leftarrow \left\{ f \in \tilde{F} \mid \text{LM}(f) \notin \text{LM}(F) \right\}$ 

4: return  $(\tilde{F}^+, F)$ 

#### 2.2.4. Function Symbolic Preprocessing

Function Symbolic Preprocessing, Algorithm 6, starts with a set L of tuples, each containing a monomial and a polynomial. These tuples were constructed in the main

function of the  $F_4$  Algorithm from the selected pairs. Then, the tuples are simplified by the function Simplify and, after multiplying polynomials with corresponding monomials, the results are put into the set F.

Next, the function goes through all monomials in the set F and for each monomial mlooks for some polynomial f from the Gröbner basis G such  $m = m' \cdot LM(f)$  where m' is some monomial. All such polynomials f and monomials m' are, after simplification, multiplied and put into the set F. The goal of this search is to have for every monomial in F a polynomial in F with the same leading monomial. This will ensure that all polynomials from F will be reduced for G after polynomial division by linear algebra.

#### Algorithm 6 Symbolic Preprocessing

```
Input:
    L a finite set of tuples of monomial and polynomial
    G a finite set of polynomials
    \mathcal{F} = (F_i)_{i=1,\dots,(d-1)}, where F_i is finite set of polynomials
Output:
    F a finite set of polynomials
 1: F \leftarrow \{multiply(Simplify(m, f, \mathcal{F})) \mid (m, f) \in L\}
 2: Done \leftarrow LM(F)
 3: while M(F) \neq Done do
 4:
        m an element of M(F) \setminus Done
        Done \leftarrow Done \cup \{m\}
 5:
        if m is top reducible modulo G then
 6:
            m = m' \cdot LM(f) for some f \in G and some monomial m'
 7:
            F \leftarrow F \cup \{multiply(Simplify(m', f, \mathcal{F}))\}
 8:
        end if
 9:
10: end while
11: return F
```

#### 2.2.5. Function Simplify

The function Simplify, Algorithm 7, simplifies a polynomial mf which is a product of a given monomial m and a polynomial f.

The function recursively looks for a monomial m' and a polynomial f' such that LM(m'f') = LM(mf). The polynomial f' is selected from all polynomials that have been reduced in previous iterations (sets  $\tilde{F}$ ). We select polynomial f' such that the total degree of m' is minimal.

This is done in the function Symbolic Preprocessing to insert polynomials that are mostly reduced and have a small number of monomials into the set F of polynomials to be reduced. This, of course, speeds up the following reduction.

#### 2.2.6. Selection strategy

For the speed of the  $F_4$  Algorithm, it is very important how the critical pairs from the list of all critical pairs P are selected in each iteration. This depends on the implementation of the function Sel. There are more possible selection strategies:

• The easiest implementation is to select all pairs from P. In this case we reduce all critical pairs at the same time.

#### Algorithm 7 Simplify

```
Input:
    m a monomial
    f a polynomial
    \mathcal{F} = (F_i)_{i=1,\dots,(d-1)}, where F_i is finite set of polynomials
Output:
    (m', f') a non evaluated product of a monomial and a polynomial
 1: for u \in \text{list of all divisors of } m \text{ do}
        if \exists j \ (1 \leq j \leq d) such that (uf) \in F_j then
 2:
            F_j is the Row Echelon Form of F_j
 3:
            there exists a (unique) p \in \tilde{F}_i such that LM(p) = LM(uf)
 4:
            if u \neq m then
 5:
                return Simplify(\frac{m}{u}, p, \mathcal{F})
 6:
 7:
            else
                return (1, p)
 8:
            end if
 9:
        end if
10:
11: end for
12: return (m, f)
```

- If the function Sel selects only one critical pair then the  $F_4$  Algorithm is the Buchberger Algorithm. In this case the Sel function corresponds to the selection strategy in the Buchberger Algorithm.
- The best function that Faugère has tested is to select all critical pairs with a minimal total degree. Faugère calls this strategy the *normal strategy for*  $F_4$ . Pseudocode of this function can be found as Algorithm 8.

```
Algorithm 8 Sel – The normal strategy for F_4
```

```
Input:

P a list of critical pairs

Output:

P_d a list of critical pairs

1: d \leftarrow min \{ deg(lcm(p)) \mid p \in P \}

2: P_d \leftarrow \{ p \in P \mid deg(lcm(p)) = d \}

3: return P_d
```

# **2.3.** $F_5$ Algorithm

Since in the Buchberger Algorithm or in the  $F_4$  Algorithm we spend much computation time to compute S-polynomials which will reduce to zero, the  $F_5$  Algorithm [7] by Jean-Charles Faugère was proposed to eliminate these reductions to zero. The  $F_5$  Algorithm saves computation time by removing useless critical pairs which will reduce to zero. The syzygies [4] are used to recognize useless critical pairs in advance.

There are several approaches how to use syzygies to remove useless pairs. For example the idea of [13] is to compute a basis of the module of syzygies together with the

computing of the Gröbner basis of the given polynomial system. Then a critical pair can be removed if the corresponding syzygy is a linear combination of the elements of the basis of syzygies.

The strategy of the  $F_5$  Algorithm is to consider only principal syzygies without computing the basis of the syzygies. The principal syzygy is a syzygy such that  $f_i f_j - f_j f_i = 0$  where  $f_i$  and  $f_j$  are polynomials. This restriction implies that not all useless critical pairs have to be removed so a reduction to zero can still appear later. However it was proved that if the input system is a regular sequence then there is no reduction to zero.

To show how to distinguish which pairs need not be considered, we use the example taken from [7]. Consider polynomials  $f_1$ ,  $f_2$  and  $f_3$ . Then, the principal syzygies  $f_i f_j - f_j f_i = 0$  can be written as follows:

$$u(f_2f_1 - f_1f_2) + v(f_3f_1 - f_1f_3) + w(f_2f_3 - f_3f_2) = 0 (2.4)$$

where u, v and w are arbitrary polynomials. This can be also rewritten as

$$(uf_2 + vf_3)f_1 - uf_1f_2 - vf_1f_3 + wf_2f_3 - wf_3f_2 = 0. (2.5)$$

We can see that all relations  $hf_1$  are such that h is in the ideal generated by polynomials  $f_2$  and  $f_3$ . So if we have computed Gröbner basis of the polynomials  $f_2$  and  $f_3$ , it is easy to decide which new generated polynomials can be removed. We can remove all polynomials in the form  $tf_1$  such that t is a term divisible by leading monomial of an element of the ideal generated by  $f_2$  and  $f_3$ . Therefore the  $F_5$  Algorithm is an incremental algorithm so if we have polynomials  $f_1, \ldots, f_m$  on the input we have to compute all Gröbner basis of the following ideals:  $(f_m), (f_{m-1}, f_m), \ldots, (f_1, \ldots, f_m)$  in this order.

Many reviews, implementations and modifications of the  $F_5$  Algorithm have been made. Let us emphasize some of them. The first implementation of the  $F_5$  was made by Jean-Charles Faugère himself in the language C. Then there is an implementation in Magma by A. J. M. Segers [18]. Another review and implementation in Magma was done by Till Stegers [19]. Since there is no proof of termination of the  $F_5$  Algorithm, a modification [5] has been introduced such that it always terminates.

# 3. Automatic generator

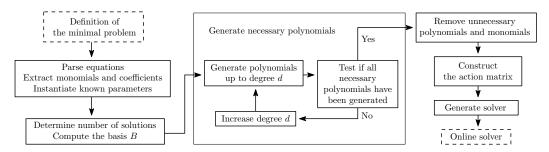
The automatic generator of Gröbner basis solvers is used to easily solve problems leading to systems of polynomial equations. These systems usually arise when solving minimal problems [17] in computer vision. Typically, these systems are not trivial so special solvers have to be designed for concrete problems to achieve efficient and numerically stable solvers. But solvers generated for concrete problems can not be easily applied for similar or new problems and therefore the automatic generator was proposed in [12]. Solvers generated by the automatic generator can be easily used to solve complex problems even by non-experts users.

The input of the automatic generator is a system of polynomial equations with a finite number of solutions and the output is a MATLAB or a Maple code that computes solutions of the given system for arbitary coefficients. One of the goals of this thesis is to improve previous implementation [12] of the automatic generator to construct more efficient and numerically stable solvers.

The newest version of the automatic genenerator implemented in MATLAB can be downloaded from [16].

### 3.1. Description of the automatic generator

In this section we would like to briefly describe the procedure for generating solvers. The procedure is based on computation of the action matrix from which solutions can be obtained. The automatic generator consists of several independent modules, see Figure 3.1. Since all these modules are independent, they can be easily improved or replaced by more efficient implementations. Next we describe each of these modules, full description can be found in [12, 10].



**Figure 3.1.** Block diagram of the automatic generator

#### 3.1.1. Definition of the minimal problem

Definitions of minimal problems to be solved are given in separate functions that are stored in the folder minimalProblems. Each of the definitions has to contain the necessary information about the minimal problem. First of all, the system of polynomial equations with symbolic variables and parameters has to be provided. Next, we have to specify the list of unknown variables and known parameters. Optionally, if we know the monomial basis B of the polynomial system in advance, we can specify it to save

some computation time. The monomial basis B is a set  $\{m \mid \overline{m}^G = m\}$  where m is a monomial and G is the Gröbner basis of the given polynomial system. At last we have to set some settings for the automatic generator. We recommend to obtain the default settings by calling the function  $gbs\_InitConfig()$  and only overwrite the settings we want to change. In the folder minimalProblem, there are some examples which are self explanatory and can be used as templates to create new minimal problem definitions.

#### 3.1.2. Equations parser, instantiating

In the next step, we have to parse the given equations, which means that we extract monomials and parameters used and obtain total degrees of the polynomials. Then, we instantiate each known parameter with a random number from  $\mathbb{Z}_p$ . We assign a unique identifier to each parameter used. The reason is that we need to track the parameters through the process of manipulating with the polynomials in order to be able to restore the process in the solver generation module.

#### 3.1.3. Monomial basis B computation

We need to know the monomials basis B to recognize when we have generated all polynomials that are necessary to build the action matrix. If the basis B was not provided within the definition of the minimal problem, we have to compute it by ourselves. Because there is no function or simple script to compute the basis in MATLAB we have to do it by calling an external software.

The easiest solution to implement was to use the Maple toolbox for MATLAB. This enables us to call Maple functions from the MATLAB environment directly. To use this option we have to set cfg.GBSolver = @gbs\_findAlgB\_maple in the settings of the automatic generator. Unfortunately it shows up that the Maple toolbox for MATLAB in not compatible with the MATLAB Symbolic Math Toolbox in versions newer than R2008 so we do not recommend to use this option nowadays, but the option is still available on older computers.

The second implemented option is to use the algebraic geometry software Macaulay2 [9]. In the folder gbsMacaulay there is a template code\_template.m2 into which we simply write the given polynomial system. This updated file is saved as code.m2 which is executed by Macaulay2 and the results are parsed back in MATLAB. To set up this option we need to install the software Macaulay2 and set cfg.GBSolver = Qgbs\_findAlgB\_macaulay in the automatic generator settings. A problem could be that the Macaulay2 is not easy to set up under the Windows OS. Therefore, the installation file of Macaulay2 is provided within the automatic generator. The only thing that has to be done is to edit the file calc.bat in the folder gbsMacaulay and follow the instructions in the file.

Because of the modularity of the generator, this part can be replaced by another function computing the monomial basis B.

The last option is to compute the basis B in advance and set it into the definition of the minimal problem.

In the end, we have to check the number of solutions of the given polynomial system. If there is a finite number of solutions we can continue with the computation.

#### 3.1.4. Polynomial generator

To be able to build the action matrix, we have to generate enough polynomials such that after their reduction we get polynomials  $q_i$  which have leading monomials from

the set  $(x_k \cdot B) \setminus B$  where  $x_k$  is a variable and all remaining monomials of  $q_i$  are from the set B. That is the reason why we had to compute the basis B in the previous step.

In this part of the automatic generator, we represent polynomials as row vectors so that systems of polynomials can be represented by matrices. This representation enables us to easily multiply polynomials with monomials only by shifting the coefficients in the vectors or to reduce the whole polynomial systems by performing the Gauss-Jordan eliminations on the corresponding matrices.

Let  $f_i \in F$  where F is a set of polynomials from the input. Let maxdeg is a maximal total degree of all polynomials  $f_i$ . At the beginning we put all polynomials  $\{mf_i \mid f_i \in F; \deg(mf_i) = \deg(f_i), \ldots, maxdeg\}$  into the matrix M, where m is a monomial. Now we perform the Gauss-Jordan elimination on the matrix M and save the result as matrix  $\tilde{M}$ . Then we check if there exists a variable  $x_k$  for which all required polynomials  $q_i$  are present in  $\tilde{M}$ . If we find such a variable we can continue with the construction of the action matrix for the variable. If not, we have to add more polynomials to the matrix M. We increment maxdeg by one and add all polynomials  $\{mf_i \mid f_i \in F; \deg(mf_i) = maxdeg\}$  to the matrix M. Then we continue with the elimination and with the checking the action matrix requirements as explained above. We repeat these steps until all required polynomials  $q_i$  are generated so the action matrix can be built. The pseudocode of this process is shown in Algorithm 9. The function Check Action Matrix Conditions from this code checks if all polynomials  $q_i$  are generated in  $\tilde{M}$  for at least one variable from the given list of variables. If such a variable is found, the function returns it, otherwise it returns an empty set.

#### **Algorithm 9** Polynomial generator – One elimination solver

#### Input:

```
F a set of polynomials variables a list of variables algB a monomial basis B
```

## Output:

M a matrix representing a set of polynomials var a variable

```
1: maxdeg \leftarrow \max\{\deg(f_i) \mid f_i \in F\}

2: M \leftarrow \{mf_i \mid f_i \in F; \deg(mf_i) = \deg(f_i), \dots, maxdeg; m \text{ is a monomial}\}

3: \tilde{M} \leftarrow \text{Reduction to a Row Echelon Form of } M

4: var \leftarrow Check \ Action \ Matrix \ Conditions(\tilde{M}, variables, algB)

5: \mathbf{while} \ var = \emptyset \ \mathbf{do}

6: maxdeg \leftarrow maxdeg + 1

7: M \leftarrow M \cup \{mf_i \mid f_i \in F; \deg(mf_i) = maxdeg; m \text{ is a monomial}\}

8: \tilde{M} \leftarrow \text{Reduction to a Row Echelon Form of } M

9: var \leftarrow Check \ Action \ Matrix \ Conditions(\tilde{M}, variables, algB)

10: \mathbf{end \ while}

11: \mathbf{return} \ (\tilde{M}, var)
```

In this whole process we need to keep track how the matrix M was built. Recall that each coefficient of the polynomials  $f_i$  has a unique identifier assigned to it in the equations parser. Because the whole matrix M contains only the polynomials  $f_i$  or their multiples with monomials, only the coefficients from the polynomials  $f_i$  appear in the matrix M. We just have to keep the positions of the coefficients. This is done by matrix  $\hat{M}$ . The matrix  $\hat{M}$  is built at the same time as the matrix M by this way. When

we put a coefficient into the matrix M, we also put the corresponding indentifier to the matrix  $\hat{M}$  at the same possition. The matrix  $\hat{M}$  enables us to recover the process of polynomial generation in the code generator module.

#### 3.1.5. Removing unnecessary polynomials and monomials

Since the polynomials were generated systematically in the previous step there may appear some polynomials which are not necessary for the constructing of the action matrix. The goal of this part of the automatic generator is to remove as many as possible polynomials which are not necessary.

We can remove a polynomial r from the matrix M if the corresponding eliminated matrix  $\tilde{M}$  still contains all required polynomials  $q_i$ . In this way we try to remove all polynomials from M.

Because the success of removing a polynomial depends on the previous removals, the number of removed polynomials depends on the ordering in which the polynomials are removed. In the automatic generator we start removing polynomials from the one with the largest leading monomial by monomial ordering used. Because it is very inefficient to remove polynomials one by one and perform each time an expensive Gauss-Jordan elimination, we can enhance the procedure by trying to remove more polynomials at the time. In the automatic generator, this heuristic is used. If we have successfully removed k polynomials, we try to remove 2k polynomials in the next step. If the removal of k polynomials have failed, we try to remove  $\frac{1}{4}k$  polynomials in the next step. The pseudocode of this removing process is shown as Algorithm 10.

Moreover we can reduce the size of the matrix M by removing unnecessary monomials. A monomial is unnecessary when its removal does not affect the building of the action matrix. We have to keep all monomials such that they are leading monomials of polynominals in the corresponding matrix  $\tilde{M}$  and all monomials that are present in the basis B. All other monomials can be removed. If we remove all such unnecessary monomials then the matrix M will have dimensions  $n \times (n+N)$  where n is the number of the polynomials in the matrix M and N is the number of solutions of the given system.

#### 3.1.6. Construction of the action matrix

This part of the automatic generator starts with the eliminated matrix  $\tilde{M}$  of polynomials and variable  $x_k$  for which all required polynomials  $q_i$  are present in  $\tilde{M}$ .

Let us describe the construction of the action matrix in an informal and practical way rather than by using the theory. If the theory is needed, it can be found in [10]. The action matrix  $M_{x_k}$  corresponding to the variable  $x_k$  is a square matrix of dimensions  $N \times N$  where N is the number of elements of the monomial basis B. Each row and column of  $M_{x_k}$  corresponds to a monomial  $b_i \in B$ . Let the monomials  $b_i$  be sorted such that if  $b_l \prec b_k$  then k < l where  $\prec$  is the monomial ordering used. We put coefficients of the polynomial  $m_i = \overline{(x_k b_i)}^F$  to the i-th row where F are polynomials corresponding to  $\tilde{M}$ . Because  $\tilde{M}$  is in a row echelon form there are two possibilities how the i-th row can be constructed:

- 1.  $x_k b_i = b_j$  for some  $b_j \in B$ . That means that  $x_k b_i$  is irreducible by F and  $m_i$  is a monomial in B. In this case we set  $M_{x_k}(i,j) = 1$  and  $M_{x_k}(i,k) = 0$  where  $k \neq j$ .
- 2.  $x_k b_i \neq b_j$  for all  $b_j \in B$ .

#### Algorithm 10 Remove unnecessary polynomials

#### Input:

```
M a matrix representing a set of polynomials variable a variable alg B a monomial basis B
```

#### **Output:**

M a matrix representing a set of polynomials

```
1: rows \leftarrow number of rows of M
 2: step \leftarrow \max\{|rows/32|, 1\}
 3: up \leftarrow 1
 4: filter \leftarrow \{1, 2, \dots, rows\}
 5: while up \leq rows do
         down \leftarrow up + step - 1
 6:
         if down > rows then
 7:
             down \leftarrow rows
 8:
             step \leftarrow down - up + 1
 9:
         end if
10:
         filter_{Old} \leftarrow filter
11:
         filter \leftarrow filter \setminus \{up, up + 1, \dots, down\}
12:
         M \leftarrow \text{Reduction to a Row Echelon Form of } M \text{ only with rows specified by } filter
13:
         v \leftarrow Check\ Action\ Matrix\ Conditions(M, variable, alg B)
14:
         if v = variable then
15:
             up \leftarrow down + 1
16:
             step \leftarrow 2step
17:
18:
         else
             if step = 1 then
19:
                 up \leftarrow up + 1
20:
             else
21:
                 step \leftarrow \max\{|step/4|, 1\}
22:
23:
24:
             filter \leftarrow filter_{Old}
         end if
25:
26: end while
27: return M only with rows specified by filter
```

In this case there is f such that  $LM(m_i) = LM(f)$  where  $f \in F$  so  $m_i = x_k b_i - f$ . Since all monomials of f except LM(f) are from B, all monomials of  $m_i$  are also from B. We put coefficient of  $m_i$  at the monomial  $b_j$  to the matrix  $M_{x_k}$  on the (i, j) position.

Now, the solutions of the given system can be easily found by computing right eigenvectors of the action matrix  $M_{x_k}$ .

#### 3.1.7. Solver generator

The last task of the automatic generator is to create a solver which will solve the given polynomial system for an arbitrary set of parameters preserving its structure. The current version of the automatic generator can generate solvers for MATLAB and Maple but new code generators can be easily added. It can be set in the min-

imal problem definition by setting cfg.exportCode which solvers will be generated. For instance to create both MATLAB and Maple solvers we set cfg.exportCode = {'matlab' 'maple'}.

To create a solver we have to restore the process of creation of the matrix M. This process is saved as the matrix  $\hat{M}$  which contains unique identifiers on the positions where the given parameters have to be put. So the matrix M can be built for each given set of parameters. Then the Gauss-Jordan elimination is called on M so we get the matrix  $\tilde{M}$ . Now the action matrix is built in the same way as above and the solutions are extracted from it. To sum up the final solver just creates the matrix M by putting parameters to the correct places. After Gauss-Jordan elimination the action matrix is built by copying some parts of rows of  $\tilde{M}$  and then the solutions are extracted by using the eigenvectors of the action matrix.

#### 3.1.8. Usage

The automatic generator is designed to be used even by non-expert users and to be easily expanded or improved.

At first, the script setpaths.m should be executed from the root directory of the automatic generator. This will add all required paths to the MATLAB environment.

Next we have to set up the definition of the minimal problem we want to solve. How the definition have to be specified is written in the section 3.1.1. All these definitions are stored in the folder minimalProblems. To generate the solver we call the function gbs\_GenerateSolver(MinimalProblem) where MinimalProblem is the name of the definition of the minimal problem, i.e. the name of the function in the folder minimalProblems. This will generate us solver solver MinimalProblem.m for the MATLAB solver and solver\_MinimalProblem.txt for the Maple solver. These solvers are stored in the folder solvers.

For example, let us present generating of a solver for the 6-point focal length problem [14]. We have defined this problem as a function sw6pt.m in the folder minimalProblems. By calling the function gbs\_GenerateSolver('sw6pt') we get solvers solver\_sw6pt.m and solver\_sw6pt.txt in the folder solvers.

### 3.2. Improvements

The bottleneck of the automatic generator is the polynomial generator module. Since the polynomials are generated systematically, the matrices in the resultant solvers are often bigger than necessary which means that the solvers are not efficient. So, many improvements of the automatic generator [12] can be done. For example, if we want to generate multiple elimination solvers as suggested in [10], the polynomial generator module have to be improved or totaly replaced. In the same way, some strategies from other algorithms, for example from the  $F_4$  Algorithm [6], can be taken over and implemented into the automatic generator. Because we are mostly working with sparse matrices in the automatic generator, Gauss-Jordan elimination for sparse matrices can be implemented to save some computation time.

#### 3.2.1. Reimplementation

The previous implementation [12] of the automatic generator was implemented in MATLAB R2008. It shows up that new versions of MATLAB are not backward compatible so the automatic generator fails when lauched in some newer version than R2008.

Our first task was to reimplement the old implementation into new version of MATLAB. All changes of the code were minor and just in the implementation level.

One important problem was that the Maple Toolbox for MATLAB in not compatible with the MATLAB Symbolic Math Toolbox anymore. One of the options in the monomials basis B module was to use the Maple toolbox to compute the basis B. Because of the new incompatibility, this option can no longer be used and the algebraic geometry software Macaulay2 [9] have to be used instead.

The new version of the automatic generator, as it is desribed in this thesis, is compatible with the version R2015 of MATLAB 64-bit and 32-bit under Windows and Unix operation systems.

#### 3.2.2. Multiple elimination solver

In the section 3.1.4, there is described a strategy how to generate polynomials for one elimination solvers. However, it may be better to create multiple elimination solvers in some cases. A multiple elimination solver is a solver where polynomials are generated systematically by multiplying already generated polynomials by monomials and reduced each time by Gauss-Jordan elimination. So the task of this section is to describe how the polynomial generator of the automatic generator can be improved to be able to generate polynomials for multiple elimination solvers.

To generate polynomials for multiple elimination solvers, we generate all polynomials up to degree d in each step and then, we perform a Gauss-Jordan elimination on them. We increase the degree d when no new polynomials can be generated by multiplying already generated polynomials by some monomials. We stop this process when all polynomials  $q_i$  are generated.

This strategy is very usefull especially when we are generating solvers for systems with many variables. The reason is that the increasing of the degree d of generated polynomials leads to a large number of new generated polynomials.

Now, let us look on the process of generating polynomials in more details. The pseudocode is shown as Algorithm 11. Let the maxdeq is the maximal total degree of all polynomials from the given system F. At the beginning, we put into the matrix  $M_1$  all polynomials up to degree maxdeg such that they are product of polynomials from F and some monomials. We get the matrix  $\tilde{M}_1$  by eliminating the matrix  $M_1$ . We check if there exists a variable for which all polynomials  $q_i$  are generated in  $M_1$ . If no such variable is found, we generate new polynomials with higher total degree. We increase the degree maxdeg which is the maximal total degree of all already generated polynomials. The variable step tells how much we want to increase the total degree in one step. We save the maxdeg + 1 from the previous iteration into the variable mindegto keep the track to which degree we have generated polynomials already. We get new matrix  $M_2$  by copying the matrix  $M_1$  and we add all polynomials with total degrees from mindeg to maxdeg to  $M_2$ . These polynomials are multiples of polynomials from  $\tilde{M}_1$  by some monomials. We save the result of the Gauss-Jordan elimination as the matrix  $M_2$ . We repeat this process until no new polynomials are added in the iteration. That situation happens when two matrices  $M_j$  and  $M_{j-1}$  have the same number of non zero rows. In this case we check if all polynomials  $q_i$  are generated for some variable. If not, we have to generate new polynomials with higher total degree. If the variable has been found, we return the last generated matrix  $M_j$  and the variable. You may notice that when we are leaving the repeat-until loop on line 13, there are two equivalent matrices  $M_j$  and  $M_{j-1}$ . These matrices have the same non zero rows and differ only with the number of zero rows. This is the reason why we can remove the matrix  $M_i$  by

#### Algorithm 11 Polynomial generator – Multiple elimination solver

```
Input:
     F a set of polynomials
     variables a list of variables
     algB a monomial basis B
     step an integer
Output:
     M a matrix representing a set of polynomials
     var a variable
 1: maxdeg \leftarrow max\{deg(f_i) \mid f_i \in F\}
 2: i \leftarrow 1
 3: M_1 \leftarrow \{mf_i \mid f_i \in F; \deg(mf_i) = \deg(f_i), \dots, maxdeg; m \text{ is a monomial}\}
 4: M_1 \leftarrow \text{Reduction to a Row Echelon Form of } M_1
 5: var \leftarrow Check\ Action\ Matrix\ Conditions(M_1, variables, alg B)
 6: while var = \emptyset do
         mindeg \leftarrow maxdeg + 1
 7:
         maxdeg \leftarrow maxdeg + step
 8:
         repeat
 9:
             j \leftarrow j + 1
10:
             M_j \leftarrow \tilde{M}_{j-1} \cup \left\{ mf_i \mid f_i \in \tilde{M}_{j-1}; \right.
11:
                   deg(m \times f_i) = mindeg, \dots, maxdeg; m \text{ is a monomial}
             \tilde{M}_i \leftarrow \text{Reduction to a Row Echelon Form of } M_i
12:
         until number of nonzero rows of M_i = number of nonzero rows of M_{i-1}
13:
14:
         j \leftarrow j-1
         var \leftarrow Check\ Action\ Matrix\ Conditions(M_i, variables, algB)
15:
16: end while
17: return (M_i, var)
```

We have to keep track how the matrices  $M_j$  were built to be able to restore the process of generation of polynomials and generate the code of the solver in the solver generator module. Therefore, we build a matrix  $\hat{M}_j$  for each matrix  $M_j$ . The first matrix  $\hat{M}_1$  is built in the same way as the matrix  $\hat{M}$  when generating one elimination solver. This matrix contains only the unique identifiers of parameters on positions were real values will be put. Because each matrix  $M_j$  is built from the matrix  $\tilde{M}_{j-1}$ , the matrix  $M_j$  contains only coefficients from the matrix  $\tilde{M}_{j-1}$ . So, when a coefficient from (m,n) position in  $\tilde{M}_{j-1}$  is put into  $M_j$  at (k,l) position, the tuple (m,n) is put at the position (k,l) of  $\hat{M}_j$ . So if we have the matrix  $\tilde{M}_{j-1}$  and the matrix  $\hat{M}_j$ , the matrix  $M_j$  can be built easily.

To enable generation of multiple elimination solvers, we assign an integer to cfg.PolynomialsGeneratorCfg.GJstep in the settings of the automatic generator. The value of the GJstep has the same meaning as the variable *step* from the Algorithm 11. E.g. by setting GJstep = 1 the generated polynomials will be eliminated each time the total degree of generated polynomials is incremented by 1. If we set GJstep = 0, the one elimination solver will be generated as described in the section 3.1.4.

#### 3.2.3. Removing redundant polynomials

We may notice that there appear matrices  $\tilde{M}_j$  with many zero rows in the process of generation polynomials for multiple elimination solvers. This is because there are many dependent rows in the corresponding matrices  $M_j$ . These zero rows in matrices  $\tilde{M}_j$  give us no new information so we want to remove as many as possible rows from the matrices  $M_j$  such that the resulting matrices  $\tilde{M}_j$  will have the same nonzero rows as before the removal and will have no zero rows.

We know that we can remove the same number of rows from  $M_j$  as it is the number of zero rows in  $\tilde{M}_j$ . But we do not know which rows we should remove. So we try to remove each row r from  $M_j$  and if the number of nonzero rows of  $\tilde{M}_j$  stays the same, the removal is successful, if not, we have to return the row r back into  $M_j$ . We end this process of removal when there is no zero row in the matrix  $\tilde{M}_j$ . Because performing the Gauss-Jordan elimination in each step of removing single row is inefficient we use the same heuristic as desribed in the section 3.1.5. For better understanding, we are providing the pseudocode of this removing as Algorithm 12.

Because this removing process removes only zero rows from the matrices  $\tilde{M}_j$  and no others rows are touched it does not influences the process of adding polynomials. Therefore, this removing process is enabled by default and can not be disabled by any option in the automatic generator settings.

#### 3.2.4. Matrix partitioning

In the automatic generator, we are dealing with matrices that are mostly sparse, so some efficient techniques can be used to work with them. This will, of course, result in generation of more efficient and numerically stable solvers.

In this section, we focus on how to speed up the Gauss-Jordan elimination of sparse matrices. We use the technique proposed in [11]. We observe that by permuting the rows and the columns of sparse matrices they can be transformed into matrices with block-diagonal structure known as singly-bordered block-diagonal (SBBD) form. Each diagonal block of the SBBD matrices forms an independent problem and therefore, it can be independently eliminated. This speeds up the process of Gauss-Jordan elimination because eliminating more smaller matrices is faster than eliminating one big matrix. If we divide the matrix into k independent blocks that contain comparable number of entries, the speed up is approximately  $n^3 \to k \cdot \left(\frac{n}{k}\right)^3$ . Moreover, the permutation matrices that transform matrices to the SBBD forms are precomputed during the solver generation process and therefore, the resultant solver is working already with matrices in the SBBD forms and does not have to spend time by computing the permutation matrices again.

Let us say we want to eliminate matrix M. First of all, we have to remove columns that correspond to monomials from the basis B from the matrix because these columns should not be permuted and eliminated. Then, we need to compute the permutation matrix. To compute the permutation matrices, we use the state of the art hypergraph partitioning tool PaToH [3] with settings to divide the matrix into two independent blocks. We do the permutations of rows a columns and get two diagonal blocks  $M_{11}$  and  $M_{22}$  and coupling columns that can not be assigned to any of the diagonal blocks. Next, we perform two independent eliminations of the blocks  $M_{11}$  and  $M_{22}$  and permute all rows of M to get the identity matrix in the left top corner. However, we get non-eliminated submatrix in the right bottom corner of M. After eliminating this submatrix, the rows above this submatrix are still not eliminated. If this is the last elimination in

#### Algorithm 12 Remove redundant polynomials

#### Input:

M a matrix representing a set of polynomials

 $\tilde{M}$  a matrix in a Row Echelon Form representing a set of polynomials

#### Output:

M a matrix representing a set of polynomials

M a matrix in a Row Echelon Form representing a set of polynomials

```
1: toRemove \leftarrow number of zero rows of M
 2: nonZero \leftarrow number of non zero rows of \tilde{M}
 3: rows \leftarrow number of rows of M
 4: step \leftarrow \max\{|toRemove/4|, 1\}
 5: up \leftarrow 1
 6: filter \leftarrow \{1, 2, \dots, rows\}
 7: while toRemove \neq 0 do
        down \leftarrow up + step - 1
        if down > rows then
 9:
10:
             down \leftarrow rows
             step \leftarrow down - up + 1
11:
        end if
12:
         filter_{Old} \leftarrow filter
13:
         filter \leftarrow filter \setminus \{up, up + 1, \dots, down\}
14:
        M \leftarrow \text{Reduction to a Row Echelon Form of } M \text{ only with rows specified by } filter
15:
        if number of non zero rows of M < nonZero then
16:
             if step = 1 then
17:
                 up \leftarrow up + 1
18:
19:
             else
                 step \leftarrow \max\{|step/4|, 1\}
20:
             end if
21:
             filter \leftarrow filter_{Old}
22:
23:
        else
             toRemove \leftarrow number of zero rows of M
24:
             up \leftarrow down + 1
25:
             step \leftarrow \min\{2step, toRemove\}
26:
        end if
27 \cdot
28: end while
29: return (M only with rows specified by filter, M)
```

the solver, we have to eliminate only the rows which we need to build the action matrix from. If this is not the last elimination in the solver, we have to eliminate them all.

**Example** Consider sparse matrix M of size  $119 \times 143$  as it is shown in Figure 3.2a. Columns 119 - 143 correspond to monomials from the monomial basis B. When we remove these columns, we get a rectangular matrix of size  $119 \times 119$  on which we execute the partitioning tool PaToH. After the permutation of the rows and columns, we get the matrix in the SBBD form which can be seen in Figure 3.2b. We get two diagonal blocks  $M_{11}$  of size  $71 \times 30$  and  $M_{22}$  of size  $48 \times 38$  and 51 coupling columns which can not be assigned to any diagonal block. After two independent Gauss-Jordan eliminations of the blocks  $M_{11}$  and  $M_{22}$  and row permutation, we get the identity matrix in the left

#### 3. Automatic generator

top corner, see Figure 3.2c. Now, we have to perform the Gauss-Jordan elimination on submatrix of size  $51 \times 75$  in the right bottom corner. After this, we get matrix as it is shown in Figure 3.2d. Now, uneliminated submatrix of size  $68 \times 75$  lefts in the right top corner of the matrix. If this is the last elimination in the solver, we do not have to eliminate it whole, but it is sufficient to eliminate only the rows we need to the constructing of the action matrix. If this is not the last elimination in the solver, we have to eliminate the whole remaining uneliminated submatrix.

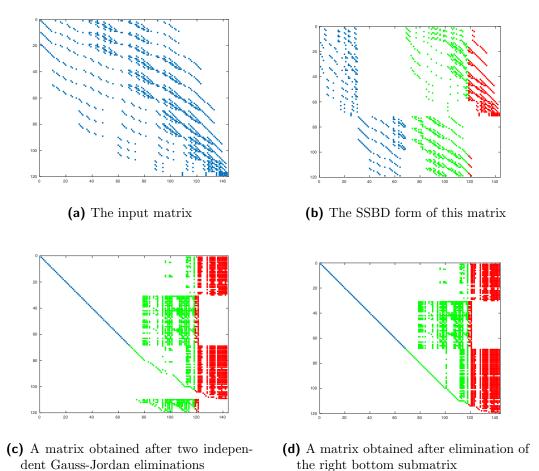


Figure 3.2. Example of the process of eliminating matrix using matrix partitioning. Colors: blue – two independent diagonal blocks, green – coupling columns, red – columns corresponding to the basis B

Since matrix partitioning does not have to be efficient for all minimal problems, it can be easily enabled or disabled in the automatic generator. By setting cfg.matrix-Partitioning = 'all' matrix partitioning is used to all Gauss-Jordan eliminations in the solver. If cfg.matrixPartitioning = 'last' is set, matrix partitioning is used only to the last Gauss-Jordan elimination in the solver. Matrix partitioning can be totally disabled by setting cfg.matrixPartitioning = 'none'. At last, we want to warn that the tool PaToH is not available under Windows OS so the matrix partitioning can not be used under this system.

#### **3.2.5. F4** strategy

Because polynomials are generated systematically in the polynomial generator module, as it is described in section 3.1.4, many of them are supefluous and therefore, need not be generated. Since the automatic generator consists of independent modules, we can replace the polynomial generator module by some better implementation.

We have described some of the state of the art techniques how polynomial systems can be solved in the chapter 2. Therefore, we can take over some of them and implement them into the automatic generator. In the section 2.2, we have understood the  $F_4$  [6] Algorithm so we have implemented this technique into the automatic generator.

#### Implementation in Maple

Before we could start implementing the  $F_4$  strategy into the automatic generator, we had to deeply understand the  $F_4$  algorithm. Therefore, we have implemented the  $F_4$  Algorithm in Maple according to [6].

We have choosen the software Maple because, there is implemented the  $F_4$  Algorithm by J. Ch. Faugère himself in the Groebner package. Therefore, we are able to compare our implementation with the implementation by Faugère.

Our implementation of the  $F_4$  algorithm is available at http://cmp.felk.cvut.cz/~trutmpav/bachelor-thesis/F4 and it is divided into the same functions as described in [6] and we have described in the section 2.2. The Gröbner basis generated by the  $F_4$  Algorithm are not reduced Gröbner basis and therefore, we add the reduction of Gröbner basis to get the reduced Gröbner basis to the end of the algorithm. This enables us to easily compare the results with results computed by the Faugère's implementation.

The main function is named  $F_4$  in our implementation and it is called F4(F, Se1, ordering) where F is a set of input polynomials, Se1 is a function which selects critical pairs as descibed in the section 2.2.6 and ordering is a monomial ordering. Output of this function is a Gröbner basis. Our implementation prints information as the number of pairs, number of selected pairs and their total degree and sizes of matrices during the computation, see an example of the output in Figure 3.3. The implementation by Faugère can be called by Groebner[Basis] (F, ordering, method=fgb) where the parameters F and ordering have the same meanings as the parameters in our implementation. We can force the Faugère's implementation to print some information about the process by setting infolevel[GroebnerBasis] := 5, see an example of the output in Figure 3.4. Therefore, we can compare, not only the results, but the sizes of the matrices, too. There are many choices in the  $F_4$  Algorithm which can be differently implemented. For example, the list of divisors may be sorted differently in the function Simplify. Therefore, the sizes of the matrices may differ in our implementation and in Faugère's implementation, as you can see in the example below.

**Example** We show both implementations on the cyclic 4 problem taken from [6]. You can see the input and the output of our implementation in Figure 3.3 and of Faugère's implementation in Figure 3.4.

#### Integration into the automatic generator

Implementation of the  $F_4$  Algorithm in the automatic generator is the same as J. Ch. Faugère has described in [6] and we did in section 2.2. The only difference is that we need to track how the polynomials are constructed to be able to reconstruct the process in the solver generator module. In the  $F_4$  Algorithm, to be concrete, in the function

```
> F4([a*b*c*d - 1, a*b*c + a*b*d + a*c*d + b*c*d, a*b + b*c + a*d + c*d, a + b + c + d],
     SelMinDeg, tdeg(a, b, c, d));
        0.00: |P| = 3, Selected 1 of degree 2
        0.00: Symbolic Preprocessing: |L| = 2, |Fd| = 0
        0.01: Reducing matrix of size 3x7
        0.02: Reduction finished
        0.02: |P| = 2, Selected 1 of degree 3
        0.02: Symbolic Preprocessing: |L| = 2, |Fd| = 3
        0.03: Reducing matrix of size 4x9
        0.03: Reduction finished
        0.03: |P| = 2, Selected 2 of degree 4
        0.03: Symbolic Preprocessing: |L| = 4, |Fd| = 7
        0.04: Reducing matrix of size 8x12
        0.05: Reduction finished
        0.05: |P| = 2, Selected 2 of degree 5
        0.05: Symbolic Preprocessing: |L| = 4, |Fd| = 15
        0.06: Reducing matrix of size 6x12
        0.06: Reduction finished
        0.06: |P| = 3, Selected 3 of degree 6
        0.07: Symbolic Preprocessing: |L| = 6, |Fd| = 21
        0.08: Reducing matrix of size 10x14
        0.08: Reduction finished
        0.09: |P| = 2, Selected 2 of degree 7
        0.09: Symbolic Preprocessing: |L| = 4, |Fd| = 31
        0.10: Reducing matrix of size 7x10
        0.10: Reduction finished
        0.10: Computing reduced Groebner basis
                                     b^2 + 2bd + d^2
                                      a+b+c+d
                                  bc^2 + c^2d - bd^2 - d^3
                                    bd^4 + d^5 - b - d
                                  c^3d^2 + c^2d^3 - c - d
                               d^4c^2 + bc - bd + cd - 2d^2
```

**Figure 3.3.** The input and the output of the cyclic 4 problem using our implementation of  $F_4$  Algorithm in Maple

 $-1 + bcd^2 + c^2d^2 - bd^3 + cd^3 - d^4$ 

Symbolic Preprocessing, we are constructing matrices  $F_i$  from the selected pairs and polynomials from Gröbner basis G that are multiplied by a monomial. Polynomials that are constructed from the selected critical pairs are from G and multiplied by a monomial too. But polynomials in G are just the input polynomials or polynomials from  $\tilde{F}_i$  from all previous iterations. All these polynomials that are added into  $F_i$  are simplified by the function Simplify. That means that such a polynomial may be replaced by another polynomial taken from  $\tilde{F}_i$ . To sum up, the matrix  $F_i$  is built from multiples

```
> infolevel[GroebnerBasis] := 5:
> Groebner[Basis]([a*b*c*d - 1, a*b*c + a*b*d + a*c*d + b*c*d, a*b + b*c + a*d + c*d,
                  a + b + c + d, tdeg(a, b, c, d), method=fgb);
-> FGb
domain: rat_int_cof
Set offset primes to 0/20000
Lin Bk ignored [NEW lib]/S:0 -> 8/
 [2](3x7)100%/
 [3] (5x10) 100%/
 [4] (8x12)100%/
 [5] (6x12) 100%/
 [6] (11x16)100%/
 [7](8x11)0.0%/100%/
Mingbasis2
 (7x24)100%/restore Z1 Copy 80.00 for 24/30 exposants
  SWAP Z1/2 Memory usage (estimate): 0.000
   7 polynomials, 30 terms
    total time:
                  0.014 sec
                                      a+b+c+d
                                      b^2 + 2bd + d^2
                                  bc^2 + c^2d - bd^2 - d^3
                           -1 + bcd^2 + c^2d^2 - bd^3 + cd^3 - d^4
                                    bd^4 + d^5 - b - d
                                   c^3d^2 + c^2d^3 - c - d
                                d^4c^2 + bc - bd + cd - 2d^2
```

**Figure 3.4.** The input and the output of the cyclic 4 problem using Faugère's implementation of  $F_4$  Algorithm in Maple

of the input polynomials or polynomials from  $\tilde{F}_{1,\dots,i-1}$  and monomials. So to track the process of building the matrices  $F_i$ , we just have to keep the track of which polynomials from which matrices  $\tilde{F}_i$  (we can look at the input polynomials as if it is a matrix  $\tilde{F}_0$ ) are multiplied with which monomials. In the end, we have to recover how the matrix G was build, but this is the same case as reconstructing a matrix  $F_i$  because the Gröbner basis G consists only of the input polynomials or polynomials from matrices  $\tilde{F}_i$ .

Unnecessary and redundant polynomials can be, of course, removed from the matrices  $F_i$  as we have described in sections 3.1.5 and 3.2.3. Moreover, if nothing is added from the matrix  $\tilde{F}_i$  to the Gröbner basis G, i.e. the matrix  $\tilde{F}_i^+$  is empty, the matrix  $F_i$  can be removed totally. Therefore, there will be no reduction to zero in the generated solver so much computation time can be saved. We are still working with sparse matrices so the matrix partitioning which is described in section 3.2.4 can be used to speed up the Gauss-Jordan eliminations.

By default, this strategy is disabled in the automatic generator and the systematic polynomial generator as described in sections 3.1.4 and 3.2.2 is used. This is done by setting cfg.PolynomialsGenerator = 'systematic'. To enable the polynomial gen-

erator using the  $F_4$  strategy, set cfg.PolynomialsGenerator = 'F4' in the automatic generator settings.

The key function in the  $F_4$  strategy is the Sel function which is descibed in section 2.2.6. While different Sel functions can be efficient for different minimal problems, it can be easily changed which function will be used. To use the normal strategy, set cfg.PolynomialsGeneratorCfg.Sel = @F4\_SelNormal and to select only one critical pair each time, set cfg.PolynomialsGeneratorCfg.Sel = @F4\_SelFirst in the automatic generator settings. This second function emulates the Buchberger Algorithm. Both this functions are stored in the folder generator/F4 and everybody can implement and use his own Sel function.

#### 3.3. Benchmark

In the automatic generator as presented in this thesis, there are many different methods which can be used to generate the solver. Efficiency of the generated solver depends on the choosen method. It shows up, that different methods are efficient for different minimal problems. Therefore, we need a tool which will generate more solvers for a choosen minimal problem. Each solver will be generated with different method used. In the end, the tool will compare the generated solvers so we will be able to choose which solver we will use in the application. The tool have to compare the solver in many aspects because in different applications we have different requirements. For example, in one application, we need solvers which are very fast, but in another, we may prefere solvers that are slower, but more numerically stable. Therefore, in this section, we present such a tool which we call the benchmark of the automatic generator.

#### 3.3.1. Structure

Now, we will briefly describe the structure of the benchmark of the automatic generator. The structure is shown in Figure 3.5 and consists of independent blocks which can be easily modified and replaced.

At the beginning, we have to load the definition of the minimal problem for which we want to generate the solver. These definitions are the same as described in the section 3.1.1.

Next, we load the benchmark templates. A benchmark template is a set of settings of the automatic generator that will be used to generate a solver. For example, if we want to compare one elimination and multiple elimination solvers, we will have two benchmark templates. In the first one, the settings will be set to generate one elimination solver, in the second one, the generation of multiple elimination solver will be set. We store a set of benchmark templates which are related to each other in one MATLAB function. In the example above, we will have both of the templates in one function called  $bench\_elimination$ . Another example can be, that benchmark templates for benchmarking the polynomial generation method, i.e. if the  $F_4$  method is better that the systematic method, will be in one function called  $bench\_polynomialGenerator$ . We keep all these functions in the folder benchmark in the automatic generator.

When the benchmark templates are loaded, we use the automatic generator to generate the solver for each benchmark template. Then, we run each solver on each instance of parameters and we write down the time spent by the computation. These instances of parameters may be specified by user or if they are not provided, they are randomly generated. The random generator of parameters assign to each parameter random

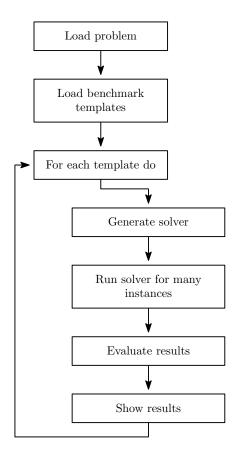


Figure 3.5. Block diagram of the benchmark of the automatic generator

number from the normal distribution with mean set to zero and standard deviation to one.

The main part of the benchmark is the evaluation of the results. It is left to the user to implement the evaluation function. This function gets the set of instanced parameters, results from the solver and the correct results, if they are provided with the set of instanced parameters. These correct results may be precomputed in other software only to check the correctness of the results from the generated solvers. The evaluation function returns the set of errors, e.g. how much the results from the solver differ from the correct values. If the evaluation function is not provided by the user, the default evaluation function is used. This function gets the set of given polynomials, substitutes the unknowns by computed solutions and evaluates each polynomial. In the best case, we should get zero for each polynomial. But we obtain non zero values. Absolute values of these errors depend on the numerical stability of the solver so we are comparing the numerical stability of the generated solvers.

Finally, we have to show the results in some well-arranged way. The presentation of the results is dependent on which data we want to show. Therefore, we leave the implementation of this part to the user. If the user do not specify the function which presents the results, the default presentation function is used. In this function, we get  $\log_{10}$  of absolute values of the errors and show them as histogram. To be able to compute the  $\log_{10}$  of the errors we have to remove all zero values from the set of errors. We also print the percentage of zero errors amongs all errors.

#### 3.3.2. Usage

The benchmark of the automatic generator is designed to be used easily, but also to be easily modified to give us the results we want to know. The benchmark of minimal problem minimalProblem is called by the command gbs\_Benchmark(minimalProblem, benchmark, inputData, correctOutput, validationFunction, renderFunction) where the first two arguments are compulsory and the others are optional. The parameter benchmark is a function handler to the function which provides the set of benchmark templates. The inputData is a set of parameters on which we want to test the generated solvers. If this argument is not provided, random generated parameters are used. The correctOutput is a set of expected results of the solvers. These correct outputs are used to compare the numerical stability of the solvers. The function handler validationFunction is a handler to a function which evalutes the results and returns the set of errors. The default validationFunction substitutes computed solutions into the given polynomials and evalutes them. The function handler renderFunction is a handler to a function which shows the results. If this function is not provided, the default function is used. This default function renders histogram of  $\log_{10}$  of absolute values of errors.

For clarity, we provide an example. We want to benchmark solvers for the 6-point focal length problem [14] and we are considering one elimination and multi elimination solvers. We have no real data to test these solvers and therefore, we have no expected results. We want to use the default validation function and the default function for presenting the results of the benchmark. In this case we call the benchmark with these parameters gbs\_Benchmark('sw6pt', @bench\_elimination). From the shown results, we decide which solver we will use in the application.

# 4. Experiments

To show the usefulness of our improvements of the automatic generator, we have compared several solvers of some minimal problems using different methods in the automatic generator. We have used the benchmark of the automatic generator to generate the solvers and to compare the results because this tool is designed to it perfectly.

We have divided the experiments into three parts. In each part, we are comparing easily comparable methods on which the speed up of the new implemented method can be straightforwardly seen. In the first section, we are comparing one elimination solvers against multiple elimination solvers. In the second part, we are comparing solvers without the matrix partitioning used, with matrix partitioning used only to the last elimation and solvers with matrix partitioning used to all eliminations in the solver. In the last section, we are comparing solvers with different strategies of polynomial generation used. One solver is generated by the systematical generator and the second one is using the  $F_4$  strategy.

To be able to compare the solvers, we have measured the time of computing the solutions from each set of parameters for each solver. In the tables below we have picked up the maximal and minimal value and median of the times for each solver. To be able to compare the numerical stability of the solvers, we have substituted back each computed solution into the given equations and wrote down the results as errors. We have removed the errors equal to zero and compute the  $\log_{10}$  of absolute values of errors. We have presented these values as histogram for each solver.

We have choosen the 9-point relative pose different radial distortion problem [15] for the testing. This problem consists of four polynomial equations in four unknowns. The number of all parameters is 63. The definition of this minimal problem can be found under the name ku9pt in the folder minimalProblems in the automatic generator. To generate the solvers, we have used the default settings of the automatic generator obtained by calling of the function gbs\_InitConfig if not specified differently. We have tested the generated solvers on randomly generated data, but the data remained the same within each experiment. Each solver was tried on 1 000 instances of parameters. All test were performed on Intel Xeon E5-2630 2.30 GHz based computer. The MATLAB R2014a 64-bit was used to the tests.

# 4.1. Multiple elimination solver

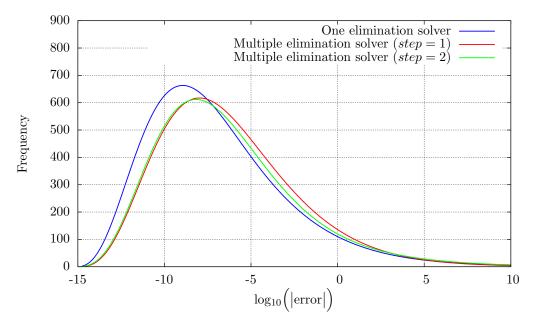
In this part, we are comparing one elimination solver with multiple elimination solvers. We have generated one solver according to the description in the section 3.1.4. This first solver consist only of one elimination in the end. The second and the third solvers have been generated as explained in the section 3.2.2. The second solver has been generated with the variable *step* set to 1 therefore, an elimination is performed always when the maximal total degree of the generated polynomials is increased by 1. This second solver consist of four Gauss-Jordan eliminations. The third solver has been generated with the variable *step* set to 2. This means, that an elimination is performed when the maximal total degree of the generated polynomials is increased by 2. This

#### 4. Experiments

solver consists of two eliminations.

We have used the benchmark templates specified in the function bench\_elimination from the folder benchmark in the automatic generator. All other settings have remained default.

The computing times of these solvers are in the Table 4.1. The numerical stability of the solvers is shown in the Figure 4.1 as histogram of  $\log_{10}$  of absolute values of errors.



**Figure 4.1.** Histogram of  $\log_{10}$  of absolute values of errors for one elimination and multiple elimination solvers.

	One elimination	Multiple elimination	Multiple elimination
	solver	<b>solver</b> $(step = 1)$	$\mathbf{solver}\ (step=2)$
minimal	2.159 s	1.044 s	1.060 s
median	2.930  s	$2.097 \mathrm{\ s}$	$2.015 \mathrm{\ s}$
maximal	3.129  s	$2.301 \; s$	$2.444 \mathrm{\ s}$

**Table 4.1.** Computing times of one and multiple elimination solvers.

You can see that the numerical stability of the multiple elimination solvers is slightly worse than the numerical stability of the one elimination solver. In the contrary, the multi elimination solvers are approximately 1.5 times faster than the one elimination solver. Interesting is that the second and third solvers are equivalently fast, but one of them consists of four eliminations and the second one only of two eliminations. Therefore, we can not say that more eliminations lead to faster solvers. So, it is important to run the benchmark to find the optimal number of eliminations for each minimal problem and then choose the best solver for the application.

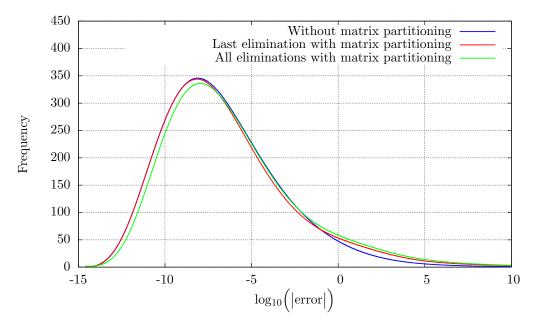
### 4.2. Matrix partitioning

In this section, we compare the solvers using matrix partitioning as described in the section 3.2.4. For comparing, we have used multiple elimination solvers as described in the section 3.2.2. We have set the variable *step* to one so the generated solvers have

four eliminations. The first solver has been generated without matrix partitioning. In the second one, matrix partitioning was used only for the last elimination and the third solver has been generated with matrix partitioning for all four eliminations used.

The benchmark templates used for this comparison are specified in the function bench\_matrixPartitioning in the folder benchmark in the automatic generator. The variable step was set to 1 and all other settings have remained default.

The computing times of these tree solvers are in the Table 4.2 and the numerical stability is shown as histogram of  $\log_{10}$  of absolute values of errors in the Figure 4.2.



**Figure 4.2.** Histogram of  $\log_{10}$  of absolute values of errors for solver without matrix partitioning, solver with matrix partitioning only for the last elimination and solver with matrix partitioning for all eliminations in the solver.

	Without matrix	Matrix partitioning	Matrix paritioning
	partitioning	for the last elimination	for all eliminations
minimal	1.353 s	$0.956 \; \mathrm{s}$	0.794  s
median	2.185  s	$1.071 \; s$	$0.932 \mathrm{\ s}$
maximal	2.604  s	$1.189 \mathrm{\ s}$	$2.273 \mathrm{\ s}$

**Table 4.2.** Computing times of solver without matrix partitioning, of solver with matrix partitioning used for the last elimination and of solver with matrix partitioning used for all eliminations in the solver.

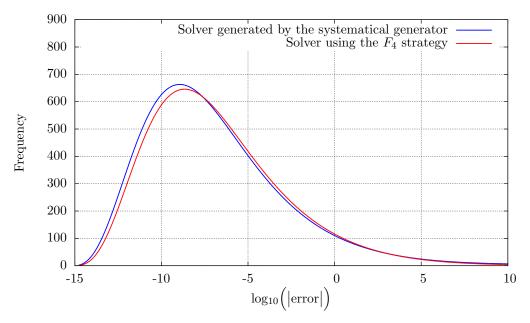
We can see that the numerical stability remained the same for all three solvers. The solver with matrix partitioning for the last elimination used is approximately twice faster than solver without matrix partitioning. There is no big difference between the solver with matrix partitioning for the last elimination used and the solver with matrix partitioning used for all eliminations in the solver. The reason is, that when eliminating not the last matrix in the solver, we have to eliminate all the coupling columns. We do not have to do this when it is the last elimination. Therefore, the speed up of elimination of not the last matrix is not so big as the speed up of the last elimination.

#### 4.3. $F_4$ strategy

At last, we have compared the solver generated by the systematical polynomial generator with solver generated with using the  $F_4$  strategy. The first solver has been generated according to the description in the section 3.1.4 using the systematical polynomial generator. The second one uses the  $F_4$  strategy described in the section 3.2.5.

We have used the benchmark templates that are defined in the function bench\_poly-nomialGenerator which is stored in the folder benchmark in the automatic generator. All other settings have remained default.

The numerical stability of both solvers are shown in the Figure 4.3 as histogram of  $\log_{10}$  of absolute values of errors. The computing times are in the Table 4.3.



**Figure 4.3.** Histogram of  $\log_{10}$  of absolute values of errors for solver generated by the systematical generator and for solver using the  $F_4$  strategy.

	Systematical	$F_4$ strategy
	generator used	$\mathbf{used}$
minimal	2.213 s	$0.615 \; { m s}$
median	2.895  s	$0.713 \mathrm{\ s}$
maximal	$3.320 \mathrm{\ s}$	$1.161 \mathrm{\ s}$

**Table 4.3.** Computing times of solver generated by the systematical generator and of the solver using the  $F_4$  strategy.

From the results, we can see that the numerical stability has remained the same for both solvers. The solver which uses the  $F_4$  strategy is about 4 times faster that the solver generated by the systematical polynomial generator.

#### 5. Conclusion

In this work, we have focused on how to solve systems of polynomial equations fast and how to automatically generate efficient solvers for these systems.

In the first part, we have reviewed the state of the art algorithms for computing Gröbner basis of polynomial systems. We have described the Buchberger Algorithm [2], then we have explained the  $F_4$  Algorithm [6] in details and in the end, we have pointed out the main features of the  $F_5$  Algorithm [7].

In the second part, the automatic generator [12] of minimal problem solvers has been presented. This tool enables us to easily generate solvers of systems of polynomial equations which arises when solving minimal problems in computer vision. We have described the process of generation of solvers in details. Then, we have suggested some improvements of the automatic generator and we have implemented them. For example, we have presented an improvement which allows us to generate multiple elimination solvers which are usually better for systems of polynomial equations in many unknowns. We have also shown that the solvers can be sped up when Gauss-Jordan eliminations for sparse matrices are used. Next, we have taken over a strategy from the  $F_4$  Algorithm [6] and we have implemented it into the automatic generator. For better understanding, we have implemented the  $F_4$  Algorithm [6] in Maple first. The description of this implementation is provided in this section, too. In the end, we have presented the benchmark of the automatic generator. This tool helps us to decide which generated solver is better for our application.

In the end, we have taken the 9-point relative pose different radial distortion problem [15] and compared solvers generated with different methods for this problem on set of randomly generated data. We have shown that solvers generated with the new implemented methods may be faster than solvers generated by the old implementation of the automatic generator. We have register the most visible speed up when the  $F_4$  strategy is used. In this case, the solver using the  $F_4$  strategy is four times faster that the solver generated by the systematical generator for the 9-point relative pose problem [15].

# A. Contents of the enclosed CD

/				
Aut	omaticGenerator	. folder with the automatic generator		
, <u> </u> 1	benchmark	.folder contaning benchmark tem-		
		plates		
.	bench_elimination.m	-		
	bench_matrixPartitioning.m	that are used in the experiment 4.1		
'	bench_matrixrartitioning.m	that are used in the experiment 4.2		
	bench_polynomialGenerator.m	-		
	- 1 7	that are used in the experiment 4.3		
	gbsMacaulay			
		for calling Macaulay2 [9]		
.	<del>-</del> :			
_ [	generator	. folder with main functions of the au-		
		tomatic generator		
.	<u> </u>			
r	minimalProblems	. folder with minimal problem defini-		
		tions		
.	ku9pt.m	definition of the 9-point relative		
		pose different radial distortion prob-		
	sw5pt.m	lem [15]		
'	Swopu.m	pose problem [?]		
.	sw6pt.m			
	-	pose problem [14]		
.	test.m	•		
		tions in two unknowns)		
	prerequisites			
		Macalulay2 [9] and PaToH [3]		
	<del>'-</del> :			
<del>   </del>	solvers	_		
	installation.txt	stored installation how to		
	license.txt			
	setpahs.m			
		to the MATLAB environment		
F4	E4 m	implementation of the E. Almerithm		
<del> </del>	F4.mw	Implementation of the $F_4$ Algorithm [6] in Maple		
thesis [0] III Maple				
	thesis.pdf	digital copy of this thesis		

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