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# COMPUTING ROOTS OF POLYNOMIALS OVER NUMBER FIELDS USING COMPLEX EMBEDDINGS

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ABSTRACT. We explore a fairly generic method to compute roots of polynomials over number fields through complex embeddings. Our main contribution is to show how to use a structure of a relative extension to decode in a subfield. Additionally we describe several heuristic options to improve practical efficiency. We provide experimental data from our implementation and compare our methods to the state of the art algorithm implemented in PARI/GP.

## 1. INTRODUCTION

In the past few years, the rise of cryptographic protocols using the structure of number fields has driven attention to computational number theory. Indeed, in order to obtain efficient implementations of such cryptosystems as well as testing their security, fast computations over number fields are required. In particular, number field units and  $S$ -units are the main objects used to retrieve short generators of principal ideals [16, 2, 25] and short elements of general ideal lattices [15, 32, 6, 7] respectively. Their computation commonly requires a saturation process [35, 17, 10], which includes a necessary step of  $n$ -th roots computations. This task is a special case of solving a polynomial equation with coefficients in a number field.

We are therefore interested in the following problem. Given  $K$  a number field and  $f(T) \in K[T]$ , find the roots of  $f(T)$  in  $K$  i.e. find  $Z_K(f) = \{x \in K \mid f(x) = 0\}$ . As a matter of fact we will only consider the cases such that all the roots can be expressed as integral combinations of a known basis of  $K$  i.e.

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such that:  $\exists (b_i)_i \in K^n, \forall x \in Z_K(f), \exists (x_i)_i \in \mathbb{Z}^n \mid x = x_1 b_1 + \dots + x_n b_n$ . We believe that we are not too restrictive as this is often the case in standard computations, such as  $S$ -units computations or the Number Field Sieve [22, 41].

The state of the art algorithm to solve this problem – that we will call the *algebraic* or  *$p$ -adic* method – is implemented in softwares such as MAGMA or PARI/GP [11, 31]. It follows the ideas used to factorise polynomials over number fields developed for example in [3, 17, 36, 37] and uses finite places. More precisely, the procedure is as follows:

- (1) pick a prime ideal  $\mathfrak{p}$  which defines an isomorphism between  $\mathcal{O}_K/\mathfrak{p}$  and  $\mathbb{F}_{p^s}$  for some prime number  $p \in \mathbb{Z}$ ;
- (2) compute the roots modulo  $\mathfrak{p}$  in  $\mathbb{F}_{p^s}[X]$ ;
- (3) lift these to elements to  $K$  modulo a power  $\mathfrak{p}^k$ ;
- (4) if  $\mathfrak{p}^k$  is large enough compared to the size of the solutions and it is given by a LLL-reduced basis then we can recover a solution  $x$  given  $y \equiv x \pmod{\mathfrak{p}^k}$ .

In practice, the method requires three main computational operations:

- (1) compute first an “approximation” of the solutions;
- (2) compute a reduced basis of a lattice;
- (3) retrieve the solutions from the approximation using the lattice.

The third step – called the *decoding* phase – can be seen as solving instances of the Bounded Distance Decoding (BDD) problem.

However, it is possible to use complex embeddings instead of prime ideals. Indeed, it is known at least since the seminal paper introducing the LLL algorithm [23] how to use lattices in order to find short integral relations between algebraic numbers or find the irreducible polynomial of a number field element. As an example, assume one knows approximations of real numbers  $\alpha_1, \dots, \alpha_n$ . Now define the embedding

$$\begin{aligned} \mathbb{Z}^n &\longrightarrow \mathbb{R}^{n+1} \\ (\lambda_i)_{i \in \llbracket 1, n \rrbracket} &\longmapsto \left( C \sum_{i=1}^n \lambda_i \alpha_i, \lambda_1, \dots, \lambda_n \right) \end{aligned}$$

which gives a lattice of  $\mathbb{R}^{n+1}$  represented by the row matrix

$$\begin{bmatrix} C\alpha_1 & 1 & 0 & \dots & 0 \\ C\alpha_2 & 0 & 1 & \ddots & \vdots \\ \vdots & \vdots & \ddots & \ddots & 0 \\ C\alpha_n & 0 & \dots & 0 & 1 \end{bmatrix}.$$

Here  $C$  is a large coefficient used to ensure the shortness of the solution. Reducing this basis would allow us to retrieve a short vector  $(\lambda_i)_{i \in \llbracket 1, n \rrbracket}$  such that  $C \sum_{i=1}^n \lambda_i \alpha_i$  is small. Note that this strategy can be seen as solving an instance of the Shortest Vector Problem (SVP). Similarly it is common to use LLL algorithm to solve knapsack problems. Assume we are given  $(\alpha_1, \dots, \alpha_n) \in \mathbb{Z}^n$  and  $S \in \mathbb{Z}$ , and that we want to find a short vector  $(\lambda_1, \dots, \lambda_n)$  such that  $\sum_{i=1}^n \lambda_i \alpha_i = S$ . Then one can consider the similar

matrix

$$\begin{bmatrix} -C\alpha_1 & 1 & 0 & \dots & 0 \\ -C\alpha_2 & 0 & 1 & \ddots & \vdots \\ \vdots & \vdots & \ddots & \ddots & 0 \\ -C\alpha_n & 0 & \dots & 0 & 1 \\ CS & 0 & \dots & \dots & 0 \end{bmatrix}.$$

Again a LLL reduction would yield a small solution of the linear equation.

**Our work.** One can combine both approaches to retrieve the coefficient of  $x \in \mathbb{Z}[\mathcal{B}]$  from an embedding  $\sigma_i(x)$ , where  $\mathcal{B} = (b_1, \dots, b_n)$  is a  $\mathbb{Q}$ -basis of  $K$ . This is done through solving an instance of BDD. For all  $j \in \llbracket 1, n \rrbracket$ , denote by  $\beta_j$  an approximation of  $\sigma_i(b_j)$  and  $S$  an approximation of  $\sigma_i(x)$ . Now consider  $(x_i)_{i \in \llbracket 1, n \rrbracket} \in \mathbb{Z}^n$  the coefficients of  $x$  in the basis  $\mathcal{B}$ . Then one can expect  $S - \sum_{i=1}^n x_i \beta_i$  to be close to 0. Thus solving a BDD instance with input the row matrix

$$\begin{bmatrix} -\beta_1 & 1 & 0 & \dots & 0 \\ -\beta_2 & 0 & 1 & \ddots & \vdots \\ \vdots & \vdots & \ddots & \ddots & 0 \\ -\beta_n & 0 & \dots & 0 & 1 \end{bmatrix}$$

and target vector  $[S, 0, \dots, 0]$  is expected to yield the solution  $[\epsilon, x_1, \dots, x_n]$  if approximations are computed with high enough precision.

In this article, we study this last method and apply it to the problem of retrieving polynomial roots in number fields. In particular, under the conjecture that the lattices we consider follow the Gaussian heuristic, we establish a *heuristic* precision for which one is ensured to retrieve an element, see [Theorem 3.17](#). We note that a similar study has been done in [\[33\]](#) when considering the version solving a SVP instead of a BDD. Following their analysis, we provide in [Theorem 3.11](#) a *proven* precision ensuring correctness of the decoding, yet larger than the heuristic one.

Additionally, one of our main contribution is to show how to take advantage of a number field extension  $L/K$  to replace the decoding of an element in  $L$  by the decoding of  $[L : K]$  elements in  $K$ . We will call it the *relative method*. When applied to the problem of solving polynomial equations, applying this method comes with the cost of searching through a set of size  $d^{[L:K]}$ , where  $d$  is the degree of the equation considered.

In order to improve the *practical* efficiency of our algorithms, we develop several heuristic strategies and observations allowing for substantial speed-ups.

- We describe a modification of the LLL algorithm – called **SpecLLL** – which improves its running time when  $\mathcal{B}$  is of the form  $(1, \alpha, \dots, \alpha^{n-1})$ : we pre-reduce each new basis vector of index  $i + 1$  using the relation used to size-reduce the previous vector of index  $i$  with respect to the lattice generated by the  $(i - 1)$ -th first rows, see [Algorithm 9](#). Note that this modification was inspired by [\[34\]](#), which is a version of LLL using the specific structure of ideal lattices. Asymptotically, we gain between 20% and 50% in the execution time, depending on which implementation of LLL is tested, as shown in [Figure 2](#) and [3](#).
- We provide a tighter evaluation of the precision required to retrieve the coefficients of a targetted element – see [Section 5.2](#).

- We use a early abort strategy in the decoding phase (with Babai’s nearest plane algorithm [1]) to avoid unnecessary computations, see Algorithm 10. This is particularly impactful when using our relative method.

Using our implementation in GP [31]<sup>1</sup>, we study the generic behaviour of our algorithms, how our improvements impact their running time, and compare them to the function `nroots` implemented in PARI/GP [31]. In most of our experiments we consider number fields defined by polynomials whose coefficients are randomly drawn in a given segment. Respectively to this definition of “random number fields”, we are then able to study the *average behaviour* of the aforementioned algorithms.

Concerning the absolute method we can observe the following facts.

- We first remark that our absolute method is *more stable* than the  $p$ -adic algorithm implemented in PARI/GP. Indeed, there can be as much as 10% of fields over which the execution time of `nroots` explodes. From experiments, these are fields for which `nroots` do not find a suitable inert prime. Consequently, our algorithm in its certified version has similar running time *in average*, depending on the parameters, despite being slower than `nroots` at its best.
- It is more influenced than PARI/GP `nroots` function by several parameters : the dimension  $[K : \mathbb{Q}]$  and the size of the roots. This is mostly linked to the fact that the lattices we use to decode take longer to be reduced than powers of prime ideals as in `nroots`. Our heuristic improvements have significant impact on the time efficiency of our algorithm (10 times faster than the certified version in dimension 100 for example, see Figure 5).

Then when we consider relative extensions  $L/K$  and our relative method we have the following main observations:

- the heuristic observations allow us to obtain a more efficient algorithm as well in the relative case: between 10 and 100 times faster for relatively small parameters chosen for our experiments, see Figure 7;
- when the relative degree  $[L : K]$  and the degree of the equation are relatively small, our relative algorithm can offer significant speed-ups compared to the absolute one: again for small parameters, our relative algorithm can already be up to 10 times faster than the absolute one, see Figure 8 and Figure 9 for example.

Finally we study specific families of number fields, namely cyclotomic fields and real Kummer extensions of the form  $\mathbb{Q}(\sqrt[m_1]{m_1}, \dots, \sqrt[m_r]{m_r})$  with  $(m_1, \dots, m_r) \in \mathbb{N}^r$  and  $p$  a prime integer. We can observe the following facts:

- over cyclotomic fields, our relative method applied to the extension  $K_m/K_m^+$  – it has relative degree  $[K_m : K_m^+] = 2$  – does not outperforms PARI/GP `nroots` function, which highlights the fact that cyclotomic fields are “good” fields for this method;
- over real Kummer extensions, which are “bad” fields for `nroots`, the situation is completely different, and our relative algorithm can be up to 500 times faster than PARI/GP `nroots` function, see Section 6.3.4.

**Future work.** First we plan to study the possibility of using the structure of relative extensions within the  $p$ -adic method. If we were to be successful, it might be adapted to general global fields as well, which are a natural generalisation of number fields [5].

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<sup>1</sup>Code publicly available at [https://github.com/AndLesav/nf\\_polynomial\\_roots](https://github.com/AndLesav/nf_polynomial_roots)

Since the main bottleneck of our method is the reduction of lattices to decode, we plan to improve SpecLLL. One direction would be to follow the strategy described in [21] for knapsack-like lattices which could lead to smaller running times. Then when there is no subfield, it is possible to adapt the strategy by using several embeddings. Intuitively, adding relations should allow us to decode more easily.

**Outline.** The rest of this article is as follows. We give necessary recalls on number fields and Euclidean lattices in Section 2. In Sections 3 and 4 we describe the generic methods considered. Then Section 5 is used to detail the improvements that we developed to speed-up the running time of our algorithms. Finally, we provide experimental results obtained from our implementation in Section 6. We also give additional data in Appendix B.

## 2. RECALLS ON NUMBER FIELDS AND LATTICES

In this section, we recall some notions and facts that we will use in the rest of the article. First we will quickly mention number fields, then describe Euclidean lattices and reduction algorithms.

**Notations 2.1.** The inner product is denoted by  $(\cdot | \cdot)$ . Given  $x \in \mathbb{R}$ , we denote by  $\{x\}$  its fractional part  $x - \lfloor x \rfloor$ .

Algebraic closures of number field extensions considered will be designated by  $\Omega$ . Given a ring morphism  $\sigma : A \rightarrow B$  and  $a \in A$ ,  $a^\sigma$  can be used to design  $\sigma(a)$ . This extends to any object  $f$  which can be identified to a vector with coefficients in  $A$ , such as polynomials in  $A[X]$ . Additionally, if  $A \subset B$  and  $f(X) \in A[X]$  we write  $Z_B(f)$  the set of roots of  $f(X)$  in  $B$ .

Given two ordered sets  $A$  and  $B$  we will denote by  $A \otimes B$  the tensor product of  $A$  and  $B$  (when it makes sense), i.e.  $A \otimes B = \{ab, b \in B \mid a \in A\}$ . If  $A$  and  $B$  are not ordered  $A \otimes B$  will be the collection of all the products  $ab$  such that  $a \in A$  and  $b \in B$ . Finally, we also use this notation for the tensor product of vectors and matrices.

Given a matrix  $M$ , we will denote by  $M[i]$  its  $i$ -th row and  $M^T$  its transpose. Given a vector  $v$  we will write  $s(v)$  for  $\log_2 \|v\|_\infty$ . This notation extend to any object that can be expressed as a vector, such as polynomials.

**2.1. Number fields.** We refer the reader to [4, 12, 13, 28, 38] for anything related to number fields and computational number theory.

**Definition 2.2.** A *number field*  $K$  is a field which is a finite extension of  $\mathbb{Q}$ , i.e. a finite dimensional  $\mathbb{Q}$ -vector space.

**Notation 2.3.** Given an extension  $L/K$  of number fields, we will call the dimension of  $L$  over  $K$  the *degree* of  $L/K$ . It will denoted by  $[L : K]$ .

**Proposition 2.4.** *Let  $L/K$  be an extension of number fields. Then there is an irreducible polynomial  $P(X) \in K[X]$  such that*

$$L \cong \frac{K[X]}{(P(X))}.$$

and  $[L : K] = \deg P(X)$ . Moreover  $P(X)$  has  $[L : K]$  distinct roots in an algebraic closure  $\Omega$  of  $K$  containing  $L$ . These roots define  $[L : K]$  distinct  $K$  – isomorphisms of  $L$  into  $\Omega$ . If  $\alpha$  is such a root,

then the corresponding isomorphism  $\sigma_\alpha$  is the following,

$$\begin{aligned} \sigma_\alpha : \frac{K[X]}{(P(X))} &\longrightarrow K[\alpha] \subset \Omega \\ \sum_{i=0}^{[L:K]-1} c_i X^i &\longmapsto \sum_{i=0}^{[L:K]-1} c_i \alpha^i. \end{aligned}$$

**Notation 2.5.** We call any such polynomial a *defining polynomial of  $L$  over  $K$* , and use denote it by  $P_K(X)$ .

An element  $\sigma_\alpha \in \text{Hom}(K, \mathbb{C})$  is a field embedding of  $K$  into  $\mathbb{C}$  corresponding a root  $\alpha$  of  $P_K(X)$  in  $\mathbb{C}$ . There are  $r_1$  real embeddings and  $r_2$  pairs of (strictly) complex embeddings. The two elements of a given pair are conjugates one from each other. It is usual to write  $\sigma_1, \dots, \sigma_{r_1}$  the real embeddings and to consider that  $\sigma_{j+r_2} = \overline{\sigma_j}$  for all  $j \in \llbracket r_1 + 1, r_1 + r_2 \rrbracket$ .

**Definition 2.6** (Minkowski embedding). Given a number field  $K$  defined by a degree  $n$  irreducible polynomial  $P(X)$ , the *canonical embedding* or *Minkowski embedding* is defined as

$$(2.1) \quad \begin{aligned} \sigma_K : K &\longrightarrow \mathbb{R}^{r_1} \times \mathbb{C}^{r_2} \cong \mathbb{R}^n \\ x &\longmapsto (\sigma_i(x))_{i \in \llbracket 1, r_1 + r_2 \rrbracket}. \end{aligned}$$

Then  $K$  can be seen as embedded in  $\mathbb{R}^n$ . We will relax this classical definition by considering any algebraic closure  $\Omega$  instead of  $\mathbb{C}$ . This way  $\sigma_K$  defines an embedding of  $K$  into  $\Omega$ . Then one can define a relative version of the Minkowski embedding: for a field extension  $L/K$  we define

$$\sigma_{L/K} : x \longmapsto (\sigma(x))_{\sigma \in \text{Hom}_K(L, \Omega)}.$$

We call this map the *Minkowski embedding relative to  $L/K$* .

**2.2. Euclidean lattices.** We refer the reader interested in more in-depth presentations on Euclidean lattices to [27, 14].

*Gram-Schmidt orthogonalisation.* An important and classical computation in linear algebra is the Gram-Schmidt orthogonalisation (GSO). It allows transforming a free family of a vector space (with a scalar product) into in an orthogonal family. The naive algorithm can be found in Algorithm 1.

---

**Algorithm 1** GSO

---

**Require:** A free family  $\mathcal{B} = \{b_1, \dots, b_r\}$  of  $\mathbb{R}^n$

**Ensure:** A family  $\tilde{\mathcal{B}} = \{\tilde{b}_1, \dots, \tilde{b}_n\}$  with  $\tilde{b}_1 = b_1$  and  $(\tilde{b}_i \mid \tilde{b}_j) = 0$  for  $i \neq j$

- 1:  $\tilde{\mathcal{B}} \leftarrow \mathcal{B}$
  - 2: **for**  $i = 1$  to  $r$  **do**
  - 3:    $\tilde{b}_i \leftarrow b_i$
  - 4:   **for**  $j = 1$  to  $i - 1$  **do**
  - 5:      $\tilde{b}_i \leftarrow \tilde{b}_i - \frac{(\tilde{b}_i \mid \tilde{b}_j)}{(\tilde{b}_j \mid \tilde{b}_j)} \tilde{b}_j$
  - 6:   **end for**
  - 7: **end for**
  - 8: **return**  $\tilde{\mathcal{B}}$
- 

**Definition 2.7** (GSO). Given a basis  $\mathcal{B}$ , we will call the basis  $\tilde{\mathcal{B}}$  outputted by Algorithm 1 *the GSO* of  $\mathcal{B}$ .

For the GSO computation, one can follow a QR decomposition as described in [29] for example which gives an algebraic complexity of  $O(n^3)$ .

**Definition and first properties.**

**Definition 2.8.** A *Euclidean lattice* is a discrete subgroup of  $\mathbb{R}^n$  where  $n$  is a positive integer. We say a lattice is an *integral* lattice when it is a subgroup of  $\mathbb{Z}^n$  or *rational* when it is in  $\mathbb{Q}^n$ . A *basis* of a lattice  $\mathcal{L}$  is a basis of  $\mathcal{L}$  as a  $\mathbb{Z}$ -module. The cardinal of said basis is called the *rank* of the lattice.

**Notation 2.9.** Given a matrix  $B$  we will write  $\mathcal{L}(B)$  the lattice generated by its row vectors.

**Definition 2.10.** Consider a Euclidean lattice  $\mathcal{L}$  with basis matrix  $B$ . Then the *determinant* of  $\mathcal{L}$ , denoted by  $\det(\mathcal{L})$ , is the value  $\sqrt{\det(BB^\top)}$ .

**Notation 2.11.** The determinant of a lattice  $\mathcal{L}$  is also called its *volume* because it is the volume of the fundamental domain defined by the vectors of one of its bases. Thus it is also written  $\text{vol}(\mathcal{L})$ .

One of the most important problem over lattices is to compute short vectors. In particular, an important quantity is the length of one of the shortest non-zero vector. It is usually denoted by  $\lambda_1(\mathcal{L})$ . We can give a minimal bound of  $\lambda_1$  using the Gram-Schmidt orthogonalisation of a basis we have at hand.

**Theorem 2.12.** Let  $B = (b_i)_{i \in \llbracket 1, r \rrbracket}$  be a basis of a lattice  $\mathcal{L}$ . Then  $\lambda_1(\mathcal{L}) \geq \min\{\|\tilde{b}_i\| \mid i \in \llbracket 1, r \rrbracket\}$ .

Then one can analyse further the geometry to obtain approximate values on  $\lambda_1$ .

**Heuristic 1** (Gaussian heuristic). Given a lattice  $\mathcal{L}$  of rank  $r$ , an approximate value for  $\lambda_1(\mathcal{L})$  is

$$(2.2) \quad \lambda_1(\mathcal{L})_{\text{gauss}} = \sqrt{\frac{r}{2\pi e}} \times \sqrt[r]{\text{vol}(\mathcal{L})}.$$

In our context, the problem that we will be interested in is the Bounded Distance Decoding with its approximate version.

**Definition 2.13** (**BDD**: Bounded Distance Decoding). Given a basis  $B$  of a lattice  $\mathcal{L}$  and a point  $x$  such that  $d(x, \mathcal{L}) < \lambda_1(B)/2$ , find the lattice vector  $v \in \mathcal{L}$  closest to  $x$ .

**Definition 2.14** (**BDD $_\gamma$** :  $\gamma$ -Approximate Bounded Distance Decoding). Given a basis  $B$  of a lattice  $\mathcal{L}$ , a point  $x$  and a approximation factor  $\gamma$  ensuring  $d(x, \mathcal{L}) < \gamma\lambda_1(B)$  find the lattice vector  $v \in \mathcal{L}$  closest to  $x$ .

In practice, one considers **BDD $_\gamma$**  for  $\gamma < \frac{1}{2}$ . This ensures that there is only one lattice vector  $v$  satisfying  $d(x, v) < \gamma\lambda_1(B)$ .

**Algorithms for lattices.** Intuitively if a basis  $\mathcal{B}$  is composed by vectors globally more orthogonal to each other than the vectors of another basis matrix  $\mathcal{B}'$  are, then the vectors of  $\mathcal{B}$  will be globally shorter than the vectors of  $\mathcal{B}'$ . Therefore, given a basis it is natural to orthogonalise it as much as possible to solve problems on lattices. Since it likely results in a basis with shorter vectors, we call such a process a *reduction process* or *reduction algorithm*.

First let us define a weak notion of basis reduction due to Hermite [19].



**Definition 2.15** (Size-reduce). A basis  $\mathcal{B} = (b_1, \dots, b_r)$  of a lattice is said to be *size-reduced* if its GSO satisfies the following condition:  $\forall i \in \llbracket 1, r \rrbracket, \forall 1 \leq j < i, |\mu_{i,j}| \leq \frac{1}{2}$ .

One can find the simple algorithm computing a size-reduced basis in Algorithm 2, which is essentially an approximation of the GSO algorithm (Alg. 1).

---

**Algorithm 2** SizeReduce

---

**Require:** A free family  $\mathcal{B} = \{b_1, \dots, b_k\}$  of  $\mathcal{L}$  such that  $|\mu_{i,j}| \leq 1/2$  for all  $i \neq j$ , an element  $b \notin \mathcal{L}$ .

**Ensure:** An element  $b_{k+1}$  such that  $b_{k+1} \equiv b \pmod{\mathcal{L}}$  and  $|\mu'_{k+1,i}| \leq 1/2$  for  $i \neq k+1$

- 1:  $b_{k+1} \leftarrow b$
  - 2:  $\tilde{\mathcal{B}}, G \leftarrow \text{GSO}(\mathcal{B})$
  - 3: **for**  $i = k$  to 1 **do**
  - 4:    $b_{k+1} \leftarrow b_{k+1} - \left\lfloor \frac{(b_{k+1} | \tilde{b}_i)}{\|\tilde{b}_i\|^2} \right\rfloor b_j$
  - 5:   Update  $G$
  - 6: **end for**
  - 7: **return**  $\mathcal{B} \cup \{b_{k+1}\}, G$
- 

A fairly natural way of reducing a basis would be to follow Algorithm 1, and replace all coefficients  $\mu_{i,j}$  by their closest integers. This way one could obtain a basis close to the GSO of the starting matrix, i.e. which is size-reduced. In order to obtain a polynomial time algorithm outputting a basis which is proven to be reduced for varying dimension, one has to introduce a new reduction condition. This is the result of the ground breaking work by A. K. Lenstra, H. W. Lenstra and L. Lovász in [23]. They define a new notion of *reduced basis*, that is then called *LLL-reduced*.

**Definition 2.16.** Consider a lattice  $\mathcal{L}$  defined by a basis  $\mathcal{B} = (b_1, \dots, b_r)$  and  $\delta \in ]\frac{1}{4}, 1[$ . Then  $\mathcal{B}$  is called LLL-reduced with parameter  $\delta$  (or  $\delta$ -LLL reduced) if it satisfies the following conditions:

- (1) It is *size-reduced*:  $\forall i \in \llbracket 1, r \rrbracket, \forall 1 \leq j < i, |\mu_{i,j}| \leq \frac{1}{2}$ ;
- (2) It satisfies the *Lovász conditions*:  $\forall i \in \llbracket 2, r \rrbracket, \delta \|\tilde{b}_{i-1}\| \leq \|\tilde{b}_i + \mu_{i,i-1} \tilde{b}_{i-1}\| = \|\tilde{b}_i\|^2 + \mu_{i,i-1}^2 \|\tilde{b}_{i-1}\|^2$ .

The LLL algorithm shown in Algorithm 3 essentially consists of applying **SizeReduce** to new vector basis incrementally, verify if Lovász condition is true and continue if so. Otherwise we swap the two last vectors and reduce again.

**Theorem 2.17.** Consider  $\mathcal{L}$  a lattice of rank  $r$ ,  $\mathcal{B} = (b_1, \dots, b_r)$  a  $\delta$ -LLL reduced basis of  $\mathcal{L}$  for  $\delta = 3/4$ , and  $\tilde{\mathcal{B}}$  the GSO of  $\mathcal{B}$ . Then the following properties are true.

- (1)  $\forall i \in \llbracket 1, r \rrbracket, \forall 1 \leq j \leq i, \|b_j\| \leq 2^{(i-1)/2} \|\tilde{b}_i\|$
- (2) For any  $x \in \mathcal{L} \setminus \{0\}$ ,  $\|b_1\| \leq 2^{(r-1)/2} \|x\|$ .

Theorem 2.17 shows that a LLL reduced basis has good properties. It provides upper bounds related to the norms of the basis vectors. In particular, one can remark that the norm of shortest basis vectors cannot be too large compared to  $\lambda_1(\mathcal{L})$ . Indeed, the second point shows that the LLL algorithm solves in deterministic polynomial time  $\text{SVP}_\gamma$ , for  $\gamma = 2^{(r-1)/2}$ . Many improvements and versions were developed since the original version of LLL. Among many others, one can consider the use of floating-point arithmetic [29].

---

**Algorithm 3** LLL

---

**Require:**  $\mathcal{B}$ , a basis of  $\mathcal{L}$  of rank  $r$ , and a constant  $\delta \in [\frac{1}{4}, 1[$ .**Ensure:**  $\mathcal{B}'$ , a  $\delta$ -LLL reduced basis of  $\mathcal{L}$ .

```

1:  $i \leftarrow 2$ 
2:  $\mathcal{B}' \leftarrow \{b_1\}$ 
3: while  $i \leq r$  do
4:    $\mathcal{B}' \leftarrow (b'_1, \dots, b'_{i-1})$ 
5:    $\mathcal{B}', G \leftarrow \text{SizeReduce}(\mathcal{B}', b_i)$ 
6:   if Lovász condition is satisfied for  $\delta, i$  then  $i \leftarrow i + 1$ 
7:   else  $\text{Swap}(b'_i, b'_{i-1})$  and  $i \leftarrow \max\{2, i - 1\}$ 
8:   end if
9: end while
10: return  $\mathcal{B}$ 

```

---

**Theorem 2.18** (Complexity of LLL [30]). *Consider  $\mathcal{L} = \mathcal{L}(\mathcal{B}) \subseteq \mathbb{R}^n$  a lattice of rank  $r$ , and  $M = \max\{\log_2 \|b\|_2 \mid b \in \mathcal{B}\}$ . Then for input  $\mathcal{B}$ , one can compute a LLL-reduced basis of  $\mathcal{L}$  in time complexity*

$$(2.3) \quad O(r^2 n(r + M)MM(r)).$$

**Remark 2.19.** When considering a full-rank lattice, i.e.  $r = n$ , we get an algebraic complexity  $O(n^4 M + n^3 M^2)$ .

Babai's nearest plane algorithm. In order to solve BDD instances, one can use an algorithm due to [1], and called the *nearest plane algorithm*. It is an inductive technique, described in Algorithm 4, which is essentially a process of size-reduction (Alg. 2)

---

**Algorithm 4** Babai's Nearest Plane Algorithm – NearestPlane

---

**Require:**  $t \in \mathbb{R}^n$ ,  $\mathcal{B} = (b_1, \dots, b_r)$  a basis of a lattice  $\mathcal{L}$ ,  $\tilde{\mathcal{B}}$  the GSO of  $\mathcal{B}$ **Ensure:**  $v \in \mathcal{L}$  a close vector of  $t$ 

```

1:  $v \leftarrow t$ 
2: for  $i = r$  down to 1 do
3:    $v \leftarrow v - \left\lfloor \frac{(v|\tilde{b}_i)}{\|\tilde{b}_i\|^2} \right\rfloor \tilde{b}_i$  ▷ Make  $t$  more orthogonal to  $b_i$ 
4: end for
5: return  $v - t$ 

```

---

Again the output depends on the quality of the basis given as input. We can prove that the output is not too far from the target given the basis is LLL-reduced.

**Proposition 2.20.** *Consider a rank  $r$  lattice  $\mathcal{L}$ , given by a basis  $\mathcal{B}$ . Then for input  $t$  and  $\mathcal{B}$ , Algorithm 4 outputs  $v \in \mathcal{L}$  lying in  $t + \mathcal{F}(\tilde{\mathcal{B}})$ . Thus it solves  $BDD_\gamma$  for  $\gamma \leq \frac{1}{2} \min\{\|\tilde{b}_i\| \mid i \in [1, r]\}$ .*

Following [30], the algebraic complexity of NearestPlane is at most  $O(n^3 + n^2 M)$ .

## 3. ABSOLUTE METHOD

In this section, we study the methods to recover number field elements from complex embeddings, and how to apply them for solving polynomial equations.

**Definition 3.1.** If  $x$  is in  $\mathbb{R}$ , the *approximation of  $x$  up to  $l$ -bits* for  $l \in \mathbb{N}$  is the integer  $\lfloor 2^l x \rfloor$ . If  $x \in \mathbb{C}$  then its *approximation up to  $l$ -bits* is  $\lfloor 2^l \Re(x) \rfloor + i \lfloor 2^l \Im(x) \rfloor$ . In both cases it is denoted by  $\lfloor x \rfloor_l$ .

**Remark 3.2.** We commonly identify a complex number  $x$  with the pair of real numbers  $(\Re(x), \Im(x))$ . We extend this identification to approximations.

**3.1. Recovering elements from complex embeddings.** First we describe how we compute the coefficients of an element given one of its complex embeddings.

**Definition 3.3.** Consider a number field  $K$ ,  $\mathcal{B}$  a  $\mathbb{Q}$ -basis of  $K$ ,  $\sigma \in \text{Hom}(K, \mathbb{C})$ , and  $l \in \mathbb{N}$ . We call a *basis lattice of  $K$  up to precision  $l$  relative to  $\mathcal{B}$  and  $\sigma$*  and denote by  $\mathcal{L}(\mathcal{B}, \sigma, l)$  the lattice generated by the matrix  $B(\mathcal{B}, \sigma, l)$  defined as follows:

$$(3.1) \quad B(\mathcal{B}, \sigma, l) = \begin{bmatrix} -\lfloor \sigma(b_1) \rfloor_l & C & 0 & \dots & 0 \\ -\lfloor \sigma(b_2) \rfloor_l & 0 & C & \ddots & \vdots \\ \vdots & \vdots & \ddots & \ddots & 0 \\ -\lfloor \sigma(b_n) \rfloor_l & 0 & \dots & 0 & C \end{bmatrix}$$

The matrix  $B(\mathcal{B}, \sigma, l)$  is called the *basis matrix* of  $\mathcal{L}(\mathcal{B}, \sigma, l)$ .

**Remark 3.4.** • When there is no ambiguity regarding  $K$ ,  $\mathcal{B}$  or  $\sigma$ , we will simply denote  $\mathcal{L}(\mathcal{B}, \sigma, l)$  by  $\mathcal{L}_l$  and call it the *lattice basis of  $K$  up to precision  $l$* . Similarly the matrix  $B(\mathcal{B}, \sigma, l)$  will be written  $B_l$ .

- The constant  $C$  is used to ensure the validity of the decoding, and accelerates the reduction algorithm on  $B_l$ . It will be determined later.
- If the embedding  $\sigma$  is not a real embedding, then the first column of the matrix in Equation 3.1 is in fact two columns, containing respectively the real and imaginary parts of the corresponding complex numbers.

**Notation 3.5.** If  $\sigma(K) \subset \mathbb{R}$ , we denote by  $\sigma(\mathcal{B})_l$  the first column vector of a basis matrix  $B(\mathcal{B}, \sigma, l)$ . If  $\sigma(K) \not\subset \mathbb{R}$ , we write  $\sigma(\mathcal{B})_l = [\Re(\sigma(\mathcal{B})_l), \Im(\sigma(\mathcal{B})_l)]$  for the matrix composed of the two first column vectors of  $B(\mathcal{B}, \sigma, l)$ .

As before, consider  $K$  a number field,  $\mathcal{B} = (b_1, \dots, b_n)$  a  $\mathbb{Q}$ -basis of  $K$ , and  $x \in K$  such that there is  $(x_i)_{i \in [1, n]} \in \mathbb{Z}^n$  with  $x = x_1 b_1 + \dots + x_n b_n$ . Now assume that  $\lfloor \sigma(x) \rfloor_l$  is known for some  $\sigma \in \text{Hom}(K, \mathbb{C})$  and  $l \in \mathbb{N}$ . Then one can use Babai's nearest plane algorithm (see Algorithm 4) using  $\mathcal{L}(\mathcal{B}, \sigma, l)$  and  $t = (\lfloor \sigma(x) \rfloor_l, 0, \dots, 0)$ . It is also better to reduce  $B(\mathcal{B}, \sigma, l)$  with LLL first, and use Algorithm 4 with input this reduced basis and  $t$ . This leads to Algorithm 5.

**Notation 3.6.** Given  $K$  a number field,  $\mathcal{B}$  a  $\mathbb{Q}$ -basis of  $K$ ,  $\sigma \in \text{Hom}(K, \mathbb{C})$  and a precision  $l$ , we denote by  $L(\mathcal{B}, \sigma, l)$  the LLL-reduced basis of  $B(\mathcal{B}, \sigma, l)$ . Similarly, we will write  $L_l$  when there is no ambiguity.

The output of Algorithm 5 is a vector of coefficients which are expected to be the coefficients of  $x$  in the basis  $\mathcal{B}$ . The correctness of the outcome depends on the parameters chosen, i.e. the precision  $l$  and the constant  $C$ . We will determine a lower bound on the precision in Section 3.2, then under a reasonable conjecture on lattices  $\mathcal{L}_l$  we will exhibit a slightly better lower bound in Section 3.3.

Given  $x \in \mathcal{O}_K$ , if one denote by  $t_x$  the target vector  $[\lfloor 2^l x \rfloor, 0, \dots, 0]$  we wish to evaluate a precision  $l$  for which  $e_x := \text{NearestPlane}(\mathcal{L}_l, t_x)$  is equal to  $[\epsilon, Cx_1, \dots, Cx_n]$ . It is known that Babai's nearest plane

**Algorithm 5** TestDecode

**Require:** An integer  $l$ , the matrix  $L_l$  of a reduced basis of  $\mathcal{L}(\mathcal{B}, \sigma, l)$ ,  $[\sigma(x)]_l$  for some  $x \in K$ .

**Ensure:** A candidate  $y = (y_1, \dots, y_n)$  for the vector of coefficients of  $x$  expressed in  $\mathcal{B}$ .

- 1:  $r \leftarrow [K_\sigma : \mathbb{R}]$   $\triangleright r = 1$  if  $\sigma(K) \subset \mathbb{R}$ , and 2 otherwise
- 2:  $t \leftarrow ([\sigma_i(x)]_l, 0, \dots, 0) \in \mathbb{Z}^{n+r}$
- 3:  $v \leftarrow \text{NearestPlane}(t, L_l)$   $\triangleright$  (Alg. 4)
- 4: **return**  $[v_{r+1}, \dots, v_{n+r+1}]/C$

algorithm outputs a vector with norm smaller than half the norm of the shortest vector of the GSO. Thus, obtaining a lower bound on the first minimum of  $\mathcal{L}_l$  will allow us to obtain a bound on the norm of the output of  $\text{NearestPlane}(\mathcal{L}_l, t_g)$ . Coupled with an upper bound on the norm of  $e_g$ , this will give a condition for a correct decoding, corresponding to

$$(3.2) \quad \|e_g\|_2 \leq \frac{\lambda_1(\mathcal{L}_l)}{2^n}.$$

**Proposition 3.7.** *Consider  $K$  a number field,  $\mathcal{L}_l = \mathcal{L}(\mathcal{B}, \sigma, l)$  a basis matrix of  $K$ , and  $x \in \mathbb{Z}[\mathcal{B}]$ . Let  $m$  be an integer such that  $m = 1$  if  $\sigma(K) \subset \mathbb{R}$  and  $m = 2$  otherwise. Then Algorithm 5 outputs the correct vector of coefficients  $(x_1, \dots, x_n)$  of  $x$  in  $\mathcal{B}$  if the following holds:*

$$(3.3) \quad \frac{2^{m-1}(1 + \|x\|_1)^2}{4} + C^2 \|x\|_2^2 \leq \frac{\lambda_1(\mathcal{L}_l)^2}{2^{2n}}.$$

*Proof.* In TestDecode we wish to solve the BDD with respect to the lattice  $\mathcal{L}(\mathcal{B}, \sigma, l)$  and target vector  $t = ([\sigma(x)]_l, 0, \dots, 0)$ . As Proposition 2.20 states, the output is known to be correct if the distance between the target and the lattice is smaller than  $\frac{1}{2} \min \{\|\tilde{b}_i\|_2 \mid i \in [1, r]\}$ . The lattice vector  $v$  which is expected to be the closest to  $t$  is  $v = (\sum_{i=1}^n x_i [\sigma(b_i)]_l, -Cx_1, \dots, -Cx_n)$ . The error vector  $e = t - v$  is then equal to

$$\left( [\sigma(x)]_l - \sum_{i=1}^n x_i [\sigma(b_i)]_l, Cx_1, \dots, Cx_n \right).$$

Now let us consider the case where  $m = 1$  and look at the first coordinate of  $e$ . First write  $(\eta, \epsilon_1, \dots, \epsilon_n) \in [-\frac{1}{2}, \frac{1}{2}]^{n+1}$  the vector of errors due to the approximations, i.e.  $[\sigma(x)]_l = 2^l \sigma(x) + \eta$  and for all  $i \in [1, n]$ ,  $[\sigma(b_i)]_l = 2^l \sigma(b_i) + \epsilon_i$ . Then we have

$$e_1 = 2^l \sigma(x) + \eta - \sum_{i=1}^n 2^l x_i \sigma(b_i) + x_i \epsilon_i = \eta - \sum_{i=1}^n x_i \epsilon_i$$

which gives  $e_1^2 \leq (1 + \sum_{i=1}^n |x_i|)^2 / 4 = (1 + \|x\|_1)^2 / 4$ . If  $m = 2$ , or equivalently  $\sigma(K) \not\subset \mathbb{R}$ , one needs to consider the real and imaginary parts. Thus we get  $e \in \mathbb{R}^{n+2}$  and  $(\eta, \epsilon_1, \dots, \epsilon_n) \in [-\frac{1}{2}, \frac{1}{2}]^{n+2}$  with  $\eta = (\eta_1, \eta_2)$ . Following the previous analysis we obtain  $e_1^2 \leq (1 + \sum_{i=1}^n |x_i|)^2 \times 2/4 = (1 + \|x\|_1)^2 / 2$ . Thus we obtain the following upper bound for  $\|e\|_2^2$ :

$$\|e\|_2^2 \leq \frac{2^{m-1}}{4} (1 + \|x\|_1)^2 + C^2 \|x\|_2^2.$$

Finally remark that Theorem 2.17 tells us that a LLL-reduced basis  $(b_1, \dots, b_n)$  of a lattice  $\mathcal{L}$  (with  $\delta = \frac{3}{4}$ ) satisfies  $\|\tilde{b}_i\|_2^2 \geq \|\tilde{b}_1\|_2^2 / 2^{n-1} \geq \lambda_1(\mathcal{L}) / 2^{n-1}$ .  $\square$

**Notation 3.8.** Given  $x \in \mathbb{Z}[\mathcal{B}]$  and  $C \in \mathbb{N}^*$ , we will write  $m_x$  for the upper bound of  $\|e_x\|_2^2$ , i.e. such that  $m_x^2 := \frac{2^{m-1}(1 + \|x\|_1)^2}{4} + C^2 \|x\|_2^2$

Now if one has a lower bound for  $\lambda_1(\mathcal{L}_l)$  depending on the parameters, one can deduce a condition for the correctness of the output of `TestDecode`. In particular it is possible to obtain a lower bound of the precision  $l$  for which the computation is correct.

**3.2. Proven precision and correctness.** Let us determine for which precision [Algorithm 5](#) outputs the correct vector of coefficients.

**Notations 3.9.** Consider a number field  $K$  and  $\mathcal{B}$  a  $\mathbb{Q}$ -basis of  $K$ . Let us write  $D_{\mathcal{B}}$  for a denominator of  $\mathcal{B}$ , i.e. such that for all  $i \in \llbracket 1, n \rrbracket$ ,  $D_{\mathcal{B}}b_i \in \mathbb{Z}[X]$  where  $K \cong \mathbb{Q}[X]/(P_K(X))$ . Additionally we will denote  $\max_{i \in \llbracket 1, n \rrbracket} T_2(b_i)$  by  $m(\mathcal{B}, T_2)$ .

**Lemma 3.10.** *Consider  $K$  a number field,  $\mathcal{L}_l = \mathcal{L}(\mathcal{B}, \sigma, l)$  a basis matrix of  $K$ , and  $x \in \mathbb{Z}[\mathcal{B}]$ . For any integral vector  $\lambda \in B(0, 2^n m_x / C)$  we get the following lower bound for  $\|\lambda B_l\|_2$*

$$(3.4) \quad \|\lambda B_l\|_2 \geq \frac{2^l C^{n-1}}{n\sqrt{n}^n 2^{n(n-1)} m_x^{n-1} m(\mathcal{B}, T_2)^{n-1} D_{\mathcal{B}}^n (1 + \|P_K\|_2)^{2n-1}}.$$

*Proof.* We will focus on the case where  $\sigma$  is a real embedding, the proof in the case where  $\sigma(K) \not\subset \mathbb{R}$  being almost identical. We follow the proof of [\[33, Lemma A.7\]](#), where similar lattices are considered. A given  $\lambda \in \mathbb{Z}^n$  generates the vector of  $\mathcal{L}_l$  of the form

$$\left[ -\sum_{i=1}^r \lambda_i (2^l b_i^\sigma + \epsilon_{b_i}), C\lambda_1, \dots, C\lambda_n \right].$$

If we write  $s(\lambda, l) := \sum_{i=1}^n \lambda_i (2^l b_i + \epsilon_{b_i})$  then  $|s(\lambda, l)| \geq 2^l \left| \sum_{i=1}^n \lambda_i b_i^\sigma \right| - \sum_{i=1}^n \frac{|\lambda_i|}{2} \geq 2^l |\lambda^\sigma| - \frac{\|\lambda\|_1}{2}$ , where  $\lambda$  is seen as an element of  $K$ . Then we can write  $\|\lambda B_l(\mathcal{B})\|_2 \geq \|\lambda B_l(\mathcal{B})\|_1 / \sqrt{n} \geq 2^l |\lambda^\sigma| / \sqrt{n}$ . Let us find a lower bound for  $|\lambda^\sigma|$ . If  $\mu = D_{\mathcal{B}}\lambda$  then from [\[33, Lemma A.8\]](#) we have

$$(3.5) \quad |\mu^\sigma| \geq \frac{1}{n(1 + \|P_K\|_2^{2n-1}) T_2(\mu)^{n-1}}.$$

Then  $T_2(\mu) = T_2(D_{\mathcal{B}} \sum_{i=1}^n \lambda_i b_i) \leq D_{\mathcal{B}} \sum_{i=1}^n |\lambda_i| T_2(b_i) \leq D_{\mathcal{B}} m(\mathcal{B}, T_2) \|\lambda\|_1$ , so [Equation \(3.5\)](#) gives

$$|\lambda^\sigma| \geq \frac{1}{n \|\lambda\|_1^{n-1} D_{\mathcal{B}}^n m(\mathcal{B}, T_2)^{n-1} (1 + \|P_K\|_2)^{2n-1}}.$$

Finally, recalling that we consider  $\lambda \in B(0, 2^n m_x / C)$  and that  $\|\lambda\|_1 \leq \sqrt{n} \|\lambda\|_2$  we get the following lower bound for  $\lambda_1(\mathcal{L}_l)$

$$\lambda_1(\mathcal{L}_l) \geq \frac{2^l C^{n-1}}{n\sqrt{n}^n 2^{n(n-1)} m_x^{n-1} D_{\mathcal{B}}^n m(\mathcal{B}, T_2)^{n-1} (1 + \|P_K\|_2)^{2n-1}}.$$

□

**Theorem 3.11** (Proven precision for correctness). *Consider  $K$  a number field,  $\mathcal{B}$  a  $\mathbb{Q}$ -basis of  $K$  and  $x \in \mathbb{Z}[\mathcal{B}]$ . Additionally fix  $\sigma \in \text{Hom}(K, \mathbb{C})$ ,  $C \in \mathbb{N}$ , and  $m \in \mathbb{N}$  such that  $m = 1$  if  $\sigma(K) \subset \mathbb{R}$  and  $m = 2$  otherwise. If  $l \in \mathbb{N}$  with  $l \geq 1$  satisfies*

$$(3.6) \quad \begin{aligned} l \geq & n^2 + (n+1) \log_2 C + n \left( \log_2 \|x\|_2 + \frac{3 \log_2(n)}{2} + \log_2 D_{\mathcal{B}} + 2 \right) \\ & + (n-1) \log_2 m(\mathcal{B}, T_2) + (2n-1) \log_2 \|P_K\|_2 \end{aligned}$$

*then [Algorithm 5](#) outputs the vector of coefficients of  $x$  in the basis  $\mathcal{B}$ . In particular, the coefficients of  $x$  given one of its embedding can be recovered in time polynomial with respect to the size of the entries.*

*Proof.* Using Equations (3.2) and (3.5) we can conclude that the decoding is correct if

$$2^n m_x \leq \frac{2^l C^{n-1}}{n\sqrt{n} 2^{n(n-1)} m_g^{n-1} D_{\mathcal{B}}^n m(\mathcal{B}, T_2)^{n-1} (1 + \|P_K\|_2)^{2n-1}}$$

which is equivalent to

$$2^l \geq \frac{2^n m_x^n n \sqrt{n} 2^{n(n-1)} D_{\mathcal{B}}^n m(\mathcal{B}, T_2)^{n-1} (1 + \|P_K\|_2)^{2n-1}}{C^{n-1}}$$

which is verified as soon as

$$l \geq n^2 + n(\log_2(m_x) + \log_2(n)/2 + \log_2 D_{\mathcal{B}}) + (n-1)(\log_2 m(\mathcal{B}, T_2) - \log_2(C)) + (2n-1) \log_2(1 + \|P_K\|_2).$$

Now we can replace  $m_x$  by its value. However let us first rewrite it for simplicity, using  $\|x\|_2$  only.

Since  $\|x\|_1 \leq \sqrt{n} \|x\|_2$ , one can safely replace  $m_x^2$  by  $\frac{2^{m-1}(1+\sqrt{n}\|x\|_2)^2}{4} + C^2 \|x\|_2^2 = \|x\|_2^2 (C^2 + \frac{2^{m-1}}{4}(n + 2\sqrt{n}/\|x\|_2 + 1/\|x\|_2^2)) \leq \|x\|_2^2 (C^2 + 2n)$ . Replacing  $m_x^2$  by this last value we get that  $e_g$  is equal to some  $[\epsilon, Cg_1, \dots, Cg_n]$  as soon as

$$l \geq n^2 + n \left( \log_2 \|x\|_2 + \frac{3 \log_2(n)}{2} + \log_2(C) + \log_2 D_{\mathcal{B}} + 2 \right) + (n-1) \log_2 m(\mathcal{B}, T_2) + (2n-1) \log_2 \|P_K\|_2 + \log_2 C,$$

which gives the claimed inequality.  $\square$

**Remark 3.12.** If  $\mathcal{B}$  is an integral basis of  $\mathcal{O}_K$  then one can bound  $D_{\mathcal{B}}$  by  $n^n \sqrt{n} \|P_K\|_2^{2(n-1)}$  [26]. Additionally if  $\mathcal{B}$  is LLL-reduced for the  $T_2$ -norm, we have that  $m(\mathcal{B}, T_2) \leq D_K^{O(1)}$  [33] so Equation (3.6) becomes

$$l \geq n^2 + n \left( \log_2 \|g\|_1 + \frac{\log_2(n)}{2} + (n+1/2) \log_2 n \right) + (n-1)O(1) \log_2 D_K + (n+1)(2n-1) \log_2 \|P_K\|_2 + \log_2 C.$$

In the case where  $\mathcal{B}$  is a power basis, i.e.  $\mathcal{B} = (1, \alpha, \dots, \alpha^{n-1})$  we have  $D_{\mathcal{B}} \leq D_{\alpha}^{n-1}$  and  $m(\mathcal{B}, T_2) = \sqrt{n} \max_{\sigma} |\alpha^{\sigma}|^{n-1}$  so Equation (3.6) can be replaced by

$$l \geq n^2 + (n+1) \log_2 C + n \left( \log_2 \|x\|_2 + \frac{3 \log_2(n)}{2} + (n-1) \log_2 D_{\alpha} + 2 \right) + (n-1)^2 \log_2 \max_{\sigma} |\alpha^{\sigma}| + (n-1) \frac{\log_2 n}{2} + (2n-1) \log_2 \|P_K\|_2.$$

**3.3. Heuristic precision and correctness.** Let us now determine a lower bound under a conjecture regarding lattices of the form  $\mathcal{L}_l$  we handle. First we determine the determinant of said lattices.

**Lemma 3.13** (Matrix determinant lemma [18]). *Let  $A$  be a ring,  $M \in M_n(A)$  be an invertible matrix and  $U, V \in M_{n,m}(A)$ . Then the following is true:*

$$(3.7) \quad \det(M + UV^T) = (\text{Id}_m + V^T M^{-1} U) \det(M).$$

**Lemma 3.14.** *Let  $K$  be a number field and  $\mathcal{L}_l = \mathcal{L}(\mathcal{B}, \sigma, l)$  be a basis lattice of  $K$ . Also let  $m$  be an integer equal to 1 if  $\sigma$  is real, and 2 otherwise. Then*

$$(3.8) \quad \text{vol}(\mathcal{L}_l)^2 = C^{2n} \det \left( \text{Id}_m + \frac{1}{C^2} \sigma(\mathcal{B})_l^T \sigma(\mathcal{B})_l \right)$$

*Proof.* By definition  $\text{vol}(\mathcal{L}_l)^2$  is the determinant of the matrix  $B(\mathcal{B}, \sigma, l) B(\mathcal{B}, \sigma, l)^T = C^2 \text{Id}_n + \sigma(\mathcal{B})_l \times \sigma(\mathcal{B})_l^T$ . Then Lemma 3.13 gives the claimed identity.  $\square$

One can deduce from Lemma 3.14 a lower bound on the volume of  $\mathcal{L}(\mathcal{B}, \sigma, l)$  which depends on the precision  $l$  and the size of  $\sigma(\mathcal{B})$ .

**Notation 3.15.** Consider a number field  $K$ ,  $\mathcal{B}$  a  $\mathbb{Q}$ -basis of  $K$ ,  $\sigma \in \text{Hom}(K, \mathbb{C})$  and  $l \in \mathbb{N}$ . Let us define the value  $\Delta(\mathcal{B}, \sigma, l)$ . If  $\sigma(K) \subset \mathbb{R}$  we set  $\Delta(\mathcal{B}, \sigma, l) = \|\sigma(\mathcal{B})\|_2^2 - \|\sigma(\mathcal{B})\|_1 / 2^l$ , and if  $\sigma(K) \not\subset \mathbb{R}$  we set  $\Delta(\mathcal{B}, \sigma, l) = \|\Re(\sigma(\mathcal{B}))\|_2^2 + \|\Im(\sigma(\mathcal{B}))\|_2^2 - \|\Re(\sigma(\mathcal{B}))\|_1 / 2^l - \|\Im(\sigma(\mathcal{B}))\|_1 / 2^l$ .

**Proposition 3.16.** *Let  $K$  be a number field and  $\mathcal{L}_l = \mathcal{L}(\mathcal{B}, \sigma, l)$  be a basis lattice of  $K$ . Also let  $m$  be an integer equal to 1 if  $\sigma$  is real, and 2 otherwise. Then the following is true,*

$$(3.9) \quad \text{vol}(\mathcal{L}_l)^2 \geq C^{2n} \left( 1 + \frac{2^{2l}}{C^2} \Delta(\mathcal{B}, \sigma, l) \right).$$

*Proof.* From Equation 3.8 we can write for  $m = 1$

$$\text{vol}(\mathcal{L}_l)^2 = C^{2n} \left( 1 + \frac{1}{C^2} \|\sigma(\mathcal{B})_l\|_2^2 \right) = C^{2n} \left( 1 + \frac{1}{C^2} \sum_{i=1}^n [\sigma(b_i)]_l^2 \right).$$

Then for each  $i \in \llbracket 1, n \rrbracket$ , there is  $\epsilon_i \in [-\frac{1}{2}, \frac{1}{2}]$  such that  $[\sigma(b_i)]_l = 2^l \sigma(b_i) + \epsilon_i$ . Thus we obtain

$$\begin{aligned} \text{vol}(\mathcal{L}_l)^2 &= C^{2n} \left( 1 + \frac{1}{C^2} \sum_{i=1}^n (2^l \sigma(b_i) + \epsilon_i)^2 \right) \\ &= C^{2n} \left( 1 + \frac{2^{2l}}{C^2} \sum_{i=1}^n (\sigma(b_i)^2 + 2\sigma(b_i) \frac{\epsilon_i}{2^l} + \frac{\epsilon_i^2}{2^{2l}}) \right). \end{aligned}$$

Since for all  $i \in \llbracket 1, n \rrbracket$ , one has  $|\epsilon_i| \leq \frac{1}{2}$ , we obtain the inequality

$$\text{vol}(\mathcal{L}_l)^2 \geq C^{2n} \left( 1 + \frac{2^{2l}}{C^2} \sum_{i=1}^n (\sigma(b_i)^2 - 2 \times \frac{1}{2} \times \frac{|\sigma(b_i)|}{2^l}) \right).$$

If  $\sigma(K) \not\subset \mathbb{R}$  then from Equation 3.8 we have  $\text{vol}(\mathcal{L}_l)^2 = C^{2n} \det(\text{Id}_2 + \frac{1}{C^2} \sigma(\mathcal{B})_l^\top \sigma(\mathcal{B})_l)$ . If we write  $M = \text{Id}_2 + \frac{1}{C^2} \sigma(\mathcal{B})_l^\top \sigma(\mathcal{B})_l$ , then

$$M = \begin{bmatrix} 1 + \frac{\|\Re(\sigma(\mathcal{B})_l)\|_2^2}{C^2} & \frac{(\Re(\sigma(\mathcal{B})_l) \mid \Im(\sigma(\mathcal{B})_l))}{C^2} \\ \frac{(\Re(\sigma(\mathcal{B})_l) \mid \Im(\sigma(\mathcal{B})_l))}{C^2} & 1 + \frac{\|\Im(\sigma(\mathcal{B})_l)\|_2^2}{C^2} \end{bmatrix}$$

so we get

$$\begin{aligned} \det(M) &= 1 + \frac{\|\Re(\sigma(\mathcal{B})_l)\|_2^2}{C^2} + \frac{\|\Im(\sigma(\mathcal{B})_l)\|_2^2}{C^2} + \frac{\|\Re(\sigma(\mathcal{B})_l)\|_2^2 \|\Im(\sigma(\mathcal{B})_l)\|_2^2}{C^4} \\ &\quad - \frac{(\Re(\sigma(\mathcal{B})_l) \mid \Im(\sigma(\mathcal{B})_l))^2}{C^4}. \end{aligned}$$

By Cauchy-Schwarz inequality we have

$$\frac{\|\Re(\sigma(\mathcal{B})_l)\|_2^2 \|\Im(\sigma(\mathcal{B})_l)\|_2^2}{C^4} - \frac{(\Re(\sigma(\mathcal{B})_l) \mid \Im(\sigma(\mathcal{B})_l))^2}{C^4} \geq 0,$$

therefore

$$\text{vol}(\mathcal{L}_l)^2 \geq C^{2n} \left( 1 + \frac{\|\Re(\sigma(\mathcal{B})_l)\|_2^2}{C^2} + \frac{\|\Im(\sigma(\mathcal{B})_l)\|_2^2}{C^2} \right).$$

Following the same reasoning that we did for the real case, we can conclude that

$$\frac{\|\Re(\sigma(\mathcal{B})_l)\|_2^2}{C^2} \geq \frac{2^{2l}}{C^2} \left( \|\Re(\sigma(\mathcal{B}))\|_2^2 - \frac{\|\Re(\sigma(\mathcal{B}))\|_1}{2^l} \right)$$

and

$$\frac{\|\Im(\sigma(\mathcal{B}))_l\|_2^2}{C^2} \geq \frac{2^{2l}}{C^2} \left( \|\Im(\sigma(\mathcal{B}))\|_2^2 - \frac{\|\Im(\sigma(\mathcal{B}))\|_1}{2^l} \right),$$

which gives the desired result.  $\square$

It is now possible to certify the correctness of the decoding. The Gaussian heuristic provides an *estimation* of  $\lambda_1(\mathcal{L}_l)$ , which can be used to obtain a *heuristic* condition for the correctness of the algorithm, as in Theorem 3.17.

**Theorem 3.17** (Correctness of decoding). *Consider  $K$  a number field,  $\mathcal{B}$  a  $\mathbb{Q}$ -basis of  $K$  and  $x \in \mathbb{Z}[\mathcal{B}]$ . Additionally fix  $\sigma \in \text{Hom}(K, \mathbb{C})$ ,  $C \in \mathbb{N}$ , and  $m \in \mathbb{N}$  such that  $m = 1$  if  $\sigma(K) \subset \mathbb{R}$  and  $m = 2$  otherwise. If  $l \in \mathbb{N}$  with  $l \geq 1$  satisfies*

$$(3.10) \quad l \geq n^2 + n \log_2 \|x\|_2 + \frac{n}{2} \log_2(2\pi e(\frac{1}{n} + 2)) + \log_2(C) - \frac{\log_2 \Delta(\mathcal{B}, \sigma, l)}{2}$$

*then Algorithm 5 outputs the vector of coefficients of  $x$  in the basis  $\mathcal{B}$ , if  $\mathcal{L}(\mathcal{B}, \sigma, l)$  satisfies the Gaussian heuristic (eq. 2.2).*

*Proof.* Let us fix  $l \in \mathbb{N}$ . The Gaussian heuristic states  $\lambda_1(\mathcal{L}) \sim \sqrt{\frac{n}{2\pi e}} \det \mathcal{L}^{1/n}$ . Then one can combine Equation (3.3) to obtain a conditional inequality. In order to simplify the expression, we will replace  $m_x^2$  by  $\|x\|_2^2 (C^2 + 2n)$  as explained in the proof of Theorem 3.11. This leads to the following inequality:

$$\|x\|_2^2 (C^2 + 2n) \leq \frac{1}{2^{2n}} \left( \frac{n}{2\pi e} C^2 \sqrt[n]{\left(1 + \frac{2^{2l}}{C^2} (\Delta(\mathcal{B}, \sigma, l))\right)} \right).$$

We can obtain the condition

$$(3.11) \quad \|x\|_2^2 (C^2 + 2n) \leq \frac{n}{2\pi e} C^2 \sqrt[n]{\left(\frac{2^{2l}}{C^2} (\Delta(\mathcal{B}, \sigma, l))\right)}.$$

Under the logarithmic map, the left side of Equation (3.11) gives  $2 \log_2 \|x\|_2 + 2 \log_2(C) + \log_2(1 + \frac{2n}{C^2})$  while the right side becomes  $\log_2(\frac{n}{2\pi e}) - 2n + \frac{1}{n} \times (2(n-1) \log_2(C) + 2l + \log_2(\Delta(\mathcal{B}, \sigma, l)))$ , and combining them allows us to retrieve the claimed condition.  $\square$

One can remark that Theorem 3.17 gives a better condition than the theoretical one given by Theorem 3.11.

**Remark 3.18.** As a matter of fact, correctness of decoding is ensured by Equation (3.10) as long as  $\mathcal{L}_l$  satisfies  $\lambda_1(\mathcal{L}_l) \geq \lambda_1(\mathcal{L}_l)_{\text{gauss}}$ .

In order to estimate the correctness of the value given by Theorem 3.17, we verified if the Gaussian heuristic holds for lattices  $\mathcal{L}_l$ . We computed the average value of the quotient  $\lambda_1(\mathcal{L}_l)/\lambda_1(\mathcal{L}_l)_{\text{gauss}}$  for increasing values of  $l$ , over random number fields  $K$  with fixed degrees. Here “random” means defined by integral polynomials whose coefficients are drawn uniformly in  $[-2^s, 2^s]$  for randomly chosen  $1 \leq s \leq 10$ . The results can be found in Figure 1. One can see that  $\lambda_1(\mathcal{L}_l)$  is larger than the value predicted by the Gaussian heuristic. This tends to show that asymptotically, one can safely consider that Theorem 3.17 provides a correct value certifying the correctness of the output of TestDecode. Note that running this kind of experiment for large dimensions (say larger than 120) would be intractable. Indeed, even approximating  $\lambda_1(\mathcal{L})$  is a hard problem.

**Heuristic 2** (Gaussian heuristic for bases lattices). *Consider  $K$  a number field,  $\mathcal{B}$  a  $\mathbb{Q}$ -basis of  $K$ ,  $\sigma \in \text{Hom}(K, \mathbb{C})$  and  $l \in \mathbb{N}$ . Then  $\lambda_1(\mathcal{L}(\mathcal{B}, \sigma, l)) \geq \lambda_1(\mathcal{L}_l)_{\text{gauss}}$ .*



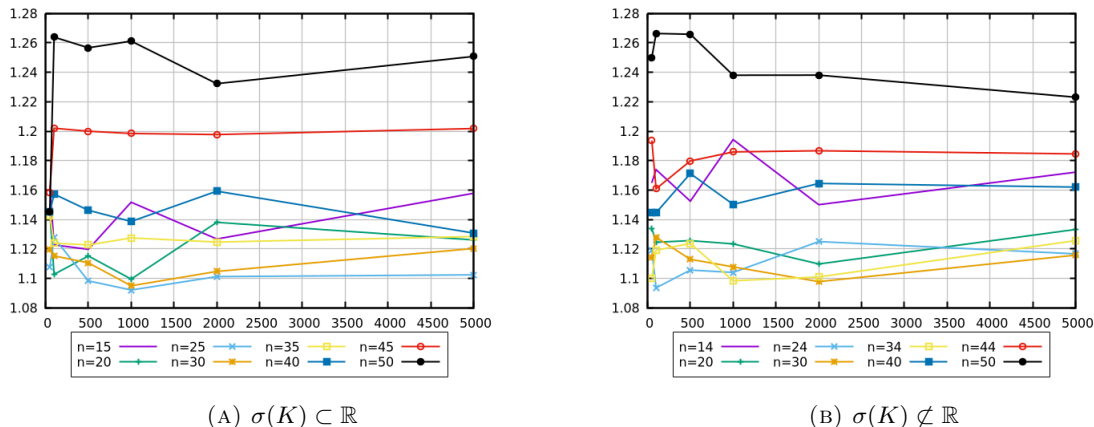


FIGURE 1. Average value of  $\lambda_1(\mathcal{L}_l)/\lambda_1(\mathcal{L}_l)_{\text{gauss}}$  plotted against the precision  $l$ , for several  $n = [K : \mathbb{Q}]$ .

**3.4. Computing polynomial roots.** We described which lattice we use, and the decoding method. Obviously the reduced matrix  $L(\mathcal{B}, \sigma, l)$  is computed once and used to retrieve all roots. Therefore we have the following main steps:

- (1) fix an embedding  $\sigma : K \hookrightarrow \mathbb{C}$ ;
- (2) compute a precision  $l \in \mathbb{N}$  ensuring that the decoding is correct;
- (3) compute a LLL-reduced basis  $L_l$  of the lattice generated by the basis of  $K$ ;
- (4) using  $\sigma$ , compute approximations of the roots of  $f(T)$  up to a precision  $l$ ;
- (5) use  $L_l$  to retrieve the roots of  $f(T)$  using TestDecode.

We will call `PrecisionEvaluation` the function returning the needed precision for input  $f(T) \in K[T]$ . Following Theorem 3.17, it depends on the Euclidean norms of the roots of  $f(T)$ . In order to evaluate an upper bound of these norms, we can follow [3]. We will denote by `FloatPolynomialRoots` the procedure computing the real (resp. complex roots) of a real (resp. complex) polynomial. We obtain Algorithm 6 describing the method we implemented to compute  $Z_K(f)$ .

Following the results from the previous section, one can state the following theorem.

**Theorem 3.19.** *Consider a number field  $K$ ,  $\mathcal{B}$  a  $\mathbb{Q}$ -basis of  $K$  and  $f(T) \in K[T]$  such that  $Z_K(f) \subset \mathbb{Z}[\mathcal{B}]$ . Then for input  $(K, \mathcal{B}, f(T))$ , Algorithm 6 outputs  $Z_K(f)$  in polynomial time.*

Finally we give a broad estimate of the algebraic complexity in function of the precision required to ensure correctness. We will only count the complexity linked to lattices operations, as other steps are negligible. Note as well that we express these complexities in terms of the precision at which computations are made instead of the size of the roots since the latter is unknown. Additionally this allows us to take into account different choices for `PrecisionEvaluation`.

**Proposition 3.20.** *Consider a number field  $K$ ,  $\mathcal{B}$  a  $\mathbb{Q}$ -basis of  $K$  and  $f(T) \in K[T]$  such that  $Z_K(f) \subset \mathbb{Z}[\mathcal{B}]$ . Then for input  $(K, \mathcal{B}, f(T))$ , the algebraic complexity of lattice operations of Algorithm 6 (reduction of a basis lattice  $\mathcal{L}_l$ , GSO computation and decoding) are respectively in  $O(n^4l + n^3l^2)$ ,  $O(n^3)$  and  $O(dn^2)$ , where  $n$  is the dimension  $[K : \mathbb{Q}]$ .*

*Proof.* Recall by Theorem 2.18 the algebraic complexity of LLL is in  $O(n^4M + n^3M^2)$ , where  $M$  is an upper-bound on the bit size of basis elements. A broad estimate of  $M$  for  $B(\mathcal{B}, \sigma, l)$  is  $l$ . For the GSO

**Algorithm 6** AbsoluteRoots

**Require:** A number field  $K$ ,  $f(T) \in K[T]$ ,  $\mathcal{B}$  a  $\mathbb{Q}$ -basis of  $K$  such that  $Z_K(f) \in \mathbb{Z}[\mathcal{B}]$ .

**Ensure:** The set  $Z_K(f)$

```

1:  $\sigma \leftarrow \text{ChooseEmbedding}(K)$ 
2:  $l \leftarrow \text{PrecisionEvaluation}(f(T))$ 
3:  $L_l \leftarrow \text{LLL}(B(\mathcal{B}, \sigma, l))$ 
4:  $Z \leftarrow \text{FloatPolynomialRoots}(f^\sigma, l)$ 
5:  $S \leftarrow \emptyset$ 
6: for  $z \in Z$  do
7:    $y \leftarrow \text{TestDecode}(L_l, z)$ 
8:   if  $f(y) = 0$  then
9:      $S \leftarrow S \cup \{y\}$ 
10:  end if
11: end for
12: return  $S$ 

```

computation, we can follow a QR decomposition as described in [29] for example which gives a complexity of  $O(n^3)$ . Finally Algorithm 6 makes  $d = \deg f(T)$  calls to Algorithm 4, adding up to  $O(dn^2)$ .  $\square$

## 4. RELATIVE METHOD

Let us now describe a method to recover the roots of a polynomial in an extension of number fields  $L/K$ , which is one of our main contribution.

**4.1. Decoding in a subfield.** First let us describe how one can reduce knowledge of embeddings related to the extension  $L/K$  to decodings in  $K$ . Fix  $\Omega$  an algebraic closure of  $L/K$ . We will use the relative Minkowski embedding  $\sigma_{L/K} : x \in L \mapsto (\sigma(x))_{\sigma \in \text{Hom}_K(L, \Omega)}$ . It defines a  $K$ -linear embedding of  $L \cong K^n$  into  $\Omega^n$  where  $n$  is the degree of  $L/K$ . More precisely, let us fix  $\mathcal{E} = (e_1, \dots, e_n)$  a  $K$ -basis of  $L$  and let  $x = x_1 e_1 + \dots + x_r e_n \in L$ . We assume that the action of each  $\sigma \in \text{Hom}_K(L, \Omega)$  on  $\mathcal{E}$  is known. Then  $\sigma_{L/K}$  sends  $K^n$  into  $\Omega^n$ ,

$$\begin{aligned} \sigma_{L/K} : \quad K^n &\longrightarrow \Omega^n \\ (x_i)_{i \in [1, n]} &\longmapsto \left( \sum_{i=1}^n x_i \sigma(e_i) \right)_{\sigma \in \text{Hom}_K(L, \Omega)}. \end{aligned}$$

Thus given knowledge of  $\sigma_{L/K}(x)$  one can apply  $\sigma_{L/K}^{-1}$  and obtain recover  $(x_i)_i$ . For computational purposes, we need to consider embeddings into  $\mathbb{C}^n$ . In order to do so, one needs to fix  $\tau$  an embedding of  $K$  into  $\mathbb{C}$  then consider  $\sigma_{L/K}$  as the collection of complex embeddings of  $L$  extending  $\tau$ . This leads to Algorithm 7 which retrieves coefficients of  $x$  knowing its Minkowski embedding relative to  $L/K$ .

**Lemma 4.1.** *Algorithm 7 runs in polynomial time and outputs the correct vector of coefficients provided  $l$  is large enough.*

*Proof.* The polynomial running time is clear. For any  $\sigma : L \hookrightarrow \mathbb{C}$  extending  $\tau$ , its action on  $x \in L$  can be seen as  $\sum_{i=1}^n \tau(x_i) \sigma(e_i)$ . This way,  $K$  is identified with a subfield of  $\mathbb{C}$  and  $L$  is identified with  $\tau(K)^n$ . The action of  $\sigma_{L/K}$  can be expressed as an action from  $\tau(K)^n$  into  $\mathbb{C}^n$  as a matrix in  $M_n(\mathbb{C})$ , that is  $\Sigma_{L/K}$ . In particular we have  $\Sigma_{L/K}[i] = \sigma_{L/K}(e_i)$ , thus  $XM = X \Sigma_{L/K}^{-1} = (x_1^\tau, \dots, x_n^\tau)_{[K:\mathbb{Q}]}$ . Thus knowing approximations of  $\sigma_{L/K}(x)$  and  $\Sigma_{L/K}$  up to  $l + s$ , it is possible to find the approximations of  $(\tau(x_i))_i$

**Algorithm 7** Mink2coeff

**Require:** An extension  $L/K$ , given by  $\mathcal{B}$  a basis of  $K$  and  $\mathcal{E}$  a  $K$ -basis of  $L$ ,  $M = \Sigma_{L/K}^{-1}$  the matrix of  $\sigma_{L/K}^{-1}$  expressed with respect to  $\tau(\mathcal{E})$  up to a precision  $l + s$ , the matrix  $L_l$  from a reduced basis of  $\mathcal{L}(\mathcal{B}, \tau, l)$  for some  $\tau \in \text{Hom}(K, \mathbb{C})$  and  $X = [\sigma_{L/K}(x)]_{l+s}$  for some  $x \in L/K$ .

**Ensure:** A candidate  $y = (y_1, \dots, y_N)$  for the vector of coefficients of  $x$  expressed in  $\mathcal{B} \otimes \mathcal{E}$ .

```

1:  $Y \leftarrow XM$ 
2:  $y = [ \ ]$ 
3: for  $i = 1$  to  $n$  do
4:    $y \leftarrow [y \mid \text{TestDecode}(L_l, Y_i)]$  ▷ Concatenation of row vectors
5: end for
6: return  $y$ 

```

up to  $l$ . Then by the results of [Sections 3.2](#) and [3.3](#) we know that  $\text{TestDecode}(L_l, [x_i]_l)$  is the vector of coefficients of  $x_i$  with respect to  $\mathcal{B}$  provided that  $l$  is large enough.  $\square$

**4.2. An algorithm for polynomial roots.** We can apply the previous strategy to compute polynomial roots by decoding in the subfield  $K$ . Again, we fix some objects. The extension of number fields  $L/K$  is given by a  $\mathbb{Q}$ -basis  $\mathcal{B}$  and  $\mathcal{E}$  a  $K$ -basis of  $L$ . Then consider a polynomial  $f(T) \in L[X]$  such that  $Z_L(f) \subset \mathbb{Z}[\mathcal{E} \otimes \mathcal{B}]$ . From what we described previously, in order to retrieve the coefficients of  $x \in Z_L(f)$ , one can compute  $\sigma_{L/K}$  and use `Mink2coeff` with the following main steps:

- (1) compute a precision  $l$  certifying the correctness of the computation;
- (2) compute  $L_l$  a LLL reduced basis of  $\mathcal{L}(\mathcal{B}, \tau, l)$ ;
- (3) compute  $M = \Sigma_{L/K}^{-1}$  up to precision  $l + s$  for some  $s$ ;
- (4) compute  $Z = \prod_{\sigma \in \text{Hom}_K(L, \mathbb{C})} Z_{\mathbb{C}}(f^\sigma)$  up to precision  $l + s$ ;
- (5) For each  $x \in Z$ , use `Mink2coeff` to obtain a root candidate.

This leads to [Algorithm 8](#).

**Remark 4.2.** The set  $Z$  is the cartesian product of the sets  $Z(f^\sigma)$  with  $\sigma$  in  $\text{Hom}_K(L, \mathbb{C})$ . Each of such set  $Z(f^\sigma)$  has at most  $d = \deg f$  elements. Therefore,  $Z$  is a set of at most  $d^n$  complex numbers. Moreover one cannot tell a priori if an element  $z \in Z(f^\sigma)$  is of the form  $[\sigma(x)]_{l+s}$  for some  $x \in Z_L(f)$ , or even if a vector  $(z_1, \dots, z_n) \in Z$  corresponds to a root  $x$  of  $f(T)$ . Thus one has to call `Mink2coeff`  $d^n$  times in the worst case. Even if  $f(T)$  splits in  $L$ , this leads to a search of  $d$  vectors in a set of size  $d^n$ . In order to improve slightly this search, a simple observation can be made. Let us write  $Z = \prod_{\sigma} Z_{\mathbb{C}}(f^\sigma)$  where  $\sigma$  ranges over  $\text{Hom}_K(L, \mathbb{C})$ . Then one has

$$(4.1) \quad \forall x \in Z_K(f), \forall \sigma \in \text{Hom}_K(L, \mathbb{C}), \exists! z \in Z_{\mathbb{C}}(f^\sigma) \mid z = [\sigma(x)]_l,$$

which implies that once we found a correct vector in  $Z$  we can remove from the search tree all the nodes where any of its coordinates appears. A more precise study of the average cost of the search (number of vectors of  $Z$  tested) can be found in [Appendix A](#).

**Notation 4.3.** We denote by `UpdateTree` the procedure updating the search space as described.

The absolute method `AbsoluteRoots` requires at most  $d = \deg f(T)$  decodings, while `RelativeRoots.naive` requires enumerating through  $d^n$  vectors, corresponding to  $n$  decodings each. The mere enumeration of  $d^n$  elements shows that `RelativeRoots.naive` has an exponential cost when  $n$  increases. In addition, several operations on vectors and matrices are done for each of the  $d^n$  possibilities. Thus it is quickly impractical.

**Algorithm 8** RelativeRoots\_naive

**Require:** An extension  $L/K$ , given by  $\mathcal{B}$  a basis of  $K$  and  $\mathcal{E}$  a  $K$ -basis of  $L$ , and  $f(T) \in L[T]$  such that

$$Z_L(f) \in \mathbb{Z}[\mathcal{E} \otimes \mathcal{B}]$$

**Ensure:**  $Z_L(f)$

```

1:  $l \leftarrow \text{PrecisionEvaluation}(f(T))$ 
2:  $L_l \leftarrow \text{LLL}(B(\mathcal{B}, \tau, l))$ 
3:  $M \leftarrow \Sigma_{L/K}^{-1}$  ▷ Up to precision  $l + s$ 
4:  $Z \leftarrow \prod_{\sigma \in \text{Hom}_K(L, \mathbb{C})} \text{FloatPolynomialRoots}(f^\sigma, l + s)$ 
5:  $S \leftarrow \emptyset$ 
6: for  $z \in Z$  do
7:    $y \leftarrow \text{Mink2coeff}(z, M, L_l)$  ▷ Compute the inverse and decode
8:   if  $f(y) = 0$  then
9:      $S \leftarrow S \cup \{y\}$ 
10:    $\text{UpdateTree}(Z, z)$ 
11:   end if
12: end for
13: return  $S$ 

```

However, for small  $n$ , Algorithm 8 can considerably speed-up the computation of  $Z_K(f)$ . Indeed one has to remember that an important part of the computation time is dedicated to the reduction of the lattice used to decode, as is also the case for the algebraic method. Algorithm 6 requires the reduction of a lattice of rank  $[L : \mathbb{Q}]$ , while the rank of the lattice reduced in Algorithm 8 is  $[L : \mathbb{Q}]/n$ . In addition, the precision needed to certify the computation shown in Theorem 3.17 involves the dimension. Therefore, dividing the dimension allows us to do computations at a smaller precision, which also leads to smaller coefficients in the matrix  $B_l$  which is reduced.

**Theorem 4.4.** *Consider an extension  $L/K$ ,  $\mathcal{B}$  a  $\mathbb{Q}$ -basis of  $K$ ,  $\mathcal{E}$  a  $K$ -basis of  $L$  and  $f(T) \in L[T]$  such that  $Z_L(f) \subset \mathbb{Z}[\mathcal{E} \otimes \mathcal{B}]$ . Then for input  $(L/K, \mathcal{E}, \mathcal{B}, f(T))$ , Algorithm 8 outputs  $Z_L(f)$  in polynomial time, except eventually for the enumeration due to the **for** loop.*

*Proof.* Since the required precision to ensure correctness of decoding has been proved to be polynomial with respect to entry sizes, the polynomial time complexity of each intermediate operations is clear. For the correctness, remark that for any  $x \in Z_L(f)$  and any  $\sigma \in \text{Hom}_{\mathbb{Q}}(L, \mathbb{C})$ ,  $\sigma(x)$  is a root of  $f^\sigma$ . Thus,  $\sigma_{L/K}(x) \in \prod_{\sigma \in \text{Hom}_K(L, \mathbb{C})} Z_{\mathbb{C}}(f^\sigma)$ , so for each  $x \in Z_L(f)$  there is one vector of  $Z$  in Algorithm 8 which is of the form  $([\sigma(x)]_{l+s})_\sigma = [\sigma_{L/K}(x)]_{l+s}$ . This ensures that  $S$  is the correct set at the end of Algorithm 8.  $\square$

**Proposition 4.5.** *Consider an extension  $L/K$ ,  $\mathcal{B}$  a  $\mathbb{Q}$ -basis of  $K$ ,  $\mathcal{E}$  a  $K$ -basis of  $L$  and  $f(T) \in L[T]$  such that  $Z_L(f) \subset \mathbb{Z}[\mathcal{E} \otimes \mathcal{B}]$ . Then for input  $(L/K, \mathcal{E}, \mathcal{B}, f(T))$ , the algebraic complexity of lattice operations of Algorithm 8 (reduction of a basis lattice  $\mathcal{L}_l$ , GSO computation and decoding) are respectively in  $O(n_K^4 l + n_K^3 l^2)$ ,  $O(n_K^3)$  and  $n_{L/K} d^{n_{L/K}} O(n_K^2)$ , where  $l$  is the required precision.*

*Proof.* The proof is similar to the one of Proposition 3.20, noting that the lattice used to decode is linked to  $K$ , that there are  $d^{n_{L/K}}$  elements in  $Z$  and that Mink2coeff makes  $n_{L/K}$  calls to TestDecode for each of these elements.  $\square$

**4.3. Trade-off.** Now let us express some broad trade-off between our absolute and relative methods, considering only computations linked to lattice reduction and decoding. We forget about the GSO computation since it is done only once and is negligible with respect to the lattice reduction. Consider  $L/K$  a number field extension and write  $n_L, n_K$  and  $n_{L/K}$  for the degrees  $[L : \mathbb{Q}]$ ,  $[K : \mathbb{Q}]$  and  $[L : K]$  respectively. Following [Propositions 3.20](#) and [4.5](#) we can conclude that the complexity of [Algorithm 6](#) due to lattice reduction and decoding is  $O(n_L^4 l + n_L^3 l + dn_L^2)$  while these operations amount to  $O(n_K^4 l + n_K^3 l + n_{L/K} d^{n_{L/K}} n_K^2)$  for [Algorithm 8](#). Thus, we obtain a condition for a trade-off :

$$(4.2) \quad \begin{aligned} n_K^2(n_K^2 l + n_K l + n_{L/K} d^{n_{L/K}}) &= n_L^2(n_L^2 l + n_L l + d) \\ \iff n_{L/K} d(d^{n_{L/K}-1} - n_{L/K}) &= n_K l(n_{L/K}^3(n_L + 1) - (n_K + 1)). \end{aligned}$$

Using [Equation \(5.6\)](#) allows us to obtain a broad indication on which algorithm to run. For example, fixing  $[L : K] = n_{L/K} = 2$  we obtain  $2d(d-2) \leq n_K l(15n_K + 7)$ , therefore our relative algorithm would be advantageous for  $d$  up to (at least)  $\sqrt{\frac{n_K l(15n_K + 7)}{2}}$ . Note that following both theoretical and heuristic bounds from [Theorems 3.11](#) and [3.17](#),  $l$  is larger than  $n_L^2 + n_L \log_2 \|x\|_2$  so  $\sqrt{\frac{n_K l(15n_K + 7)}{2}}$  can be replaced by

$$\sqrt{\frac{n_K(2n_K(2n_K + \log_2 \|x\|_2))(15n_K + 7)}{2}} = n_K \sqrt{(2n_K + \log_2 \|x\|_2)(15n_K + 7)}.$$

## 5. IMPROVEMENTS AND HEURISTIC OBSERVATIONS

In order to improve the naive algorithms presented in [Sections 3](#) and [4](#), several directions can be explored. The main bottleneck of our absolute method ([Alg. 6](#)) is the reduction of a basis lattice  $\mathcal{L}(\mathcal{B}, \sigma, l)$ . It can be hasten by improving the reduction process itself or by evaluating the required precision  $l$  more accurately. Our relative method potentially requires many decodings, i.e. calls to a BDD solver such as [NearestPlane \(Algorithm 4\)](#). Therefore, improving this step can speed-up considerably the running time of [Algorithm 8](#).

**5.1. Basis reduction in the case of power-bases.** The first heuristic improvement that we develop concern the reduction of  $B(\mathcal{B}, \sigma, l)$ , in the case where  $\mathcal{B}$  is a power-basis, i.e there is  $\alpha \in K$  such that  $\mathcal{B} = (1, \alpha, \dots, \alpha^{n-1})$ . The idea behind the improvement comes from the following observation. First, denote by  $B^{(i)}$  the top left  $i \times (i+1)$  submatrix of  $B$ , then assume that  $B$  is reduced up to the  $i$ -th vector, i.e. we found  $U^{(i)}$  such that  $L^{(i)} = U^{(i)} B^{(i)}$  is reduced. Then we found  $u^{(i)} \in \mathbb{Z}^i$  such that  $u^{(i)} B^{(i)} = L^{(i)}[i]$ . Consider now the  $(i+1)$ -th vector of the matrix. Then

$[0, u^{(i)}] B^{(i+1)} \approx \alpha u^{(i)} B^{(i)} = \alpha L^{(i)}[i]$ . Thus applying the transformation operated to reduce the  $i$ -th vector should pre-reduce the  $(i+1)$ -th vector. This leads to [Algorithm 9](#), which computes a LLL-reduced basis  $L_l$ . Note that pre-reduction strategies have already been mentioned or used in the literature [[39](#), [21](#)]. To the best of our knowledge these apply to general knapsack-like matrices, while our observation tries to take advantage of the regularity of the basis vectors, and can be compared to [[34](#)].

We compared [SpecLLL](#) with LLL for available implementations of [MAGMA \[11\]](#), [PARI/GP \[31\]](#) and [FPLL \[40\]](#). The results can be found in [Figure 2](#) for the real case and [Figure 3](#) for the complex one.

The first observation is that [SpecLLL](#) is asymptotically faster than LLL, regardless of the software used. Then one can remark that the ratio is decreasing when both the precision or the dimension is increasing.

[SpecLLL](#) is always more efficient than LLL when used in [PARI/GP](#) or [MAGMA](#), with a asymptotic gain of 25% and 35% respectively. For [FPLL](#), using [Algorithm 9](#) results in larger running times for small

**Algorithm 9** SpecLLL

**Require:**  $\mathcal{B}$ , a basis of a lattice  $\mathcal{L}$  of rank  $n$ .

**Ensure:**  $\mathcal{B}'$ , a LLL reduced basis of  $\mathcal{L}$ .

- 1:  $\mathcal{B}' \leftarrow \{b_1\}$
- 2: **for**  $i = 2$  to  $n$  **do**
- 3:    $b'_i \leftarrow b_i + \left\| b'_{i-1} \right\|_2^2 e_{i+1}$   $\triangleright M$  is a large constant to ensure reduction
- 4:    $b'_i \leftarrow b'_i - [0, u^{(i-1)}] B^{(i)}$   $\triangleright$  Pre-reduction using the previous vector
- 5:    $\mathcal{B}' \leftarrow \mathcal{B}' \cup \{b'_i\}$
- 6:    $\mathcal{B}' \leftarrow \text{LLL}(\mathcal{B}')$   $\triangleright$  Size-reduction
- 7:    $b'_i \leftarrow b'_i + (C - \|b'_{i-1}\|_2^2) e_{i+1}$
- 8: **end for**
- 9:  $\mathcal{B}' \leftarrow \text{LLL}(\mathcal{B}')$
- 10: **return**  $\mathcal{B}'$

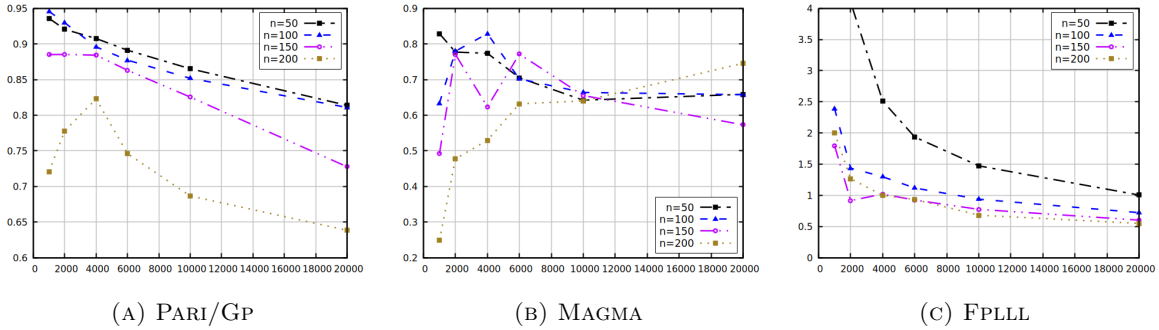


FIGURE 2. Ratios of running times of SpecLLL to LLL plotted against the precision  $l$ , for  $\sigma(K) \subset \mathbb{R}$  and several dimensions  $n$

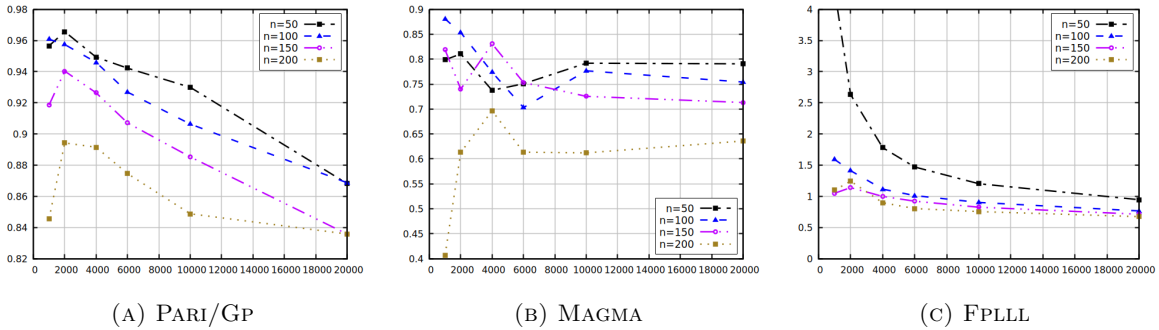


FIGURE 3. Ratios of running times of SpecLLL to LLL plotted against the precision  $l$ , for  $\sigma(K) \not\subset \mathbb{R}$  and several dimensions  $n$

precision. However, it becomes more efficient as precision increases and offers speed-ups between 25 and 50% for  $l = 20000$ , depending on the dimension.

Moreover, as expected the speed-up given by SpecLLL is larger when  $\sigma(K) \subset \mathbb{R}$ , however one still gains up to 20% when  $\sigma(K) \not\subset \mathbb{R}$ .

Note that the gain really depends on the version of LLL used. However, a better gain does not imply that the corresponding software should be used. Indeed, one should also take into account the initial

running time of LLL. For example, FPLLL is more efficient than the two other implementation. We refer the reader to the timings plots presented in [Appendix B.1](#), where are shown the corresponding timings.

Finally we verified that our modified algorithm really takes advantage of the special shape of the initial matrix. To do so, we launched SpecLLL over matrices of the shape

$$\begin{bmatrix} \beta_1 & C & 0 & \dots & 0 \\ \beta_2 & 0 & C & \ddots & \vdots \\ \vdots & \vdots & \ddots & \ddots & 0 \\ \beta_n & 0 & \dots & 0 & C \end{bmatrix},$$

where  $\beta_i$  are random integers with the same size as the entries of  $B(\mathcal{B}, \sigma, l)$ , and computed the same timings ratios. Results can be found in [Table 6](#), and one can remark that it is always close to 1, sometimes larger.

**5.2. Heuristic precision evaluation.** The second observation concern the precision needed to retrieve the roots. First we study which precision is required to retrieve an element depending on its norm, then we present an improvement concerning the evaluation of the norm of the roots.

**5.2.1. Precision from the norm.** The precision given by [Equation \(3.10\)](#) allows us to certify the correctness and the polynomial complexity of the method we described. However, it is possible to find a smaller value which is experimentally sufficient to retrieve an element.

Fix  $K$  a number field given by an irreducible polynomial  $P_K(X) \in \mathbb{Z}[X]$ . Given  $\mathcal{B} = (b_1, \dots, b_n)$  a  $\mathbb{Q}$ -basis of  $K$ , we generated random elements  $x \in \mathbb{Z}[\mathcal{B}]$  such that for all  $i \in \llbracket 1, n \rrbracket, x_i \in \llbracket -2^s, 2^s \rrbracket$  for  $s \in \{1, 25, 50, 75, 100\}$ . Then we computed the quotient  $q_t$  of the precision required to retrieve the coefficients of  $x$  by  $l_t = [K : \mathbb{Q}] \log_2 \|x\|_2$ .

We chose to test the experimental precision against  $l_t$  for several reasons. First, when  $K$  and  $\mathcal{B}$  are fixed, it is the term of [Equation 3.10](#) which is asymptotically relevant. Moreover the algebraic method of [\[3\]](#) requires the norm  $N_{K/\mathbb{Q}}(\mathfrak{p}^k)$  to be greater than a value which is essentially  $l_t$ , so that the decoding is certified. Finally, this value was also suggested by experiments when we first used this method to compute cube roots in multivariate fields [\[25\]](#). When comparing this value to [Equation \(3.10\)](#), we can remark that we gain essentially  $n(n - \log_2(n))$ .

Experimentally, the quotient  $q_t$  seems to be influenced by the size of the elements to decode and the dimension. More precisely, the maximum of the retrieved quotients in our experiments approaches 1 from above when the precision is increasing. Moreover, it seems to increase with the dimension. With these requirements in mind, we observed that it is always smaller than  $1 + \ln(n) \ln \ln(n) / (s \ln(2) + \ln(n)/2)$  – see [Figure 16](#) in appendix. Remark that the quotient in the previous value is the norm of an element with all coefficients equal to  $2^s$ .

If  $\text{Prec}$  is the precision required to retrieve the coefficients of an element  $x$ , it seems reasonable that

$$(5.1) \quad \text{Prec}(x) \leq \left(1 + \frac{\ln(n) \ln \ln(n)}{s + \ln(n)/2}\right) n \ln \|x\|_2 \leq n \ln \|x\|_2 + n \ln(n) \ln \ln(n).$$

**5.2.2. Norm of the roots.** Fix the factorisation of  $f(T)$  over  $K$  as  $f(T) = g(T)h(T)$  such  $Z_K(g) = Z_K(f)$ ,  $Z_K(h) = \emptyset$ . We can follow [\[17, 3\]](#) to bound  $\|x\|_2^2$  for  $x \in Z_K(f)$  given  $f(T)$ :

- (1) find  $B_{f, T_2}$  such  $\|\sigma_{K/\mathbb{Q}}(x)\|_2^2 \leq B_{f, T_2}$  for all  $x \in Z_K(f)$ ;
- (2) compute  $B_{\sigma_K^{-1}}$  the matrix norm of  $\sigma_K^{-1}$ , expressed relatively to the canonical basis of  $\sigma_{K/\mathbb{Q}}(K)$  and  $\mathcal{B}$ ;

(3) the value  $B_f = B_{f, T_2} B_{\sigma_K}^2$  satisfies the desired property.

This bound can be quite large compared to  $M(f) := \sup\{\|x\|_2^2 \mid x \in Z_K(f)\}$ . Without extra-information this is the best one can do. However, one can use heuristic evaluations for  $\|x\|_2^2$ , giving smaller results than  $B_f$ . Consider  $\sigma \in \text{Hom}(K, \mathbb{C})$  and  $x \in K$ . Then one has

$$|\sigma(x)|^2 = \left| \sum_{i=1}^n x_i \sigma(b_i) \right|^2 \leq \left( \sum_{i=1}^n |x_i \sigma(b_i)| \right)^2 \leq \|x\|_2^2 \|\sigma(\mathcal{B})\|_2^2.$$

Therefore for any  $x \in K$  and any  $\sigma \in \text{Hom}(K, \mathbb{C})$ ,  $\|x\|_2^2 \geq \frac{|\sigma(x)|^2}{\|\sigma(\mathcal{B})\|_2^2}$ . We define the function `HeuristicNorm` as the maximum of such quotients, as follows. For any  $\sigma \in \text{Hom}(K, \mathbb{C})$ , let  $B_{f, \sigma}$  be such that

$$(5.2) \quad \forall x \in Z_K(f), |\sigma(x)|^2 \leq B_{f, \sigma}.$$

Then we fix

$$(5.3) \quad \text{HeuristicNorm}(f) := \max \left\{ \frac{B_{f, \sigma}}{\|\sigma(\mathcal{B})\|_2^2} \mid \sigma \in \text{Hom}(K, \mathbb{C}) \right\}.$$

Even if `HeuristicNorm` gives a value which is a lower bound on  $\|x\|_2^2$  instead of an upper bound, it allows a first precision to be obtained. This evaluation is usually smaller than  $B_f$  and gives a good starting point, i.e. we do not need to increase much the precision to retrieve at least *some roots*.

We evaluated the generic difference between the proper maximal norm of  $x \in Z_K(f)$  and the value given by `HeuristicNorm`, and plotted the results in [Figure 4](#). These experimental values tend to show

$$(5.4) \quad \ln M(f) \leq \text{HeuristicNorm}(f) + \ln(n)/2.$$

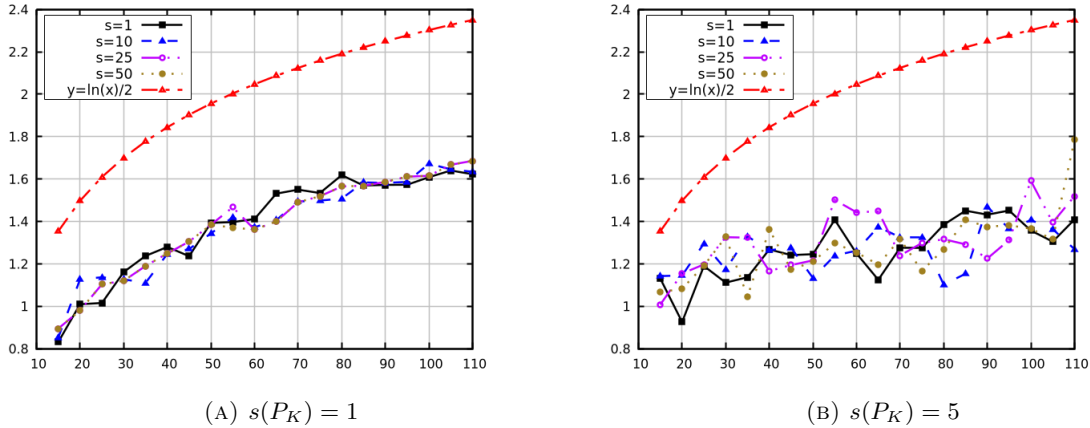


FIGURE 4. Logarithms of maximum ratios of  $M(f)$  by `HeuristicNorm`( $f$ ) plotted against  $n = [K; \mathbb{Q}]$  for several  $s$ , and  $s(P_K) \in \{1, 5\}$

*Heuristic precision.* Finally, given  $K$  and  $f(T) \in K[T]$  as above, one can compute a conjectured precision `HeuristicPrecision`( $f$ ) combining Equations (5.1) and (5.4) as

$$(5.5) \quad \text{HeuristicPrecision} := [K : \mathbb{Q}] \log_2 \text{HeuristicNorm}(f) + n \ln(n)/2 + n \ln(n) \ln \ln(n).$$



---

**Algorithm 10** Babai’s Nearest Plane Algorithm with early abort– EarlyAbortNP
 

---

**Require:**  $t \in \mathbb{R}^n$ ,  $\mathcal{B} = (b_1, \dots, b_r)$  a basis of a lattice  $\mathcal{L}$ ,  $\tilde{\mathcal{B}}$  the GSO of  $\mathcal{B}$

**Ensure:**  $v \in \mathcal{L}$  a vector close to  $t$  and  $b$  a boolean expressing whether  $t$  has been rejected or accepted.

```

1:  $w \leftarrow t$ 
2: for  $i = r$  down to 1 do
3:    $\lambda = (w \mid \tilde{b}_i) / \|\tilde{b}_i\|_2^2$ 
4:   if  $|\{\lambda\}| > 1/4$  then
5:     return  $w$ , false
6:   else
7:      $w \leftarrow w - \lfloor \lambda \rfloor b_i$ 
8:   end if
9: end for
10: return  $w$ , true

```

---

**5.3. Solution testing through finite fields.** An important step – especially when using heuristic observations – is to verify that a candidate is indeed a solution (steps 8 of both [Algorithm 6](#) and [Algorithm 8](#)). When manipulating large elements, evaluating the polynomial expression  $f(x)$  can be costly. Thus we chose to use a probabilistic method, which is summed up as follows:

- choose prime integers  $p_1, \dots, p_r$
- check whether  $f(x) = 0 \pmod{p_i}$  for all  $i \in \llbracket 1, r \rrbracket$

If the  $p_i$  are uniformly drawn as 32-bits integers and  $f(x)$  is uniformly distributed in  $\mathbb{F}_{p_i}$ , then the probability  $x$  of obtaining a false positive is  $1/2^{32r}$ , which is rapidly small. This choice is particularly useful for our relative method, since it potentially requires to test a huge amount of elements.

**5.4. Early abort in decoding.** The final observation that we use to speed-up computations consists in improving the decoding step, especially when the tested complex number  $y$  does not correspond to any solution. Here is the rationale behind our technique. If the vector  $Y$  given to `NearestPlane` correspond to a solution, then it is expected to be close to a vector of the lattice  $\mathcal{L}_l$ . However, if the vector is not a solution vector, then it should be “far” from a lattice vector. It can even be expected to be uniformly distributed modulo  $\mathcal{L}_l$ . Given a vector  $t$  and  $\mathcal{B}$  a basis, if we denote  $|(t \mid \tilde{b}_i) / \|\tilde{b}_i\|_2^2|$  by  $\lambda_i$ , then the distance between  $v$  and `NearestPlane`( $t, \mathcal{B}$ ) is  $\sum_{i=1}^r |\{\lambda_i\}| \cdot \|\tilde{b}_i\|_2^2$ . The quantity  $\{\lambda_i\} \in [-1/2, 1/2]$ , thus we distinguish between two cases : if  $|\{\lambda_i\}| \leq 1/4$  we deem  $\lambda_i$  to be acceptable and we continue the algorithm, while if  $|\{\lambda_i\}| > 1/4$  we deem  $t$  to be rejected. This gives [Algorithm 10](#).

Additionally to this strategy, we will reject a vector of complex numbers during the search of [Algorithm 8](#) as soon as one of the decoding fails.

We checked the experimental maximum number of steps done in [Algorithm 10](#) over bad vectors  $t$  of the form  $[z, 0, \dots, 0]$ , when  $z \in Z_{\mathbb{C}}(f)$  in the run of [Section 5.5](#), for several root sizes  $s_Z = \max_{x \in Z_K(f)} s(x)$ , field dimension  $[K : \mathbb{Q}]$  and coefficient sizes of the defining polynomial  $s(P_K)$ . For fields dimension between 25 and 125 and various root sizes, the maximum that we find is 7. Over all parameters this maximum does not seem to vary much and is mainly in  $\llbracket 4, 7 \rrbracket$ , see [Table 1](#) for example.

**Algorithm 11** AbsoluteRootsHeur**Require:** A number field  $K$ ,  $f(T) \in K[T]$ ,  $\mathcal{B}$  a  $\mathbb{Q}$ -basis of  $K$  such that  $Z_K(f) \in \mathbb{Z}[\mathcal{B}]$ .**Ensure:** The set  $Z_K(f)$ 

```

1:  $\sigma \leftarrow \text{ChooseEmbedding}(K)$ 
2:  $l \leftarrow \text{HeuristicPrecision}(f(T))$  ▷ Heuristic norm evaluation and precision formula
3:  $L_l \leftarrow \text{LLL}(B(\mathcal{B}, \sigma, l))$  ▷ Use of SpecLLL when  $\mathcal{B}$  is a power-basis
4:  $Z \leftarrow \text{FloatPolynomialRoots}(f^\sigma, l)$ 
5:  $S \leftarrow \emptyset$ 
6: for  $z \in Z$  do
7:    $y, b \leftarrow \text{TestDecodeHeur}(L_l, z)$  ▷ Early abort nearest plane
8:   if  $b$  and  $f(y) = 0$  then ▷ Verification through finite fields
9:      $S \leftarrow S \cup \{y\}$ 
10:  end if
11: end for
12: return  $S$ 

```

Dimension		25	50	75	100	125
$s_Z = 1$	min	1	1	1	1	1
	max	6	7	5	6	6
	average	2.7	2.4	2.5	2.2	2.8
$s_Z = 50$	min	1	2	1	1	1
	max	4	7	5	4	6
	average	2.5	2.8	2.7	2.3	2.8

TABLE 1. Number of steps before early aborts, for random targets in random number fields with  $s(P_K) = 10$ .

One can remark that the average value is always smaller than 3.

**5.5. Heuristic algorithms.** All the observations and strategies mentioned above lead to heuristic versions of the algorithms computing polynomial roots. First is the heuristic version of the decoding method `TestDecode` using `EarlyAbortNP`. We will design it by `TestDecodeHeur`. Using this version of the decoding method, together with the heuristic precision evaluation given by Equation (5.5) leads to a heuristic version of Algorithm 6 described in Algorithm 11.

Finally, regarding the relative method, we get a heuristic variant of Algorithm 7 and Algorithm 8. Remark that in Algorithm 12 (thus in Algorithm 13 as well) the use of early abort for the decoding of each coordinate respectively to  $L/K$  (step 8 of Algorithm 12) leads generally to decode only one of them when the vector of complex numbers  $X$  is not of the form  $\sigma_{L/K}(x)$  (as a matter of fact, the decoding will typically abort only after a few steps).

**5.5.1. Trade-off.** Now let us express some broad trade-off between our heuristic absolute and relative methods, again considering only computations linked to lattice reduction and decoding. Compared to the previous analysis, we need to take into account the influence of early abort. This still gives  $O(n_L^4 l + n_L^3 l + dn_L^2)$  in the worst-case for Algorithm 11 while these operations amount to  $O(n_K^4 l + n_K^3 l +$

**Algorithm 12** EarlyAbortMink2Coeff

**Require:** An extension  $L/K$ , given by  $\mathcal{B}$  a basis of  $K$  and  $\mathcal{E}$  a  $K$ -basis of  $L$ , an integer  $l$ , the matrix  $L_l$  from a reduced basis of  $\mathcal{L}(\mathcal{B}, \tau, l)$ ,  $M = \Sigma_{L/K}^{-1}$  up to a precision  $l + s$ , and  $X = [\sigma_{L/K}(x)]_{l+s}$  for some  $x \in L/K$

**Ensure:** A candidate  $y = (y_1, \dots, y_N)$  for the vector of coefficients of  $x$  expressed in  $\mathcal{B} \otimes \mathcal{E}$

```

1:  $Y \leftarrow XM$ 
2:  $y = [ \ ]$  ▷ void vector
3: for  $i = 1$  to  $n$  do
4:    $v, b \leftarrow \text{TestDecodeHeur}(L_l, Y_i)$ 
5:   if  $b = \text{false}$  then
6:     return  $X, b$ 
7:   end if
8:    $y \leftarrow [y \mid v]$ 
9: end for
10: return  $y, b$ 

```

**Algorithm 13** RelativeRootsHeur

**Require:** An extension  $L/K$ , given by  $\mathcal{B}$  a basis of  $K$  and  $\mathcal{E}$  a  $K$ -basis of  $L$ , and  $f(T) \in L[X]$  such that  $Z_L(f) \in \mathbb{Z}[\mathcal{E} \otimes \mathcal{B}]$

**Ensure:**  $Z_L(f)$

```

1:  $l \leftarrow \text{HeuristicPrecision}(f(T))$ 
2:  $L_l \leftarrow \text{LLL}(B(\mathcal{B}, \tau, l))$  ▷ Use of SpecLLL when  $\mathcal{B}$  is a power-basis
3:  $M \leftarrow \Sigma_{L/K}^{-1}$  ▷ Up to precision  $l + s$ 
4:  $Z \leftarrow \prod_{\sigma \in \text{Hom}_K(L, \mathbb{C})} \text{FloatPolynomialRoots}(f^\sigma, l + s)$ 
5:  $S \leftarrow \emptyset$ 
6: for  $z \in Z$  do
7:    $y, b \leftarrow \text{EarlyAbortMink2coeff}(z, M, L_l)$  ▷ Early abort
8:   if  $b$  and  $f(y) = 0$  then ▷ Verification through finite fields
9:      $S \leftarrow S \cup \{y\}$ 
10:     $\text{UpdateTree}(Z, z)$ 
11:  end if
12: end for
13: return  $S$ 

```

$(n_{L/K}d + d^{n_{L/K}} - d)n_K^2$ ) for Algorithm 13. This gives us the following condition for a trade-off :

$$(5.6) \quad d^{n_{L/K}} + dn_{L/K}n_K^2 - dn_L^2 - d = n_K^3 l (n_{L/K}^3 (n_L + 1) - (n_K + 1)).$$

When  $n_{L/K} = 2$  for example, this becomes  $d^2 - 2dn_K^2 - d = n_K^3 l (15n_K + 7)$ . Thus following the analysis for the naive versions of our algorithm we obtain that the relative version with early abort is advantageous compared to the absolute one for  $d$  up to

$$\sqrt{2}n_K^2 \sqrt{(n_K + 2 \log_2 \|x\|_2)(15n_K + 7)}$$

which is better than the naive trade-off by a factor of  $O(n_K)$ .

## 6. EXPERIMENTAL RESULTS

We compare in these section the practical performances of the methods described previously with the generic algebraic methods implemented in PARI/GP [31], which is the function `nroots`. Our implementation is (mainly) in GP, and is publicly available.<sup>2</sup>

Recall that we assume that  $Z_K(f) \subset \mathbb{Z}[\mathcal{B}]$ , with  $\mathcal{B} = (b_1, \dots, b_n)$  being some  $\mathbb{Q}$ -basis of  $K$ . Given  $x \in K$  we will keep denoting by  $x_1, \dots, x_n$  its coefficients relative to  $\mathcal{B}$ . In almost all our experiments, we chose  $\mathcal{B}$  to be  $\mathbb{Z}[\theta]$  with  $\theta = X \bmod P_K(X)$ , where  $P_K(X)$  is a fixed polynomial defining the field  $K$ .

Number fields are drawn randomly through the choice of their defining polynomial following a given distribution. When we consider general number fields, meaning not from a special family such as cyclotomic number fields,  $P_K(X)$  will be drawn with random coefficients in a given interval of the form  $\llbracket -2^s, 2^s \rrbracket$  with  $s \in \mathbb{N}$ . This somewhat arbitrary choice has been used in the literature [4] and allows us to study the average behaviour of our algorithms, for a well defined notion of “average”. Finally, we consider two specific classes of number fields, namely *cyclotomic fields* in Section 6.3.3 and *real Kummer fields* in Section 6.3.4, because they showcase nicely the good and bad behaviours of our algorithms (especially our relative method) and PARI/GP `nroots`.

**Remark 6.1.** We checked the number of roots retrieved by our algorithms, and all roots were retrieved by both methods (absolute and relative).

**6.1. Instability of `nroots`.** We observed experimentally that over generic number fields, `nroots` is not stable, and that fields could be divided into two groups. We call *bad fields* those number fields for which `nroots` behave poorly, and *good fields* the rest of them. In order to obtain better observations, we split them by minimising the sum of their medians, that we denote by  $m_1$  and  $m_2$  respectively. To showcase this unstable behaviour we give some statistical data regarding the running time of `nroots`, `AbsoluteRoots` and `AbsoluteRoots` in Table 2.

---

<sup>2</sup>[https://github.com/AndLesav/nf\\_polynomial\\_roots](https://github.com/AndLesav/nf_polynomial_roots)

TABLE 2. Statistical data regarding the running time of nroots, AbsoluteRoots and AbsoluteRoots over random number fields such that  $s(P_K) = 5$ .

(A) nroots				
Dimension $[K : \mathbb{Q}]$	25	50	75	100
# bad fields (%)	49	13	10	12
$m_1$ (s)	2.37	20.41	55.51	125.92
$m_2$ (s)	3.10	174.00	2189.93	16081.07
average (s)	2.76	40.08	269.67	1989.31

(B) AbsoluteRoots				
Dimension $[K : \mathbb{Q}]$	25	50	75	100
# bad fields (%)	65	77	68	67
$m_1$ (s)	1.59	39.53	318.24	1637.11
$m_2$ (s)	2.69	66.44	539.42	2646.72
average (s)	2.31	60.44	467.41	2285.76

(C) AbsoluteRootsHeur				
Dimension $[K : \mathbb{Q}]$	25	50	75	100
# bad fields (%)	57	71	72	61
$m_1$ (s)	0.42	3.86	15.95	54.84
$m_2$ (s)	0.61	5.31	21.72	67.19
average (s)	0.53	4.88	19.98	62.33

In Table 2 one can see that there are about 10% of fields for which the running time of nroots explodes. The corresponding median  $m_2$  is up to 127 times larger than the median  $m_1$  corresponding to the rest of the fields. Algorithm 6 and Algorithm 13 are more stable. For these algorithms  $m_2$  never exceeds  $2m_1$ . The instability of nroots has the following consequence. For a majority of number fields, nroots clearly outperforms AbsoluteRoots (it is more than 10 times faster for  $[K : \mathbb{Q}] = 100$ ) and this difference is growing with the dimension. However, *in average*, the timings obtained for both algorithms are similar.

*Potential explanation.* The determining parameter seems to be the possibility of retrieving an inert prime. Following the search for good primes implemented in PARI/GP, we evaluated the proportion of fields for which an inert prime will be found, see Table 3.

Dimension $[K : \mathbb{Q}]$	23	51	75	83
# bad fields (%)	5	7	17	10

TABLE 3. Proportion of fields for which no inert prime has been found

These proportions tend to follow the ones found previously, such as presented in Table 2. Additionally we recorded the running times over the two kind of fields, as showcased in Table 3.

TABLE 4. Running time of `nroots`, `AbsoluteRoots` and `AbsoluteRootsHeur` over random number fields for which inert primes can be found or not.(A) `nroots`

Dimension $[K : \mathbb{Q}]$	25	50	75
Good fields (s)	3.1	17.8	55.5
Bad fields (s)	4.9	164.1	2177.9

(B) `AbsoluteRoots`

Dimension $[K : \mathbb{Q}]$	25	50	75
Good fields (s)	2.3	52.5	437.1
Bad fields (s)	2.3	52.2	403.1

(C) `AbsoluteRootsHeur`

Dimension $[K : \mathbb{Q}]$	25	50	75
Good fields (s)	0.5	4.7	21.7
Bad fields (s)	0.5	4.7	20.7

The results presented in [Tables 2](#) and [3](#) are very similar, which confirms the fact that the determining parameter regarding the instability of `nroots` is the use of inert primes.

**Remark 6.2.** In the following, for `nroots`, we will systematically present the average value of the execution times over all fields, over “good” fields and over “bad” fields, in order to have a more complete picture of its behaviour. We will write  $m_1$  and  $m_2$  the two last quantities.

**6.2. Absolute method.** We study the impact of the different parameters of the problems which are the size of the polynomial  $P_K(X)$ , the number of roots  $|Z_K(f)|$ , and the size of the roots  $\log_2 \|x\|_2$ . We also differentiated between number fields  $K$  such that  $r_1 > 0$  or such that  $r_1 = 0$ , since the matrices used to decode do not have exactly the same sizes. However, since the results are very similar we only present here the data for  $r_1 > 0$ . Data when  $\sigma(K) \subset \mathbb{C} \setminus \mathbb{R}$  can be found [Appendix B.3](#).

**6.2.1. Choice of  $P_K(X)$ .** We study the impact of two parameters linked to  $P_K(X)$ , namely  $\deg P_K(X)$  which is the dimension of  $K$ , and the size of its coefficients. In our experiments, we fixed the parameters of the problem linked to  $f(T) \in K[T]$ . More precisely we considered  $f(T)$  such that:

- $\deg f(T) = 50$ ;
- $f(T)$  splits in  $K$ , i.e.  $|Z_K(f)| = \deg f(T)$ ;
- $\forall x \in Z_K(f), \log_2 \|x\|_\infty \leq 10$ .

Then we considered  $P_K(X) = p_0 + p_1X + \dots + p_{n-1}X^{n-1} + X^n \in \mathbb{Z}[X]$  for increasing degrees  $n$ , and several coefficient sizes  $s(P_K) = \log_2 \|P_K\|_\infty$ . More precisely for sizes  $s(P_K) \in \{1, 10\}$ , we picked polynomials  $P_K(X)$  such that  $\forall i \in \llbracket 0, n-1 \rrbracket, p_i \in \llbracket -2^{s(P_K)}, 2^{s(P_K)} \rrbracket$ , and this for  $n$  increasing. The data obtained are shown in [Figures 5](#) and [18](#).

One can remark that the time efficiency of all three methods are widely influenced by the dimension  $[K : \mathbb{Q}]$ . It is easily explained by the fact that all three methods require the computation of a LLL-reduced basis of a lattice with rank equal to  $[K : \mathbb{Q}]$ . Moreover the volume of said lattice depends also on the

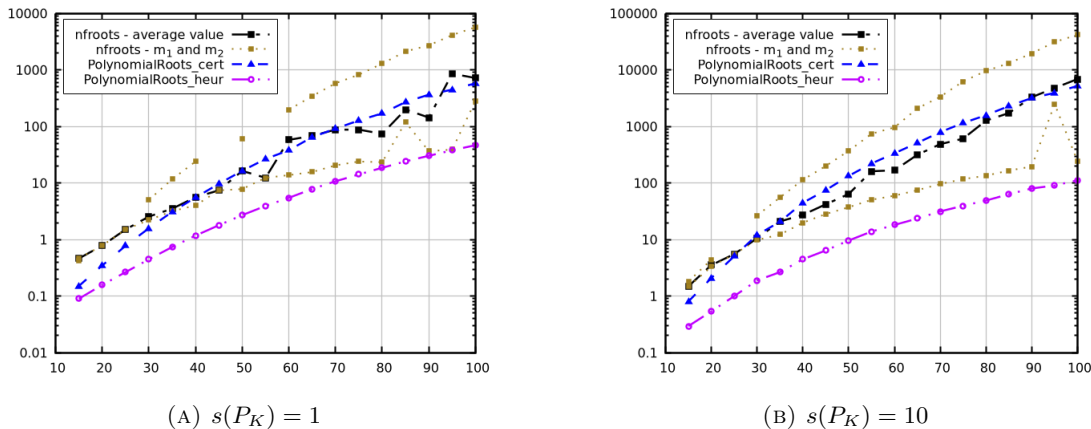


FIGURE 5. Average timings (s) of nroots, AbsoluteRoots and AbsoluteRootsHeur plotted against  $\deg P_K(X)$  for randomly generated  $P_K(X)$  such that  $r_1 > 0$ , with  $s(P_K) \in \{1, 10\}$

dimension. Remark that as observed in Table 2, the difference between  $m_1$  and  $m_2$  widens when  $[K : \mathbb{Q}]$  increases.

The parameter  $s(P_K)$ , i.e. the coefficient size of the defining polynomial of  $K$ , also influences the performances of all three algorithms. Our heuristic method seems to be slightly less impacted by this parameter.

One can see from Figure 5 and Figure 18 that AbsoluteRoots is less efficient than nroots, at least for the fixed shape of  $f(T)$ . The heuristic method AbsoluteRootsHeur – using heuristic norm evaluation and formula to compute the precision – is way more efficient than the certified version. It is also more efficient than nroots *on average*, and compete with its best running times in a number of cases. Our method seems to be more influenced by the dimension of  $K$  than the algorithm of PARI/GP, at least when the latter runs at its best.

6.2.2. *Size of the roots.* We now study how the size of the elements of  $Z_K(f)$  impacts the performance of the different algorithms. To this end we fixed the parameters linked to  $P_K(X)$  and  $f(T)$ :

- $\deg f(T) = 50$ ;
- $f(T)$  splits in  $K$ , i.e.  $|Z_K(f)| = \deg f(T)$ ;
- $s(P_K) = 1$ ;
- $\deg P_K(X) \in \{25, 50\}$ .

We did experiments for increasing size of roots. Let us denote by  $s_Z$  this size, i.e.  $\forall x \in Z_K(f), \log_2 |x_i| \in \llbracket 0, s_Z \rrbracket$ . The results can be found in Figure 19.

One can see that our algorithm is more influenced by  $s_Z$  than nroot. This is certainly due to the fact that the lattices reduced for decoding during the algorithm described in [3] are ideals, and more precisely *prime powers ideals*. These are usually easier to reduce, especially if one can use a pre-reduction such as mentioned by Belabas, or the documentation of PARI/GP [31]. Remark that the difference of running time for nroots between bad and good fields seems constant here.

6.2.3. *Partial conclusion.* From the different situations explored and the data gathered, we can conclude that the certified version of our method AbsoluteRoots is in general less efficient than the algebraic method implemented in PARI/GP. However, their running time are similar *in average* in some situations (see

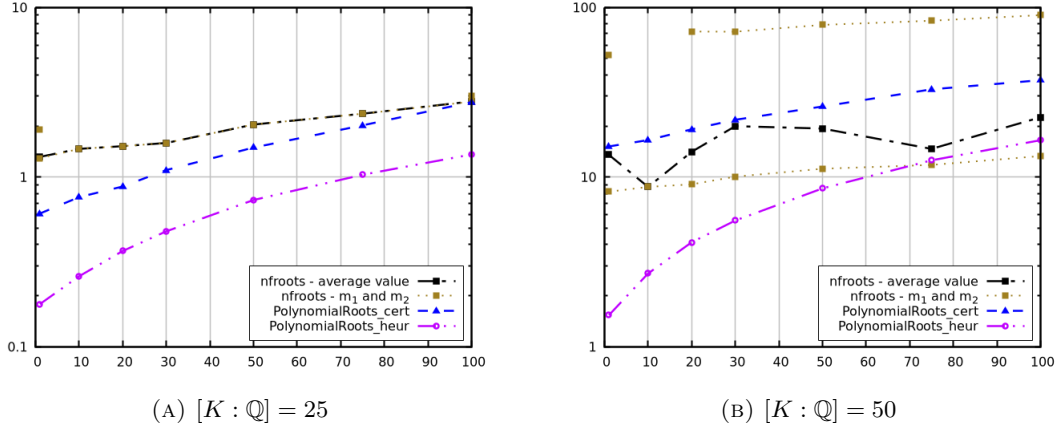


FIGURE 6. Average timings (s) of nroots, AbsoluteRoots and AbsoluteRootsHeur plotted against  $s_Z$  for randomly generated  $P_K(X)$  such that  $[K : \mathbb{Q}] = 25$  and  $[K : \mathbb{Q}] = 50$ , with  $r_1 > 0$

for instance Figure 5). The heuristic version AbsoluteRootsHeur (Alg. 11) offers a clear advantage and is more efficient than nroot, still in average. When restricted to the “good” fields for nroot, our heuristic version is still competitive for relatively small roots (crossing point at  $s_Z \sim 75$  for  $[K : \mathbb{Q}] = 50$ ).

Regarding the impact of the different parameters, our method behaves less nicely with respect to the dimension and the size of the roots. This is certainly due to the nature of the lattices to be reduced in the two different methods. The algebraic method takes fully advantage of the nature of the lattice, which is a power of a prime ideal, thus mitigating the influence of the size of the vectors and the dimension. In the best cases, it uses a pre-reduction algorithm which hasten considerably the subsequent LLL reduction [3].

**6.3. Relative extensions.** Let us now consider relative extensions  $L/K$ , and study the efficiency of Algorithm 13. First we look into the impact of our heuristic strategies. Then we compare this method to Algorithm 11. Finally, we study both algorithms together with nroots over cyclotomic fields and Kummer extensions.

Let us fix the notations. We will consider  $L/K$  together with  $P_K(X) \in \mathbb{Q}[X]$  and  $P_L(Y) \in K[Y]$  such that  $K \cong \frac{\mathbb{Q}[X]}{(P_K(X))}$  and  $L \cong \frac{K[Y]}{(P_L(Y))}$ .

**6.3.1. Impact of the heuristic strategies.** We considered several versions of the relative method : the certified one as Algorithm 8, the heuristic one as Algorithm 13, and two intermediate versions, where either the precision or the search is heuristic. We studied the impact of (a) the degree of the equation  $f(T)$ , and (b) the size  $s_Z$  of the roots. We considered extensions  $L/K$  such that  $[K : \mathbb{Q}] = 30$ ,  $[L : K] = 3$ ,  $s(P_K), s(P_L) \leq 1$ , and  $f(T)$  splits in  $K$ . Additionnally, we fixed  $s_Z = 10$  when studying  $\deg f(T)$  and  $\deg f(T) = 25$  when studying the other parameter. We focused on small parameters for these experiments, as we wished to illustrate the impact of the heuristic modifications that we described. The timings obtained can be found in Figure 7.

One can remark that both heuristic observations speed-up the computations, especially when the degree of  $f$  is increasing, as shown in Figure 7a. This leads to a speed-up up to a factor 100 for the full heuristic implementation compared to the “naive” one. Moreover from both experiments, one can



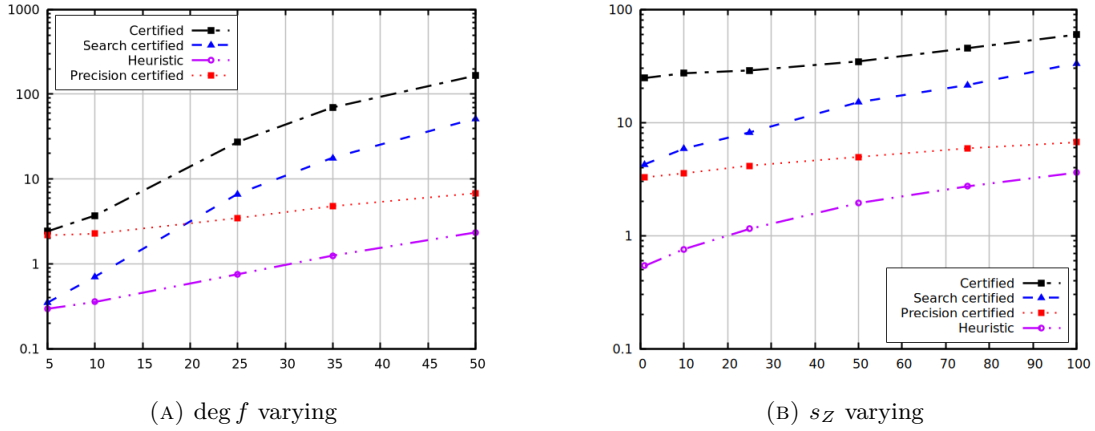


FIGURE 7. Timings (s) of certified and heuristic versions of RelativeRoots plotted against  $\deg f(T)$  (resp.  $s_Z$ ) for randomly generated  $P_K(X), P_L(Y)$

conclude that the impact of the heuristic search (using Algorithm 12) is more important than the one of the heuristic precision – at least for the small range of parameters considered.

6.3.2. *Generic number fields.* In this section, we study the performances of our most efficient procedures, i.e. the heuristic algorithms AbsoluteRootsHeur and RelativeRoots, over “generic” number fields. Here, the term “generic” refers to fields whose defining polynomials have coefficients uniformly drawn in segments of the form  $[-2^s, 2^s]$ . For practicability reasons, our experiments mainly deal with  $s = 1$ .

*Number of roots.* We first study the influence of the number of roots, as it impacts the running time of RelativeRootsHeur. We considered extensions  $L/K$  such that  $[L : \mathbb{Q}] = 90$  and  $[L : K] \in \{2, 3\}$  with:

- $s(P_K), s(P_L) \leq 1$ ;
- $\deg f(T) = 50$ ;
- $\forall x \in Z_K(f), s(x) \leq 10$ .

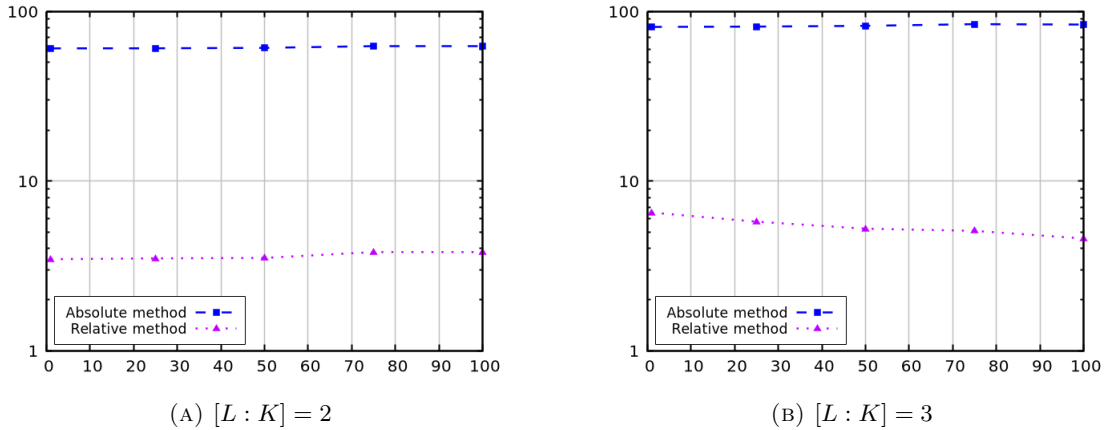


FIGURE 8. Average timings (s) of AbsoluteRoots and RelativeRoots plotted against  $q_f$  for randomly generated  $P_K(X), P_L(X)$  and  $f(T)$ , such that  $[L : \mathbb{Q}] = 90$  and  $\deg f(T) = 50$ .

The influence of the number of roots can be seen in Figure 8. The timings of the method `AbsoluteRoots` increase with the number of roots, while the ones for the relative algorithm `RelativeRoots` globally decrease, which is expected, see Appendix A. Indeed, with  $q_f$  increasing, roots are retrieved earlier and the search space is updated more quickly. One can remark on these first comparisons, that Algorithm 13 can be between 5 and 30 times faster than Algorithm 11. Moreover, as expected, Algorithm 13 is slower when  $[L : K]$  is increasing. All these observations will be verified in other experiments.

*Degree of the extension.* Let us now consider the influence of the degree  $[L : K]$  of the extension considered. We already know that the running time of `RelativeRoots` depends exponentially on this parameter. We illustrate its impact on practical computations.

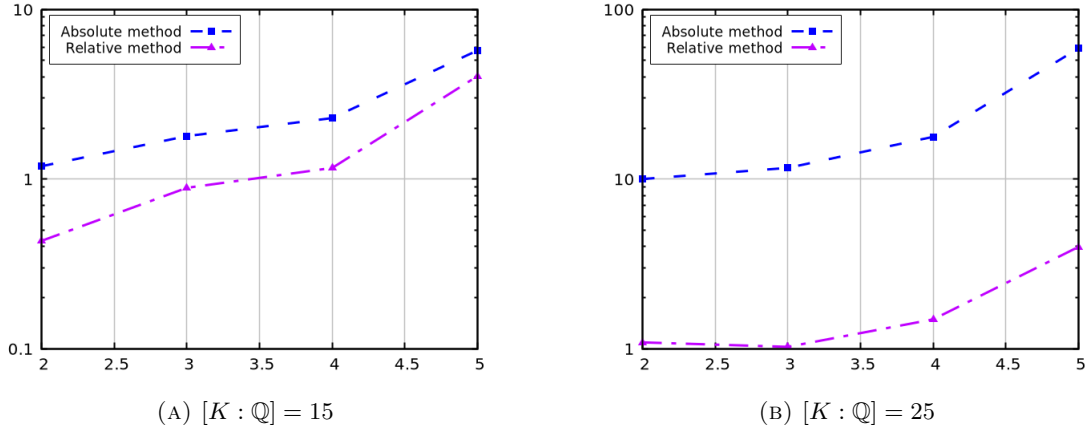


FIGURE 9. Average timings (s) of `AbsoluteRoots` and `RelativeRoots` plotted against  $[L : K]$  for randomly generated  $P_K(X)$ ,  $P_L(Y)$  and  $f(T)$ , such that  $[K : \mathbb{Q}] \in \{15, 25\}$  and  $\deg f(T) = 50$ .

As expected, the running time of Algorithm 13 is deeply impacted by an increase in the relative degree  $[L : K]$ . However, when compared to Algorithm 11, one can observe that this drawback is mitigated when considering increasing degrees  $[K : \mathbb{Q}]$ . This shows that for a constant  $[L : K]$  using our relative method should asymptotically outperform the absolute methods `AbsoluteRootsHeur`.

**6.3.3. Cyclotomic fields.** In this section we study the performances of `AbsoluteRootsHeur`, `RelativeRootsHeur` and `nroots` over cyclotomic fields. We will write  $K_m$  the cyclotomic field  $\mathbb{Q}(\zeta_m)$  with conductor  $m$ . We study these fields because they are widely used in applied mathematics such as cryptography. Moreover, since their generic defining polynomials have small norms, `nroots` is particularly efficient.

Additionally our relative method always apply be used over these fields. Indeed,  $K_m$  always has a totally real subfield  $K_m^+$  such that  $[K_m : K_m^+] = 2$ . Moreover, since  $K_m^+ = K_m^\tau$  where  $\tau$  is the complex conjugation, the search space has  $\deg f$  elements. We use this structure only in our first experiments. Results are gathered in Figure 10.

If we compare the timings obtained to previous data such as presented in Figure 5 or Figure 18, both `AbsoluteRootsHeur` and `nroots` seem more efficient than over randomly generated number fields. It is particularly the case of the latter.

The PARI/GP `nroots` is more efficient than our absolute method `AbsoluteRootsHeur` over all fields. Our relative method seems to compete with the implementation of PARI/GP [31] with the set of parameters

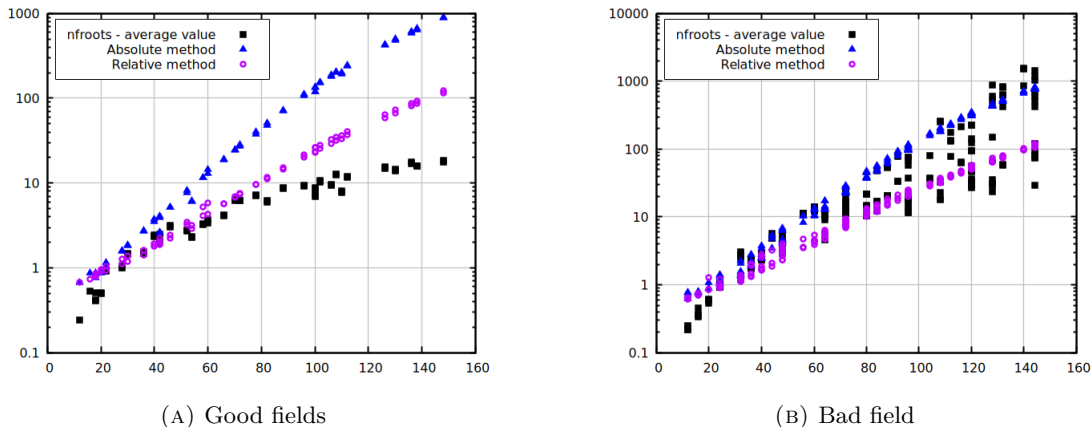


FIGURE 10. Average timings (s) of AbsoluteRootsHeur, RelativeRoots and nroots plotted against  $[K_m : \mathbb{Q}]$  for cyclotomic fields of conductor  $m$  such that  $10 \leq m \leq 500$  and  $[K_m : \mathbb{Q}] < 150$ , for  $s_Z = 50$ .

fixed in the experiments over “bad” fields, see Figure 10b. It is again more influenced by the dimension than nroots (in its best behaviour). Additionally since we know from Section 6.2 that the methods using complex embeddings have a worse behaviour with respect to the size of the roots, it seems that using only  $K_m^+$  is not enough to beat the implementation of PARI/GP over cyclotomic fields.

Note however that most of these fields do have other subfields, which can be themselves cyclotomic fields. For example, if  $m = p^r$  with  $p$  a prime integer, then writing  $m' = p^{r'}$  such that  $r' < r$  we have that  $K_{m'}$  is a subfield of  $K_m$  such that  $[K_m : K_{m'}] = p^{r-r'}$ . Using the extension  $K_m/K_{m'}$  would lead to a search space with  $d^{p^{r-r'}}$  elements. We ran some experiments over such extensions for  $p \in \{2, 3\}$ , data are gathered in Table 5. Again we chose  $s_Z = 50$  and  $\deg f(T) = 50$  (resp.  $\deg f(T) = 55$ ) for  $p = 2$  (resp.  $p = 3$ ). We retrieve similar timings and asymptotically, the choices  $p = 2, r = 2$  and  $p = 3, r = 1$  should outperform nroots. Remark that as expected AbsoluteRoots is the least efficient algorithm.

6.3.4. *Kummer extensions.* Let us now focus on real Kummer extensions, meaning such that  $L = \mathbb{Q}(\sqrt[p]{m_1}, \dots, \sqrt[p]{m_r})$ , where  $p \geq 2$  is a prime integer and  $(m_1, \dots, m_r) \in \mathbb{Q}^r$  such that  $[K : \mathbb{Q}] = p^r$ . Fix  $K = \mathbb{Q}(\sqrt[p]{m_1}, \dots, \sqrt[p]{m_{r-1}})$ . Then  $L/K$  is an extension over which one can take full advantage of RelativeRoots. Indeed  $K$  can be embedded in  $\mathbb{R}$ , so that  $\text{Hom}_K(L, \mathbb{C})$  has one real embedding and  $p - 1$  complex ones. This gives a search space with  $\deg f(T)^{(p-1)/2}$  elements. We will compare AbsoluteRoots, RelativeRoots and nroots of PARI/GP [31]. We study these fields because real Kummer extensions are all “bad” fields for the  $p$ -adic method. Recall that is mentioned by [3] that ideal lattices are usually easy to reduce – especially when one can use a pre-reduction algorithm – which occurs when the inertial degree of said ideal is large. Moreover we saw in Section 6.1 that nroots behaved badly when the search for an inert prime was unfruitful. It turns out that over real Kummer fields of the form  $\mathbb{Q}(\sqrt[p]{m_1}, \dots, \sqrt[p]{m_r})$ , the inertial degree cannot be larger than the exponent  $p$  (see for instance [24]).

The first situation that we will explore will be the same as before, i.e.  $f(T)$  is split. It will show how computations can be accelerated in good situations, and the difference between good and bad fields for nroots. Then we will study the special case where  $f(T)$  has degree  $p$ . It is the direct generalisation of  $f(T) = T^p - \alpha^p$  with  $\alpha \in L$ , which is the type of equation that are to be solved in several tasks involving

TABLE 5. Running time of nroots, AbsoluteRoots and AbsoluteRoots over cyclotomic fields of conductor  $m = p^r$  with  $p$  a prime integer.

(A)  $p = 2$

$r$	5	6	7	8
nroots	0.4	2.4	11.6	30.3
AbsoluteRoots	0.7	1.4	13.0	436.8
RelativeRoots, $[L : K] = 2$	1.4	2.0	3.8	27.9
RelativeRoots, $[L : K] = 4$	19.6	22.4	29.2	38.0

(B)  $p = 3$

$r$	3	4	5
nroots	0.5	2.8	29.7
AbsoluteRoots	0.9	6.8	1334.0
RelativeRoots, $[L : K] = 3$	3.9	7.1	33.1

(S-)units or class group [8, 17, 35]. In particular, this task arose in several practical works these past few years [2, 25, 9, 7, 10].

*Split equation.* We considered split polynomials  $f(T)$  of degree 50 for increasing size of roots  $s_Z$ , over Kummer fields of the form  $L = \mathbb{Q}(\sqrt[p]{m_1}, \dots, \sqrt[p]{m_r})$  with  $p \in \{2, 5\}$ . Moreover  $[L : \mathbb{Q}] = 128$  if  $p = 2$ ,  $[L : \mathbb{Q}] = 125$  if  $p = 5$ , and each  $m_i$  is a prime number smaller than 40. The results are shown in Figure 11.

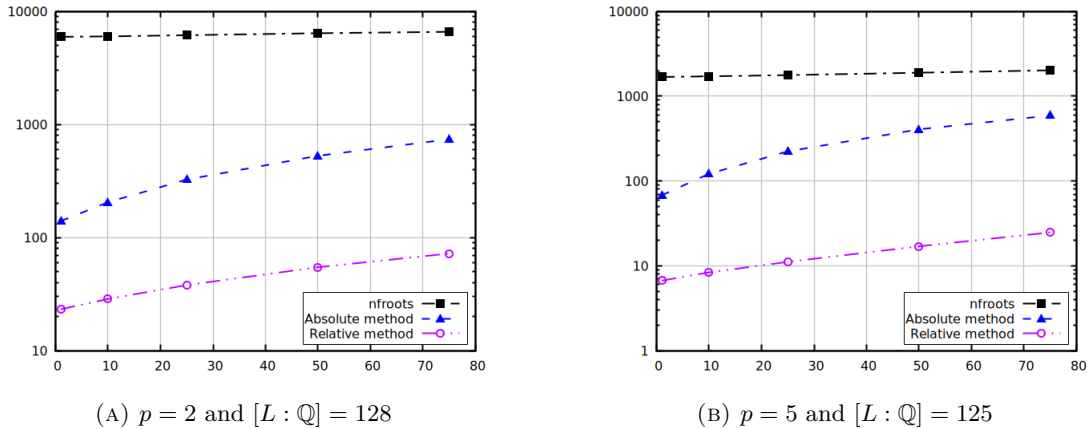


FIGURE 11. Average timings (s) of AbsoluteRootsHeur, RelativeRoots and nroots plotted against  $s_Z$  for Kummer fields of exponent  $p \in \{2, 5\}$ .

One can observe that both Algorithm 11 and Algorithm 13 outperform nroots. Our relative method is particularly efficient, and is up to 100 times faster. This contrasts with cyclotomic fields. Indeed, compare timings reported in Figure 10 and Figure 11a for which the degree of the relative extension is 2.

*Small degree equations.* As mentioned we consider here polynomials  $f(T)$  with small degrees, namely  $\deg f(T) = p$  over real Kummer fields  $\mathbb{Q}(\sqrt[p]{m_1}, \dots, \sqrt[p]{m_r})$  of exponent  $p$ . For such degrees, the search

space of Algorithm 13 is small enough to take full advantage of this method. Indeed, since the complex embeddings in  $\text{Hom}_K(L, \mathbb{C})$  are all conjugates (except one real embedding), the cardinality of the search space is  $p^{\frac{p+1}{2}}$ .

**Remark 6.3** (nroots). In this configuration, nroots does not follow the method described before. Indeed, when  $3 \deg f(T) < [L : \mathbb{Q}]$ , the implementation of PARI/GP uses Trager's method [42, 3] for factorising polynomials. We will therefore refer to it as Trager, to differentiate it from the  $p$ -adic method.

We considered Kummer fields of exponent  $p$  in  $\{2, 3, 5, 11\}$  and the different objects are drawn such that:

- each  $m_i$  is a prime number smaller than 40;
- $\deg f(T) = p$  and  $|Z_K(f)| = 1$ .

The data gathered can be found in Figure 12.

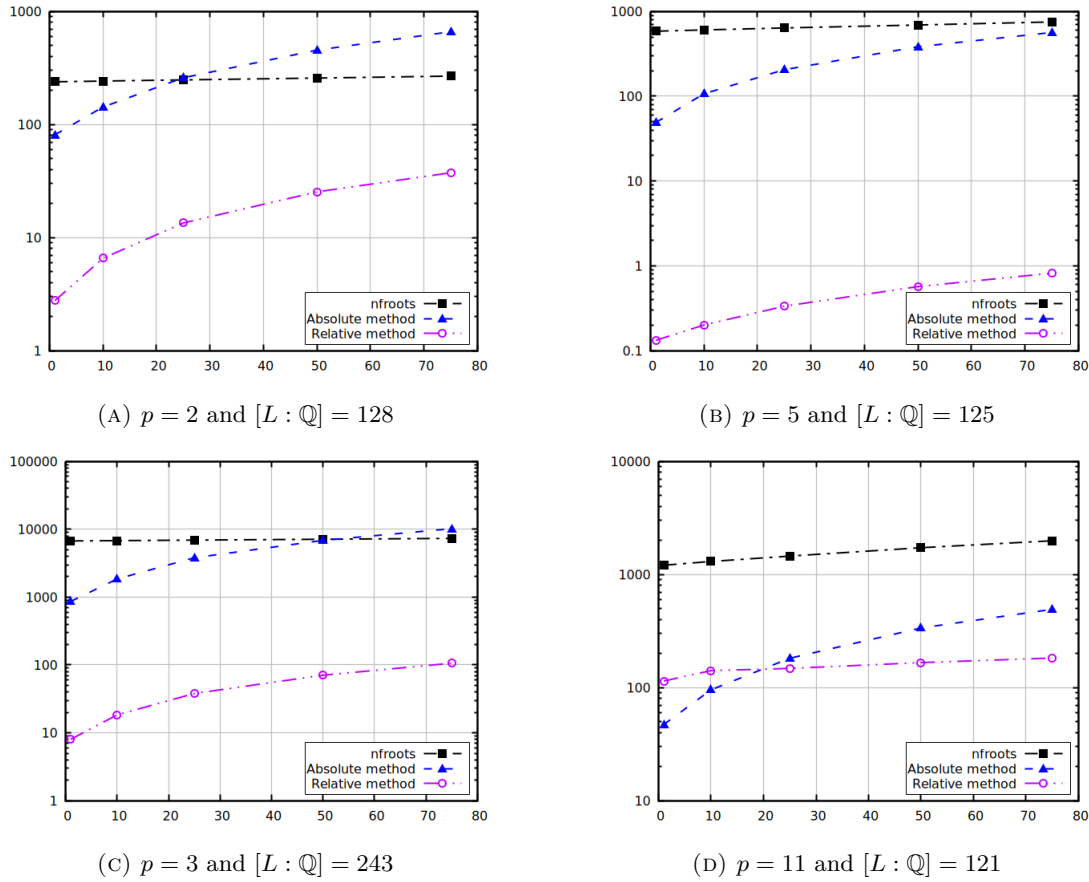


FIGURE 12. Average timings (s) of AbsoluteRoots and RelativeRoots and Trager plotted against  $s_Z$  for randomly generated Kummer fields  $L/K$  with exponent  $p \in \{2, 3, 5, 11\}$  and  $\deg f(T) = p$ .

In most cases our relative method is way more efficient than the two others. It can go up to 500 times faster for  $p = 3$  for example. The method Trager implemented in PARI/GP is always worse than both our algorithms when the roots are small, but it is stable when the size is increasing and it outperforms

Algorithm 11 for large roots and  $p \in \{2, 3, 5\}$ . Note however that a basis lattice  $\mathcal{L}_l$  can be reduced once and used to solve different equations. Thus, a batch strategy is possible with `AbsoluteRootsHeur`, which is not the case with `Trager`.

If one compares the data gathered in Figures 12a and 12b for  $p = 2$  and  $p = 5$  respectively – for which the degrees  $[L : \mathbb{Q}]$  and the size of the search space are similar – one can see that the timings for `RelativeRootsHeur` are around 10 times lower for  $p = 5$ . This is due to the fact that the dimension of the subfield  $K$  over which decodings are made is smaller in this case. These observations are confirmed with the timings gathered in Figure 12c.

Finally, one can remark in Figure 12d that the size of the search space is important, as the difference between `RelativeRootsHeur` and `AbsoluteRootsHeur` is less important for  $p = 11$ . However in this case as for the others, it seems that the size of the roots is less of a problem for the relative method than for `AbsoluteRootsHeur`.

6.3.5. *Small dimensions, small degree and large roots.* We consider a final situation to showcase the performances of the algorithms studied here, namely when  $f(T)$  has large roots and  $\deg f(T)$  is small as well as  $[K : \mathbb{Q}]$ . This situation is encountered in its extreme in the Number Field Sieve with  $\deg f(T) = 2$  and  $[K : \mathbb{Q}]$  usually smaller than 10. In our experiments, we considered  $\deg f(T) = 10$  and  $[K : \mathbb{Q}] \in \{12, 30\}$ . Data are gathered in Figure 13.

As already observed, `PARI/GP nroots` is asymptotically more efficient than the absolute method `AbsoluteRoots`. However `RelativeRoots` offers again a certain advantage compared to `nroots`, especially for larger dimensions, here  $[L : \mathbb{Q}] = 30$ . Note that `MAGMA function Roots` becomes quickly the least efficient method by a large factor.

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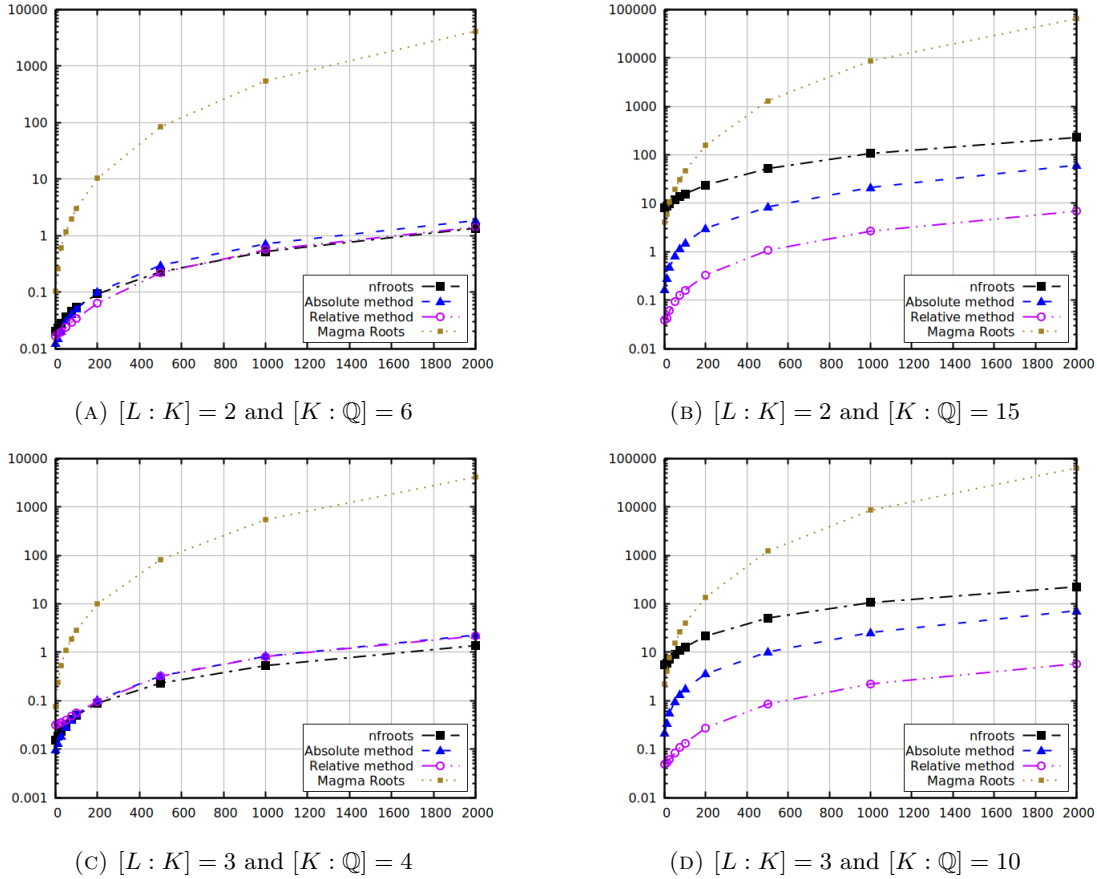


FIGURE 13. Average timings (s) of AbsoluteRoots and RelativeRoots, nroots and Magma Roots plotted against  $s_Z$  for randomly extensions  $L/K$  with  $[L : \mathbb{Q}] \in \{12, 30\}$  and  $\deg f(T) = 10$ .

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#### APPENDIX A. AVERAGE COST OF THE SEARCH

Let us study the average cost of the search phase of Algorithm 8 or Algorithm 13. In particular we will determine the average number of decodings that will occur before finding all roots. If one denotes by  $\sigma_1, \dots, \sigma_n$  the elements of  $\text{Hom}_K(L, \mathbb{C})$ , then the two algorithms can be described as a search without replacement in the set  $Z = Z_1 \times \dots \times Z_n$  where  $Z_i = Z_{\sigma_i}$ .

**Remark A.1.** A cartesian product  $S = S_1 \times \dots \times S_n$  will be ordered using the lexicographic order. This means that for all  $(x, y) \in S^2$ , if  $i(x, y) = \min\{i \in \llbracket 1, n \rrbracket \mid x_i \neq y_i\}$ , then we have  $x < y \iff x_{i(x,y)} < y_{i(x,y)}$ .

From now on, let us consider the sets  $Z_i$  as ordered sets of elements  $z_{i,j}$  with  $z_{i,j} < z_{i,j'} \iff j < j'$ . Moreover we consider that we run through  $Z$  following the lexicographic order. Assume that the state of the computation is at the state with index  $j = (j_1, \dots, j_n) \in \llbracket 1, \deg f(X) \rrbracket^n$ , and that we found a root  $x$ . We can write  $x = (z_{1,j_1}, \dots, z_{n,j_n})$ . Then the action of `UpdateTree` on  $Z$  is as follows. We mentioned that it removes  $z_{i,j_i}$  from  $Z$  for all  $i \in \llbracket 1, n \rrbracket$ . This amounts to removing  $z_{i,j_i}$  from  $Z$  for all  $i \in \llbracket 2, n \rrbracket$  and updating the index by doing  $j_1 \leftarrow j_1 + 1$  and for all  $i > 1, j_i \leftarrow 1$ .

**Definition A.2.** Let  $N, M, m$  be integers satisfying  $m \leq M \leq N$ . A random variable  $X$  taking non-negative values follows a *negative hypergeometric distribution with parameters*  $(N, M, m)$  if it satisfies

the following formula:

$$\mathbb{P}(X = k) = \frac{\binom{k+m-1}{k} \binom{N-m-k}{M-m}}{\binom{N}{m}}.$$

We will write  $X \sim NHG(N, M, m)$ .

The negative hypergeometric distribution describes exactly what we want to study. Indeed, it arises as follows. Consider a set of  $N$  elements, containing  $M$  “success elements” and  $N - m$  “fail elements”. Then if one draws uniformly in the set *without replacement* until  $m$  successes are found, then the number of failures drawn follows a negative hypergeometric distribution.

**Proposition A.3** ([20]). *Let  $N, M, m$  be integers, and  $X$  be a random variable such that  $X \sim NHG(N, M, m)$ . Then one has:*

$$\mathbb{E}[X] = m \frac{N - M}{M + 1} \quad \text{and} \quad \text{Var}[X] = m \frac{(N - M)(N + 1)}{(M + 1)(M + 2)} \left(1 - \frac{m}{M + 1}\right).$$

We are essentially interested in the average cost, so we focus on the expectation.

**Proposition A.4.** *Let  $L/K$  be an extension of number fields such that  $[L : K] = n$ . Consider  $f(X) \in L[X]$  such that  $|Z_K(f)| = s$  and write  $d = \deg f(X)$ . Then, assuming the search is naïve, the average number of “failed” decodings done in Algorithm 8 before finding one root is  $\frac{d^n + 1}{s + 1}$ . The average number of “failed” decodings before finding all the roots is  $s \frac{d^n + 1}{s + 1}$ .*

*Proof.* Let  $X_1$  be the random variable representing the number of failures before finding the first root, and  $X_s$  be the random variable of the number of failures before finding all the roots. Clearly one has  $X_1 \sim NGH(d^n, s, 1)$  and  $X_s \sim NGH(d^n, s, 1)$ , and can apply directly the formula of the expectation from Proposition A.3 to find  $\mathbb{E}[X_1]$  and  $E[X_s]$ .  $\square$

The study is slightly more complex when using `UpdateTree`. One can remark that the number of elements removed from the search space  $Z$  by `UpdateTree` depends on the index of the state. We will therefore consider the random variables corresponding to the number of failures between two found solutions.

**Notation A.5.** Let  $L/K$  be an extension of number fields such that  $[L : K] = n$ . Consider  $f(X) \in L[X]$  such that  $|Z_K(f)| = s$ . We will denote by  $x^{(1)} < \dots < x^{(s)}$  the elements of  $Z_K(f)$  ordered in  $Z$ . For each  $k \in \llbracket 1, s \rrbracket$  we will consider several random variables.

- (1)  $X^{(k)}$  is the random variable corresponding to the number of failures between the  $(k - 1)$ -th solution and the  $k$ -th solution.
- (2)  $X_1^{(k)}$  is the random variable corresponding to the number of failures in the first coordinate between the  $(k - 1)$ -th solution and the  $k$ -th solution.
- (3)  $X_2^{(k)}$  is the random variable corresponding to the number of failures occurred between the  $(k - 1)$ -th solution and the  $k$ -th solution such that  $j_1 = x_1^{(k)}$ .
- (4)  $Y^{(k)}$  is the random variable corresponding to the number of failures which occurred before the  $k$ -th solution is found.

**Lemma A.6.** *Let  $L/K$  be an extension of number fields such that  $[L : K] = n$ . Consider  $f(X) \in L[X]$  such that  $|Z_K(f)| = s$ . Then one has the following:*

$$\forall k \in \llbracket 1, s \rrbracket, Y^{(k)} = \sum_{j=1}^k \left( X_1^{(j)} (d - j + 1)^{n-1} + X_2^{(j)} \right).$$

*Proof.* Let us fix  $k \in \llbracket 1, s \rrbracket$ . Clearly one has  $Y^{(k)} = \sum_{j=1}^{(k)} X^{(j)}$ . Now let us denote by  $C_k$  the integer  $|Z_2 \times \cdots \times Z_n|$  after the  $(k-1)$ -th root and before the  $k$ -th root are found. Recall that after each new solution is found, **UpdateTree** removes one element of each  $Z_i, i > 1$ . Therefore, one obtains  $C_k = (d-k+1)^{n-1}$ . Because the search is done following the lexicographic order, it is easy to see that for each fixed  $j_1 \in \llbracket x_1^{(k-1)} + 1, x_1^{(k)} \rrbracket$  there are two possibilities. If  $j_1 < x_1^{(k)}$  then the search will run through all  $\{z_{1,j_1}\} \times Z_2 \times \cdots \times Z_n$ , which contains no solution. This leads to  $C_k$  failures. If  $j_1 = x_1^{(k)}$  then the search will run through  $\{z_{1,j_1}\} \times Z_2 \times \cdots \times Z_n$  until finding the solution. It amounts to  $X_2^{(k)}$  failures. Finally the number of  $j_1$  that are passed such that  $j_1 < x_1^{(k)}$  is  $X_1^{(k)}$ .  $\square$

**Proposition A.7.** *Let  $L/K$  be an extension of number fields such that  $[L : K] = n$ . Consider  $f(X) \in L[X]$  such that  $|Z_K(f)| = s$ . Write  $d = \deg f(X)$ . Then, for  $k \in \llbracket 1, s \rrbracket$ , the average number of “failed” decodings done in Algorithm 8 before finding  $k$  roots is*

$$(A.1) \quad \frac{2d-s+1}{2(s+1)} \sum_{j=0}^{k-1} (d-j)^{n-1} - \frac{k}{2}.$$

*Proof.* Using Lemma A.6 and by linearity of the expectation, one has:

$$\forall k \in \llbracket 1, s \rrbracket, \mathbb{E}[Y^{(k)}] = \sum_{j=1}^k \left( \mathbb{E}[X_1^{(j)}] (d-j+1)^{n-1} + \mathbb{E}[X_2^{(j)}] \right).$$

Let fix  $j \in \llbracket 1, k \rrbracket$ . Recall that  $X_2^{(j)}$  is the number of failures found during the search through the set  $\{x_1^{(j)}\} \times Z_2 \times \cdots \times Z_n$ . We know there is exactly one solution in this set. Therefore we have  $X_2^{(j)} \sim NHG((d-j+1)^{n-1}, 1, 1)$  and

$$\mathbb{E}[X_2^{(j)}] = \frac{(d-j+1)^{n-1} - 1}{2}.$$

Now let us determine  $X_1^{(j)}$ . It is the number of wrong first coordinates visited until finding  $x_1^{(j)}$ , and after finding  $x_1^{(j-1)}$ . The search is done over the set  $Z_1$  minus the elements visited before  $x_1^{(j-1)}$  included. It is a set with cardinal number

$$|Z_1| - \sum_{i=1}^{(j-1)} X_1^{(i)} - (j-1) = d-j+1 - \sum_{i=1}^{(j-1)} X_1^{(i)}$$

which contains  $s - (j-1)$  success elements. Therefore we have  $\mathbb{E}[X_1^{(j)} \mid \sum_{i=1}^{(j-1)} X_1^{(i)} = a] \sim NHG(d-j+1-a, s-(j-1), 1)$ , for all possible  $a$ . Using the law of total expectation  $\mathbb{E}[X] = \mathbb{E}[\mathbb{E}[X|Y]]$  between two variables  $X$  and  $Y$ , it is easy to see that we have

$$\mathbb{E}[X_1^{(j)}] = \frac{(d-j+1) - \sum_{i=1}^{(j-1)} \mathbb{E}[X_1^{(i)}] - (s-j+1)}{s-j+2} = \frac{(d-s) - \sum_{i=1}^{(j-1)} \mathbb{E}[X_1^{(i)}]}{s-j+2}.$$

It is possible to use this recurrence relation to obtain an expression of  $\mathbb{E}[X_1^{(j)}]$  in closed-form. As a matter of fact, we will prove that  $\mathbb{E}[X_1^{(j+1)}] = \mathbb{E}[X_1^{(j)}]$ . Indeed one has

$$\mathbb{E}[X_1^{(j+1)}] = \frac{(d-s) - \sum_{i=1}^{(j)} \mathbb{E}[X_1^{(i)}]}{s-j+1} = \frac{d-s - \mathbb{E}[X_1^{(j)}] - \sum_{i=1}^{(j-1)} \mathbb{E}[X_1^{(i)}]}{s-j+1},$$

and since

$$(s-j+2)\mathbb{E}[X_1^{(j)}] = (d-s) - \sum_{i=1}^{(j-1)} \mathbb{E}[X_1^{(i)}]$$

we obtain

$$\mathbb{E}[X_1^{(j+1)}] = \frac{d - s - \mathbb{E}[X_1^{(i)}] + (s - j + 2)\mathbb{E}[X_1^{(j)}] - (d - s)}{s - j + 1} = \mathbb{E}[X_1^{(j)}].$$

Then remark that  $X_1^{(1)} \sim NHG(d, s, 1)$ , which gives the desired result.  $\square$

**Remark A.8.** One can see that the gain of using `UpdateTree` increases with  $s$ . Indeed if  $f(X)$  has only one root, then the expected number of decodings is the same for both methods. However, if  $s = \deg f(X)$  then the expected number of decodings with the naive search is  $d \frac{d(d^{n-1}-1)}{d+1}$ , whereas it is  $\sum_{i=0}^{d-1} \frac{(d-i)^{n-1}-1}{2}$  when using `UpdateTree`.

Let us analyse more precisely the difference between the two procedures.

**Notation A.9.** Let  $L/K$  be an extension of number fields such that  $[L : K] = n$ . Consider  $f(X) \in L[X]$  such that  $|Z_K(f)| = s$ . For each  $k \in \llbracket 1, s \rrbracket$  we will consider several random variables, all related to a naïve search.

- (1)  $X_{\text{naive}}^{(k)}$  is the random variable corresponding to the number of failures between the  $(k-1)$ -th solution and the  $k$ -th solution.
- (2)  $Y_{\text{naive}}^{(k)}$  is the random variable corresponding to the number of failures which occurred before the  $k$ -th solution is found.

Finally we will denote by  $S_X^{(k)}$  the difference  $X_{\text{naive}}^{(k)} - X^{(k)}$  and by  $S_Y^{(k)}$  the difference  $Y_{\text{naive}}^{(k)} - Y^{(k)}$ .

**Proposition A.10.** *Let  $L/K$  be an extension of number fields such that  $[L : K] = n$ . Consider  $f(X) \in L[X]$  such that  $|Z_K(f)| = s$ . Write  $d = \deg f(X)$ . Then, for  $k \in \llbracket 1, s \rrbracket$  we have  $\mathbb{E}[X_{\text{naive}}^{(k)}] \geq \mathbb{E}[X^{(k)}]$ ,  $\mathbb{E}[Y_{\text{naive}}^{(k)}] \geq \mathbb{E}[Y^{(k)}]$  and the following is true:*

$$(A.2) \quad S_Y^{(k)} = k \frac{(s-1)(d^{n-1}-1)}{2(s+1)} + \frac{2d-s+1}{s+1} \sum_{i=0}^{k-1} (d^{n-1} - (d-i)^{n-1}).$$

*Proof.* First, for all  $k \in \llbracket 1, s \rrbracket$  we will denote by  $c_k$  the integer  $d - k + 1$ . Then, from the proofs of Lemma A.6 and Proposition A.7 we get

$$\mathbb{E}[X^{(k)}] = \frac{d-s}{s+1} c_k^{n-1} + \frac{c_k^{n-1}-1}{2} = \frac{2dc_k^{n-1} - (s-1)c_k^{n-1} - s - 1}{2(s+1)}.$$

Moreover  $\mathbb{E}[X_{\text{naive}}^{(k)}] = \frac{d^n - s}{s+1}$ . Therefore we obtain

$$\begin{aligned} S_X^{(k)} &= \frac{2d^n - 2s - 2dc_k^{n-1} + (s-1)c_k^{n-1} + s + 1}{2(s+1)} \\ &= \frac{2d(d^{n-1} - c_k^{n-1}) + (s-1)(c_k^{n-1} - 1)}{2(s+1)}, \end{aligned}$$

which is positive. Now let us determine how this difference evolves with  $k$ . Clearly  $c_k$  decreases when  $k$  increases, therefore so does  $\mathbb{E}[X^{(k)}]$ . Since  $\mathbb{E}[X_{\text{naive}}^{(k)}]$  is constant, we can conclude that  $S_X^{(k)}$  increases with  $k$ . Let us express  $S_Y^{(k)}$ . Given any sequence  $(u_n)_n$  let us denote  $u_{n+1} - u_n$  by  $\Delta(u, n)$ . Remark that  $S_Y^{(k)} = \sum_{i=1}^k S_X^{(i)}$  and  $S_X^{(k)} = S_X^{(1)} + \sum_{i=1}^{k-1} S_X^{(i+1)} - S_X^{(i)} = S_X^{(1)} + \sum_{i=1}^{k-1} \Delta(S_X, i)$ . Moreover one has

$$\Delta(S_X, i) = \mathbb{E}[X^{(i+1)}] - \mathbb{E}[X^{(i)}] = \frac{(2d-s+1)(-c_{i+1}^{n-1} + c_i^{n-1})}{2(s+1)}$$

which leads to

$$\begin{aligned} S_X^{(k)} &= S_X^{(1)} + \frac{2d-s+1}{s+1} \sum_{i=1}^{k-1} \Delta((-c_k^{n-1})_k, i) \\ &= \frac{(s-1)(d^{n-1}-1)}{2(s+1)} + \frac{2d-s+1}{s+1} (d^{n-1} - c_k^{n-1}). \end{aligned}$$

Finally we can write

$$\begin{aligned} S_Y^{(k)} &= \sum_{i=1}^k \frac{(s-1)(d^{n-1}-1)}{2(s+1)} + \frac{2d-s+1}{s+1} (d^{n-1} - c_i^{n-1}) \\ &= k \frac{(s-1)(d^{n-1}-1)}{2(s+1)} + \frac{2d-s+1}{s+1} \sum_{i=0}^{k-1} (d^{n-1} - (d-i)^{n-1}). \end{aligned}$$

□

## APPENDIX B. EXTRA DATA

**B.1. Different versions of LLL.** In this section we present the timings obtained when comparing the efficiency of LLL with our method SpecLLL (Alg. 9), which correspond to the ratio presented in Figures 2 and 3. One can find said timings in Figures 14 and 15.

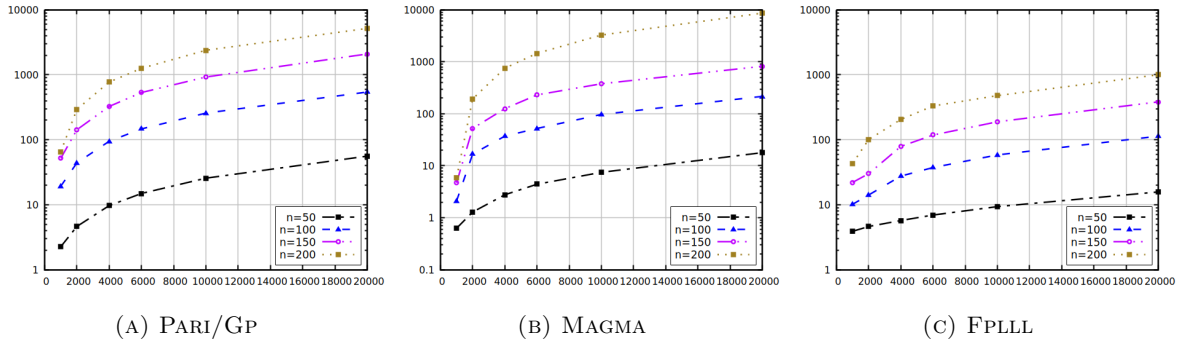
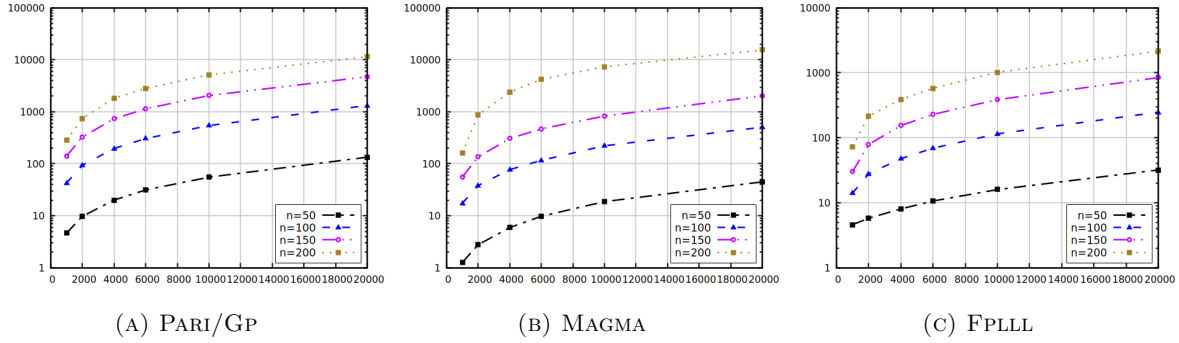


FIGURE 14. Timings (s) for  $\sigma(K) \subset \mathbb{R}$  plotted against the precision

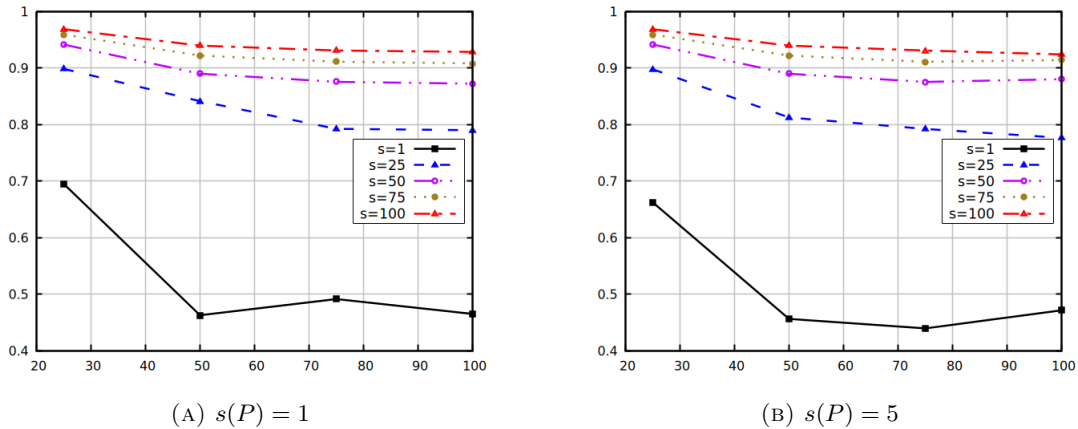
As mentioned previously, FPLLL is faster than MAGMA or PARI/GP. One can remark that the latter is around 5 times slower than FPLLL asymptotically. This gives an idea of the timings that could be obtained for all our experiments by switching from the implementation of PARI/GP to FPLLL.

Precision		1000	2000	4000	6000	10000	20000
$n = 50$	LLL	2.5	5.1	10.7	16.8	29.8	69.4
	SpecLLL	2.6	5.2	10.8	16.8	29.8	69.3
$n = 150$	LLL	74.5	166.0	371.2	639.3	1132.0	2843.6
	SpecLLL	76.1	167.6	367.7	634.9	1118.1	2841.0

TABLE 6. Timings (s) over random knapsack-like matrices

FIGURE 15. Timings (s) for  $\sigma(K) \not\subset \mathbb{R}$  plotted against the precision

**B.2. Precision evaluation.** In Figure 16, we plotted the ratio of  $\max q_t$  by  $1 + \ln(n) \ln \ln(n) / (s \ln 2 + \ln(n)/2)$ .

FIGURE 16. Maximum ratios of  $q_t$  by  $1 + \ln n \ln \ln(n) / (s \ln 2 + \ln(n)/2)$  plotted against  $n = [K : \mathbb{Q}]$ , for several  $s$ .

Note that we did similar experiments using an integral basis of  $\mathcal{O}_K$  instead of the equation basis  $(1, X, \dots, X^{n-1})$  and obtained similar results, which can be found in Figure 17. We were to smaller degrees because of the cost of computing such basis.

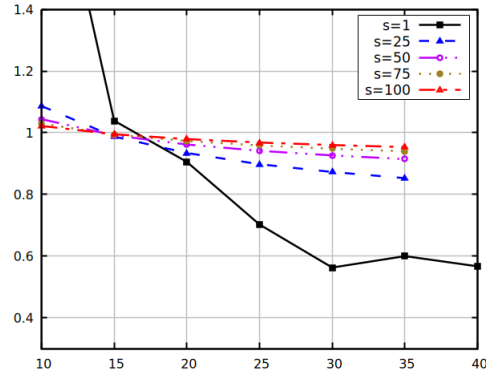
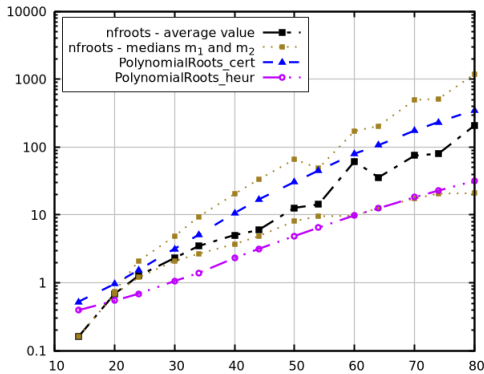


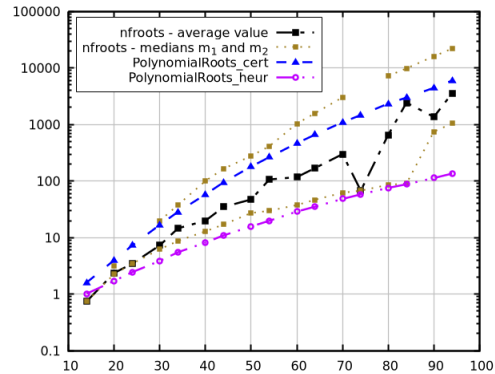
FIGURE 17. Maximum ratios of  $q_t$  by  $1 + \ln n \ln \ln(n)/(s \ln 2 + \ln(n)/2)$  plotted against  $n = [K : Q]$  for several  $s$ , where  $\mathcal{B}$  is an integral basis of  $\mathcal{O}_K$

One can see that ratio of  $q_t$  by  $1 + \ln n \ln \ln(n)/(s \ln 2 + \ln(n)/2)$  is larger than 1 only for very small dimensions, which also happens for the equation basis. In practice, the additional precision that is required does not impact significantly the running times the algorithms.

**B.3. Absolute method.** In this section we give data corresponding to experiments presented in Section 6.2, in the case where  $r_1 = 0$ , and consequently  $\sigma(K) \subset \mathbb{C} \setminus \mathbb{R}$ .



(A)  $s(P_K) = 1$



(B)  $s(P_K) = 10$

FIGURE 18. Average timings (s) of `nfroots`, `AbsoluteRoots` and `AbsoluteRootsHeur` plotted against  $\deg P_K(X)$  for randomly generated  $P_K(X)$  such that  $r_1 = 0$ , with  $s(P_K) \in \{1, 10\}$

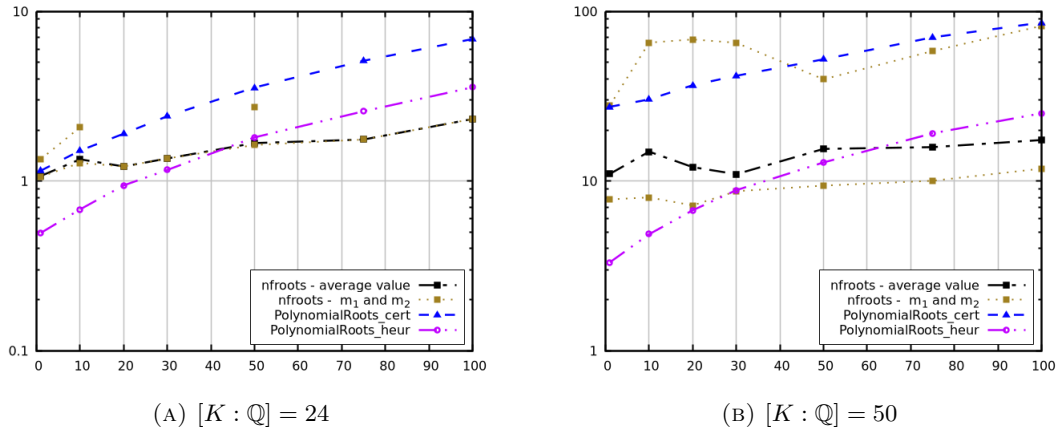


FIGURE 19. Average timings (s) of nroots, AbsoluteRoots and AbsoluteRootsHeur plotted against  $s_Z$  for randomly generated  $P_K(X)$  such that  $[K : \mathbb{Q}] = 24$  and  $[K : \mathbb{Q}] = 50$ , with  $r_1 = 0$