# **Fast Fourier Orthogonalization**

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ABSTRACT

The classical fast Fourier transform (FFT) allows to compute in quasi-linear time the product of two polynomials, in the *circular convolution ring*  $\mathbb{R}[x]/(x^d-1)$  — a task that naively requires quadratic time. Equivalently, it allows to accelerate matrix-vector products when the matrix is *circulant*.

In this work, we discover that the ideas of the FFT can be applied to speed up the orthogonalization process of matrices with circulant blocks of size  $d \times d$ . We show that, when d is composite, it is possible to proceed to the orthogonalization in an inductive way —up to an appropriate re-indexation of rows and columns. This leads to a structured Gram-Schmidt decomposition. In turn, this structured Gram-Schmidt decomposition accelerates a cornerstone lattice algorithm: the nearest plane algorithm. The complexity of both algorithms may be brought down to  $\Theta(d \log d)$ .

Our results easily extend to *cyclotomic rings*, and can be adapted to Gaussian samplers. This finds applications in lattice-based cryptography, improving the performances of trapdoor functions.

**Keywords.** Fast Fourier transform, Gram-Schmidt orthogonalization, nearest plane algorithm, lattice algorithms, lattice trapdoor functions.

# 1. INTRODUCTION

When operations involving linear algebra are to be performed over structured matrices, a natural problem is to accelerate them by exploiting the structure. The most classical example is the fast Fourier transform [2, 8], which allows to perform matrix-vector multiplication in quasilinear time

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when the matrix is circulant. This is achieved by exploiting the isomorphism between the ring of  $d \times d$  circulant matrices and the *circular convolution ring*  $\mathcal{R}_d = \mathbb{R}[x]/(x^d - 1)$ .

A widely studied and more involved question is matrix decomposition [19] for structured matrices, in particular Gram-Schmidt orthgonalization (GSO). In this work, we are interested in the GSO of *circulant matrices*, and more generally of matrices with circulant blocks. Our main motivation is to accelerate lattice algorithms for lattices that admit a basis with circulant blocks. This use case allows a helpful extra degree of freedom: one may permute rows and columns of the lattice basis since this leaves the generated lattice unchanged —up to an isometry.

As we will show, a proper re-indexation of these matrices highlights an inductive structure, with a fast Fourier flavor. This leads to accelerations of the orthogonalization process — and of the related nearest plane algorithm— down to quasilinear time and space.

# The Nearest Plane Algorithm, Lattices and Cryptography

The nearest plane algorithm [1] is a central algorithm over lattices. It allows, after precomputation of the GSO and using a quadratic number of real operations, to find a relatively close point in a lattice to an arbitrary target. It is a core subroutine of LLL [12], and can be used for error correction over analogical noisy channels. It has also found applications in lattice-based cryptography as a decryption algorithm, and a randomized variant (called discrete Gaussian sampling) [11, 5] provides secure trapdoor functions based on lattice problems. This leads to cryptosystems (attributebased encryption) with fine-grained access control, as [17, 6] to name a few.

Given a basis **B** of a lattice  $L \subset \mathbb{R}^d$  and a target vector **c**, the nearest plane algorithm finds a lattice point somewhat close to **c**. The result belongs to a fundamental domain centered in **c**, whose shape is the cuboid defined by  $\tilde{\mathbf{B}}$ , the GSO of **B** (see Figure 1). This algorithm performs  $\Theta(d^2)$  real operations. The GSO itself is required as a precomputation.

## Structured lattices in cryptography.

When it comes to practical lattice-based cryptography, a quadratic cost in the dimension is rather prohibitive considering the lattices at hand have dimensions ranging in the hundreds, or even thousands. For efficiency purposes, many cryptosystems (such as [10, 14, 15] to name a few) chose to rely on lattices with some algebraic structure, improving time and memory requirements to quasilinear in the dimension.

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This is sometimes referred as lattice-based cryptography in the *ring setting*. Technically, the chosen rings typically are *cyclotomic rings*, but those are closely related to the *convolution rings* discussed so far. The core of this optimization is the fast Fourier transform (FFT) [2, 4, 8, 18] allowing fast multiplication of polynomials. But this improvement did not apply in the case of the nearest plane algorithm or its randomized variant [11, 5]: naïve GSO do not take the algebraic structure into account.

One possible work-around [9, 21] consist of using the roundoff algorithm [1] instead of the nearest-plane algorithm. However, this simpler algorithm outputs further vectors, both in the average and worst cases, weakening those cryptosystems.

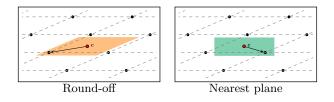


Figure 1: Round-off and nearest plane algorithms, and their associated fundamental domains.

# **Our contribution**

In this work, we discover new algorithms, obtained by crossing Cooley-Tukey's [2] fast Fourier transform algorithm together with the orthogonalization and nearest plane algorithms (not exactly the GSO, but the closely related LDL<sup>\*</sup> decomposition). Precisely, we show that, up to a re-indexation of rows and columns, the orthogonalization of matrices composed of  $d \times d$ -circulant blocks can be done in time  $\Theta(d \log d)$ when the prime factors of d are bounded. Our algorithm produces the LDL<sup>\*</sup> decomposition in a special compact format, requiring  $\Theta(d \log d)$  complex numbers to represent.

From this compact representation, the nearest plane algorithm can also be performed using  $\Theta(d \log d)$  real operations.<sup>1</sup> As a demonstration of the simplicity of our algorithms, we propose an implementation in **python** for  $d \times d$ -circulant matrices, when d is a power of 2.

### Computational model and Parallelism.

Our algorithms and their complexity are described for Random Access Machines with unitary arithmetic operations on *real numbers*, i.e. we do not study the issue related to floating-point approximations and numerical stability.

Regarding parallelism, we note that our Fast-Fourier LDL<sup>\*</sup> algorithm is almost fully parallelizable: viewed as an arithmetic circuit on real numbers, it has depth  $O(\log(n))$  and width O(n). This is unfortunately not the case of the Nearest Plane algorithm nor our Fast-Fourier variant, for which the n rounding steps are inherently sequential.

#### Techniques.

At the core of our techniques is the realization that representing elements of the convolution ring  $\mathcal{R}_d = \mathbb{R}[x]/(x^d-1)$  as circulant matrices is not the appropriate choice. To allow an induction similar to Cooley-Tukey's FFT, our representation must follow the tower of rings  $\mathbb{R} \subset \mathcal{R}_{d_1} \subset \cdots \subset \mathcal{R}_{d_{i-1}} \subset \mathcal{R}_d$ , for some chain of divisors  $1|d_1| \ldots |d_{i-1}|d$ .

Such a representation is obtained by applying the (mixedradix) digit-reversal order to the indexation of the rows and column of the circulant blocks, as pictured in Figure 2.

We show that this alternative indexation allows to represent the matrix **L** of the GSO in a factorized form: a product of  $\Theta(\log d)$  (sparse) structured matrices, each of which can be stored in space O(d). An example is given in Figure 3.

Once this hidden structure is unveiled (Theorem 1), the algorithmic implications follow quite naturally: one first easily derive an algorithm in time  $O(n \log^2 n)$  —matching previous and more general results [20]— but noting that the splitting step may be performed directly in the Fourier domain allows us to save another  $\log n$  factor. For easier algorithmic manipulations, the factorization of **L** is represented using a tree.

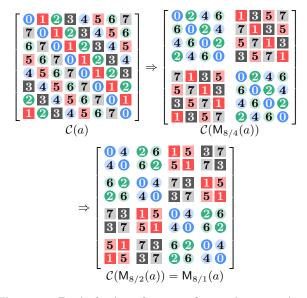


Figure 2: Re-indexing the transformation matrix of  $f \in \mathcal{R}_8 \mapsto fx \ (a = 0 + x + 2x^2 + \dots + 7x^7 \in \mathcal{R}_8)$ .

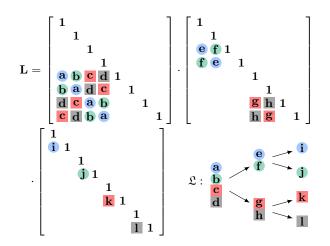


Figure 3: Factorization of L in the LDL<sup>\*</sup> decomposition of  $M_{8/1}(a)$ , and its tree representation  $\mathfrak{L}$ .

<sup>&</sup>lt;sup>1</sup>If the number  $n \times m$  of  $d \times d$  circulant blocks is not constant, they contribute a factor  $O(n^3)$  to the runtime of the fast-Fourier LDL<sup>\*</sup> algorithm,  $O(n^2)$  to its output size, and  $O(n^2)$ to the run-time of the fast-Fourier nearest plane algorithm.

# Related Works.

There exist many works related to the orthogonalization of structured bases. For Toeplitz matrices, Sweet [22] introduced an algorithm faster than the naive orthogonalization by a linear factor. Gragg [7] has shown that for Krylov bases – which are bases of the form  $\{\mathbf{b}, r(\mathbf{b}), ..., r^{d-1}(\mathbf{b})\}$  –, the Levinson recursion [13, 3] allows, when r is an isometric operator, to perform orthogonalization in time  $\Theta(d^2)$  instead of  $\Theta(d^3)$ .

There also exists superfast (running time  $O(n \log^2 n)$ ) algorithms for the orthogonalization of Toeplitz-like matrices, for example by Olshevsky and Pan [20], and those are already based on a structure-preserving induction. In this light, one may interpret our result as a Fourier-compatible version of these superfast algorithms.

The question of improving the nearest-plane algorithm for structured matrices seems less studied. As far as we know, the state of the art consist of a single result [16], applying the Levinson recursion [13] to reduce by a linear factor its *space* complexity. Alternatively, a work-around of lesser quality was proposed for the NTRU signature scheme [9, 21].

## Outline.

Section 2 introduces the mathematical tools that we will use through this paper. Section 3 presents our main result, namely the existence of a compact, factorized representation for the GSO and LDL<sup>\*</sup> decomposition, and gives a fast Fourier flavored algorithm for computing it in this form. This compact LDL<sup>\*</sup> decomposition is further exploited in Section 4, which presents a nearest plane algorithm that also has an FFT flavor. In the full version of our article<sup>2</sup>, we propose the proofs, the extension of our results to cyclotomic rings, and a **python** code fragment. The full prototype implementation in **python** —for d a power of 2— is also available online.<sup>3</sup>

# 2. PRELIMINARIES

For any ring  $\mathcal{R}$ ,  $\mathcal{R}[x]$  will denote the ring of univariate polynomials over  $\mathcal{R}$ . Scalars (which includes elements of  $\mathcal{R}$ ) will usually be noted in plain letters (such as a, b), vectors will be noted in bold letters (such as  $\mathbf{a}, \mathbf{b}$ ) and matrices will be noted in capital bold letters (such as  $\mathbf{A}, \mathbf{B}$ ). Vectors are mostly in row notation, and as a consequence vector-matrix products are done in this order unless stated otherwise.  $(a_1, \ldots, a_n)$  denotes the row vector formed of the  $a_i$ 's, whereas  $[\mathbf{a}_1, \ldots, \mathbf{a}_n]$  denotes the matrix whose rows are the  $\mathbf{a}_i$ 's.  $\mathbb{N}$  denotes the set of non-negative integers, and  $\mathbb{N}^*$  the set  $\mathbb{N}\setminus\{0\}$ . For  $i, j \in \mathbb{Z}, [[i, j]]$  will denote the set  $\{i, i + 1, \ldots, j - 1, j\}$ .

# 2.1 The Convolution Ring $\mathcal{R}_d$

DEFINITION 1. For any  $d \in \mathbb{N}^*$ , let  $\mathcal{R}_d$  denote the ring  $\mathbb{R}[x]/(x^d-1)$ , also known as circular convolution ring, or simply convolution ring.

When d is highly composite, elementary operations in  $\mathcal{R}_d$  can be performed in time  $\Theta(d \log d)$  using the fast Fourier transform [2].

We equip the ring  $\mathcal{R}_d$  with a conjugation operation as well as an inner product, making it an Hermitian inner product space. The definitions that we give also encompass other types of rings that are used in the full version of this article.

DEFINITION 2. Let  $h \in \mathbb{R}[x]$  be a monic polynomial with distinct roots over  $\mathbb{C}$ ,  $\mathcal{R} \triangleq \mathbb{R}[x]/(h(x))$  and a, b be arbitrary elements of  $\mathcal{R}$ .

- We note  $a^*$  and call (Hermitian) adjoint of a the unique element of  $\mathcal{R}$  such that for any root  $\zeta$  of h,  $a^*(\zeta) = \overline{a(\zeta)}$ , where  $\overline{\cdot}$  is the usual complex conjugation over  $\mathbb{C}$ .
- The inner product over  $\mathcal{R}$  is  $\langle a, b \rangle \stackrel{\Delta}{=} \sum_{h(\zeta)=0} a(\zeta) \cdot \overline{b(\zeta)}$ , and the associated norm is  $||a|| \stackrel{\Delta}{=} \sqrt{\langle a, a \rangle}$ .

In the particular case of convolution rings, one can check that if  $a(x) = \sum_{i=0}^{d-1} a_i x^i \in \mathcal{R}_d$ , then

$$a^{\star}(x) = a(1/x) \mod (x^d - 1) = a_0 + \sum_{i=1}^{d-1} a_i x^{d-i}.$$

The (Hermitian) adjoint  $\mathbf{B}^*$  of a matrix  $\mathbf{B} \in \mathcal{R}^{n \times m}$  is the transpose of the coefficient-wise adjoint of  $\mathbf{B}$ .

While the inner product  $\langle \cdot, \cdot \rangle$  (resp. the associated norm  $\|\cdot\|$ ) is not to be mistaken with the canonical coefficient-wise dot product  $\langle \cdot, \cdot \rangle_2$  (resp. the associated norm  $\|\cdot\|_2$ ), they are closely related. One can easily check that for any  $f = \sum_{i \in \mathbb{Z}_d} f_i x^i \in \mathcal{R}_d$ , the vector  $(f(\zeta))_{\{\zeta^d=1\}}$  can be obtained from the coefficient vector  $(f_i)_{0 \leq i < d}$  by multiplying it by the Vandermonde matrix  $\mathbf{V}_d = (\zeta_d^{ij})_{0 \leq i, j < d}$ , where  $\zeta_d$  denotes an arbitrary *d*-th primitive root of unity. The matrix  $\mathbf{V}_d$  satisfies  $\mathbf{V}_d \mathbf{V}_d^* = d \cdot \mathbf{I}_d$  and as an immediate consequence:  $\langle f, g \rangle = d \cdot \langle f, g \rangle_2$ .

DEFINITION 3. Let  $m \ge n$  and  $\mathbf{B} = {\mathbf{b}_1, \ldots, \mathbf{b}_n} \in \mathcal{R}^{n \times m}$ . We say that  $\mathbf{B}$  is full-rank (or is a basis) if for any linear combination  $\sum_{1 \le i \le n} a_i \mathbf{b}_i$  with  $a_i \in \mathcal{R}$ , we have the equivalence  $(\sum_i a_i \mathbf{b}_i = \mathbf{0}) \iff (\forall i, a_i = 0).$ 

We note that since  $\mathcal{R}$  is generally not an integral domain, a set formed of a single nonzero vector is not necessarily full-rank. In the rest of the paper, a basis will either denote a set of independent vectors  $\{\mathbf{b}_1, \dots \mathbf{b}_n\} \in (\mathcal{R}^m)^n$ , or the full-rank matrix  $\mathbf{B} \in \mathcal{R}^{n \times m}$  whose rows are the  $\mathbf{b}_i$ 's.

# **2.2** The GSO and LDL\* Decomposition

In this section,  $\mathcal{R} = \mathbb{R}[x]/(h(x))$  as in Definition 2. We first recall a few standard definitions. A matrix  $\mathbf{L} \in \mathcal{R}^{n \times n}$  is unit lower triangular if it is lower triangular and has only 1's on its diagonal.

We say that a self-adjoint matrix  $\mathbf{G} \in \mathcal{R}^{n \times n}$  is full-rank Gram (or FRG) if there exist  $m \ge n$  and a full-rank matrix  $\mathbf{B} \in \mathcal{R}^{n \times m}$  such that  $\mathbf{G} = \mathbf{B}\mathbf{B}^*$ . This generalizes the notion of positive definiteness for symmetric real matrices.

We now recall the GSO and LDL<sup>\*</sup> decomposition. The GSO decomposes any full-rank matrix as the product of a unit lower triangular matrix and an orthogonal matrix.

PROPOSITION 1. Let  $\mathbf{B} \in \mathcal{R}^{n \times m}$  be a full-rank matrix. **B** can be uniquely decomposed as

$$\mathbf{B} = \mathbf{L} \cdot \dot{\mathbf{B}},\tag{1}$$

where  $\mathbf{L}$  is unit lower triangular, and the rows of  $\tilde{\mathbf{B}}$  are pairwise orthogonal.

<sup>&</sup>lt;sup>2</sup>https://eprint.iacr.org/2015/1014

<sup>&</sup>lt;sup>3</sup>https://github.com/lducas/ffo.py

When  $\mathcal{R}$  is replaced by  $\mathbb{R}$  or a number field, Proposition 1 is standard. In our case, a proof can be found in the full version of this article.

The LDL<sup>\*</sup> decomposition writes any positive definite matrix as a product  $\mathbf{LDL}^*$ , where  $\mathbf{L} \in \mathcal{R}^{n \times n}$  is unit lower triangular with 1's on the diagonal, and  $\mathbf{D} \in \mathcal{R}^{n \times n}$  is diagonal. It is related to the GSO as for a basis **B**, there exists a unique GSO  $\mathbf{B} = \mathbf{L} \cdot \tilde{\mathbf{B}}$  and for an FRG matrix **G**, there exists a unique LDL<sup>\*</sup> decomposition  $\mathbf{G} = \mathbf{LDL}^*$ . If  $\mathbf{G} = \mathbf{BB}^*$ , then  $\mathbf{G} = \mathbf{L} \cdot (\tilde{\mathbf{B}}\tilde{\mathbf{B}}^*) \cdot \mathbf{L}^*$  is a valid LDL<sup>\*</sup> decomposition of **G**. As both decompositions are unique, the matrices **L** in both cases are actually the same. In a nutshell:

 $\mathbf{L} \cdot \tilde{\mathbf{B}}$  is the GSO of  $\mathbf{B}$ 

 $\Leftrightarrow \mathbf{L} \cdot (\mathbf{\tilde{B}}\mathbf{\tilde{B}}^{\star}) \cdot \mathbf{L}^{\star} \text{ is the LDL}^{\star} \text{ decomposition of } (\mathbf{B}\mathbf{B}^{\star}).$ 

Algorithm 1  $\mathsf{LDL}^\star_\mathcal{R}(\mathbf{G})$ 

**Require:** A full-rank Gram matrix  $\mathbf{G} = (G_{ij}) \in \mathcal{R}^{n \times n}$ . **Ensure:** The decomposition  $\mathbf{G} = \mathbf{LDL}^*$  over  $\mathcal{R}$ , where  $\mathbf{L}$ 

is unit lower triangular and **D** is diagonal. 1:  $\mathbf{L}, \mathbf{D} \leftarrow \mathbf{0}^{n \times n}$ 

1.  $\mathbf{h}, \mathbf{D} \leftarrow \mathbf{0}$ 2: for *i* from 1 to *n* do 3:  $L_{ii} \leftarrow 1$ 4:  $D_i \leftarrow G_{ii} - \sum_{j < i} L_{ij} L_{ij}^* D_j$ 5: for *j* from 1 to *i* - 1 do 6:  $L_{ij} \leftarrow \frac{1}{D_j} \left( G_{ij} - \sum_{k < j} L_{ik} L_{jk}^* D_k \right)$ 7: end for 8: end for 9: return  $((L_{ij}), \operatorname{Diag}(D_i))$ 

Algorithm 1 computes the LDL<sup>\*</sup> decomposition. When  $\mathcal{R}$  is replaced by  $\mathbb{R}$ , the decomposition is noted LDL<sup>\*</sup> rather than LDL<sup>\*</sup> and it is well-known that it terminates without encountering divisions by 0. In our case, we prove that it terminates correctly in the full version of this article.

# 2.3 Babai's Nearest Plane Algorithm

The nearest plane algorithm allows to find a lattice close to an arbitrary target in the ambient vector space. Precisely, it ensures that the difference between the target and the output lies in the fundamental parallelepiped spanned by the GSO  $\tilde{\mathbf{B}}$  of a given lattice basis  $\mathbf{B}$ , as depicted on Figure 1.

DEFINITION 4. Let  $\mathbf{B} = {\mathbf{b}_1, \dots, \mathbf{b}_n}$  be a real basis. We call fundamental parallelepiped generated by  $\mathbf{B}$  and note  $\mathcal{P}(\mathbf{B})$  the set  $\sum_{1 \leq i \leq n} \left[ -\frac{1}{2}, \frac{1}{2} \right] \mathbf{b}_j = \left[ -\frac{1}{2}, \frac{1}{2} \right]^n \cdot \mathbf{B}$ .

Algorithm 2  $\mathsf{NP}_{\mathbb{R}}(\mathbf{t}, \mathbf{L})$ 

**Require:** The decomposition  $\mathbf{B} = \mathbf{L} \cdot \tilde{\mathbf{B}}$  of  $\mathbf{B} \in \mathbb{R}^{n \times m}$ , a vector  $\mathbf{t} \in \mathbb{R}^{n}$  **Ensure:** A vector  $\mathbf{z} \in \mathbb{Z}^{n}$  such that  $(\mathbf{t} - \mathbf{z})\mathbf{B} \in \mathcal{P}(\tilde{\mathbf{B}})$ 1:  $\mathbf{z} \leftarrow \mathbf{0}$ 2: for  $j = n, \dots, 1$  do 3:  $\bar{t}_{j} \leftarrow t_{j} + \sum_{i>j} (t_{i} - z_{i})L_{ij}$ 4:  $z_{j} \leftarrow \lfloor \bar{t}_{j} \rfloor$ 5: end for 6: return  $\mathbf{z}$ 

PROPOSITION 2 (FROM [1, 12]). Algorithm 2 outputs an integer vector  $\mathbf{z}$  ( $\mathbf{zB} \in \mathcal{L}(\mathbf{B})$ ) such that  $(\mathbf{t} - \mathbf{z})\mathbf{B} \in \mathcal{P}(\tilde{\mathbf{B}})$ .

# 2.4 Coefficient Vectors and Circulant Matrices

DEFINITION 5. We define linear maps  $c : \mathcal{R}_d^m \to \mathbb{R}^{dm}$  and  $\mathcal{C} : \mathcal{R}_d^{n \times m} \to \mathbb{R}^{dn \times dm}$  as follows. For any  $a = \sum_{i \in \mathbb{Z}_d} a_i x^i \in \mathcal{R}_d$  where each  $a_i \in \mathbb{R}$ :

- 1. The coefficient vector of a is  $c(a) = (a_0, \ldots, a_{d-1})$ .
- $2. \ The \ circulant \ matrix \ of \ a \ is$

$$\mathcal{C}(a) \stackrel{\Delta}{=} \begin{bmatrix} a_{d-1}^{a_0} & a_1 & \cdots & a_{d-1} \\ a_{d-1}^{} & a_0^{} & \cdots & a_{d-2} \\ \vdots & \vdots & \ddots & \ddots & \vdots \\ a_1^{} & a_2^{} & \cdots & a_0^{} \end{bmatrix} = \begin{bmatrix} c(a) \\ c(xa) \\ \vdots \\ c(x^{d-1}a) \end{bmatrix} \in \mathbb{R}^{d \times d}.$$

3. c and C generalize to vectors and matrices in a coefficientwise manner.

PROPOSITION 3. The maps c and C satisfy the following properties:

- 1. C is an injective algebra morphism. In particular, C(a)C(b) = C(ab).
- 2.  $c(a)\mathcal{C}(b) = c(ab)$ .
- 3.  $\mathcal{C}(a)^{\star} = \mathcal{C}(a^{\star}).$

Proposition 3, item 2 illustrates the following fact: the maps c and C are complementary. Indeed, if we consider the ordered  $\mathbb{R}$ -basis  $\mathbf{B} = \{1, x, \dots, x^{d-1}\}$  of  $\mathcal{R}_d$ , then c is the map that writes any  $a \in \mathcal{R}_d$  as its unique decomposition according to  $\mathbf{B}$ . Similarly, for the same basis,  $\mathcal{C}(a)$  is the transformation matrix of the injective map  $f \in \mathcal{R}_d \mapsto fa$ . This is the intrinsic reason that makes Proposition 3, item 2 true.

In Section 2.5, we will define maps V, M which will be complementary in the same way as c, C are.

# 2.5 Linearization Operators

In this section, we introduce two partial linearization operators, a vectorial one denoted by  $V_{d/d'}$  (or V), and a matricial one denoted by  $M_{d/d'}$  (or M).

They are very similar to c, C. Indeed, for any d'|d and  $a \in \mathcal{R}_d, V_{d/d'}(a)$  will write a as a vector of the  $\mathcal{R}_{d'}$ -module  $\mathcal{R}_{d'}^{d/d'}$ , and  $\mathsf{M}_{d/d'}(a)$  will be the matrix associated to the injective map  $f \in \mathcal{R}_d \mapsto fa$ . They are in fact generalizations of the maps from Section 2.4:

- 1.  $\mathcal{R}_d$  may not only be seen as a *d*-dimensional  $\mathbb{R}$ -vector space, but also as a d/d'-dimensional  $\mathcal{R}_{d'}$ -module for any d'|d.
- 2. When writing  $\mathcal{R}_d$  as a *d*-dimensional  $\mathbb{R}$ -vector space, we will not use  $\{1, x, \ldots, x^{d-1}\}$  as our canonical basis, but a permutation of this basis

DEFINITION 6. Let  $d \in \mathbb{N}^*$  be a product of h (not necessarily distinct) primes. Let gpd(d) denote the greatest proper divisor of d. When clear from context, we also note h the number of prime divisors of d (counted with multiplicity),  $d_h \stackrel{\Delta}{=} d$  and for  $i \in [\![1,h]\!]$ ,  $d_{i-1}gpd(d_i)$  and  $k_i \stackrel{\Delta}{=} d_i/d_{i-1}$ , so that  $1 = d_0|d_1| \dots |d_h = d$  and  $\prod_{j \leq i} k_j = d_i$ .

The  $d_i$ 's defined in Definition 6 form a tower of proper divisors of d. For any composite d, there exist multiple towers of proper divisors: per example, 1|6, 1|2|6 and 1|3|6 for d = 6. In this paper, each time we mention a tower of proper divisors of d it will refer to the unique one induced by Definition 6 (e.g. for d = 6, that tower is 1|3|6).

DEFINITION 7. Let  $d, d' \in \mathbb{N}^*$  such that d' is in the tower of proper divisors of d, and let k = d/d'. We denote by x the indeterminate of the polynomial ring  $\mathcal{R}_d = \mathbb{R}[x]/(x^d-1)$  and by  $y = x^k$  the indeterminate of  $\mathcal{R}_{d'} = \mathbb{R}[y]/(y^{d'} - 1)$ . We define the partial linearization  $V_{d/d'} : \mathcal{R}_d^m \to \mathcal{R}_{d'}^{km}$  recursively as follows:

- 1. For d = d' = 1,  $V_{d/d'}$  is the identity.
- 2. For  $d' = \operatorname{gpd}(d)$  and a single element  $a \in \mathcal{R}_d$ , let  $a = \sum_{i \in \mathbb{Z}_{k}} x^{i} a_{i}(y)$  where  $a_{i} \in \mathcal{R}_{d'}$  for each *i*. Then

$$\mathsf{V}_{d/d'}(a) \stackrel{\Delta}{=} (a_0, \dots, a_{k-1})$$

In other words,  $V_{d/d'}(a) \in \mathcal{R}_{d'}^k$  is the row vector whose

- coefficients are the  $(a_i)_{i \in \mathbb{Z}_k}$ . 3. For a vector  $\mathbf{v} \in \mathcal{R}_d^m$ ,  $V_{d/d'}(\mathbf{v}) \in \mathcal{R}_{d'}^{km}$  is the componentwise applications of  $V_{d/d'}$ .
- 4. For d''|d'|d and any vector  $\mathbf{v} \in \mathcal{R}_d^m$ ,

$$\mathsf{V}_{d/d''}(\mathbf{v}) \stackrel{\Delta}{=} \mathsf{V}_{d'/d''} \circ \mathsf{V}_{d/d'}(\mathbf{v}) \in \mathcal{R}_{d''}^{(d/d'')m}$$

When d is clear from context, we simply note  $V_{d/d'} = V_{/d'}$ .

#### Interpretation.

In practice, an element  $a \in \mathcal{R}_d$  is represented by a vector of d real elements corresponding to the d coefficients of a. In this context, the operator V simply permutes coefficients. As highlighted by Figure 4, when  $d = 2^{h}$  is a power of two,  $V_{d/1}$  permutes the coefficients according to the bit-reversal order<sup>4</sup>, which appears in the radix-2 fast Fourier transform (FFT). More generally, one can show that for an arbitrary d,  $V_{d/1}$  permutes the coefficient according to the general mixedradix digit reversal order, which appears in the mixed-radix Cooley-Tukey FFT [2].

$$\begin{array}{c} (1) \ \begin{array}{c} 2 \ \begin{array}{c} 3 \ 4 \ 5 \ 6 \ 7 \end{array} \\ c(a) \end{array} \Rightarrow \begin{array}{c} (1) \ 2 \ \begin{array}{c} 4 \ 6 \ 1 \ 3 \ 5 \ 7 \end{array} \\ c(V_{/4}(a)) \end{array} \\ \end{array} \\ \Rightarrow \begin{array}{c} (1) \ 4 \ \begin{array}{c} 2 \ 6 \ 1 \ 5 \ 8 \ 7 \end{array} \\ c(V_{/2}(a)) = V_{/1}(a) \end{array} \end{array}$$

#### Figure 4: Partial vectorial linearizations.

We now move to the matrix representation M compatible with V.

DEFINITION 8. Following the notations of Definition 7, we define the operator  $\mathsf{M}_{d/d'}: \mathcal{R}_d^{n \times m} \to \mathcal{R}_{d'}^{kn \times km}$  as follows: 1. For d = d' = 1,  $M_{d/d'}$  is the identity.

2. For  $d' = \operatorname{gpd}(d), k = d/d'$  and a single element a = $\sum_{i\in\mathbb{Z}_k} x^i a_i(y)$  where each  $a_i \in \mathcal{R}_{d'}$ ,  $\mathsf{M}_{d/d'}(a)$  is the following matrix of  $\in \mathcal{R}_{d'}^{k \times k}$ :

$$\begin{bmatrix} y_{a_{k-1}}^{a_0} & a_1 & \cdots & a_{k-1} \\ y_{a_{k-1}}^{a_k} & a_0 & \cdots & a_{k-2} \\ \vdots & \vdots & \ddots & \ddots & \vdots \\ y_{a_1} & y_{a_2}^{a_k} & \cdots & a_0 \end{bmatrix} = \begin{bmatrix} \mathsf{V}_{d/d'}(a) \\ \mathsf{V}_{d/d'}(x^k a) \\ \vdots \\ \mathsf{V}_{d/d'}(x^{(d'-1)k} a) \end{bmatrix}.$$

In particular, if d is prime, then  $M_{d/1}(a) \in \mathbb{R}^{d \times d}$  is exactly the circulant matrix  $\mathcal{C}(a)$ .

3. For a vector  $\mathbf{v} \in \mathcal{R}_d^m$  or a matrix  $\mathbf{B} \in \mathcal{R}_d^{n \times m}$ ,  $\mathsf{M}_{d/d'}(\mathbf{v}) \in \mathcal{R}_{d'}^{k \times km}$  and  $\mathsf{M}_{d/d'}(\mathbf{B}) \in \mathcal{R}_{d'}^{kn \times km}$  are the componentwise applications of  $M_{d/d'}$ .

<sup>4</sup>https://oeis.org/A030109

4. For d''|d'|d and any element  $a \in \mathcal{R}_d$ ,

$$\mathsf{M}_{d/d''}(a) \stackrel{\Delta}{=} \mathsf{M}_{d'/d''} \circ \mathsf{M}_{d/d'}(a) \in \mathcal{R}_{d''}^{(d/d'') \times (d/d'')}.$$

When d is clear from context, we simply note  $M_{d/d'} = M_{/d'}$ .

#### Interpretation.

As for the operator V, the steps of applying M are depicted in Section 1 Figure 2. The linearization  $M_{d/d'}(a)$  writes the transformation matrix of the map  $f \in \mathcal{R}_d \mapsto fa$  using the same basis as the partial linearization  $V_{d/d'}$ . As a result, both operators are compatible:  $\forall a, b \in \mathcal{R}_d, V_{d/d'}(ab) =$  $V_{d/d'}(a) \cdot M_{d/d'}(b)$ . More properties follows.

PROPOSITION 4. Let  $d \in \mathbb{N}^*$ , a, b (resp.  $\mathbf{a}, \mathbf{b}$ , resp.  $\mathbf{A}, \mathbf{B}$ ) be arbitrary scalars (resp. vectors, resp. matrices) over  $\mathcal{R}_d$ , and d'|d. To be concise, we note  $V \stackrel{\Delta}{=} V_{d/d'}$  and  $M \stackrel{\Delta}{=} M_{d/d'}$ . The maps V and M satisfy the following properties:

1. M is an injective algebra morphism, and in particular:

$$\mathsf{M}(\mathbf{A} \cdot \mathbf{B}) = \mathsf{M}(\mathbf{A}) \cdot \mathsf{M}(\mathbf{B}).$$

- 2. V is an injective linear map.
- 3.  $V(ab) = V(a) \cdot M(b)$ .
- 4. V is an isometry:

$$\langle \mathsf{V}(\mathbf{a}), \mathsf{V}(\mathbf{b}) \rangle_2 = \langle \mathbf{a}, \mathbf{b} \rangle_2.$$

5. **B** is full-rank if and only if  $M(\mathbf{B})$  is full-rank.

The proofs are rather straightforward to check from the definitions, and can be found in the full version of this article.

### *Computing* V and M in the Fourier Domain.

The operators V and M that we defined can be computed very efficiently when an element  $a \in \mathcal{R}_d$  is represented by its coefficients but also when represented in the Fourier domain. In the first case, it is obvious that they can both be performed in time  $\Theta(d)$  as they (symbolically) permute coefficients of a.

If a is represented in FFT form —that is, by the vector  $(a(\zeta_d^i))_{i \in \mathbb{Z}_d}$  in  $\mathbb{C}^d$ — then computing  $\mathsf{V}_{d/d'}(a)$  and  $\mathsf{M}_{d/d'}(a)$  in FFT form can naively be done in time  $\Theta(d \log d)$  by computing its inverse FFT, permuting its coefficients, and computing d/d' FFT's over  $\mathcal{R}_{d'}$ . However, it can be done in time  $\Theta(d)$ , as it is a single step - also known as *butterfly* - of the original fast Fourier transform. This is formalized in Lemma 1, a reformulation of a simple lemma that is at the heart of Cooley and Tukey's FFT.

LEMMA 1 ([2], ADAPTED). Let  $d \ge 2, d' = \text{gpd}(d)$  and k = d/d'. Let  $\mathsf{V} \stackrel{\Delta}{=} \mathsf{V}_{d/d'}, \mathsf{M} \stackrel{\Delta}{=} \mathsf{M}_{d/d'}$ . For any  $a \in \mathcal{R}_d$  (resp.  $\mathbf{a} \in \mathsf{V}(\mathcal{R}_d), \ resp. \ \mathbf{A} \in \mathsf{M}(\mathcal{R}_d)):$ 

• V(a),  $V^{-1}(a)$ ,  $M^{-1}(A)$  can be computed in FFT form in time  $\Theta(kd)$ .

• M(a) can be computed in FFT form in time  $\Theta(k^2d)$ .

PROOF. Let  $a \in \mathcal{R}_d$  be uniquely written  $a = \sum_{i \in \mathbb{Z}_k} x^i a_i(x^k)$ , where each  $a_i \in \mathcal{R}_{d'}$ . Cooley and Tukey show in [2] (equations 7, 8) that we can switch from the FFT of a to the FFT of all the  $a_i$ 's (and conversely) in time  $\Theta(kd)$ . As the  $a_i$ 's are the coefficients of V(a) and M(a), the result follows.

In Sections 3 and 4, we will use the operators V and M to speed up the orthogonalization and nearest plane algorithms. The core idea is that these operators allow to "batch" orthogonalization operations, resulting in a  $\Theta(d/\log(d))$  speedup.

# 3. FAST FOURIER LDL DECOMPOSITION

This section presents our main result. We present the existence of a compact representation in Section 3.1, and then derive a fast algorithm to compute it in Section 3.2.

## 3.1 A Compact LDL\* Decomposition

THEOREM 1. Let  $d \in \mathbb{N}$  and  $1 = d_0|d_1| \dots |d_h| = d$  be a tower of proper divisors of d. Let  $\mathbf{b} \in \mathcal{R}_d^m$  such that  $M_{d/1}(\mathbf{b})$ is full-rank. There exists a GSO of  $M_{d/1}(\mathbf{b})$  as follows:

$$\mathsf{M}_{d/1}(\mathbf{b}) = \left(\prod_{i=0}^{h-1} \mathsf{M}_{d_i/1}(\mathbf{L}_i)\right) \cdot \tilde{\mathbf{B}}_0$$

where  $\tilde{\mathbf{B}}_0 \in \mathbb{R}^{d \times dm}$  is orthogonal, and each  $\mathbf{L}_i \in \mathcal{R}_{d_i}^{(d/d_i) \times (d/d_i)}$ is a block-diagonal matrix with unit lower triangular matrices of  $\mathcal{R}_{d_i}^{(d_{i+1}/d_i) \times (d_{i+1}/d_i)}$  as its  $d/d_{i+1}$  diagonal blocks.<sup>5</sup>

As an example, the matrix **L** of the GSO of  $M_{8/1}(a)$  for some  $a \in \mathcal{R}_8$  is depicted in Section 1 Figure 3.

PROOF. If d is prime, the theorem is trivial as it is exactly the GSO. We suppose that d is composite and that the theorem is true for any  $\mathcal{R}_i$  with i < d. By Proposition 4, item 5, the matrix  $\mathbf{B}_{h-1} \stackrel{\Delta}{=} \mathsf{M}_{d/d_{h-1}}(\mathbf{b})$  is full-rank too. Using the classical GSO, we can therefore decompose it as  $\mathbf{B}_{h-1} = \mathbf{L}_{h-1}\tilde{\mathbf{B}}$ , where  $\mathbf{L}_{h-1} \in \mathcal{R}_{d_{h-1}}^{k_h \times k_h}$ ,  $\tilde{\mathbf{B}} \in \mathcal{R}_{d_{h-1}}^{k_h \times mk_h}$ and  $k_h \stackrel{\Delta}{=} d/\operatorname{gpd}(d)$ .  $\mathbf{L}_{h-1}$  is unit lower triangular and  $\tilde{\mathbf{B}}$  is orthogonal. Noting  $\tilde{\mathbf{B}} = [\mathbf{b}_1, \dots, \mathbf{b}_{k_h}]$ , all the  $\mathbf{b}_j$ 's are pairwise orthogonal and each  $\mathsf{M}_{/1}(\mathbf{b}_j)$  is full-rank. By inductive hypothesis, they can be decomposed as follows:

$$\forall j \in \llbracket 1, k_h \rrbracket, \mathsf{M}_{d_{h-1}/1}(\mathbf{b}_j) = \left(\prod_{i=0}^{h-2} \mathsf{M}_{d_i/1}(\mathbf{L}_{i,j})\right) \cdot \tilde{\mathbf{B}}_j, \quad (2)$$

where each  $\tilde{\mathbf{B}}_{j} \in \mathbb{R}^{d_{h-1} \times md_{h-1}}$  is full-rank orthogonal and for i < h-1, each  $\mathbf{L}_{i,j} \in \mathcal{R}_{d_{i}}^{(d_{h-1}/d_{i}) \times (d_{h-1}/d_{i})}$  is a block-diagonal matrix with unit lower triangular matrices of  $\mathcal{R}_{d_{i}}^{(d_{i+1}/d_{i}) \times (d_{i+1}/d_{i})}$  as its  $d_{h-1}/d_{i+1}$  diagonal blocks. To be concise, we now note  $\mathsf{M} \triangleq \mathsf{M}_{d_{h-1}/1}$  and  $\mathsf{V} \triangleq \mathsf{V}_{d_{h-1}/1}$ . We have:

$$\begin{split} \mathcal{A}_{d/1}(\mathbf{b}) &= \mathsf{M}(\mathbf{L}_{h-1}) \cdot \mathsf{M}(\tilde{\mathbf{B}}_{h-1}) \\ &= \mathsf{M}(\mathbf{L}_{h-1}) \cdot \mathsf{M}[\mathbf{b}_1, \dots, \mathbf{b}_{k_h}] \\ &= \mathsf{M}(\mathbf{L}_{h-1}) \cdot [\mathsf{M}(\mathbf{b}_1), \dots, \mathsf{M}(\mathbf{b}_{k_h})] \\ &= \mathsf{M}(\mathbf{L}_{h-1}) \cdot \operatorname{Diag}\left(\prod_{i=0}^{h-2} \mathsf{M}_{d_i/1}(\mathbf{L}_{i,j})\right) \cdot \tilde{\mathbf{B}}_0 \\ &= \mathsf{M}(\mathbf{L}_{h-1}) \cdot \left(\prod_{i=0}^{h-2} \mathsf{M}_{d_i/1}(\mathbf{L}_i)\right) \cdot \tilde{\mathbf{B}}_0 \\ &= \left(\prod_{i=0}^{h-1} \mathsf{M}_{d_i/1}(\mathbf{L}_i)\right) \cdot \tilde{\mathbf{B}}_0, \end{split}$$

where  $\tilde{\mathbf{B}}_0 = [\tilde{\mathbf{B}}_1, \ldots, \tilde{\mathbf{B}}_{k_h}]$ . The first equality simply uses the fact that M is a ring homomorphism (Proposition 4, item 1). The second and third ones are immediate from the definitions. The fourth one uses the inductive hypothesis (equation 2) on each  $\mathbf{b}_j$  and take  $\tilde{\mathbf{B}}_0 \triangleq [\tilde{\mathbf{B}}_1, \ldots, \tilde{\mathbf{B}}_{k_h}]$ . In the fifth equality, we take  $\mathbf{L}_i \triangleq \text{Diag}(\mathbf{L}_{i,1}, \ldots, \mathbf{L}_{i,k_h})$  and just need to check that  $\tilde{\mathbf{B}}_0$  and  $\mathbf{L}$  are as stated by the theorem:

- Since each  $\mathbf{L}_{i,j}$  is block diagonal with  $d_{h-1}/d_{i+1}$  unit lower triangular diagonal blocks,  $\mathbf{L}_i$  is block diagonal with  $k_h(d_{h-1}/d_{i+1}) = d/d_{i+1}$  unit lower triangular diagonal blocks.
- We also need to show that  $\mathbf{B}_0$  is orthogonal. Each submatrix  $\tilde{\mathbf{B}}_j$  of  $\tilde{\mathbf{B}}_0$  is the orthogonalization of  $\mathsf{M}(\mathbf{b}_j)$  by induction hypothesis. Therefore, for two distinct rows  $\mathbf{u}, \mathbf{v}$  of  $\tilde{\mathbf{B}}_0$ :
  - If they belong to the same submatrix  $\mathbf{B}_j$ , they are orthogonal by induction hypothesis.
  - Suppose they belong to different submatrices:  $\mathbf{u} \in \tilde{\mathbf{B}}_j, \mathbf{v} \in \tilde{\mathbf{B}}_\ell$  and  $j \neq \ell$ . Then  $\mathbf{u}$  (resp.  $\mathbf{v}$ ) is a linear combination of rows of  $\mathsf{M}(\mathbf{b}_j)$  (resp.  $\mathsf{M}(\mathbf{b}_\ell)$ ):  $\mathbf{v} = \mathbf{a}_j \cdot \mathsf{M}(\mathbf{b}_j)$  and  $\mathbf{v} = \mathbf{a}_\ell \cdot \mathsf{M}(\mathbf{b}_j)$  for some  $\mathbf{a}_j, \mathbf{a}_\ell$  in  $\mathbb{R}^{d_{h-1}}$ . Noting  $a_j = \mathsf{V}^{-1}(\mathbf{a}_j)$  and  $a_\ell = \mathsf{V}^{-1}(\mathbf{a}_\ell)$ :

$$\begin{aligned} \langle \mathbf{u}, \mathbf{v} \rangle_2 &= \langle \mathsf{V}(a_j) \mathsf{M}(\mathbf{b}_j), \mathsf{V}(a_\ell) \mathsf{M}(\mathbf{b}_\ell) \rangle_2 \\ &= \langle \mathsf{V}(a_j \mathbf{b}_j), \mathsf{V}(a_\ell \mathbf{b}_\ell) \rangle_2 \\ &= \langle a_j \mathbf{b}_j, a_\ell \mathbf{b}_\ell \rangle_2 = 0 \end{aligned}$$

Where the second equality comes from Proposition 4, item 3, the third one from the fact that V is a scaled isometry (Proposition 4, item 4) and the fourth one from  $\mathbf{b}_i, \mathbf{b}_\ell$  being orthogonal.

# Therefore $\tilde{\mathbf{B}}_0$ is orthogonal.

The theorem we stated gives the GSO of  $M_{d/1}(\mathbf{b})$  for a vector  $\mathbf{b} \in \mathcal{R}_d^m$ , but can be easily generalized from a vector  $\mathbf{b}$  to a matrix  $\mathbf{B}$ , and also yields a compact LDL<sup>\*</sup> decomposition.

COROLLARY 1. Let  $d \in \mathbb{N}$  and  $1 = d_0|d_1| \dots |d_h = d$  be a tower of proper divisors of d. Let  $\mathbf{B} \in \mathcal{R}_d^{n \times m}$  be a full-rank matrix. There exist h + 1 matrices  $(\mathbf{L}_i)_{0 \leq i \leq h}$  such that:

- $\mathbf{L}_h \in \mathcal{R}_d^{n \times n}$  is unit lower triangular.
- For each i < h,  $\mathbf{L}_i \in \mathcal{R}_{d_i}^{n(d/d_i) \times n(d/d_i)}$  is a blockdiagonal matrix whose  $n(d/d_{i+1})$  diagonal blocks are unit lower triangular matrices of  $\mathcal{R}_{d_i}^{(d_{i+1}/d_i) \times (d_{i+1}/d_i)}$ .

Furthermore, if we note  $\mathbf{L} = \left(\prod_{i=0}^{h} \mathsf{M}_{d_i/1}(\mathbf{L}_i)\right)$  and

- $\tilde{\mathbf{B}}_0 \stackrel{\Delta}{=} \mathbf{L}^{-1} \cdot \mathsf{M}_{d/1}(\mathbf{B}), \text{ then:}$ 
  - 1. The GSO of  $M_{d/1}(\mathbf{B})$  is  $M_{d/1}(\mathbf{B}) = \mathbf{L} \cdot \ddot{\mathbf{B}}_0$ .
  - 2. The LDL<sup>\*</sup> decomposition of  $M_{d/1}(\mathbf{BB}^*)$  is

$$\mathsf{M}_{d/1}(\mathbf{B}) = \mathbf{L} \cdot (\mathbf{B}_0 \mathbf{B}_0^t) \cdot \mathbf{L}^t$$

PROOF. We have  $\mathbf{B} = \mathbf{L}_{h}\mathbf{B}'$ , where  $\mathbf{L}_{h}$  is given by either the GSO or LDL<sup>\*</sup> decomposition algorithm.  $\mathbf{B}' = {\mathbf{b}'_{1}, \ldots, \mathbf{b}'_{n}}$  is orthogonal. Applying Theorem 1 to each row vector  $\mathbf{b}'_{j}$  of  $\mathbf{B}'$  yields *n* decompositions  $(\mathbf{L}_{i,j})_{0 \leq i < h}$  and *n* orthogonal matrices  $\tilde{\mathbf{B}}_{j}$ , each spanning the same space as  $\mathbf{B}_{j} \triangleq M_{d/1}(\mathbf{b}'_{j})$ . Taking  $\mathbf{L}_{i} \triangleq \text{Diag}(\mathbf{L}_{i,j})$  and  $\tilde{\mathbf{B}}_{0} \triangleq [\tilde{\mathbf{B}}_{1}, \ldots, \tilde{\mathbf{B}}_{n}]$  yields the GSO.

The LDL<sup>\*</sup> decomposition is then given "for free" by its equivalence with the GSO, and indeed, one can check that since  $\tilde{\mathbf{B}}_0$  is orthogonal,  $\tilde{\mathbf{B}}_0 \tilde{\mathbf{B}}_0^t$  is diagonal.

Theorem 1 and Corollary 1 state that for any full-rank matrix  $\mathbf{B} \in \mathcal{R}_d^{n \times m}$ , the **L** matrix in the GSO (resp. LDL<sup>\*</sup> decomposition) of  $\mathsf{M}_{d/1}(\mathbf{B})$  (resp.  $\mathsf{M}_{d/1}(\mathbf{BB}^*)$ ) can be represented in a factorized form, where each of the factors  $\mathbf{L}_i$  is a sparse, block-diagonal matrix.

Ν

<sup>&</sup>lt;sup>5</sup>Indexed products are to be read  $\prod_{i=0}^{k} \alpha_i = \alpha_k \alpha_{k-1} \cdots \alpha_0$ .

# 3.2 A Fast Algorithm for the Compact LDL\* Decomposition

Theorem 1 and Corollary 1 are constructive: more precisely, their proofs give a fast algorithm for computing a compact factorized form of  $\mathbf{L}$  quickly. Algorithm 3 computes a compact LDL<sup>\*</sup> decomposition in the form of the tree  $\mathfrak{L}$ , which nodes are labeled by structured matrices of various sizes. We note that this decomposition depends of the tower of proper divisors chosen. Algorithm 3 uses the unique one induced by Definition 6.

Algorithm 3 ffLDL $^{\star}_{\mathcal{R}_d}(\mathbf{G})$ 

**Require:** A full-rank Gram matrix  $\mathbf{G} \in \mathcal{R}_d^{n \times n}$  **Ensure:** The compact LDL\* decomposition of  $\mathbf{G}$  in FFT form 1:  $(\mathbf{L}, \mathbf{D}) \leftarrow \mathsf{LDL}_{\mathcal{R}_d}^*(\mathbf{G})$ 2: if d = 1 then 3: return  $(\mathbf{L}, \mathbf{D})$ 4: end if 5:  $d' \leftarrow \operatorname{gpd}(d)$ 6: for  $i = 1, \ldots, n$  do 7:  $\mathfrak{L}_i \leftarrow \operatorname{ffLDL}_{\mathcal{R}_{d'}}^*(\mathsf{M}_{d/d'}(\mathbf{D}_{ii}))$ 8: end for 9: return  $(\mathbf{L}, (\mathfrak{L}_i)_{1 \leq i \leq n})$ 

Algorithm 3 computes a "fast Fourier LDL<sup>\*</sup>", instead of the "fast Fourier GSO" hinted at in the proof of Corollary 1. The reason why we favor this approach is because it allows a complexity gain. This gain can already be observed in the classic versions of the aforementioned algorithms. Indeed, consider the **L** in the GSO of  $\mathbf{B} \in \mathcal{R}_d^{2\times m}$ , which is exactly the **L** in the LDL<sup>\*</sup> decomposition of  $\mathbf{BB}^* \in \mathcal{R}_d^{2\times 2}$ . Computing it with the LDL<sup>\*</sup> algorithm is then  $\Theta(m)$  times faster than with the Gram-Schmidt process. The same phenomenon happens with their recursive variants.

LEMMA 2. Let  $d \in \mathbb{N}$  and  $1 = d_0|d_1| \dots |d_h| = d$  be the tower of proper divisors of d given by the successive gpd's, and for  $i \in [\![1,h]\!]$ , let  $k_i \stackrel{\Delta}{=} d_i/d_{i-1}$ . Let  $\mathbf{G} \in \mathcal{R}_d^{n \times n}$  be a fullrank Gram matrix. Then Algorithm 3 computes the LDL<sup>\*</sup> decomposition tree of  $\mathbf{G}$  in FFT form in time

$$\Theta(n^2 d \log d) + \Theta(n^3 d) + \Theta(nd) \sum_{1 \le i \le h} k_i^2.$$

In particular, if all the  $k_i$  are bounded by a small constant k, then the complexity of Algorithm 3 is upper bounded by  $\Theta(n^3d + n^2d \log d)$ .

The proof of Lemma 2 can be found in the full version of this article. We note that Algorithm 3 is parallelizable to up to d processes: step 1 relies on operations on polynomials which are parallelizable and all the iterations of step 7 are independent.

## 4. FAST FOURIER NEAREST PLANE

In this section, we show how to exploit further the compact form of the LDL<sup>\*</sup> decomposition to have a fast Fourier variant of the nearest plane algorithm. It outputs vectors of the same quality (*ie* as close to the target vector) as its classical iterative counterpart Algorithm 2, but runs  $\tilde{\Theta}(d)$  times faster.

DEFINITION 9. Let  $\mathcal{Z}_d$  denote the ring  $\mathbb{Z}[x]/(x^d-1)$  of elements of  $\mathcal{R}_d$  with integer coefficients.

#### Algorithm 4 ffNP $_{\mathcal{R}_d}(\mathbf{t}, \mathfrak{L})$

- **Require:**  $\mathbf{t} \in \mathcal{R}_d^n$ , a precomputed tree  $\mathfrak{L}$ , (implicitly) a matrix  $\mathbf{B} \in \mathcal{R}_d^{n \times m}$  such that  $\mathfrak{L}$  is the compact LDL<sup>\*</sup> decomposition tree of  $\mathbf{BB}^*$ .
- **Ensure:**  $\mathbf{z} \in \mathbb{Z}_d^n$  such that  $V((\mathbf{z} \mathbf{t})\mathbf{B}) \in \mathcal{P}(\tilde{\mathbf{B}}_0)$ , where  $\tilde{\mathbf{B}}_0$  is the orthogonalization of  $M(\mathbf{B})$ .
- 1: if d = 1 then 2:  $(\mathbf{L}, \mathbf{D}) \leftarrow \mathfrak{L}$ 3: return  $\mathsf{NP}_{\mathbb{R}}(\mathbf{L}, \mathbf{t})$ 4: end if 5:  $(\mathbf{L}, (\mathfrak{L}_i)_{1 \leq j \leq n}) \leftarrow \mathfrak{L}$ 6:  $d' \leftarrow \operatorname{gpd}(d)$ 7: for  $j = n, \dots, 1$  do 8:  $\overline{\mathbf{t}}_j \leftarrow \mathbf{t}_j + \sum_{i>j} (\mathbf{t}_i - \mathbf{z}_i) L_{ij}$ 9:  $\mathbf{z}_j \leftarrow \bigvee_{d/d'}^{-1} [\operatorname{ffNP}_{\mathcal{R}_{d'}}(\bigvee_{d/d'}(\overline{\mathbf{t}}_j), \mathfrak{L}_j)]$ 10: end for 11: return  $\mathbf{z} = (\mathbf{z}_i, \dots, \mathbf{z}_n)$

LEMMA 3. Let  $\mathbf{B} = {\mathbf{b}_1, \dots, \mathbf{b}_n} \in \mathcal{R}_d^{n \times m}$  and  $\tilde{\mathbf{B}} = {\tilde{\mathbf{b}}_1, \dots, \tilde{\mathbf{b}}_n}$  be its Gram-Schmidt orthogonalization in  $\mathcal{R}_d$ . The vectors  $\mathbf{z}$  and  $\bar{\mathbf{t}} = (\bar{\mathbf{t}}_1, \dots, \bar{\mathbf{t}}_n)$  in Algorithm 4 satisfy

$$(\mathbf{z} - \mathbf{t}) \cdot \mathbf{B} = (\mathbf{z} - \overline{\mathbf{t}}) \cdot \widetilde{\mathbf{B}}.$$

PROOF. We recall that for each  $i \in [\![1,n]\!]$ ,  $\tilde{\mathbf{b}}_i = \mathbf{b}_i - \sum_{j < i} L_{ij} \tilde{\mathbf{b}}_j$ . We have:

$$\begin{aligned} (\mathbf{z} - \overline{\mathbf{t}}) \cdot \tilde{\mathbf{B}} &= \sum_{1 \leq j \leq n} (\mathbf{z}_j - \overline{\mathbf{t}}_j) \tilde{\mathbf{b}}_j \\ &= \sum_{1 \leq j \leq n} \left[ (\mathbf{z}_j - \mathbf{t}_j) + \sum_{i > j} L_{ij} (\mathbf{z}_i - \mathbf{t}_i) \right] \tilde{\mathbf{b}}_j \\ &= \sum_{1 \leq i \leq j \leq n} (\mathbf{z}_i - \mathbf{t}_i) L_{ij} \tilde{\mathbf{b}}_j \\ &= \sum_{1 \leq i \leq n} (\mathbf{z}_i - \mathbf{t}_i) \mathbf{b}_i \\ &= (\mathbf{z} - \mathbf{t}) \cdot \mathbf{B}. \end{aligned}$$

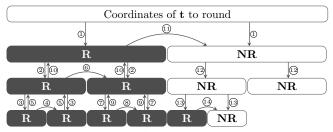
The first and last equalities are trivial, the second one replaces the  $\mathbf{\bar{t}}_j$ 's by their definitions, the third one just simplifies the sum and the fourth one is another way of saying that  $\mathbf{L} \cdot \mathbf{\tilde{B}} = \mathbf{B}$ .

THEOREM 2. Let  $\mathsf{M} \triangleq \mathsf{M}_{d/1}$  and  $\mathsf{V} \triangleq \mathsf{V}_{d/1}$ . Algorithm 4 outputs  $\mathbf{z} \in \mathcal{Z}^n$  such that  $\mathsf{V}((\mathbf{z} - \mathbf{t})\mathbf{B}) \in \mathcal{P}(\tilde{\mathbf{B}}_0)$ , where  $\tilde{\mathbf{B}}_0$  is the orthogonalization of  $\mathsf{M}(\mathbf{B})$  over  $\mathbb{R}$ .

PROOF. The result is trivially true if d = 1. We prove it in the general case. By definition, each subtree  $\mathfrak{L}_j$  is the LDL<sup>\*</sup> decomposition tree of  $\mathsf{M}_{d/\operatorname{gpd}(d)}(\tilde{\mathbf{b}}_j)$ . By induction hypothesis, we therefore know that  $\mathsf{V}((\mathbf{z}_j - \bar{\mathbf{t}}_j)\tilde{\mathbf{b}}_j) \in \mathcal{P}(\tilde{\mathbf{B}}_j)$ , where  $\tilde{\mathbf{B}}_j$  is the orthogonalization of  $\mathbf{B}_j \triangleq \mathsf{M}(\tilde{\mathbf{b}}_j)$ . From Lemma 3, we have

$$(\mathbf{z} - \mathbf{t})\mathbf{B} = \sum_{j=1,...,n} (\mathbf{z}_j - \overline{\mathbf{t}}_j) \cdot \tilde{\mathbf{b}}_j$$

so  $V((\mathbf{z} - \mathbf{t})\mathbf{B}) \in \mathcal{P}([\tilde{\mathbf{B}}_1, \dots, \tilde{\mathbf{B}}_n])$ . Now, from the proof of Corollary 1, we know that  $\tilde{\mathbf{B}}_0 = [\tilde{\mathbf{B}}_1, \dots, \tilde{\mathbf{B}}_n]$  is actually the orthogonalization of  $M(\mathbf{B})$ , which concludes the proof.  $\Box$ 



An ongoing execution of Algorithm 4 for a single  $t \in \mathcal{R}_8$ .

- 1. Arrows are labeled in their order of execution. Downward  $(\downarrow \downarrow)$  (resp. upward  $(\uparrow \uparrow)$ ) arrows correspond to computing  $V_{d/d'}(\bar{\mathbf{t}}_j)$  (resp.  $V_{d/d'}^{-1}(\ldots)$ ) at step 9. Transverse  $(\curvearrowright)$  arrows correspond to step 8.
- 2. Cells labeled with **R** (as in **R**ounded) correspond to already completed subcalls of Algorithm 4, as opposed to those labeled with **NR** (as in **Not R**ounded).

# Figure 5: High-level execution of the fast Fourier nearest plane algorithm

Unlike the fast Fourier transform, Algorithm 4 is not fully parallelizable, due to step 8 ( $\bigcirc$  arrows in Figure 5). However, its complexity in d is  $\Theta(d \log d)$ : informally, this is because each arrow  $\downarrow, \uparrow$  or  $\bigcirc$  has a linear complexity in the size of the cells it connects. A more precise statement follows.

LEMMA 4. Let  $d \in \mathbb{N}$  and  $1 = d_0|d_1| \dots |d_h = d$  be be the tower of proper divisors of d given by the successive gpds, and for  $i \in [\![1,h]\!]$ , let  $k_i \stackrel{\Delta}{=} d_i/d_{i-1}$ . Let  $\mathbf{B} \in \mathcal{R}_d^{n \times m}$  and  $\mathfrak{L}$  be its  $LDL^*$  decomposition tree. The complexity of Algorithm 4 is given by:

$$\Theta(nd\log d) + \Theta(n^2d) + \Theta(nd) \sum_{1 \le i \le h} k_i^2.$$

In particular, if all the  $k_i$  are bounded by a constant, then the complexity of Algorithm 4 is  $\Theta(n^2d + nd\log d)$ .

The proof of Lemma 4 can be found in the full version of this article.

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