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## Lattice Problems and Rounding Algorithms

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### 1 Introduction

In the previous lecture, we established that any lattice  $L$  admits a basis  $\mathbf{B}$ , and furthermore that each basis induces a parallelepiped  $\mathcal{P}(\mathbf{B})$  that is a tiling for  $L$ . We will first show that this tiling can be performed efficiently by an algorithm. This, in turns allows to solve computational problems (such as the approximate closest vector problem) up to a certain approximation factor that depends on the basis. This problem is of theoretical interest in Complexity Theory, but also has very practical applications (error correction, quantization, lossy compression).

We will then show that the same basis  $\mathbf{B}$  induces another parallelepiped that is tiling for  $L$ , namely  $\mathcal{P}(\mathbf{B}^*)$  where  $\mathbf{B}^*$  is the Gram-Schmidt orthogonalization of  $\mathbf{B}$ , and that this is true despite  $\mathbf{B}^*$  not being a basis for  $L$  —an example of this tiling can be viewed in many brick walls, especially in Leiden... An efficient algorithm for the approximate shortest vector problem can also be designed based on this tiling.

The above motivates the rich area of research called basis reduction, which will be studied in the following lectures. Motivating questions include: given that a lattice in 2 or more dimensions can be represented by infinitely many bases, how should you choose one for a specific purpose? How “good” of a basis is guaranteed to exist? Can we find a “good” basis with an algorithm? How costly are those algorithms?

We will finish the lecture with the notion of the Voronoi cell of a lattice, which is in many senses a geometrically optimal tiling of the  $\mathbb{R}$ -span of the lattice. But it is also computationally costly to deal with. In particular, we will relate its quality for the closest vector problem (CVP) directly to the properties of the lattice.

#### 1.1 Notation

As hinted above, the Euclidean norm (a.k.a. the  $\ell_2$  norm) will play a special role in this lecture, and in most lectures involving lattice reduction. We will (do our best to<sup>1</sup>) always specify this restriction by using the notation  $\|\cdot\|_2$ , while  $\|\cdot\|$  denotes an arbitrary norm. The associated inner product is denoted  $\langle \cdot, \cdot \rangle$ . When considering metric properties of a lattice, like the minimal distance or the covering radius, the use of the Euclidean metric will be denote in parentheses in the exponent, e.g.  $\lambda_1^{(2)}(L)$ ,  $\mu^{(2)}(L)$ .

**Algorithms and Complexity for the Mathematician.** This course is written for students with a mathematical background, but the topic really lies at the intersection between mathematics and computer science, and in particular complexity theory. We will do our best to provide the necessary notion of complexity theory as we go, in a progressive manner. This section is an overview of the mindset, ideas, and terminology that we will encounter, which will be made more precise later on.

A **computational problem** is the task of an algorithm: given an **input** guaranteed to have a certain property, the algorithm must produce an **output** with prescribed properties. An algorithm is said to be **correct** for a task if it always terminates and fits the requirement stated by the task.

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<sup>1</sup>This goes against the habits of the lecturer, so mistakes are likely. Please notify such typos.

When describing an algorithm, it will often be listed together with its **specification**, that is, the problem that it solves. Note that an algorithm can be correct for several tasks, and it will often be the case that we give an algorithm with a tight specification, only to show later that it is correct for a less specific task. For example, we will soon see that there is an algorithm (Algorithm 2) to tile the space  $\mathbb{R}^n$  using  $\mathcal{P}(\mathbf{B})$  for some basis  $\mathbf{B}$  of  $L$ . This is a tight specification of the algorithm: the specification dictates a unique valid output for each input. We will prove that this algorithm is correct (Lemma 4) and, as a corollary that it solves the **closest vector problem** up to a certain radius (Corollary 5).

Computational problems and algorithms *are* formal objects, and not some alien object interacting with mathematics. They have several equivalent formal foundations (Church, Turing), though algorithms are often described using "pseudo-code" for convenience and readability. Pseudo-code is no less formal than theorems and proofs as usually written in textbooks and mathematical papers; it is "pseudo" in the sense that it is not machine-readable (in the same way that proofs are not machine-readable unless written in a language like Coq or Lean).

The **complexity** of an algorithm is a measure of how many operations it performs. This is typically given in number of bit operations, but this is sometimes inconvenient. As an intermediate step, we may just count the number of arithmetic operations in an algorithm: the gap between both methods of counting depends on how large the numbers in the computations are. Take Algorithm 1 as an example.

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**Algorithm 1:** Addition Algorithm

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**Input** : An integer  $x$  and an integer  $y$ .

**Output:** An integer  $z$  such that  $z = x + y$

**return**  $x + y$

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The number of arithmetic operations in this algorithm is 1, because the algorithm just adds two numbers and outputs the answer. But the bit complexity is larger, because a computer will add two integers the same way we all learn to in school: add digits in one column, and if the value is greater than ten, *carry the one* to the next columns and continue. In a computer this is all done in base two, so the number of **bit operations** is bounded above by the number of bits required to write the two numbers in binary.

The complexity can be given explicitly in terms of the input, such as the dimension  $n$  of the input matrix  $\mathbf{B} \in \mathbb{Q}^{n \times n}$ . But when we merely say that an algorithm is quadratic, or **polynomial time** without specifying in which variable, it is implicitly meant as a function of the **bitsize of the input**. Explicitly, an algorithm runs in polynomial time if there exists constants  $C, d$  such that for any valid input of bitsize  $K$ , the algorithm outputs a correct answer after no more than  $CK^d$  bit operations. This is more often abbreviated by saying that the running time is  $O(K^d)$ . For an integer output  $x \in \mathbb{Z}$ , the bitsize of the input is  $\log_2 |x| + O(1)$ , while a rational  $x/y \in \mathbb{Q}$  can be represented using two integers  $x$  and  $y$ , and an  $n \times k$  matrix can be represented using  $nk$  inputs, etc.

Not only can we write theorems about algorithms (correctness, complexity) and quantify over classes of algorithms, but we can also formally relate computational problems to one another. In particular, complexity theory is often interested in showing that a computational problem  $A$  is "easier" or "not much harder" than a computational problem  $B$ . Or, one can say equivalently that computational problem  $A$  is **reducible** to computational problem  $B$ . This means that one

can solve an instance of problem  $A$  by (cheaply) transforming it to be solvable in one or more instances of problem  $B$ . When designing such a **reduction**, the unknown algorithm for solving  $B$  is called an **oracle** for  $B$ : we only assume that it solves  $B$ , but we do not know how it does it.

These type of statements are also central in cryptography. We will want to convincingly argue for or against the security of a cryptosystem, which is the practical (im)possibility of breaking it. A cryptosystem might be complex; it might be interactive between the sender and receiver; or there may even be many parties involved. But by reducing its (formally defined) security to a simple and non-interactive computational problem, we can better focus the effort of attempting attacks (cryptanalysis) and be reassured about its security.<sup>2</sup>

## 2 Lattice Computational Problems

There are many (many!) lattice problem variants that are of interest, and we should not list them all here in full detail. We focus on the most important ones.

**DEFINITION 1 (EXACT SVP AND  $\alpha$ -SVP)** *The approximate Shortest Vector Problem with approximation factor  $\alpha \geq 1$ , denoted  $\alpha$ -SVP is defined as follows:*

- Given as input the basis  $\mathbf{B}$  of a lattice  $L$
- Output a short non-zero vector  $\mathbf{v} \in L \setminus \{\mathbf{0}\}$ , satisfying  $\|\mathbf{v}\| \leq \alpha \cdot \lambda_1(L)$ .

The Exact version of the problem, denoted SVP, is the special case  $\alpha = 1$ .

Note that  $\lambda_1(L)$  is not necessarily known, which can be inconvenient. In particular we cannot even efficiently verify that a given solution is indeed correct. For this reason, some variations of these problems are sometime considered, such as  $\alpha$ -Hermite SVP, or  $\alpha$ -HSVP. Here, shortness is defined relative to the determinant:  $\|\mathbf{v}\| \leq \alpha \cdot \det(L)^{1/k}$  where  $k$  is the rank of  $L$ . The factor  $\det(L)^{1/k}$  makes the problem invariant by rescaling  $L \mapsto s \cdot L$  for  $s > 0$ . Or it might be sometimes convenient to simply state an absolute bound for the desired vector (AbsSVP).

Note further that, even for  $\alpha = 1$ , the solution is never unique since  $\|-\mathbf{v}\| = \|\mathbf{v}\|$ . But even up to flipping the sign, there may be many solutions. Take for example the lattice

$$L = \{(x_1, \dots, x_n) \in \mathbb{R}^n : x_i \in \mathbb{Z} \ \forall i\}.$$

This lattice has  $2n$  shortest vectors. Meanwhile, in dimension  $n = 24$  there is a lattice with 196560 shortest vectors. Therefore we should always speak of *a* shortest vector rather than *the* shortest vector.

**DEFINITION 2 (EXACT CVP AND  $\alpha$ -CVP)** *The approximate Closest Vector Problem with approximation factor  $\alpha \geq 1$ , denoted  $\alpha$ -CVP is defined as follows:*

- Given as input the basis  $\mathbf{B}$  of a lattice  $L \subset \mathbb{R}^n$  and a target  $\mathbf{t} \in \mathbb{R}^n$
- Output a vector  $\mathbf{v} \in L$  satisfying  $\|\mathbf{v} - \mathbf{t}\| \leq \alpha \cdot d(\mathbf{t}, L)$

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<sup>2</sup>And yet, the saying is that “Cryptographers seldom sleep well at night”. In particular, if  $P = NP$ , then cryptography as we know it would entirely collapse. In fact, cryptography relies on the stronger and more specific assumption than  $P \neq NP$ .

where  $d(\mathbf{t}, L) := \min_{\mathbf{x} \in L} \|\mathbf{x} - \mathbf{t}\|$  denotes the distance to the lattice. The Exact version of the problem, denoted CVP, is the special case  $\alpha = 1$ .

For the same reasons as for SVP, variants of this problem where the norm is bounded in absolute terms is also convenient (AbsCVP). Lastly, we define a restriction of the Exact CVP problem, where this time we are also given a guarantee that the solution is particularly close to the lattice, and therefore, unique.

**DEFINITION 3 ( $\alpha$ -BDD)** *The Bounded Distance Decoding Problem with approximation factor  $\alpha \leq 1/2$ , denoted  $\alpha$ -BDD is defined as follows:*

- Given as input the basis  $\mathbf{B}$  of a lattice  $L \subset \mathbb{R}^n$  and a target  $\mathbf{t} \in \mathbb{R}^n$  such that  $d(\mathbf{t}, L) < \alpha \lambda_1(L)$
- Output the unique closest vector  $\mathbf{v} \in L$  satisfying  $\|\mathbf{v} - \mathbf{t}\| < \alpha \lambda_1(T)$

where  $d(\mathbf{t}, L) := \min_{\mathbf{x} \in L} \|\mathbf{x} - \mathbf{t}\|$  denotes the distance to the lattice.

Once again, an absolute version (AbsBDD) will also be convenient to consider.

**Close Vectors and Tiling.** Many algorithms of interest for CVP and BDD are related to  $L$ -tilings. That is, for a subset  $T \subseteq \text{Span}_{\mathbb{R}}(L)$  that is  $L$ -tiling, an algorithm is said to *solve  $T$ -tiling* if, when given an input basis  $\mathbf{B}$  and target  $\mathbf{t}$ , the algorithm outputs a CVP/BDD vector  $\mathbf{v} \in L$  such that the error  $\mathbf{e} = \mathbf{v} - \mathbf{t}$  is in  $T$ .

In this case, we can state that the algorithm solves CVP and BDD up to radii that depend on the following geometric properties of  $T$ . For any bounded body  $T \subset \mathbb{R}^n$ , we define its inner radius  $\nu(T)$  and its outer radius  $\mu(T)$  by

$$\nu(T) := \sup\{r \in \mathbb{R} \mid r \cdot \mathfrak{B} \subset T\}, \quad \mu(T) := \inf\{r \in \mathbb{R} \mid T \subset r \cdot \mathfrak{B}\}. \quad (1)$$

Alternatively, the above may be re-written as:

$$\nu(T) := \inf_{\mathbf{x} \in \mathbb{R}^n \setminus T} \|\mathbf{x}\|, \quad \mu(T) = \sup_{\mathbf{x} \in T} \|\mathbf{x}\|. \quad (2)$$

If a close vector algorithm is associated with a tiling  $T$ , then it correctly solves:

- AbsBDD up to a radius  $\nu(T)$
- AbsCVP up to a radius  $\mu(T)$
- $\alpha$ -CVP up to an approximation factor  $\alpha = \mu(T)/\nu(T)$ .

Note that the inner and outer radius of tilings are bounded by the metric properties of the lattice: for any tiling  $T$  of  $L$ , it must be that  $\mu(T) \geq \mu(L)$  and  $\nu(T) \leq \lambda_1(L)/2$ .

### 3 The Simple Rounding algorithm and the $\mathcal{P}(\mathbf{B})$ Tiling

The Simple Rounding algorithm merely rounds the coordinate of the target vector when it is expressed in terms of the basis  $\mathbf{B}$ . One can understand the algorithm as a reduction to solving CVP in the trivial lattice  $\mathbb{Z}^n$ : multiply by  $\mathbf{B}^{-1}$  which sends  $L$  to  $\mathbb{Z}^n$ , solve CVP on the lattice  $\mathbb{Z}^n$  by coordinate-wise rounding to the nearest integer, and then translate the solution back to  $L$  by multiplying it by  $\mathbf{B}$ .

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**Algorithm 2:** SimpleRounding( $\mathbf{B}, \mathbf{t}$ ): Simple Rounding Algorithm

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**Input** : A basis  $\mathbf{B} \in \mathbb{Q}^{n \times n}$  of a full rank lattice  $L$ , a target  $\mathbf{t} \in \text{Span}_{\mathbb{R}}(L)$ .

**Output:**  $\mathbf{v} \in L$  such that  $\mathbf{e} = \mathbf{t} - \mathbf{v} \in \mathcal{P}(\mathbf{B})$

$\mathbf{t}' \leftarrow \mathbf{B}^{-1} \cdot \mathbf{t}$

$\mathbf{v}' \leftarrow (\lfloor t'_i \rfloor)_{i \in \{1 \dots n\}}$

$\mathbf{v} \leftarrow \mathbf{B} \cdot \mathbf{v}'$

**return**  $\mathbf{v}$

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LEMMA 4 *Algorithm SimpleRounding is correct and runs in polynomial time.*

PROOF: Let us start with correctness. Because  $L$  is full-rank, its basis  $B$  is invertible, so the first line of the algorithm is well-defined. By construction,  $\mathbf{v}'$  belongs to  $\mathbb{Z}^n$ , and because  $\mathbf{B}$  is a basis of  $L$ ,  $\mathbf{v} = \mathbf{B}\mathbf{v}'$  belongs to  $L$ . Now define  $\mathbf{e}' := \mathbf{t}' - \mathbf{v}'$ , which by construction belongs to  $[-1/2, 1/2]^n$ . Note that  $\mathbf{e} = \mathbf{B} \cdot \mathbf{e}'$ , so by definition of  $\mathcal{P}(\mathbf{B}) = \mathbf{B} \cdot [-1/2, 1/2]^n$  we have  $\mathbf{e} \in \mathcal{P}(\mathbf{B})$ .

For proving polynomial time complexity, the main issue is inversion of the matrix  $\mathbf{B}^{-1}$ . The Gaussian elimination process requires  $O(n^3)$  arithmetic operation, but one must also show that the numerators and denominators of the rational numbers that occur in the intermediate steps remain small enough.  $\square$

COROLLARY 5 *For  $\mathbf{B}$  a basis of  $L$ , the algorithm SimpleRounding( $\mathbf{B}, \cdot$ ) solves  $\mu(\mathcal{P}(\mathbf{B}))$ -AbsCVP and  $\nu(\mathcal{P}(\mathbf{B}))$ -AbsBDD.*

**Inner and Outer Radii of  $\mathcal{P}(\mathbf{B})$ .** The outer radius  $\mu(\mathcal{P}(\mathbf{B}))$  admits a natural upper bound by direct application of the triangle inequality:

$$\mu(\mathcal{P}(\mathbf{B})) \leq \frac{1}{2} \sum_{i=1}^n \|\mathbf{b}_i\|. \quad (3)$$

In the particular case of the Euclidean norm, this bound is never reached, that is  $\mu^{(2)}(\mathcal{P}(\mathbf{B})) < \frac{1}{2} \sum_{i=1}^n \|\mathbf{b}_i\|_2$  because an equality instance of the triangle inequality occurs when vectors are collinear, but the vectors of a basis can never be collinear. However, one can get arbitrarily close to it, for example

$$\mathbf{B} = \begin{pmatrix} 1 & 1 & 1 & \dots & 1 \\ 0 & \varepsilon & 0 & \dots & 0 \\ 0 & 0 & \varepsilon & \dots & 0 \\ \vdots & & & \ddots & \\ 0 & 0 & 0 & \dots & \varepsilon \end{pmatrix} \quad \text{gives} \quad \frac{\mu^{(2)}(\mathcal{P}(\mathbf{B}))}{\frac{1}{2} \sum_{i=1}^n \|\mathbf{b}_i\|_2} \geq \frac{n}{n\sqrt{1+\varepsilon^2}} \rightarrow 1 \text{ as } \varepsilon \rightarrow 0, \quad (4)$$

On the contrary, the same example with any  $\varepsilon \in (-1, 1)$  shows that this upper bound can be reached for the  $\ell_\infty$  norm.

Similarly, one would like a lower bound on the inner radius  $\nu(\mathcal{P}(\mathbf{B}))$ . This turns out to be more challenging for general norms. However, it can be exactly determined for the Euclidean norm. Let us start with a characterization of the parallelepiped  $\mathcal{P}(\mathbf{B})$  which allows us to describe a parallelepiped by a system of linear inequalities rather than by a spanning basis.

LEMMA 6 *Let  $\mathbf{B} \in \text{GL}_n(\mathbb{R})$  and  $\mathbf{C} = (\mathbf{B}^{-1})^\top$  be its inverse-transpose. Then,*

$$\mathcal{P}(\mathbf{B}) = \left\{ \mathbf{x} \in \mathbb{R}^n \mid \mathbf{c}_i^\top \cdot \mathbf{x} \in [1/2, 1/2] \text{ for all } i \in \{1, \dots, n\} \right\}.$$

PROOF: Recall that  $\mathcal{P}(\mathbf{B}) = \mathbf{B} \cdot [-1/2, 1/2]^n$  and rewrite the condition  $\mathbf{x} \in \mathcal{P}(\mathbf{B})$  as  $\mathbf{B}^{-1} \cdot \mathbf{x} \in [-1/2, 1/2]^n$ . Each coordinate  $y_i$  of  $\mathbf{y} = \mathbf{B}^{-1}\mathbf{x}$  writes  $\mathbf{c}_i^\top \cdot \mathbf{x} \in [-1/2, 1/2]$  for all  $i \in \{1, \dots, n\}$ .  $\square$   
 From there, we obtain an explicit

LEMMA 7 *Let  $\mathbf{B} \in \text{GL}_n(\mathbb{R})$  and  $\mathbf{C} = (\mathbf{B}^{-1})^\top$  be it's inverse-transpose. Then*

$$v^{(2)}(\mathcal{P}(\mathbf{B})) = \min_i \frac{1}{2 \cdot \|\mathbf{c}_i\|_2}.$$

PROOF: Recall that  $v(T) := \inf_{\mathbf{x} \in \mathbb{R}^n \setminus T} \|\mathbf{x}\|$ . Note that for the canonical inner product  $\langle \cdot, \cdot \rangle$  of  $\mathbb{R}^n$ ,  $\mathbf{c}_i^\top \cdot \mathbf{x} = \langle \mathbf{c}_i, \mathbf{x} \rangle$ . By Cauchy-Schwarz and the above characterization, it holds that any  $\mathbf{x}$  of norm strictly less than  $\min_i \frac{1}{2 \cdot \|\mathbf{c}_i\|_2}$  belongs to  $\mathcal{P}(\mathbf{B})$ . On the contrary, the vector  $\mathbf{x} = \frac{\mathbf{c}_i}{2\|\mathbf{c}_i\|_2^2}$  satisfies  $\mathbf{c}_i^\top \cdot \mathbf{x} = 1/2$  and therefore does not belong to  $\mathcal{P}(\mathbf{B})$ , and this vector has norm  $\frac{1}{2 \cdot \|\mathbf{c}_i\|_2}$ . Minimizing over  $i$  yields the claim.  $\square$