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Height map compression techniques

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Abstract: The goal of this thesis is to design a suitable method for lossy compression of heightmap terrain data. This method should accept blocks of float samples of dimensions $2^n \times 2^n$ as an input, for which it should be able to perform progressive mip-maps (progressive lower-resolution representations) decompression. For every mip-map, it should keep the reconstructed data within a certain maximum absolute per-sample error bound in the unit of meters adjustable by the user. Given these constraints, it should be as efficient as possible. Our method is inspired by the second generation of progressive wavelet-based compression scheme modified to satisfy the maximum-error constraint. We simplified this scheme by factoring out unnecessary computations in order to improve the efficiency. Our method can compress a 256×256 block in about 30 ms and decompress it in about 1 ms. Thanks to these attributes, the method can be used in a real-time planet renderer. It achieves the compression ratio of 37:1 on the whole Earth 90m/sample terrain dataset transformed and separated into square blocks, while respecting the maximum error of 5m.

Keywords: heightmap, lossy, compression, mip-map, guaranteed maximum error bound

Dedication.

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Introduction

A planetary renderer based on real (non-procedural) data has to work with huge amounts of information. This includes its storage and distribution to both the developers working on the application and the users of it. Data compression can be very beneficial in this situation - it can greatly reduce all the storage demands. This results in decreased network and disk utilization - the data needed to transmit via network becomes smaller, so the transmission is quicker. Similarly for the rendering, the data which has to be fetched from disk into RAM is smaller which decreases the transfer time, often being one of the major real-time performance bottlenecks.

In this kind of renderer, terrain information forms a significant subset of the utilized data. This is basically the information about the height of the terrain at all places of the planet. To keep the size of the data reasonable, the terrain height is usually discretely sampled in a regular square grid of some density (eg. 90m per sample). Each one of these height samples is of the same common form - usually a float number (4 bytes), the accuracy of which is sufficient enough for usual planets (eg. Earth). This thesis focuses on compression of this kind of data.

Classical image compression techniques (png, jpeg) [18, 14] are not suitable for this purpose, because they are designed to compress four 1-byte channels per pixel. They benefit from patterns and similarities of the channel values among adjacent pixels. Putting a 4-byte float height value into four channels directly does not result in anything meaningful, from which these methods might benefit - even with a slight change of the height value, all the four channels, into which it has been put, can change significantly. This is why they do not achieve good compression ratios when used this way. Of course, this might be sorted out by a more sophisticated translation of the height value into the channels, but this approach would still not be able to control the maximum absolute height error of the compressed data which is one of the attributes our compression should have. Even though JPEG 2000 offers a similar feature, it is only able to control the per-channel maximum error, which is impossible to translate into the height error, as a single height value would be decomposed into four channels. Thus, heightmap compression requires a different approach - the heights have to be handled directly as float values and similarity or closeness of adjacent height values has to be exploited.

Generally, every application designed to render huge scenes must implement some kind of multiresolution (LOD-ing) approach in order to keep the rendering frame rates reasonable. LOD is the abbreviation of level of detail - degradation of quality of the displayed data with the growing distance in order to optimize the rendering. A planetary renderer certainly belongs to this group of applications, so, among the other things, it must somehow ensure that the further the terrain is from the camera, the less detailed it is. With the growing distance, the amount of observable detail decreases anyway, so displaying it would be just an unnecessary performance burden.

Such LOD-ing approach is almost always implemented by a multiresolution tree hierarchy of nodes [16]. Usually, all these nodes have the same form - triangles, squares, multipolygons, etc. The leaf (bottom-level) nodes of this hierarchy

are the partitions of the original terrain data and the other non-leaf nodes are gradual simplifications of the leaf nodes - the higher the level a node belongs to (the further it is from the leaf level), the coarser representation of data it presents. This hierarchy gives a very good basis for the subsequent multiresolution rendering of scene - given a camera position, we determine which nodes will be rendered on the basis of their distance from the camera, using an appropriate metrics. This choice forms a cut in the hierarchy, with its lowest point at the finest node - the one closest to the camera. The height of this cut increases with growing distance from the camera.

To prepare a terrain LOD hierarchy, several transformations of the input terrain data are required - projecting it from the geographic coordinates into the space of the LOD hierarchy followed by splitting it into its basic partitions forming the leaf nodes of the hierarchy, which are then iteratively grouped into coarser parent nodes in order to build the whole multiresolution hierarchy. The nodes of this hierarchy from subsequent levels might either be stored completely independently from each other or dependently - for the children of a certain node, we remember just the data which is required to reconstruct the children from their parent. The independent approach introduces significant data redundancy - every non-leaf level of the hierarchy stores independently again the same terrain as the following finer level, just in a coarser form. This results in the greater size of the resulting data hierarchy prepared for rendering. On the other hand, this approach makes the rendering quicker - in order to fetch a certain node, we do not have to fetch the data of all its ancestors. In practice, both approaches are used. The application, into which this compression method should be plugged, uses the first approach - every node of its LOD hierarchy is a square, stored completely independently from the others. The method designed by us should be able to compress such a square, not interfering with the LOD hierarchy of the application at all - not trying to remove the redundancy caused by independent storage of its nodes in any way, because, as it has been said, it is beneficial in some aspects.

More precisely, the aim of this thesis is to review the existing heightmap compression methods and either find or come up with a method which is able to compress a regular square of float terrain samples the dimension of which is 2^n as efficiently as possible, while enabling subsequent real-time progressive decompression of its data from the coarsest to the finest mip-map. Mip-mapping is a standard method for multiresolution representations of textures. Mip-maps are gradually smaller and coarser squares with the dimensions $2^{n-1..0}$ representing the same data [11]. Usually, every pixel inside a certain mip-map is the average of the four corresponding children pixel inside the previous larger mip-map. This averaging can be performed in the domain of colors, but in our case, it should be performed in the domain of heights. The decompression should be as fast as possible. We should explore the possibility to perform the decompression on GPU mainly in order to save RAM. We should keep in mind that from among the mentioned optimization objectives, the resulting size of compressed data is the most important one. The maximum absolute per-sample deviation of the compressed data must be controllable by the user. No rendering of data has to be handled, as it is supposed that the renderer using this method will handle this.

The mip-maps which the decompression should be able to provide are in fact

a form of level of detail technique too, but it must be made clear that the ability to provide simplified multiresolution representations of every square of the LOD hierarchy is just a tiny part of the multiresolution rendering pipeline, a LODing terrain engine can certainly not be build solely on this ability. The application using the desired method should be able to traverse its multiresolution hierarchy of square nodes on itself in order to render a scene. After selecting which squares should be displayed (the lower the distance, the more detailed the displayed square), the application then should decide for each of the squares which mip-map of it will be displayed. Thus, the mip-maps present a less significant LOD concept inside the greater LOD concept - the multiresolution squares hierarchy. The mip-map selection can be based on the screen-space area of the square in order to reduce the terrain aliasing. For example, when looking at a certain terrain square from a side, a coarser mip-map of it should be chosen than in case we look at it from the top.

In Chapter 1, we list the methods which contain heightmap compression and form a conclusion to get inspired by the compression used inside one of them named C-BDAM in order to build our own method. In Chapter 2, we briefly describe the core component of compression in many of the mentioned papers, including C-BDAM - wavelet transform. Following this, we compare the basic wavelet approaches to the compression inside C-BDAM and our method. Lastly, we briefly describe the basic methods of lossless compression which are often utilized in wavelet compression. In Chapter 3, we briefly describe the basic outline of our method. In Chapter 4, we describe the details of the method. In Chapter 5, we compare the core algorithm of this method to the algorithm of C-BDAM. We present the results in Chap 6. In Chapter 7, we discuss these results along with the possibility how the decompression could be put onto GPU in order to spare RAM. So far, we have not found a way how this could be done.

1. Related works

A survey paper by Pajarola et al. [16] summarizes the best known multiresolution terrain rendering methods. All these methods also handle the rendering supported by their own LOD-ing hierarchies. We only looked at those of them which contain data compression as a sub-task. Some of them are designed to render just a flat area [10, 15, 7, 12], others are able to render the whole planet [3, 9]. In the rest of this chapter, we will briefly describe the methods which contain terrain data compression

C-BDAM¹ [10] and P-BDAM² [3] by Gobbetti et al. turned out to be the most promising ones after comparing with the rest of the methods. Both these methods handle the LOD rendering too and perform wavelet-based [2] data compression in the refinement of a node of their LOD hierarchy. Once the values of a certain node are known, they are used to predict the values of its children as accurately as possible. After that, the differences between these predictions and the real values are computed. These are called residuals. With the help of them, the real values can be restored with absolute accuracy. However, the residuals are then quantized to achieve better compression ratio which means that the compression is lossy. Then, they are losslessly compressed by an entropy codec [19, 13]. Both these methods are able to compute the residuals in the way which ensures that the error of the reconstructed data is kept within a maximum error bound adjustable by the user in every node of their LOD hierarchy. This can be achieved by a slight modification of the second-generation wavelet lifting scheme [20]. C-BDAM is designed to render just a flat portion of terrain, whereas P-BDAM is just C-BDAM modified to be able to render a whole planet. These modifications do not include any improvements to the efficiency of the compression, so, from our point of view, it is sufficient to know just C-BDAM from these two methods. What makes C-BDAM the most interesting are two aspects: the outstanding compression ratio achieved and the ability to respect a certain user-set maximum per-sample error bound (the ratio of 64:1 on the whole planet - 58GB with 90m resolution - with 16m maximum height deviation).

Wavelet tiled pyramids by Ricardo Olanda [15] is another method for rendering a flat portion of terrain. This method contains data compression based on the same principle - the residuals needed to reconstruct the children of a square node of the terrain LOD hierarchy are compressed. The computation of residuals is based on the wavelet-based JPEG2000 standard. This method is not able to reconstruct the data within a certain maximum-error bound which makes it almost not interesting to us at all. Due to this fact, we cannot judge its compression ratios, as they are not connected with any maximum-error bound. Besides, the visual artifacts between adjacent nodes of different LODs are not handled by its rendering pipeline, but it is not our concern anyway.

Apart from that, we found several methods dedicated to compressing triangulated terrain representations [7, 12, 9]. Two of them [7, 12] perform decompression on GPU and are also able to keep the compressed data within a certain maximum deviation adjustable by the user. However, they add more complexity to

¹Compressed Batched Dynamic Adaptive Meshes

²Planet Sized Batched Dynamic Adaptive Meshes

the task as they allow for greater irregularity of their input data. Unlike the previous methods which work with regularly distributed height samples, the triangle meshes these methods compress can be irregular which introduces the need to handle more spatial information. Due to this, the compression ratios they achieve are all way lower than the ones of the previous methods. Our method is expected to work with just regularly distributed height samples - square mip-maps, so we do not see as much potential in getting inspired by compression methods which have to handle much more complicated data which results in worse efficiency and greater complexity.

HFPaC [8] by Durdevic and Tartalja is the only method we found which focuses solely on heightmap compression, which looked promising at first. Moreover, it is GPU-based which can bring greater performance. However, it has several severe limitations. It is near-lossless, but the maximum deviation of the reconstructed data cannot be controlled by the user. Quite to the contrary, it varies on the basis of both the characteristics of the input and the settings of its internal parameters, such as the size of its compression block. Even though the authors of the method observed that this deviation is relatively small most of the time (up to 2 meters), this is undoubtedly not exactly what we want, as our method should be able to support any arbitrary maximum error set by the user. In addition, this method is only able to provide exactly three layers of resolution of the compressed data, but our method should be able to provide a larger number of mip-maps which varies depending on the dimension of the input. It would not be easy to tailor this method to our needs, because it is really complicated. Additional modifications made to it could result in its even greater complexity and possible lack of efficiency compared to the other simpler available methods which are already able to respect a maximum deviation adjustable by the user and to perform multi-level progressive decompression.

Because terrain exhibits some fractal characteristics [1], we also read a survey about fractal compression [17] by Saupe et al. However, such methods are quite complicated and designed to compress the whole data at once without any possibility of progressive decompression. More precisely, it might be possible to introduce the progression there, but to our best knowledge, no one has found the way yet and we did not see any straightforward option to do this either. Additionally, the aspect of data this family of methods benefits from the most is its regularity or at least certain repeating. This can be best exploited when compressing huge amounts of data. However, our method should be able to compress and decompress just one small square of height samples completely independently from the others and we cannot benefit from much repetition on such a small area. All in all, with respect to the planned usage of the method, the path of fractal compression seemed not very perspective, but it might be interesting for further research.

At the end, we decided to rather stick to the more promising options - the methods which support multiresolution approach by definition. When designing our method, we decided to start off the ideas behind the compression inside C-BDAM. The main reasons are that C-BDAM achieves the best compression ratios and is able to respect a certain maximum-error bound adjustable by the user which is exactly what we needed, too. As we already mentioned in the beginning, the size of the compressed data was the most important constraint,

which is why we chose to follow the best compression ratio. The compression inside C-BDAM is relatively simple which makes it easy to understand and also fast when implemented. However, C-BDAM contains features which we did not need, so we omitted them and just isolated the compression and tailored it for our needs.

C-BDAM performs the compression inside the refinement of a node of its LOD hierarchy. Due to the request that our method should be able to decompress the mip-maps progressively, it seemed convenient to put the compression in our method inside the refinement of a mip-map. This way, only the residuals needed to reconstruct the following finer mip-map would have to be fetched in order to perform a single step of refinement. So, instead of a LOD node in C-BDAM, we put a mip-map in our method. Additionally, we significantly simplified the compression equations performed in C-BDAM in order to increase the efficiency and speed of our method, while still being able to satisfy the required maximum absolute error bound constraint. However, let it be repeated again that a LOD node in C-BDAM is actually the basic component of its multiresolution rendering hierarchy, whereas a mip-map in our method not, it is only a side component of the LOD hierarchy of the application into which our method should be plugged. The main component of the multiresolution rendering hierarchy of this application is the square which should be compressed by this method, so our method lies outside this hierarchy. However, it is not hard to utilize the compression in C-BDAM inside our method, because the principle remains the same in both cases - we refine a set of height samples to a more detailed one.

2. The preliminaries

This chapter consists of three sections. In the first one (2.1), we will briefly and formally describe the main principle and usage of second generation wavelet transformation methods which are relevant for this thesis. In the second one (2.2), we will compare the compression inside C-BDAM and our method to these methods. Even though C-BDAM is based on the same principle, it differs from these methods a bit, so we will describe the basic differences. Then we will perform the same basic comparison with our proposed method. Our method differs from the described wavelet scheme a bit more, see Section 4.2 for the details.

2.1 The introduction to second-generation wavelets

Basically, there are two generations of wavelets. The first generation uses dilated and translated wavelet function [2] for computation. The second one uses filter banks to perform high-pass and low-pass filtering [4]. The computational equivalency of these two approaches has been proven [6]. For this work, the second generation of discrete wavelet transform methods is most relevant, so we will briefly describe it in this section in order to give the reader an idea of the wavelet concept which is referred to in many places of this thesis. We will base this description on 1-dimensional input data. We will also briefly describe how the transition of the principle to 2-dimensional data, such as images, can be performed.

Every method of this generation consists of just several subsequent applications of lifting onto the input. The lifting is the basic step of the method. It splits the set of its input signal samples into two parts - low-pass (the low frequency information) and high-pass (residuals, the high frequency information). The lifting is firstly applied on the input set of signal samples and then is recursively applied to the low-pass part produced by the previous iteration until the length of the latest low-pass part is 1. In order to make this recursion possible, the count of samples of the original input of the method must be a certain power of two. If the length of the input is 2^n , the method performs n iterations of lifting. The described successive application of lifting on smaller and smaller input is called the bottom-up pass. We can imagine this as building a pyramid of low-pass outputs the first tier of which is the input itself and every following higher tier is the low-pass output of lifting applied to the tier right below. Every tier is half the width of the previous one and after the bottom-up pass, the highest tier has the width of 1 (Fig. 2.1).

After this bottom-up pass, we can perform the inverse top-down pass. This pass does not know how the produced pyramid looks, it only knows its highest tier, sized 1. However, it is supposed to be able to progressively reconstruct the whole pyramid from the top to the bottom, only utilizing the knowledge of the high-pass information (Fig. 2.1). Producing a certain tier from the previous one is called the reconstruction which is the exact opposite of lifting.

Now, let us briefly describe how this principle can be extended to 2-dimensional image data. Basically, this extension introduces just one change to the described principle - when the lifting is applied to an image with dimension 2^n , it produces four equally-sized outputs with dimension 2^{n-1} : one low-pass part and three high-

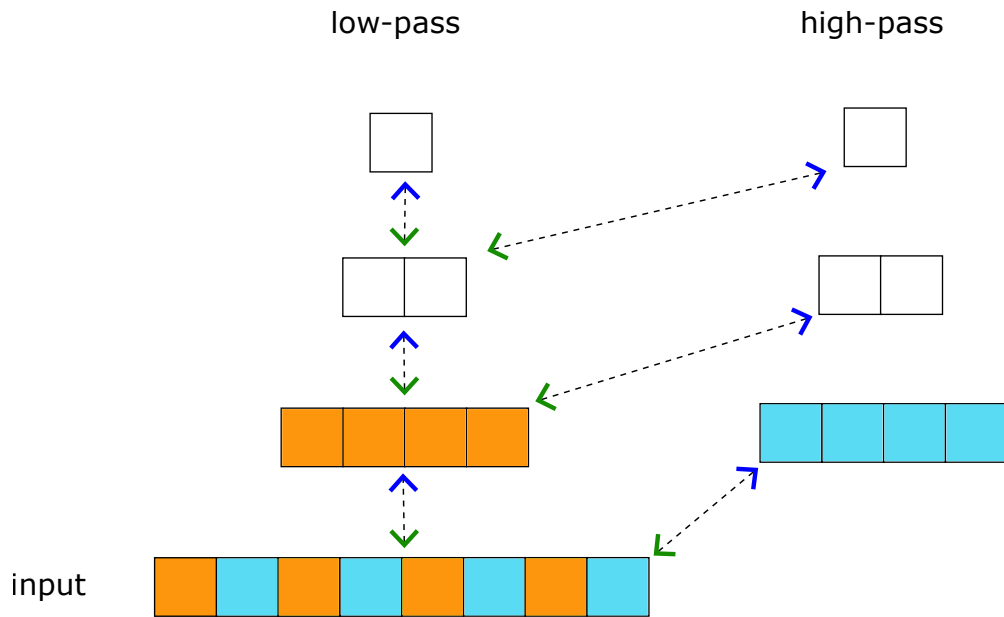


Figure 2.1: An example of recursive wavelet decomposition of 1-dimensional data of samples. On the left side, there is a pyramid of low-pass tiers, the bottom tier of which is the input itself. On the right side, there is the corresponding pyramid of high-pass tiers. The dashed lines represent both the decomposition (lifting) inside the first bottom-up pass (when following the blue arrows) and the inverse compositions (reconstruction) inside the second top-down pass (when following the green arrows). On the input samples, it is also marked which samples form the smaller low-pass part (the orange ones) and the smaller high-pass part (the blue ones). The orange ones are the even ones, as the indexing of samples is supposed to start at zero.

pass parts (Fig. 2.2). Each of these high-pass parts contains different information - for example, the first one might contain vertical edges, the second one horizontal edges and the third one diagonal edges. During the subsequent reconstruction, all these four parts are used to produce the larger low-pass part with twice better resolution.

What makes these decompositions and backward compositions useful is the fact that the top-down pass just needs the high-pass information (residuals) to fully reconstruct the input. This information tends to be sparse and input-dependent - the smoother the input, the less high-pass information it contains. If we compress it well, we can save much storage space. Thus, if we want to store a set of samples the count of which is a power of two in as little space as possible, we will not store the samples directly, but we will store just the compressed residuals produced by the successive iterations of lifting applied to the input. If we are not required to accurately reconstruct the input, we can even decimate (quantize) the residuals. Because this information often contains just details, its careful decimation does not deform the reconstruction much and ensures better compression ratio. One more interesting fact is that the residuals bound to lower tiers of the pyramid carry finer details than those bound to the higher ones. Thanks to this, the more-detailed (larger) sets of residuals can be compressed more aggressively than the less-detailed (smaller) ones. This is called progressive compression and it is used for example in JPEG standard [18].

In the following lines, we will describe the lifting and reconstruction steps in 1-dimensional domain more formally. Let us say that the lifting is given the input samples x_k . It splits them into the even ones: $x_{2k} = x_e$ and the odd ones: $x_{2k+1} = x_o$. This splitting is not yet based on any frequency properties of the samples, it is based just on their order. However, these two sets of samples will subsequently be modified, so that the even ones will contain the low-pass information and the odd ones will become the residuals - the high-pass information. This will be performed with the help of two operators: the prediction operator P and the update operator U . P will be used to produce the residuals d from x_o and U will be used to produce the low-pass part s from x_e .

Up to this point, just the common properties of the second-generation methods have been described. Now will come the differences between them. The only thing they differ in is the way they perform lifting and reconstruction. The way the lifting step is performed clearly determines the way how the reconstruction is performed, as the reconstruction must be the exact inverse of lifting. The lifting step varies in the order in which the operators P and U are applied. According to this, the methods can be split into two main groups - the prediction-first ones and the update-first ones.

In the prediction-first methods, the prediction is applied first:

$$\begin{aligned} d &= x_o - P(x_e) \\ s &= x_e + U(d) \end{aligned}$$

The reconstruction must be the exact inverse:

$$\begin{aligned} x_e &= s - U(d) \\ x_o &= d + P(x_e) \end{aligned}$$

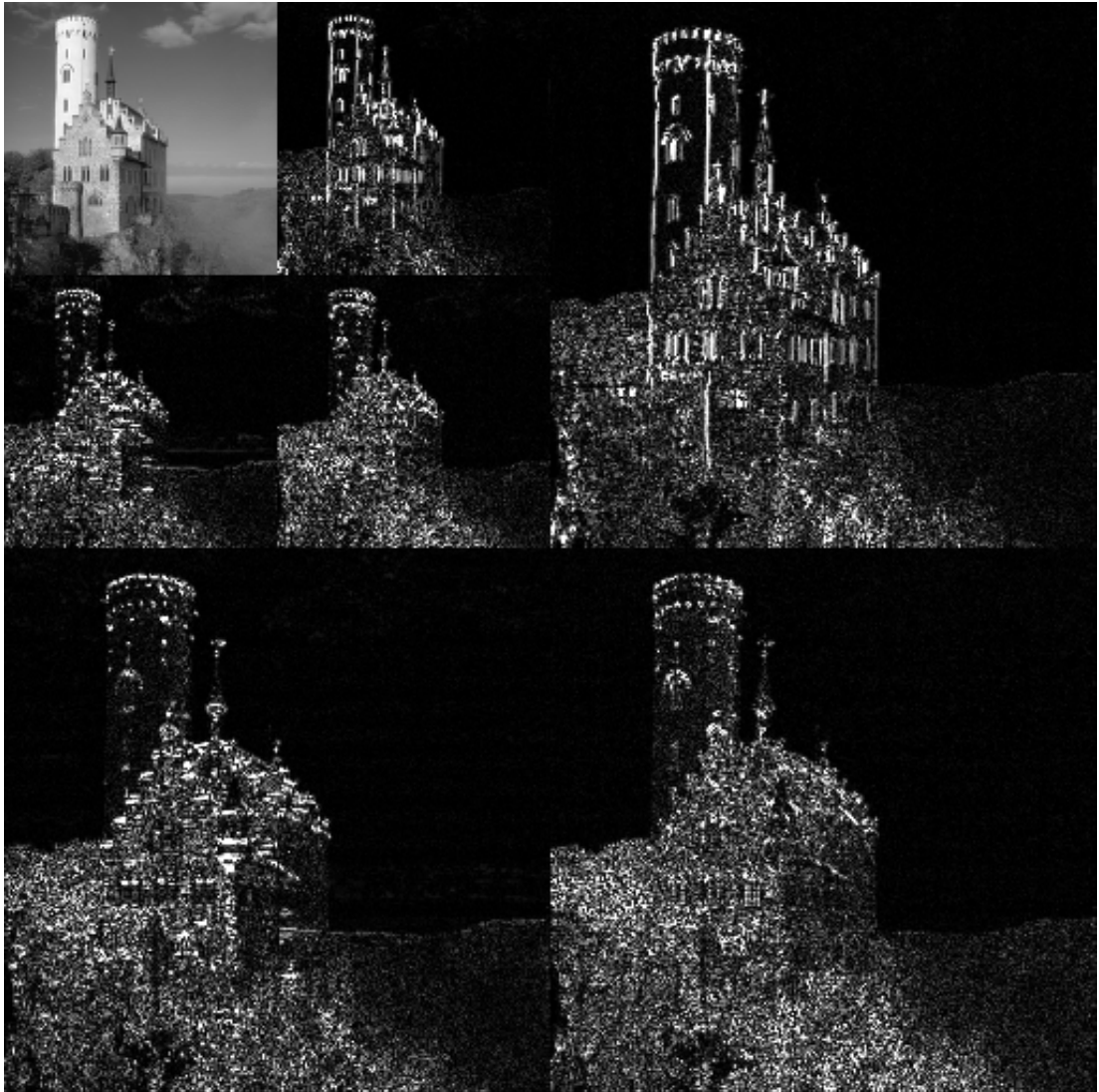


Figure 2.2: An example of recursive wavelet decomposition of 2-dimensional black-and-white image. The sub-image at the top-left corner stands for the low-pass part after two iterations of lifting applied to the input. The original input is the same as this sub-image, except for four times better resolution. The neighbors of this low-pass part are its corresponding high-pass parts. The top-right one of them contains vertical edges, the bottom-left contains horizontal edges and the bottom-right one contains diagonal edges. From these four top-left sub-images, twice as detailed low-pass part can be reconstructed. You can imagine this reconstruction as replacing these four sub-images with the twice more detailed top-left sub-image. After this, we can merge it with the remaining three largest high-pass parts to obtain the original input.

Source: Wikimedia Commons [5]

In the update-first methods, the update operator is applied first:

$$\begin{aligned}s &= x_e + U(x_o) \\ d &= x_o - P(s)\end{aligned}$$

Here is how the reconstruction looks then:

$$\begin{aligned}x_o &= d + P(s) \\ x_e &= s - U(x_o)\end{aligned}$$

2.2 Comparing the wavelets to C-BDAM and our method

In this section, we will briefly describe how the compression inside C-BDAM and our method differ from the basic second-generation wavelet scheme. Even though C-BDAM works with 2D data, its lifting still has just two outputs, unlike in the case of images. It is like it because of the spatial arrangement of its height samples. In one iteration of lifting, a half of samples is omitted (Fig. 2.3). Its lifting is a slight variation of the update-first approach. The main difference is that the input to the first update is not only x_o , but the whole x . In addition, the computation of s is not the summation of the product of x_e and U anymore, because inside $U(x)$, x_e is multiplied:

$$\begin{aligned}s &= U(x) \\ d &= x_o - P(s)\end{aligned}$$

The inverse reconstruction is then:

$$\begin{aligned}x_o &= d + P(s) \\ x_e &= U^{-1}(x)\end{aligned}$$

Moreover, the samples x are regularly distributed in the plane, so the splitting into x_o and x_e no longer depends on the indices of the samples, but on their positions instead (Fig. 2.3). Nevertheless, this is just a formal difference which has no effect on the computation. The size of x_o and x_e is still half the size of x which is crucial to keep the original form of 1D lifting.

Note that if the residuals d were simply quantized after lifting and used in the reconstructions inside the second top-down pass, each step of the reconstruction would increase the maximum absolute deviation from the original low-pass values produced in the first bottom-up pass. To ensure that the reconstructed values are within the maximum-error bound from their corresponding values produced in the first pass at each tier, the residuals computed in the first pass are slightly corrected according to the actual values in another additional top-down pass. What is basically done is that we compare each reconstructed value with its target value. When it is too far from the target value, its residual is shifted by a multiple of the quantization interval, so that the resulting value is close enough to the target value. Thanks to the selection of the quantization interval which

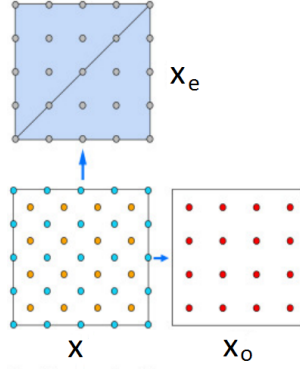


Figure 2.3: Lifting in C-BDAM - the samples x are split into the even ones (x_e) which will become low-pass (s) and the odd ones (x_o) which will become high-pass (d)

Source: C-BDAM [10] (edited)

respects the set maximum-error bound, it is always possible to find such a shift. The following reconstruction (decompression) is again the same top-down pass, except for the fact that just the corrected residuals are read, decompressed and used to progressively reconstruct the data.

The method proposed in this thesis is update-first and uses the whole x as the input to U , but has several differences: the size of x_e and thus s is not half the size of x , but one fourth of it instead, just like in the 2D example of images, as each four neighboring pixels of x are collapsed into one inside s (Fig. 4.2). Additionally, the lifting is not complete, because the prediction operator is not applied there and the computation of residuals is not performed there either. The correct residuals which already ensure the satisfaction of the maximum-error bound constraint are computed directly in the second top-down pass, also utilizing the prediction operator. Similarly to C-BDAM, the computations inside this pass are identical to the reconstruction of the data, except for the fact that during the reconstruction, the residuals are not computed anymore. Additionally, the prediction operator is applied multiple times in one step of the reconstruction which is explained in Chapter 4. The rationale behind all these differences is explained in Chapter. 5.

2.3 Lossless compression

Lossless compression decreases the data size without any information loss. It is often used to compress quantized residuals - the output of wavelet transform. In this section, we briefly describe two main approaches to such compression - Huffman coding and dictionary compression. We do this because both these approaches are used in Zlib - a publicly available lossless compression library which we use in our method to losslessly compress the quantized residuals. We chose it because it achieved the best compression ratios on our data from among all publicly available lossless compression libraries. It is also quite fast. It firstly applies LZ77 to the data and then Huffman coding. Both these approaches basically profit from the redundancy of data, but in different ways.

2.3.1 LZ77 compression

This method is able to detect duplicate subsequences inside the input which are relatively close to each other. It is very suitable for binary data. It reads the input sequentially during the compression. During this, it keeps track of the last N characters. We call this track the sliding window, with N being its length. In practice, N is relatively large - 32KB in Zlib.

We start at the beginning of the input stream, with an empty window. Then, being at a certain position, we examine the data after the window for the longest continuous sequence beginning right after window which can also be found somewhere inside the window. When we have found it, we just write a triple (D, L, C) to the output, with D being the distance of the start of the matching sequence found inside the window to the end of the window, L being the length of the sequence and C being the character in the stream right after this sequence. Of course, it might happen that no such sequence is found - especially at the very start - then both D and L will be 0. Note that L might be greater than D - in this case, when we reach the end of the window while trying to find the longest match, we start again D steps back from the window's end. After outputting the triple (D, L, C) , we move the sliding window $L + 1$ steps forward. Note that this procedure works even for the initial state - the empty sliding window. In this case, $(0, 0)$ is outputted, followed by the first character of the input stream. In this particular situation, $(0, 0)$ can even be omitted, as it has to always be written at the beginning. Then, we repeat the same matching of sequences again, as long as there is something to read. When we have reached the end of the stream, we write a special end character at the place of C .

The decompression then works inversely as follows - it starts with an empty sliding window and then it starts reading all the following triples (D, L, C) . After reading each triple, it moves D steps back from the end of the sliding window and copies L following characters from the current position to the end of the decompressed stream. If L is greater than D , it goes back D steps from the window's end, or equally, it starts traversing the output stream right after the window's end, as in both cases, the same characters will be found (the first character which has been written after the window was the one located D steps back from the window's end). After writing all L characters to the decompressed stream, it writes C to the output stream, pushes the sliding window forward by $L + 1$ steps and continues again with the following triple. When it encounters the special end character, it reads no more triples.

2.3.2 Huffman coding

Huffman coding is a form of prefix coding. It is usually used to compress data composed of characters. It might not be very suitable for arbitrary binary data, as it would most probably have to partition them into certain characters (bytes) in order to work. Thus, if the binary data is actually composed of float numbers or something else distinct from characters, Huffman coding might not work very well and LZ77 is definitely a better choice. However, in Zlib, Huffman coding is used to compress the data produced by LZ77 - the triples consisting of two natural numbers and one character - which can be easily interpreted as characters.

This algorithm starts by assigning a certain weight to each of the characters

contained in the data. The more often the character occurs, the larger its weight should be. Then, according to these weights, it builds a binary Huffman tree, with the characters being the leaf nodes of it. From each non-leaf node of the tree, two edges go - they are marked as "0" and "1". The marks of the edges through which we have to traverse from the root of the tree to a certain character form the code of the character in the compressed stream. Ideally, the greater the weight of a certain character, the closer to the root it should be, so the shorter its code is.

With the tree built, we can easily compress the data. However, it remains a question how exactly. The main advantage of the coding should be that the more frequent characters have shorter codes, so we definitely would not profit from putting each code regularly into one or more bytes. But how to represent distinct lengths then? Should we somehow store the code's length before the code itself? Would not it be quite an overkill?

Zlib comes with a good solution - it constructs the tree always in such way that no two characters have the same code length. Thus, for the subsequent decoding it is sufficient to store just the lengths of the codes. This is a great solution, but did we need the theory of Huffman coding here? This approach is practically equivalent to substituting each character with a number. The most frequent character is substituted by 1, the second one by 2, etc.

3. The outline of the method

In this chapter, we will briefly describe how the compression works. Basically, two consecutive passes are performed on the input heightmap the bottom-up pass followed by the top-down pass (Fig. 3.1). These passes are analogic to the passes of the described second-generation wavelet methods.

The first bottom-up pass computes the target mip-maps - from the largest one to the smallest one. Those will be the mip-maps, against which the accuracy of reconstruction will be measured. The largest mip-map is the input itself. Every smaller mip-map is computed from the previous larger one. These mip-maps will be used just as a reference and will not be stored in any way.

The second top-down pass then directly constructs the decompressed mip-maps. During this construction, it computes, compresses and stores the information needed to subsequently reconstruct these mip-maps. These mip-maps are constructed from the smallest one to the largest one with respect to the target mip-maps in order to ensure that the maximum deviation of every decompressed mip-map from its corresponding target mip-map is within the maximum error bound set by the user. The smallest of these mip-maps is just the suitably quantized sole value of the corresponding target mip-map. It is directly stored as first. After that, none of the following decompressed mip-maps are stored directly. Each following larger decompressed mip-map is firstly predicted from its previous decompressed mip-map. These predictions are then subtracted from the corresponding same-sized target mip-map to obtain the residuals. These residuals are then quantized in such way that the maximum deviation of the current decompressed mip-map is still respected. Then, they are losslessly compressed and stored (Fig. 3.2). Thus, for all the mip-maps except the smallest one, we store just losslessly encoded quantized residuals which are sufficient to reconstruct them in the same way - for each mip-map, we add its decompressed residuals to the predictions from the previous smaller decompressed mip-map (Fig. 3.3).

More formally, the first pass is given the input square block of float height samples \mathbf{L}_n sized $2^n \times 2^n$ and produces n mip-maps $\mathbf{L}_{n-1..0}$ from it, one by one. The dimension of \mathbf{L}_i is half the dimension of \mathbf{L}_{i+1} . Just like it is common in mip-maps \mathbf{L}_i is computed from \mathbf{L}_{i+1} by straightforward averaging of pixels - see the details in the following chapter.

The second top-down pass has already $\mathbf{L}_{0..n}$ available and computes $\mathbf{L}_{0..n}^\bullet$ - the decompressed mip-maps. The dimension of \mathbf{L}_i and \mathbf{L}_i^\bullet is the same. The computation ensures that the maximum absolute deviation between their corresponding samples is not greater than D - the parameter set by the user. This will be denoted by:

$$\maxdev(\mathbf{L}_i, \mathbf{L}_i^\bullet) \leq D,$$

where

$$\maxdev(A, B) = \arg \max_{x,y} |A[x][y] - B[x][y]|$$

We will achieve this with the help of the uniform quantizer Q_D the quantization step of which is set to the maximum value which still respects this error bound:

$$\maxdev(Q_D(x), x) \leq D,$$

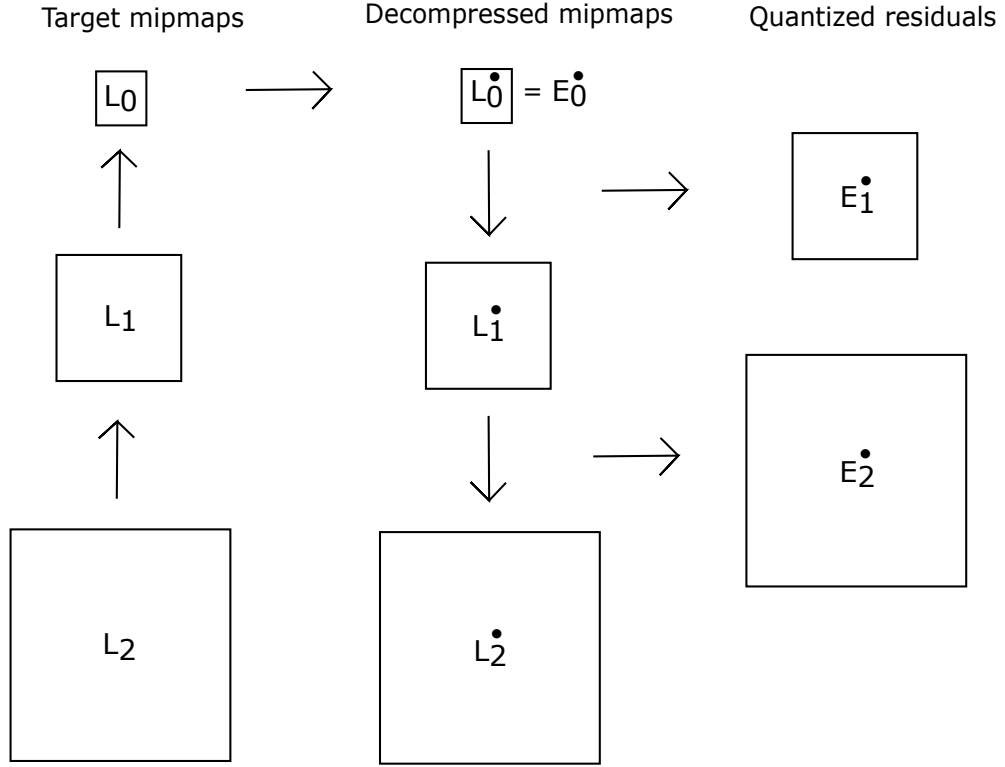


Figure 3.1: The main schema of compression of the method. On the left, the target mip-maps \mathbf{L}_n are constructed in the first bottom-up pass. On the right, the decompressed mip-maps \mathbf{L}_n^\bullet are constructed with respect to the target mip-maps in the second top-bottom pass. Each constructed mip-map is predicted from the previous one and subtracted from the neighbor target mip-map to obtain the residuals, which are then quantized to \mathbf{E}_n^\bullet and compressed.

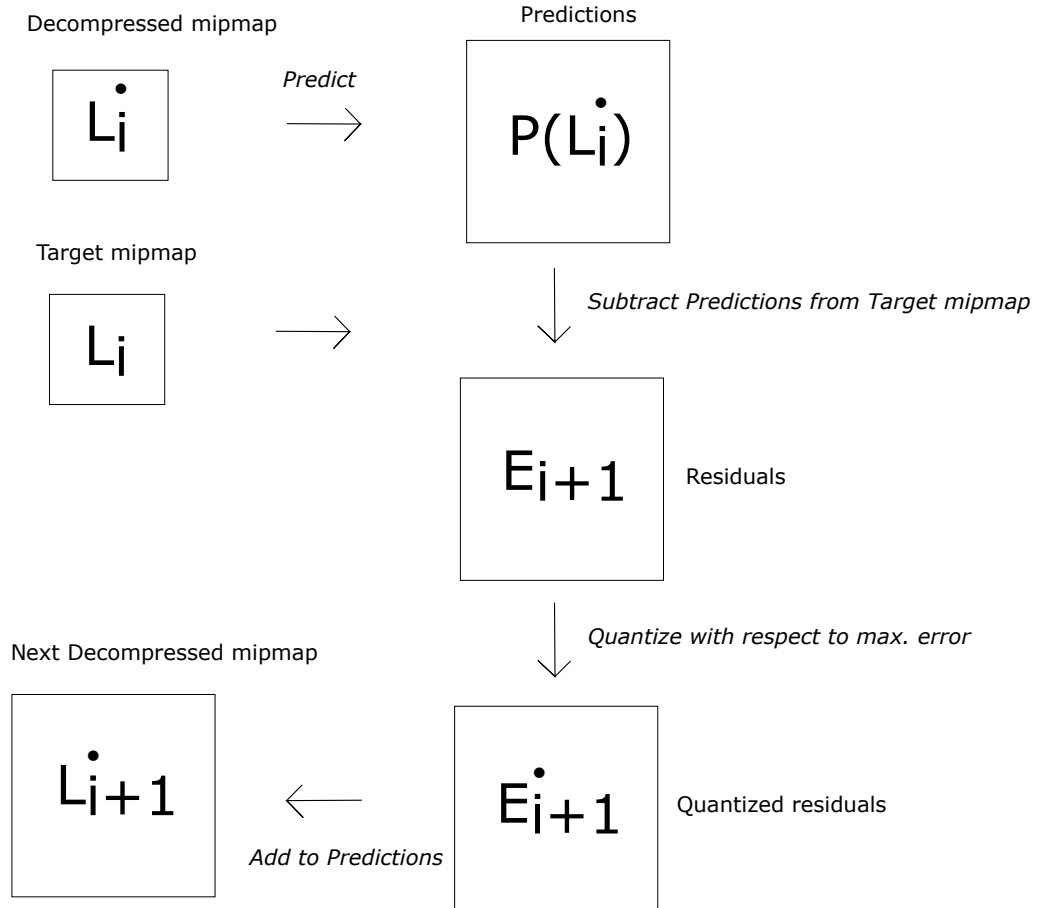


Figure 3.2: The diagram of construction of larger decompressed mip-map L_{i+1}^\bullet from the smaller L_i^\bullet during the off-line compression. Only the quantized residuals E_{i+1}^\bullet are losslessly compressed and stored. They are sufficient to reconstruct L_{i+1}^\bullet from L_i^\bullet .

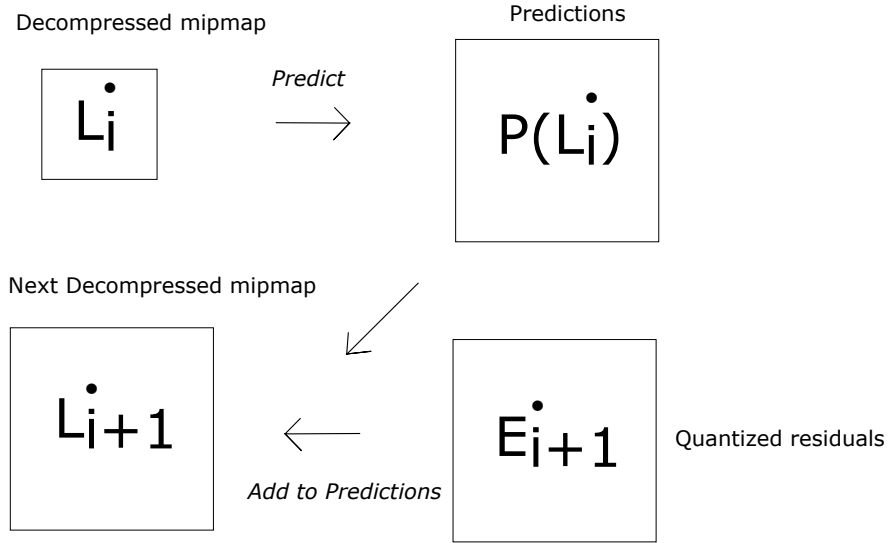


Figure 3.3: The diagram of construction of larger decompressed mip-map L_{i+1}^\bullet from the smaller L_i^\bullet during the real-time decompression. The quantized residuals E_{i+1}^\bullet are read and decompressed. Then, they are added to the predictions from the smaller mip-map.

where x is an arbitrary float sample or block of samples. The quantizing step of this quantizer is $2D - 1$ in case $D \geq 0.5$ and $2D$ otherwise.

As we already mentioned, L_o^\bullet is just the quantized sole value of L_o :

$$L_o^\bullet = Q_D(L_o)$$

Thanks to the fact that the quantizer respects the maximum-error bound D , $\maxdev(L_o, L_o^\bullet) \leq D$.

Then, the values of every following L_{i+1}^\bullet are predicted from the values of the previous L_i^\bullet . The raw differences between the target values and the predicted values are denoted as E_{i+1} (the residuals). With the help of them and the predictions from L_i^\bullet , we would be able to accurately reconstruct the target mip-map L_{i+1} . However, these residuals are then quantized with the uniform quantizer Q_D to E_{i+1}^\bullet . With the help of the quantized residuals, we are no longer able to accurately reconstruct L_{i+1} , but thanks to the fact that the used quantizer keeps the maximum absolute error within the bound D , we can guarantee that the reconstructed L_{i+1}^\bullet will satisfy the maximum-error constraint: $\maxdev(L_{i+1}^\bullet, L_{i+1}) \leq D$. Here is how we construct L_{i+1}^\bullet :

$$\begin{aligned} E_{i+1} &= L_{i+1} - P(L_i^\bullet) \\ E_{i+1}^\bullet &= Q_D(E_{i+1}) \\ L_{i+1}^\bullet &= P(L_i^\bullet) + E_{i+1}^\bullet \end{aligned} \tag{3.1}$$

Thanks to the fact that the residuals E_{i+1} are computed with respect to the target mip-map L_{i+1} , the maximum-error constraint is satisfied, no matter what values are in L_i^\bullet and what the prediction operator P looks like. At the end, the quantized residuals $E_{o..n}^\bullet$ are losslessly compressed with the help of Zlib and stored ($E_o^\bullet = L_o^\bullet$). The order of their storage is from E_o^\bullet to E_n^\bullet , so that progressive decompression is possible. The more accurate P is, the smaller the residuals are, thus the higher the compression ratio is. The details of the prediction operator used in this method are described in the following chapter. The higher D is, the less entropy there is among the residuals, thus the higher the compression ratio is, but the lower the reconstruction quality is.

The real-time decompression then just reads the stored quantized residuals and decompresses them with the help of the same lossless codec. Thanks to the fact that the residuals of a smaller mip-map are stored before the residuals of a larger mip-map, progressive decompression of mip-maps $L_{o..n}^\bullet$ is possible, utilizing the same principle of producing predictions from the previous reconstructed mip-map and adding residuals to them (eq. 3.1). Of course, in order for this to work, the prediction operator must be identical to the one used in the compression.

4. Details of the method

In this chapter, the method is described in more detail. Unlike the previous outline, this description should be sufficient enough for the reader to implement this method by themselves. Section 4.1 details the construction of the mip-maps inside the first bottom-up pass and what alternative constructions we also considered. In Section 4.2, we explain what is the form of P - the prediction operator - and how exactly it is applied in the second top-down pass in order to compute the residuals needed to reconstruct a finer mip-map from the coarser one and also perform the reconstruction with the help of these residuals. Note that this method does not use an update operator, see Chapter 5 for the explanation and details.

4.1 Bottom-up pass

As we already said, in the first bottom-up pass, we just construct the target mip-maps one by one, from the largest one - the input itself - to the smallest one, sized 1. At each step, we construct a smaller mip-map from the last constructed one. The dimension of the new mip-map is half the dimension of the last one, in other words, it is half as detailed. Generally, we can build the new mip-map by any form of averaging of pixels of the larger mip-map. In the previous chapter, we explained that the maximum absolute error of the reconstruction is not dependent on how the mip-maps look, as long as they contain valid values (no infinities, NaNs). However, the appearance of the mip-maps affects the compression ratio. The less different are the heights between the neighboring mip-maps, the lower the residuals of the transition from the smaller one to the larger one are, thus the higher the compression ratio is. Additionally, as mentioned in the introduction, inside the renderer in which the method has been applied, the mip-maps of a certain terrain square are carefully selected, so that aliasing is minimized. This decision is based on the area of the square projected to the screen. This means that while looking at a certain square from above, its finest mip-map is displayed and during a fixed-radius circular traversal around it up to the point when we look at it from a side, we will be gradually displaying coarser (less detailed) mip-maps of the square. This means that if the mip-maps are significantly different from each other, disturbing visual artifacts might occur during this traversal. The best way how to minimize these artifacts is to use the simplest form of averaging of heights when producing a lower-resolution mip-map where the height at every pixel of the smaller mip-map \mathbf{L}_i will be the average of the heights of the four corresponding neighboring pixels inside the larger mip-map \mathbf{L}_{i+1} :

$$\mathbf{L}_i[m][n] = \frac{\sum_{om=0}^1 \sum_{on=0}^1 \mathbf{L}_{i+1}[2m+om][2n+on]}{4} \quad (4.1)$$

For a comparison, in transition to a coarser LOD in C-BDAM, a different form of heights averaging is utilized. It properly conforms to the standard lifting scheme - it uses the update operator to produce the coarser LOD. This averaging is even parametrized by one coefficient named subsampling weight the value of

which can span from 0 to 1. To the contrary, our method does not use the update operator there which is explained in Chapter 5. However, we tried to use a similar averaging of pixels inspired by the one performed in the update operator in C-BDAM. With the subsampling weight well set, we achieved a slightly better compression ratio, but the mentioned visual artifacts were more disturbing. Thus, in order to minimize the visual artifacts, as they really affect the user experience, we decided to use the described averaging scheme (eq. 4.1, Fig. 4.1).

4.2 Top-down pass

This pass is performed in the offline compression after the first bottom-up pass, in which case it computes the residuals $\mathbf{E}_{0..n}^\bullet$ needed to completely progressively reconstruct the compressed multiresolution representation of the input \mathbf{L}_n . During the following real-time decompression, this is the only pass which is performed, with the only difference that it no longer computes the residuals, but it just reads them and uses them to reconstruct the data. Let us describe it in detail, so that it becomes clear how this pass is implemented.

In the previous chapter describing the outline of the method, we claimed that we construct a larger decompressed mip-map \mathbf{L}_{i+1}^\bullet from the smaller \mathbf{L}_i^\bullet in just one step (eq. 3.1). This is not exactly true, it is a simplification which we made for several reasons: to give the reader a simple high-level idea of the method, to make the fact that $\maxdev(\mathbf{L}_{i+1}^\bullet, \mathbf{L}_{i+1}) < D$ easier to see and to make it clear that the way the target mip-maps $\mathbf{L}_{n-1..0}$ look has no effect on this constraint. However, the truth is that to get \mathbf{L}_{i+1}^\bullet from \mathbf{L}_i^\bullet , the prediction operator P is applied consecutively three times. Its form is different at each of these applications which reflects the fact that before each application, different height values are known - the later the application, the more values are known. After each of these applications, the residuals are computed during the compression and added to the already computed values during both the compression and decompression. Nevertheless, the two main principles which ensure that the maximum error bound between \mathbf{L}_{i+1} and \mathbf{L}_{i+1}^\bullet is kept remain unchanged - during the compression, the residuals are still computed against the target mip-map \mathbf{L}_{i+1} after each application of P and all predictions are made from the reconstructed values which ensures that both the compression and the decompression add the residuals to the same values. Hence, let us explain how these three steps are performed.

When constructing the larger compressed mip-map \mathbf{L}_{i+1}^\bullet from \mathbf{L}_i^\bullet , we can imagine it as every pixel p from \mathbf{L}_i^\bullet being substituted by four pixels a, b, c, d in \mathbf{L}_{i+1}^\bullet as depicted in Fig. 4.2. This substitution is the exact inverse of the one performed in the first bottom-up pass described in the previous section ???. We will apply the prediction operator and subsequent residuals computation and addition three times in order to compute the values of the four pixels and the residuals needed to reconstruct them from p during the decompression.

In the first of the three steps, we compute the pixels labeled a . To predict them from their corresponding p pixels inside \mathbf{L}_i^\bullet , we use a simple prediction operator $P_a(\mathbf{L}_i^\bullet) = p$. We compute the residuals \mathbf{E}_a and \mathbf{E}_a^\bullet with respect to the target value a_t in \mathbf{L}_{i+1} and then assign a the final value a^\bullet (eq. 4.2, recall that Q_D is a uniform quantizer respecting the maximum deviation D : $\maxdev(v, Q_D(v)) < D$). It is clear that $\maxdev(a^\bullet, a_t) \leq D$. The explanation would be the same as in

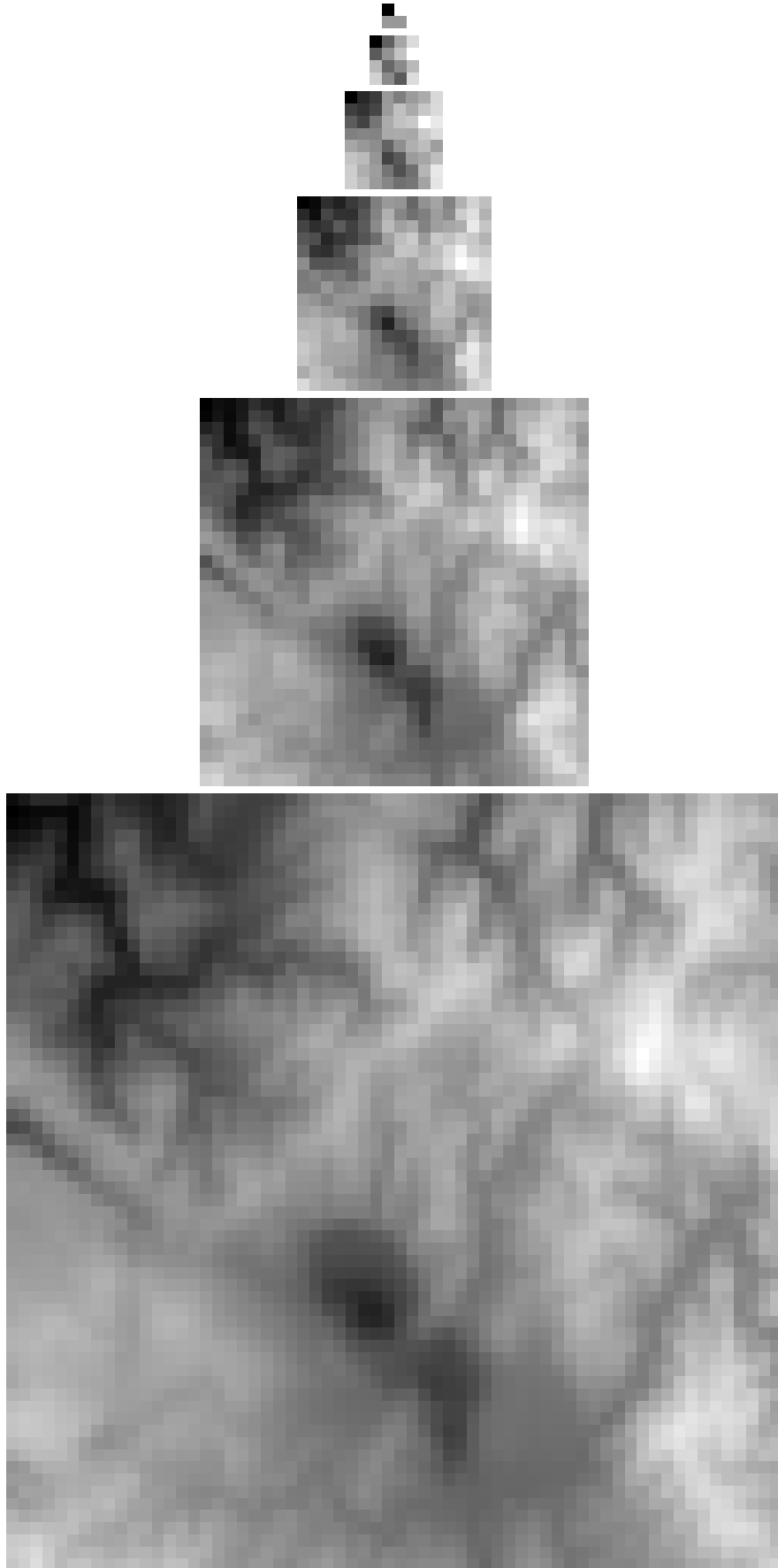


Figure 4.1: An example of the averaging used to produce mip-maps; each mip-map is produced from the previous larger one by averaging of each four neighboring pixels. The brighter a pixel, the higher point it represents.



Figure 4.2: Substituting the pixel p in \mathbf{L}_i^\bullet with four pixels in \mathbf{L}_{i+1}^\bullet



Figure 4.3: The prediction operator of b - $P_b(\mathbf{L}_{i+1}^\bullet)$ - averages the compressed heights at all the displayed $a\bullet$.

Chapter 3.

$$\begin{aligned}
 \mathbf{E}_a &= a_t - p \\
 \mathbf{E}_a^\bullet &= Q_D(\mathbf{E}_a) \\
 a\bullet &= p + \mathbf{E}_a^\bullet
 \end{aligned} \tag{4.2}$$

In the second of the three steps, we compute the pixels labeled b . We do not predict them from \mathbf{L}_i^\bullet anymore, but from the already available pixels $a\bullet$ inside \mathbf{L}_{i+1}^\bullet . The prediction operator P_b used for this has now the form of a straight-oriented Neville interpolating filter of order 2. All it does is that when it is requested to predict the height at some pixel, it just averages the heights of its certain four neighboring pixels as depicted in Fig. 4.3. It is easy to see that as long as it is requested to predict the height at pixels b , it always averages only the already known $a\bullet$ pixels. This is the same prediction operator as the one used in C-BDAM to predict the heights of the samples located at the border of a LOD hierarchy node. Once the predictions of b pixels are known, we perform an analogic computation of residuals \mathbf{E}_b and their quantizations \mathbf{E}_b^\bullet , again with respect to the corresponding target values b_t in \mathbf{L}_{i+1} . Finally, we assign b its final value $b\bullet$ (eq. 4.3).

$$\begin{aligned}
 \mathbf{E}_b &= b_t - P_b(\mathbf{L}_{i+1}^\bullet) \\
 \mathbf{E}_b^\bullet &= Q_D(\mathbf{E}_b) \\
 b\bullet &= P_b(\mathbf{L}_{i+1}^\bullet) + \mathbf{E}_b^\bullet
 \end{aligned} \tag{4.3}$$

The reason why C-BDAM uses the order 2 Neville interpolating filter at the borders is that thanks to the way the samples are organized inside a node of

its LOD hierarchy, the filter does not pick the samples behind the node’s border. We can view the mip-map in our method as an analogy to the node in C-BDAM. However, the spatial organization of mip-map samples in our method differs from the organization of samples inside a LOD node in C-BDAM, so unlike in C-BDAM, in our method it might happen that this interpolating filter comes out of the underlying mip-map. We handle this by only including the valid interior values in the resulting average and completely ignoring the imaginary values behind the mip-map borders. Thus, when computing a certain prediction, we count how many times the filter has hit the interior of mip-map and divide the sum of the valid interior heights with this number of hits. Most of the times, the number of hits will be 4, but it will be 2 at the borders and just 1 at the corners. This way, it is always ensured that the filter does not make any data up, unlike the possible alternative of some mirror extension of data behind the borders. ??? Comparison with mirror extension ???

C-BDAM uses the larger order 4 Neville interpolating filter to predict the heights at the interior of a node of its LOD hierarchy. This filter covers larger area - it samples twelve points instead of four. In addition, it does not compute the average of these points, but their weighted sum. Just like in the case of simple averaging, the sum of the weights is 1. The difference is that the four closest points have a certain positive weight, whereas the remaining eight further points have a different negative weight, the absolute value of which is lower than the first weight (Fig. ???). The property with the lower absolute value indicates that the points which are further affect the result less. The fact that their weights are negative basically means that the valleys and hills are predicted better (Fig. ???).

Unlike C-BDAM, this method uses the smaller order 2 filter even for the interior samples. Let us explain the reasons why we decided to do so. The first reason is the increase of speed. The order 2 filter only averages four values, whereas the order 4 filter averages 12 values. Moreover, the subsequent averaging performed by the order 2 filter can easily be cached during the horizontal traversal which is an additional reduction of the computation overhead (Fig. ???). We also tried using the order 4 filter with various weights settings for the interior values, too. This slightly increased the compression ratio - probably because this filter is better at predicting hills and valleys - but worsened the quality of compression by producing more significant artifacts near smooth terrain’s borders (Fig. 4.4) and sharp terrain transitions (Fig. 4.5). The most probable cause of this is that the predictions made by the order 4 filter tend to differ from the neighboring heights more. This emphasises the artifacts.

Generally, the reason why these artifacts occur is that as long as the predictions are close enough to the target mip-map and their quantized residuals are equal to zero, the compressed values might remain above/under the terrain for a long time, but only until one prediction gets a bit further from the target terrain. As soon as it happens, its associated residual will be quantized to a certain non-zero value which will result in the reconstructed value being flipped to the opposite side of the real terrain which produces a visual artifact. It is not a coincidence that this often occurs near a sharp change in the terrain. The predictions produced by the averaging filter get a bit different from the adjacent ones near this change, because at these places, the filter reaches out to the area behind the change (Fig. 4.6). This difference might then cause the difference in

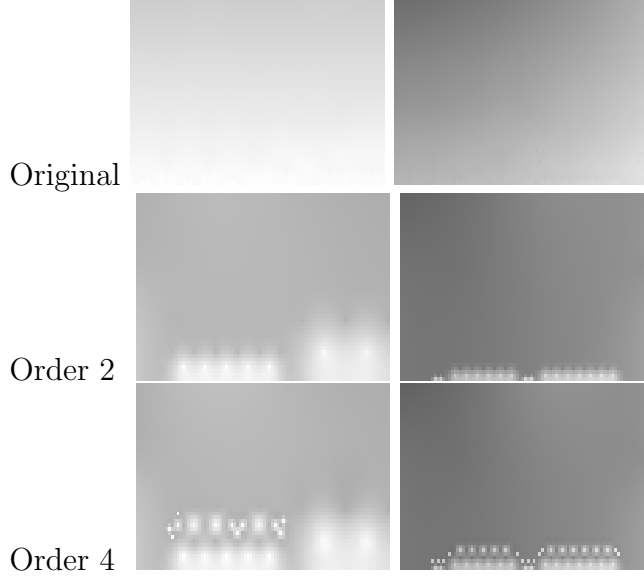


Figure 4.4: Two examples of the difference between artifacts caused by order 2 and order 4 filters near smooth terrain’s border - in the first row there are the target heightmaps, in the second, there are the same heightmaps compressed using the order 2 filter, in the third row, the heightmaps compressed with the order 4 filter.

residuals - the quantized residuals further from this change might be all zeroes, whereas the residual near this change not, causing a spike to occur. This spike will then get propagated to the following compressed mip-map levels. The only thing that is guaranteed is that the maximum error bound is still satisfied. ???(musim overit, clipping je nasadeny len v Bohemke, ale je to lepsi napad ako mirroring, tusim, ze to aj zlepsilo kompresny pomer, no nezistoval som pri nom, ake su artefakty)The clipping performed by the predicting filter near the mip-map borders creates the effect similar to a sharp terrain change, too, in a bit different way - by the sole fact that the terrain behind the border no longer follows its trend up to the border (rising, for example), but is practically mirrored behind the border (following the example, falling), because instead of reaching out to the non-existing values out of the mip-map, the existing ones are used.???

In the last of the three steps, we compute the pixels labeled c . We predict them from the already available compressed values of both $a\bullet$ and $b\bullet$ inside \mathbf{L}_{i+1}^\bullet . The prediction operator P_c used for it has now the form of a diagonally-oriented Neville interpolating filter of order 2. It is the same as the previously applied P_b , except for its different orientation - relatively to P_b , it is rotated by 45 degrees and averages the 4-connected neighbors of the point of application (Fig. 4.7). Once the predictions of heights at c pixels are known, we perform an analogic computation of residuals \mathbf{E}_c and their quantizations \mathbf{E}_c^\bullet , again with respect to the corresponding target values c_t inside \mathbf{L}_{i+1} . At last, we assign every pixel c its final value $c\bullet$ (eq. 4.4).

$$\begin{aligned}\mathbf{E}_c &= c_t - P_c(\mathbf{L}_{i+1}^\bullet) \\ \mathbf{E}_c^\bullet &= Q_D(\mathbf{E}_c)\end{aligned}$$

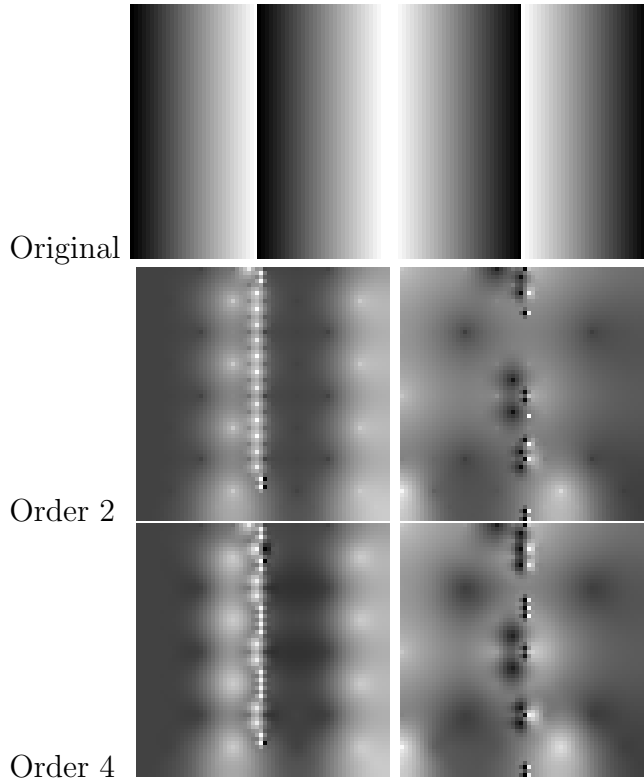


Figure 4.5: Two examples of the difference between artifacts caused by order 2 and order 4 filters near a sharp terrain change - in the first row there are the target heightmaps, in the second row, the same heightmaps compressed using the order 2 filter, in the third row, the heightmaps compressed with the order 4 filter. The span of the values in the original images is from 0 to 16 and the maximum absolute deviation (D) of compression is set to 9.

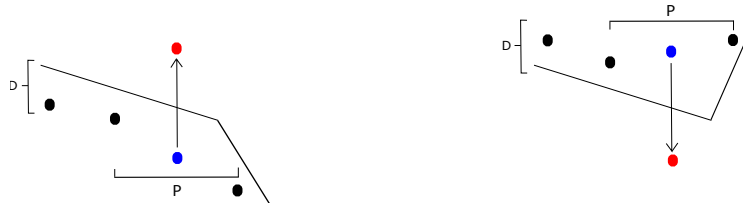


Figure 4.6: Two illustrations of how artifacts can occur near sharp terrain changes - the black dots stand for the predictions which are still within the maximum-error bound D from the target terrain, the blue dots represent the predictions which are just slightly further from the terrain, because their filters span to the area behind the change. Due to the fact that a uniform quantizer with the step of $2D - 1$ is applied to the residuals, the residuals added to the blue predictions will cause them to be shifted by $2D - 1$ to the top (the image on the left) or to the bottom (the image on the right), creating sharp peaks in the reconstructed values - the artifacts.

	•	
•	c	•
	•	

Figure 4.7: The prediction operator of c - $P_c(\mathbf{L}_{i+1}^\bullet)$ - averages the compressed heights at all the pixels marked with a dot - •. Both a_\bullet and b_\bullet are among these pixels.

		$b_1 \bullet$		$b_2 \bullet$	
	$a_1 \bullet$		$a_2 \bullet$		$a_2 \bullet$
$b_1 \bullet$		$b_1 \bullet$		$b_2 \bullet$	
	$a_3 \bullet$		$a_4 \bullet$		$a_4 \bullet$
		$b_1 \bullet$		$b_2 \bullet$	

Figure 4.8: (!!!aktualizuj) Handling of border cases in the computation of $P_c(\mathbf{L}_{i+1}^\bullet)$ - the red line represents the border.

$$c_\bullet = P_c(\mathbf{L}_{i+1}^\bullet) + \mathbf{E}_c^\bullet \quad (4.4)$$

Just like in the case of P_b , P_c can reach out behind the mip-map borders, too. We handled these situations exactly the same way - we only average the heights which are inside the mip-map (Fig. 4.8). Similarly to the predictions of b pixels, we use the order 2 filter to predict the heights at all c pixels - even the interior ones. The subsequent predictions of neighboring c pixels can again be cached to spare some computations. However, the traversal with P_c must now be diagonal in order to make such caching possible (Fig. ???).

Now, let us sum up what we have already done in a few sentences. We have performed all the three subsequent applications of the prediction operator in different forms - P_a , P_b and P_c on the already computed compressed heights in order to obtain the predictions of the yet unknown heights. After each of these applications, we calculated the differences between the predictions and the target values located at the target mip-map at the same places, obtaining the raw residuals \mathbf{E}_a , \mathbf{E}_b and \mathbf{E}_c . Then we quantized these residuals with Q_D - the uniform quantizer respecting the maximum deviation D which ensures that when the quantized residuals are added back to the predictions, all these summations will be within the deviation D from the corresponding target heights. We called the quantized residuals \mathbf{E}_a^\bullet , \mathbf{E}_b^\bullet and \mathbf{E}_c^\bullet . Together, they form \mathbf{E}_{i+1}^\bullet - all the residuals required to reconstruct the larger compressed mip-map \mathbf{L}_{i+1}^\bullet from the previous compressed \mathbf{L}_i^\bullet .

Finally, with all the quantized residuals computed, we losslessly compress them and store them. We firstly compress and store $\mathbf{E}_0^\bullet = Q_D(\mathbf{L}_0)$, then \mathbf{E}_1^\bullet , up to

E_n^\bullet . Thanks to this organization, when we want to run-time decompress any L_i^\bullet , we will be required to read just the starting continuous block of the compressed data $E_{0..i}^\bullet$. This is called the progressive decompression. The decompression itself is performed in a similar way. The only difference is that the quantized residuals are no longer calculated, but just read and decompress. Thus, with L_i^\bullet available, we obtain L_{i+1}^\bullet by substituting every pixel labeled p inside L_i^\bullet by four neighboring pixels labeled a, b, c in L_{i+1}^\bullet (Fig. 4.2), the heights of which will then be computed in three steps. At each of these steps, we will just predict the heights of the pixels with the relevant prediction operator. This will be followed only by adding the read and decompressed residuals to the predictions (the last lines of eq. 4.2, 4.3, 4.4).

The lossless compression of residuals is performed in two steps - packing followed by lossless compression by Zlib. During packing, we divide all the residuals with their quantization interval. Then, we shift them, so that all the values are non-negative. Then, we crop the number of bits of each residual to the number of bits of the largest value. Obviously, we also store the information needed to perform the inverse of packing in the reconstruction - the quantization interval and the upper-shift amount. Only after this packing are the residuals losslessly encoded by Zlib. The packing proved to be useful, as it additionally increased the compression ratio. The decompression is just the inverse - lossless decompression by Zlib followed by the unpacking - expanding the residuals back into floats, shifting them down by the shifting amount and multiplying them with the quantization interval.

5. Functional comparison to C-BDAM and wavelets

In this chapter, with the details of this method described, we will compare it with C-BDAM in more detail. The main difference which we already mentioned is that unlike in C-BDAM where the whole rendering pipeline is contained, our method only focuses on compression. However, we can compare how lifting is performed in these two methods. From the point of view of lifting, a mip-map level in our method is analogic to a node of LOD hierarchy of C-BDAM and the per-mip-map maximum absolute error bound constraint which our method has to keep is analogic to the same per-LOD-node constraint in C-BDAM. At the end of Section 2.2, we already mentioned that when constructing a coarser LOD node from the finer one, C-BDAM omits a half of the samples of the finer node. To the contrary, our method omits three fourths of the samples of the finer mip-map level when constructing a coarser one. With respect to the spatial arrangement of the samples, this transition is equivalent to two fine-to-coarser transitions in C-BDAM (nieco viac)(Fig. 2.3). In the first transition, the pixels labeled as b are removed and in the second one, the pixels labeled as c are removed, as depicted in Fig. 4.2. However, this equality remains only spatial, not computational.

As we already said in Section 2.2, it cannot be really claimed that our method performs the lifting - during the first bottom-top pass, it does not use any prediction or update operator. However, while slightly simplifying the reality, the averaging of four neighboring pixels described in the beginning of Section ?? which is used to construct \mathbf{L}_i from \mathbf{L}_{i+1} can be viewed as an analogy of applying the update operator of lifting in C-CDAM. The crucial difference is that in our method, the lifting is not complete as no prediction operator is applied there to produce the high-pass part - it computes no residuals yet. In C-BDAM, the lifting is complete - a prediction operator is applied there too in order to calculate intermediate quantized residuals. However, reconstructing the data from the coarsest LOD to the finest one in the subsequent top-bottom pass using just these residuals inside the inverse lifting equations would not ensure the required per-sample satisfaction of the maximum absolute error bound, even if these residuals have been quantized with a uniform quantizer respecting this error bound. The reason is that the calculations inside the proper lifting are much more intricate. After each step of the reconstruction, the maximum deviation from the target LODs would increase and would become uncontrollable.

This is exactly why C-BDAM corrects the intermediate residuals against the heights inside the target LODs computed in the first bottom-top pass. This correction is done in another top-bottom pass. It can be simply described as follows - we just reconstruct a certain finer LOD node from its coarser parent and compare it to its corresponding target LOD node produced in the first bottom-top pass. In the places where the reconstruction differs from the target more than the maximum-deviation constraint allows, the corresponding residuals are shifted within the quantization buckets, so that this constraint becomes satisfied. Because C-BDAM uses the same quantizer as our method - the uniform one, set to satisfy the maximum-error bound - such a shift is possible to find. To find

the correct number of quantization buckets by which a residual should be shifted, some intricate computations must be performed, including division, which is undoubtedly a large performance hit. These computations are straightforwardly derived from the lifting equations.

After studying all these equations, we saw an opportunity for simplification there - once it is required to perform a top-bottom pass to correct the intermediate residuals in order to satisfy the maximum error bound constraint, we did not really see it as necessary to calculate any temporary residuals inside the lifting of the bottom-top pass. This is the reason why we use just an analogy of the update operator in the first bottom-top pass to produce the target mip-maps and do not compute any high-pass information (residuals) there yet, and so do not utilize any prediction operator there. We just let the suitable values of residuals be computed in the following top-bottom pass. These values directly satisfy the maximum-deviation constraint and thanks to the fact that we do not need to ensure this with respect to any intricate lifting scheme, it becomes very simple, if not trivial, to compute them (sec. ??). Inside one reconstruction step of the following top-bottom pass, all we do is predict the yet unknown heights in the finer LOD with as much accuracy as possible. These predictions, however, are not linked to the previously performed bottom-top pass, because they have not been applied there at all. Thus, even though the prediction operator is applied three times inside one reconstruction step, we do not have to pay any attention for the equations resulting from it to be exactly inverse to the ones performed in the bottom-top pass. Then we compute the final residuals directly with respect to the target values calculated in the first bottom-top pass. This is undoubtedly a significant deviation from both C-BDAM and the standard second-generation wavelet scheme.

All in all, the way the residuals are computed in this method is a great simplification of the approach used in C-BDAM. Our approach does not even conform to the wavelet scheme of second generation - the lifting is incomplete and the reconstruction is not the inverse of lifting. However, our opinion is that without the per-level residuals correction in the subsequent top-bottom pass, it makes sense to respect this wavelet scheme, because it ensures computational equivalence with the wavelets of the first generation. But as soon there comes the need to correct the residuals at each level, we think that it starts to make no sense to still conform to this scheme, as these corrections destroy this equivalence at a glance - once a quantized residual is manually shifted, so that the resulting value gets closer to the target data, it can no longer be claimed that any of the subsequent reconstruction steps is the exact inverse to the corresponding lifting performed in the first bottom-top pass. Additionally, due to the deviation of C-BDAM from the normal update-first wavelet scheme of second generation which has already been discussed in Section 2.2, we question whether it is still computationally equivalent with the wavelets of the first generation even if it were not for the residuals quantization or cropping. These are the reasons why we suppose that we can optimize the computations performed inside the second top-bottom pass without any cost. With respect to all the discussed matters, this method should be called wavelet-inspired rather than wavelet-based.

6. Results

We have plugged this method into the real-time planet renderer which we already mentioned in the introduction. With this method, we compressed and then real-time viewed the whole-Earth height data with 90m span between height samples (SRTM). The total size of this dataset is 58 GB. Due to the data redundancy of the LOD hierarchy of the renderer which is caused by the fact that it stores all LODs of terrain completely separately, the total size of the SRTM dataset converted to this hierarchy was 260GB. When we applied this compression method to every independent square node of this LOD hierarchy, we managed to compress this data down to 7GB with the maximum deviation of the compression set to 5m. This yield the compression ratio of 37:1. However, if we take the size of the original SRTM dataset as the original size, the compression ratio will drop to just about 8:1. It needs to be said that the first ratio is solely the credit of our method, whereas the second one is influenced by both the way the renderer handles the data and our method.

For a comparison, C-BDAM achieved the compression ratio of 64:1 on the same 90m-resolution SRTM dataset with the maximum error bound set to 16m. As we already mentioned, C-BDAM contains its own LOD rendering hierarchy without any redundancy - a finer LOD is constructed from the coarser one which is where the compression takes place. Thanks to this, there is no data redundancy in the LOD hierarchy of C-BDAM, so the compression ratio of this method was evaluated with respect to the original size of the dataset which is 58 GB. The final size of the compressed data prepared for rendering in C-BDAM was just 870MB.

The most accurate comparison of our method with C-BDAM we could perform was compress the SRTM dataset prepared for rendering in the mentioned application by our method with the maximum deviation set to 16m. We did so, the final size of the data was ??? GB which yield the compression ratio of ??? when compared to the size of the dataset prepared for the renderer (260 GB) and the compression ratio of ??? when compared to the original size of the dataset (58 GB). Again, the first compression ratio is only the credit of our method, whereas the second one is also influenced by the way the data is handled by the renderer (the LODs redundancy, overlaps, etc.). However, in terms of functionality, C-BDAM is not analogic to our method, because our method cannot handle the rendering. It is only analogic to the renderer with our method plugged in - only this way, both compression and rendering are achieved. For this reason, the second compression ratio is the most relevant one for comparison, even though the implementation of the renderer is completely out of the scope of this thesis. Even though the compression ratio resulting from the usage of this renderer together with our method is worse than the one achieved by C-BDAM, the redundancy of the LOD hierarchy of the renderer has some benefits which have already been mentioned in the introduction - mainly much faster access to the data. Whereas this renderer is able to render a scene viewed from quite close to the ground quite quickly because it only needs to fetch the relevant LODs which are stored indepentendly, this is not the case of C-BDAM - to display such a scene, it needs to gradually reconstruct every ancestor LOD node of the nodes displayed in the scene which undoubtedly takes more time. However, we did not

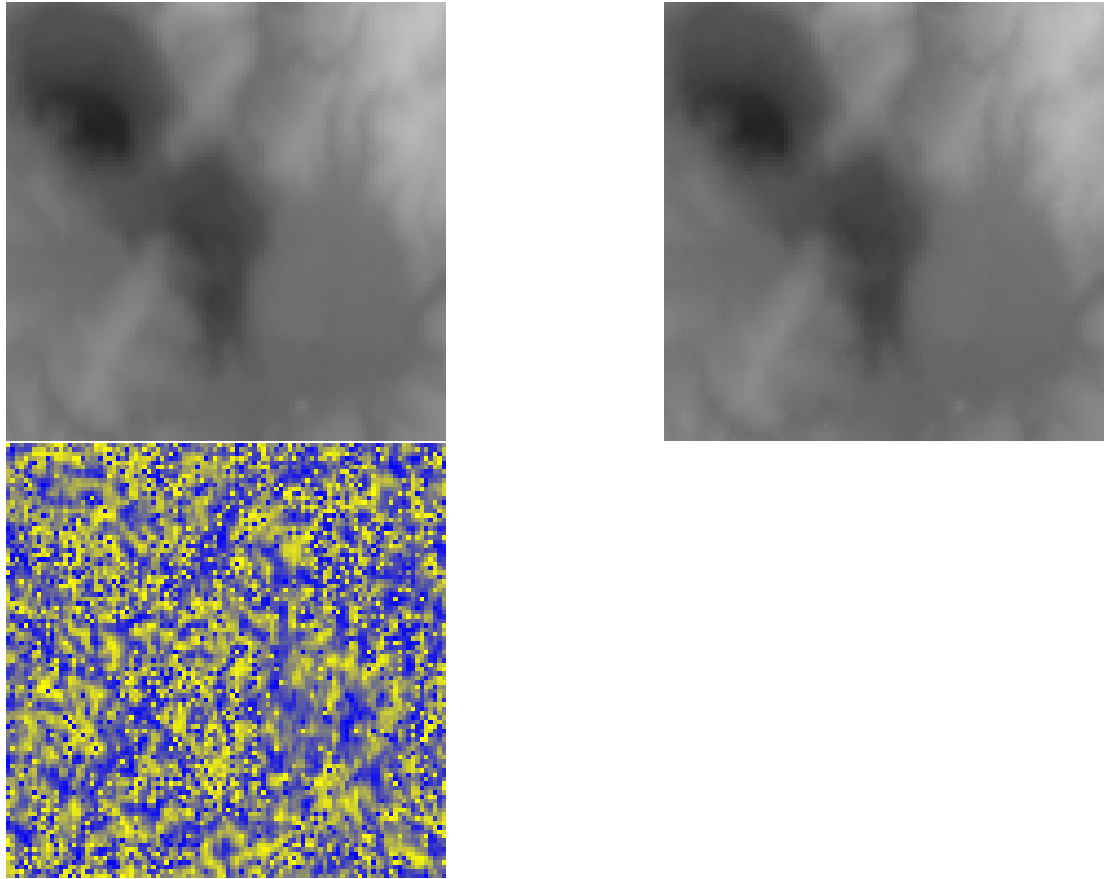


Figure 6.1: From the top to the bottom - the original terrain, the same terrain compressed with the maximum deviation of 5m, the difference between these two. The brighter the color, the greater the value. In the difference image, the yellow color means 4.5m, whereas the blue color means -4.5m.

measure these times.

Fig. 6.1 shows a part of a heightmap compressed by this method, along with the differences from the original.

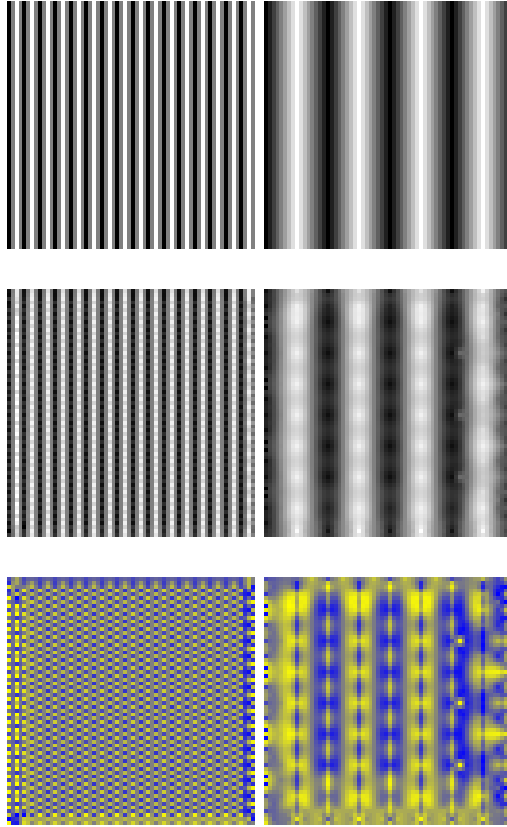


Figure 6.2: Two synthetic test images of size 64x64, each one containing spiky terrain with the heights ranging from -16 to 16. On the left, the longitude of spikes is 4, on the right, it is 16. From the top to the bottom - the original, compressed with the maximum deviation of 5, the difference between these two. The brighter the color, the greater the value. In the difference image, the yellow color means 4.5, whereas the blue color means -4.5.

7. Conclusion

The aim of this thesis was to design a heightmap compression plugin into an existing real-time planet renderer. This plugin should be able to compress a regular square block of height samples and progressively decompress it in the real-time, from the smallest mip-map to the largest one. Apart from this, the plugin should not in any way interfere with the rendering pipeline of the application. After summarizing the available literature, we decided to write our own method inspired by the most promising compression approach used inside C-BDAM, the method for rendering compressed flat terrain. Our method turned out to be suitable for the purpose. The decompression of one 256x256 square which is the size of square used in the application takes only about 1ms. The compression of it takes about 30ms which is still bearable, as it does not have to be performed in the real time. The compression ratios of our method are comparable to the ones of C-BDAM which provides the best compression ratios among the studied methods.

We should also investigate the possibility to perform the decompression on GPU in order to spare RAM. However, the main priority was still to minimize the size of the compressed data which is why we followed the compression ratio in the first place. With respect to it, we designed our method and only after this we analyzed whether and how its decompression could be put onto GPU. We saw two places where the decompression could already switch to GPU:

1. before the decompression of residuals
2. right after the decompression of residuals

The first possibility would imply that the decompression performed by Zlib would have to be put onto GPU. As we already described in Sec. ???, Zlib contains both entropy and dictionary compression, so this would be really complicated. At the end, we did not have time to try to do it.

The second possibility means less spared RAM, but still some, because, as we described in Sec. 4.2, before the residuals are losslessly compressed they are packed - divided by their quantization interval, shifted, so they are all positive and then skewed to the number of bits of the largest value. So, right after their decompression by Zlib, they are still a bit compressed. Theoretically, their unpacking could be done on GPU. Then, the subsequent reconstruction of a larger mip-map from the smaller one would have to be performed on GPU, too. The ability to do this surely already depends on the implementation of the renderer. This is where we encountered a problem.

The renderer uses just one vertex buffer per each square. Then, it sends a mip-map of the square to GPU as a texture and according to this heights in this texture, it performs displacement of vertices of the buffer. Thus, if the unpacking of residuals and the progressive mip-maps reconstruction were all to be performed on GPU, we would no longer send any textures to GPU, but we would just send the unpacked residuals there which is the only remaining way how to spare RAM in our case. The problem is that when constructing a certain mip-map, we must have the previous one at hand. However, in the current implementation of the renderer which does not use CUDA, we have no possibility to create and maintain an array of values. The second possibility would be to directly utilize

the information in vertices - their location. To determine the height of a certain vertex, we would have to average the suitably backward transformed heights of its neighbors (Sec. 4.2). This would have to be done in vertex shader. However, this is impossible, as in vertex shader, we cannot access neighboring vertices of the current vertex. Due to all these factors, we finally decided not to implement any part of decompression on GPU.

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List of Tables

List of Abbreviations

Attachments