Similarity Based Deduplication with Small Data Chunks

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Abstract. Large backup and restore systems may have a petabyte or more data in their repository. Such systems are often compressed by means of deduplication techniques, that partition the input text into chunks and store recurring chunks only once. One of the approaches is to use hashing methods to store fingerprints for each data chunk, detecting identical chunks with very low probability for collisions. As alternative, it has been suggested to use similarity instead of identity based searches, which allows the definition of much larger chunks. This implies that the data structure needed to store the fingerprints is much smaller, so that such a system may be more scalable than systems built on the first approach.

This paper deals with an extension of the second approach to systems in which it is still preferred to use small chunks. We describe the design choices made during the development of what we call an approximate hash function, serving as the basic tool of the new suggested deduplication system and report on extensive tests performed on a variety of large input files.

Keywords: approximate hash scheme, deduplication, compression

1 Introduction and Motivation

Huge amounts of data have to be processed daily and the current trend suggests that these amounts will continue being ever-increasing in the foreseeable future. An efficient way to alleviate the problem is by using deduplication: large parts of the available data is copied again and again and forwarded without any change; the idea underlying a deduplication system is to locate repeated data and store only its first occurrence. Subsequent copies are replaced by pointers to the stored occurrence, which significantly reduces the storage requirements if the data is indeed repetitive [2].

Several approaches have been proposed to solve the problem, each concentrating on another aspect of the input characteristics. One of the approaches, based on hashing, can be schematically described as follows [11,13,14].

The available data is partitioned into parts called chunks $C_i$. These chunks can be of fixed or variable size, and the (average) size of a chunk can be small, say 4–8 KB, up to quite large, say, about 16 MB. A cryptographically strong hash function $h$ is applied to these chunks, meaning that if $h(C_i) = h(C_j)$, it can be assumed, with
very low error probability, that the chunks $C_i$ and $C_j$ are identical. The set $S$ of different hash values, along with pointers to the corresponding chunks, is kept in a data structure $D$ allowing fast access and easy update, typically a hash table or a B-tree. For each new chunk to be treated, its hash value is searched for in $D$, and if it appears there, one may assume that the given chunk is a duplicate. It is thus not stored again, rather, it is replaced by a pointer to its earlier occurrence. If the hash value is not in $D$, the given chunk is considered new, so it is stored and its hash value is adjointed to the set $S$.

The suggested methods mainly differ in the way they define the chunk boundaries, and in the suggested size of the chunks. The chunk size may indeed have a major impact on the performance: if it is too small, the number of different chunks may be so large as to jeopardize the whole approach, because the data structure $D$ might not fit into RAM, so the system might not be scalable. On the other hand, if the chunk size is chosen too large, the probability of getting identical chunks decreases: many instances of chunks might exist, that could have been deduplicated had the chunk size been chosen smaller, but which, for the larger chunk size, have to be kept.

A possible solution to this chunk size dilemma has been suggested in [1]. The main idea there is to look for similar rather than identical chunks. If such a similar chunk is located, only the difference is recorded, which is generally much smaller than a full chunk. This allows the use of much larger chunks than in identity based systems.

For many applications, such as data backups and archiving, data is more fine-grained, and much better deduplication can be performed if one can use significantly smaller chunks. A simple generalization of the above system in which the chunk size would be reduced from 16 MB to 8 KB, that is, by a factor of 2000, without changing anything else in the design, would imply a 2000 fold increase of the size of the index, from 4 GB to about 8 TB. This cannot be assumed to fit into RAM in the near future. Moreover, keeping the definition of the notion of similarity and reducing the size of the chunks will lead to an increased number of collisions, which may invalidate the approach altogether.

The idea of the current work is to implement the required similarity by what we call an approximate hash scheme. This is an extension of the notion of locality-sensitive hashing introduced in [8]. The basic idea is that such an approximate hash function is not sensitive to “small” changes within the chunk, and yet behaves like other hash functions as far as the close to uniform distribution of its values is concerned. As a consequence, one can handle the set of approximate hash values as is usually done in hash applications (using a hash table, or storing the values in a B-Tree), but detect also similar, and not only identical chunks. If a given chunk undergoes a more extended, but still minor, update, its new hash value might be close to the original one, which suggests that in the case of a miss, the values stored in the vicinity of the given element in the hash table should be checked. Such vicinity searches are useless in a regular hash approach.

An approximate hash could be defined by a property that reminds the definition of a continuous function: let $A$ and $B$ be data chunks of fixed size, and let $d(x, y)$ be some distance function to be defined on the set of chunks; a hash function $ah$ will be called an $\varepsilon$-approximate hash if

$$\exists \delta > 0 \quad d(A, B) < \delta \implies |ah(A) - ah(B)| < \varepsilon.$$  

Note the difference with the common continuity definition, in which we would have $\forall \varepsilon \exists \delta$, implying that we can get function values as close as wanted ($\varepsilon$ can tend to 0)
if we start from close enough arguments. In our case, it would be exaggerated to impose such a property, and we can relax it to find two bounds $\delta$ and $\varepsilon$ such that if the distance between chunks is bounded by the first, then the distance between the hash values of these chunks is bounded by the second, for reasonably chosen small values of $\varepsilon$.

Actually, even this definition could be too restrictive, and we should allow a small number of exceptions for certain extreme chunks. This leads to a probabilistic version of the above definition: a hash function $ah$ will be called an $\varepsilon$-approximate hash with probability $p$ if

$$\exists \delta > 0 \quad d(A, B) < \delta \rightarrow Pr (|ah(A) - ah(B)| > \varepsilon) < 1 - p,$$

where the probability is taken over a uniform selection of the possible chunks $A$ and $B$.

There are several possibilities to define the distance function $d$. A simple solution would be the Hamming distance, defined either on bits (number of 1 bits in $A \oplus B$) or on characters (number of differing characters), but this requires the chunks to be of the same length. A more significant, yet more involved, function could be the edit distance: the minimal number of single character insert, delete and substitute operations needed to transform $A$ into $B$.

The challenge is now to find such a function $ah$, giving a tradeoff between how well it can be adapted to reflect the approximate nature described above, and how long it takes to evaluate it. We should still bear in mind that one of the most basic requirements of a hash function is that it should not require too much CPU time.

The general algorithm for storing the repository will then be as follows. The number $k$ of bits in the signature will be chosen in advance, and a hash table $H$ with $2^k$ entries will be used as basic data structure. During the building process, each chunk $C$ will be assigned its approximate hash value $ah(C)$, and the index, or address, of the chunk will be stored at $H[ah(C)]$, the entry in $H$ indexed by the hash value of the chunk. If the location in the table is not free, it is overwritten. This may happen in case the new chunk is identical or very similar to a previously encountered chunk, in which case we prefer to store the address of the more recent chunk for potential later reference; but a collision may also be the result of two completely different chunks hashing to the same value, and then the pointer to the older chunk that has been overwritten will be lost.

In the next section, we describe the details leading to the design of the approximate hash function, and then report on extensive tests in Section 3, showing a noticeable improvement of the suggested method over identity based deduplication with small data chunks.

## 2 An approximate hashing function

Once it has been decided to base the system on approximate hashes according to the ideas stated above, the problem reduces to devising an appropriate function. This is the main thrust of the present suggestion.

Classical hashing functions have been studied for decades and many good solutions are known. The major challenge in the design of an approximate hash function is finding the right balance between the following three competing criteria:
– Uniformity: the function should yield a distribution of values as close as possible to uniform, so as to minimize the number of collisions (false matches);
– Simplicity: the function should be easy and fast to calculate;
– Sensitivity: small changes in the chunk should not, or only slightly, affect the corresponding approximate hash value.

The first two are properties that are common to all hashing functions, the third one, sensitivity, is proper to the approximate version suggested herein. For standard hashing schemes, just the contrary is required: even very small changes in the chunk should lead to extensive changes in the hash value, otherwise the uniformity would be hurt. Some works dealing with similarity rather than identity can be found in [4,5,10]. Our approach is different and will be described below.

The value produced by a hash function is, in a certain sense, a summarization of the information contained in the data on which the function has been applied. This is reminiscent of similar functions, like the Fourier Transform with its many applications, or the Discrete Cosine Transform, used in JPEG image compression. Such transforms allow to recode the information of the given data into a different form, which may be more useful for certain applications, for example, being more compressible. Similarly, we would like to recode compactly much of the information contained within a given data chunk under the constraint that this recoding should be immune to small fluctuations.

This lead to the decision of using the distribution of the various characters that appear in the data as the basis for the suggested approximate hash. The data will be partitioned into relatively small chunks \( C \) of fixed or variable length, with (average) size of about 8–16 K. Each such chunk will be analyzed as to the distribution of the bytes forming it and their frequencies. We define the sequence of different bytes, ordered by their frequency of occurrence in the chunk, as the c-spectrum of \( C \), and the corresponding sequence of frequencies as the f-spectrum of \( C \). In addition, we consider also the sequence of different byte pairs, ordered by their frequency of occurrence in the chunk, and call it the p-spectrum of \( C \). The suggested approximate hash function \( ah(C) \) will be a combination of certain elements of these spectra. The reasoning behind the decision of relying on these distributions is that on the one hand, they usually behave like fingerprints, and it will be rare that essentially different chunks will exhibit the same distributions, but on the other hand, small perturbations in the data will often have no, or just a minor, impact on the corresponding spectra. This is the goal we wish to achieve in designing an approximate hash.

The size of the hash values will be fixed in advance, so as to exploit the space of the allocated hash table. For example, one could decide that the table will have about 4 billion entries, which corresponds to a hash value of 32 bits. A much larger hash value using \( k > 32 \) bits could be prohibitive, since the corresponding hash table would then have \( 2^k \) entries. On the other hand, a small value of \( k \) limits the number of elements of the spectra that can be chosen to be a part of the definition of the signature. To overcome this limitation, the chosen elements of the spectra, and more precisely, only a part of their bits, will be arranged appropriately by shifting them to the desired positions, and all these bit strings will be XORed. By using different indents for the different elements, the final value will not only depend on each of the building blocks, but also on their internal order. Figure [Fig.1] is a schematic representation of a possible layout of these elements. The columns correspond to the 32 bit positions, and each
rectangle stands for one of the elements, with the upper elements corresponding to the c-spectrum, the lowest elements corresponding to the f-spectrum, and the element in the middle corresponding to the p-spectrum, as detailed below. As can be seen, each bit position of the final signature is influenced by several elements.

We do not claim that the suggested layout is the best possible, not even for the sample data on which it has been tested. Rather, it is brought as an illustration of the ideas leading to its design. The specific values of the various parameters (lengths and shifts) shown in this example have been set empirically by iterating experiments to locally optimize the performance.

2.1 Using elements of the c-spectrum

Let $a_1, a_2, \ldots, a_n$ be the sequence of different bytes in the chunk, or, more precisely, the numerical value of these bytes, ordered by non-increasing frequency in the chunk. Ties are broken by sorting bytes with identical frequency by their numerical value. Let $f_1, f_2, \ldots, f_n$ be, respectively, the corresponding frequencies. The number $n$ of different bytes in the chunk can vary between 1 (for chunks of identical bytes, like all zeroes or blanks) and $|C|$, the size of the chunk. As this size is mostly much larger than the maximum numerical value of a byte, one may assume that $1 \leq n \leq 256$.

A first attempt would be to consider each byte on its own as one of the building blocks of Figure 1, but this might result in a function that is too sensitive to noise. It will often happen that frequencies of certain bytes may be equal or very close. In such a case, a small perturbation might change the order of the bytes and yield a completely different hash value, contrarily to our goal of the approximate hash function being immune to small changes. To circumvent this problem, the $a_i$ will be partitioned into blocks, gathering several bytes together and treating them symmetrically. The representation of all the elements in a block will be aligned with the same offset and will be XORed together, so that the internal order within the blocks may be arbitrary, since the XOR operation is commutative.
The sizes of the blocks should not be fixed in advance, but depend on the values themselves. Consider the sizes $d_i$ of the gaps between the frequencies, $d_i = f_i - f_{i+1}$, for $i = 1, \ldots, n - 1$. The boundaries between the blocks should be chosen according to the largest gaps, however, sorting according to $d_i$ alone would strongly bias the definition of the gaps towards inducing blocks with single elements, since the largest gaps will tend to occur between the largest values. We should therefore normalize the size of the gaps by dividing by an appropriate weight. We chose harmonic weights $\frac{1}{i}$ for $i \geq 1$ according to Zipf’s law [15]. The gaps are therefore sorted with respect to $i \times d_i = i \times (f_i - f_{i+1})$, which has the advantage of requiring only integer arithmetic.

For a given parameter $\ell$, the $\ell - 1$ gaps with largest weights are chosen and the $\ell$ sets of consecutive elements delimited by the beginning of the sequence, these $\ell - 1$ gaps, and the end of the sequence, are defined as the blocks. Figure 2 is a schematic representation of an example partition into blocks with $\ell = 8$. The squares represent elements $a_i$, the arrows stand for weighted gaps $i (f_i - f_{i+1})$, and the numbers under the arrows are the indices of the weighted gaps in non-increasing order. In this example, the induced blocks would consist of 3, 1, 3, 2, 4, . . . bytes, respectively.

The number of elements forming the last block is limited, if necessary, to include at most a predetermined number of bytes, say 10, otherwise the speed of calculation could be hurt, and spurious bytes that appear possibly only once or twice in the chunk would have too strong of an influence. For the same reason, there are also lower bounds on the number of occurrences of a byte to be considered at all (for example, 15) and on the size $d_i$ of a gap (for example, 5). If after these adjustments, the number of blocks in a given chunk is smaller than the selected value of $\ell$, a different layout will be chosen that is adapted to the given number of blocks. In any case, one has to prepare layouts also for the possibility of having any number of blocks between 1 and $\ell$, since certain extreme chunks may contain only a small number of different bytes.

Each block taken from the c-spectrum will be represented by a string of 8 bits, using the full representation of the corresponding bytes. The strings are depicted in Figure 1 as white rectangles. Each of these rectangles is shifted as indicated in Figure 1, where they are listed in order top down. That is, the first rectangle is shifted by 24 bits to the left (in fact, to get it left justified in the 32-bit layout), the next rectangle is shifted 21 bits, then 18, 15, 12, 9, 6 and 3 bits.

2.2 Using elements of the f-spectrum

The elements of the f-spectrum are incorporated into the signature independently from the partition into blocks of the corresponding bytes. For each frequency value, which can be an integer between 1 and $|C|$, consider first its standard binary representation (say, in 16 bits), and extend this string by 8 additional zeros to the right. We thus assign to each frequency $f_i$ a 24-bit string $F_i$, for example, if $f_i = 5$, then $F_i = 00000000 00000101 00000000$. We further define $D_i$ as the substring of
$F_i$ of length $m$ bits, for some predetermined small integer $m$, starting at the position immediately following the most significant 1-bit, for our case 00000000 000010100000000, the bits forming $D_i$ for $m = 3$ appear bold faced. To give another example with a value of more than 8 bits, consider $f_i = 759; 0000001011011100000000$ then displays both $F_i$ and $D_i$. In the example of Figure 1, the elements included in the layout are the $D_i$, and the size $m$ of all the elements is chosen as 3 bits. We experimented also with other values of $m$, from 1 to 8, but got better results with $m = 3$.

The 8 bit padding allows values of $m$ up to 8. The offsets of the chosen elements are as indicated, this time bottom up, with the largest frequency being depicted as the lowest element in the figure: 0, 0, 0, 1, 1, 2, 2, 3, 3, 4, 4, 5, 5, 6, 6 and 6. The idea behind this choice of bits is to select those with highest variability so as to get a broad spread of values, but to ignore, for the larger frequencies, the lowest bits, which are those most influenced by small fluctuations.

### 2.3 Using elements of the p-spectrum

Though much of the information of a chunk is already contained in the c- and f-spectrum, we decided to adjoin also elements of the p-spectrum and got empirical evidence that this improved the performance. When the maximal number $\ell$ of blocks could be used, a single element based on the p-spectrum was sufficient. The corresponding rectangle, depicted in the center of Figure 1, is of length 12 bits and will be placed left-justified in the layout. It is defined as follows: order the pairs by non-increasing frequencies and consider those indexed 5, 6, 7, 8 and 9 in this ordering. The reason for not choosing the most frequent pairs as we did for the individual bytes is that their distribution is much more biased, with the pairs (0,0) and (255,255) appearing as the most frequent in an overwhelming majority of the cases we tested. On the other side, there was already a great variability in the pairs in positions 5 to 9.

For each of the 5 pairs, the following bitstring is constructed. Given the 2 bytes $A = a_7 \cdots a_0$ and $B = b_7 \cdots b_0$, we rotate $A$ cyclically to the left by 3 bits, and $B$ cyclically to the right by 3 bits. The bytes are aligned so that the rightmost 4 bits of $A$ overlap with the leftmost 4 bits of $B$, and then the strings are XORed. Formally, the 12 resulting bits are now

$$a_4,a_5,a_2,a_1,a_0 \oplus b_2,a_7 \oplus b_1,a_6 \oplus b_0,a_5 \oplus b_7,b_6,b_5,b_4,b_3,$$

where the $\oplus$ operator stands for bitwise XOR. Note that the most and least significant bits of both $A$ and $B$ are in the overlapping part, so if their distribution is biased, they have an additional chance to correct the bias by the additional XOR. This is important for special cases, for example, when the chunk only contains printable text. The representation of all the bytes would then start with the same one to three bits, which could have a negative effect on the uniformity we seek.

### 2.4 Putting it all together

Finally, all the elements of the layout are XORed, yielding a 32 bit string, representing a number between 0 and $2^{32} - 1$ that will act as the hash value of the given chunk $C$.

The geometry of the layout of the signature has been chosen on purpose as given in Figure 1, with the most frequent bytes being placed left-justified, thereby influencing the most significant (highest) bits, and the lowest elements of the f-spectrum appearing in the area influencing the least significant (lowest) bits. The intention was
that in case of small fluctuations in the frequencies, the order of the most frequent
characters might remain the same, so only some low order bits would change, yield-
ing just a small difference in the signature values. Minor changes affecting even lower
frequencies may go undetected, either because the corresponding frequencies are not
among those chosen, or because the change is in the lower order bits that are not
recorded in the signature.

3 Experimental Results

We performed a series of tests to assess the usefulness of the approach. A first concern
was to verify that the proposed approximate hash indeed spreads its values evenly.
Once this has been confirmed, we have to check that this uniformity does not come at
the price of sensitivity, as it would for a standard hashing scheme. We thus checked the
impact of the signature scheme in some artificial perturbation and clustering tests,
described below. Finally, we bring examples of applying the whole deduplication
process in comparison with an identity based approach.

As testbed, a subset of an Exchange database (EXC) of about 27 GB has been
chosen, as well as the entire operating system of one of our computers (OS), a file of
about 5 GB. The first set of tests was done with chunks of fixed length 8 K. These
tests were then repeated for variable length sized chunks, the boundary of a chunk
being defined by applying a simple Rabin-Karp rolling hash on the \( d \) rightmost bytes
of the chunk under consideration. If this hash value equals some predefined constant
\( c \), the chunk is truncated after these \( d \) bytes; otherwise, the following byte is adjoined
and the test with the rolling hash is repeated. In the test, \( d = 25 \), \( c = 2718 \) and the
hash function is \( RK(x) = x \mod P \), where \( P = 2^{48} - 257 \) is a prime number. To
avoid extreme values for the chunk lengths, a lower limit of 2 K and an upper limit
of 64 K has been imposed. The average size of a chunk was around 12 K on our test
databases.

\[
\begin{array}{|c|c|c|c|c|}
\hline
\text{EXC} & 3,300,000 & 0.5050 & 1.0\% & 0.2991 & 3.6\% & 1.0033 \\
\text{OS} & 594,969 & 0.5085 & 1.7\% & 0.2858 & -2\% & 1.0996 \\
\hline
\end{array}
\]

Table 1. Some statistics on the test databases and signatures

3.1 Uniformity

Table 1 summarizes some statistics about the test databases, the number of 8K blocks,
the average signature value (normalized to the \([0,1]\) range), the standard deviation
of these normalized values, as well as the deviation from the expected results for a
uniform distribution. As can be seen, the values are very close to the expected ones.
On the EXC database, the chunk containing only zeros appeared 1756 times, but
beside the corresponding signature, all the others appeared mostly only once, some
appeared twice, etc. No signature appeared more than 45 times. The last column of
Table 1 gives the average number of occurrences for each signature.

For a more precise evaluation, inspecting each individual bit, the graph of Figure 3
shows the probability of getting a 1-bit in each of the 32 positions of the signatures.
Bit distribution on example data

Figure 3. Bit distribution on example data

3.2 Perturbation tests

We now turn to observing the properties of the signature when introducing perturbations. Recall that the challenge was to reconcile two contradicting demands: on the one hand, the function is required, similarly to usual hash functions, to spread its values as much as possible, so as to minimize the number of collisions; on the other hand, we want small perturbations to yield only slight differences, if at all, in the corresponding signature values, a property one explicitly prohibits for classical hashing.

To simulate real life changes, the modified bytes did not get a random value, but rather another randomly chosen byte from within the same chunk was copied into the location to be modified. Thus the perturbations were introduced as follows: a random position \( i \) between 1 and \( |C| \), the size of the chunk, was chosen, and the character from position \( |C| - i + 1 \) was copied to position \( i \), overwriting the current one. The idea was to change the chunk slightly, but without introducing any new characters that are not already present in the chunk. Obviously, there is a small chance that this “perturbation” is in fact void (when overwriting a character by itself), but the corresponding probability is small enough so as not to bias the overall statistics. The signature function was then applied to the modified chunk and compared to the signature of the original chunk. In many cases, we got the same signature, meaning that changing a single byte in the chunk did not change the function, contrarily to what would be expected from a real hash function.

The above perturbation procedure has then been repeated, and the signature was reevaluated after 2, 3, \( \ldots \), 10, 20, 30, \( \ldots \), 100, 110 byte changes. The changes were cumulative, that is, each test added one (or 10) more perturbations. Table 1 is a sample of some consecutive lines of the corresponding table.

One could define the distance between two signature values as the absolute value of their difference, reflecting the intention of the design of the signature layout to yield changes in the low order bits of the signature as result of small changes in the chunk. However, in the intended application to a deduplication system, one cannot afford too many search attempts in the vicinity of the hash value. More precisely, suppose...
Table 2. Hamming distance with original chunk after 1, 2, ..., 110 perturbations

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A chunk $C$ is given. We would evaluate $ah(C)$ and check whether there is a pointer to a chunk $D$ at the address $H[ah(C)]$ in the hash table. If so and indeed $D = C$, the newly arrived chunk $C$ can be deduplicated by pointing to $D$. But if $D \neq C$, the intention was to look for pointers to chunks identical to $C$ at the neighboring locations of the hash table. But each trial is costly, so the number of trials will have to be restricted. It might possibly only be reasonable to check at $H[ah(C)]$, $H[ah(C) + 1]$ and $H[ah(C) - 1]$. In that case, we can as well restrict ourselves to the Hamming distance between signatures, i.e., the number of differing bits in their standard binary representation, rather than their arithmetic difference. These are the values displayed in Table 2.

The first column gives the original signature (as a decimal number) before applying any perturbation, then in the column headed $i$ is the Hamming distance between the original and the new signatures when $i$ perturbations have been applied. Note that this distance is not always an increasing function of the number of perturbations, indicating that there might be quite a few cases in which the signature tends to “correct itself” when there are many changes; however, the overall trend is clearly increasing, as can be seen in the graphs below.

The plot in Figure 4 shows the average number of changed bits as a function of the number of perturbations, for both the Exchange and the OS databases. The average Hamming distance was between 0.3 and 5 to 6. There are slight differences between the databases, but the trend is the same.

Note that the distances, at least for the small number of perturbations, are quite low, and very often even zero, meaning that very small changes often yield the same signature as before. This is in sharp contrast with regular hashing schemes, for which the corresponding graph is expected to be a straight line (that is, independent of the number of changes as long as this number is $> 0$) at the level of about 16 (that is, about half of the bits are expected to change).

To verify this fact, we devised a control experiment in which a regular hash function (modulo $P = 2^{32} - 17 = 4, 294, 967, 279$), was applied the same blocks, and then...
performed the same perturbation tests as for our function, recording the Hamming distance between the original and perturbed signatures. Note that \( P \) is a prime number (actually the largest one fitting into our unsigned 32 bit signature). As expected, the number of changed bits was indeed around 16, as can be seen on the plot in Figure 5, more precisely, the average values were in the range from 15.988 to 16.018. Figure 5 also displays again the graph for the OS database, for comparison.

For our function, even if there are more than 100 bytes changed, this implies, at the average, only a change in about 5 to 6 bits. The resulting signature might thus be very different (depending on the position of those 6 changed bits), but the change is clearly not as radical as if a regular hash had be applied. In any case, this is just a noteworthy observation as in the intended application, there is no intention to look for similar chunks so far away.

To get a feeling on how far one can insert perturbations without yet changing the signature value, consider the plot in Figure 5, giving for each number \( i \) of perturbations, the probability of getting a non-zero value in the column headed \( i \) of the perturbation table. The plots are again given for both the Exchange and the OS databases. The probability of a non-zero value for a single perturbation is just about 0.06, and we see a clear ascending trend, reaching probability about 0.8 for
more than 100 changes. For the control test with the real hashing function, we again got practically always non-zero values, more precisely, the probability for getting a non-zero for each of the columns was between 0.99986 and 1.00000.

3.3 Clustering

After testing that the proposed signature indeed has the properties required from an approximate hash function, that is: it gives a uniform spread, yet preserves locality in the sense that similar blocks give similar signatures, we turn now to a more general clustering test, which in fact checks the transpose of the above implication, that similar signatures also imply similar blocks.

For each of the tested databases, N centroid chunks have been chosen (we used \(N = 11\)), so that they were mutually not similar. This is achieved by choosing the chunks in a random sequence, and checking for each new candidate that it is different enough from all the preceding chosen chunks in the sequence. \(X\) and \(Y\) are said to be different enough if \(LD(X, Y) \geq T\), where \(T\) is some independently chosen threshold (we used 1000), and \(LD\) defines the Levenshtein distance \([12]\).

Each of the centroids is then used to generate a number \(M\) of perturbation chunks (we used \(M = 10\)), which are obtained by either changing a predetermined number \(K\) of bytes of the map to a random value, or by copying to each of these \(K\) bytes the value of another, randomly chosen, byte value from within the same chunk. The number of perturbations \(K\) has been chosen to vary from 2 to 1024, doubling in each step. Finally, the approximate hashing is applied to each of the generated chunks, and the whole set of \(N \cdot M\) signatures is then sent to a clustering procedure, which partitions the set of signatures, and thereby the set of corresponding chunks, into subset of similar chunks. The number of hits, that is, correctly assigned correlations between a generated chunk and its generating centroid, is recorded as a function of the number \(K\) of perturbations.

Three different alternatives have been considered to perform the clustering: the hierarchical Tree-method (repeatedly choosing the pair of closest chunks among the set of remaining subsets and dynamically updating the sets), K-means (minimizing the within-cluster sum of squares) \([9]\), and simply checking the distance from every generated chunk to each of the centroids and choosing that with minimal distance. The results were similar, with the first method consistently giving slightly better performance.
Each experiment was repeated 10 times and the values averaged. The results for our test databases of the hit ratio as function of the number of perturbations are displayed in Figure 7. As can be seen, the success rate is indeed decreasing with increasing $K$, and for a small number of perturbations, the number of successful assignments may be as large as 95%.

### 3.4 Comparison of similarity with identity

As has been mentioned earlier, the ultimate aim of these hash based systems is to perform deduplication. One approach is to use standard hashing, even with graphically strong hash functions that reduce the probability of false alarms to almost zero, but can thereby detect only identical chunks. The alternative suggested in this paper is the *approximate hash*, which could be able of locating also similar and not necessarily identical data chunks.

It might not be possible to quantify the relative improvement caused by shifting from a system based on identity to one based on similarity: the results will be extremely data dependent, based on the nature of the data and its repetitiveness. It obviously makes no sense to simulate the system’s behavior on random data, as is done for many other applications, since truly random data is not compressible. On the other hand, also compressed files cannot be compressed even further, but they may be able to take advantage of deduplication, for example when several copies of such a file appear in the database.

We therefore decided, by lack of what could be agreed on as being “typical” data, to test the performance in tests on publicly available files and report the results just as examples, without claiming that these results are representative. Indeed, on different input data, the figures could be higher or lower, depending on the data at hand.

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*Table 3. Comparing identity with similarity based systems*
The files we chose were images of virtual machines obtainable from the web, a sample of which is presented in Table 3. The sizes are given in MB, and the columns headed identity and similarity list the corresponding compression ratios. The compression ratio is defined as the size of the original file divided by the size of the compressed file. For identity, we used the SHA1 secure hash function \([7]\) and second and subsequent copies of identical chunks were removed. For similarity, we used our approximate hashing scheme, and in case a similar chunk has been found, the new chunk was delta-encoded using vcdiff \([3]\). Fixed length chunks of size 8 K have been used for both parts of the experiment. The final column lists the relative gain, in percent, of using similarity instead of identity.

Table 4 gives a more specific insight in the distribution of where the matching chunks have been located by our system. We checked first at \(H[ah(C)]\), and if this entry did not contain a pointer to \(C\), we also checked \(H[ah(C) \pm i]\), for \(i = 1, 2, \ldots, 5\). On our example data, in the overwhelming majority of cases among those where the chunk could indeed be deduplicated, the pointer was found at \(H[ah(C)]\) itself. But in 18% of the cases, it was found nearby. As could be expected, the probability of locating the chunk decreases with the distance from \(ah(C)\), but interestingly, the decrease is not monotonic: the values for \(\pm 4\) are larger than for \(\pm 3\). Clearly, this is due to the fact that a difference of 4 means that only one bit is different in the signature, while for a difference of 3, there are two differing bits.

### 4 Conclusion

We have presented the main ideas leading to the design of a similarity rather than identity based deduplication system working with relatively small data chunks. Similarity has been explored earlier in this context \([1]\), but the performance depended critically on the fact that the chunk size could be chosen large enough, in the MB range, which reduced the size of the required data structures. The current work is a first attempt to adapt the similarity approach also to systems in which a more fine grained resolution is required, with data chunks typically in the KB range.

The tests we performed suggest that the proposed approximate hash function indeed combines quite contradicting properties, like uniformity and sensitivity as required, though this can only be empirically tested on chosen examples, and not quantitatively checked in controlled statistical experiments. The scalability of the system will obviously depend on the amount of duplicate data it contains.
References