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# Revisiting and Improving Algorithms for the 3XOR Problem

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**Abstract.** The 3SUM problem is a well-known problem in computer science and many geometric problems have been reduced to it. Here, we study the 3XOR variant which is more cryptologically relevant. In this problem, the attacker is given black-box access to three random functions F, G and H. She has to find three inputs x, y and z such that  $F(x) \oplus G(y) \oplus H(z) = 0$ . The 3XOR problem is a difficult case of the more-general k-list birthday problem.

Wagner's celebrated k-list birthday algorithm and the ones that it inspired work by querying the functions more than strictly necessary from an information-theoretic point of view. This gives them some leeway to target a solution of a specific form, at the expense of processing a huge amount of data.

We restrict our attention to solving the 3XOR problem for which the total number of queries to F is minimal. If F is an n-bit random function, we want to solve the problem with roughly  $\mathcal{O}\left(2^{n/3}\right)$  queries. In this setting, the folklore quadratic algorithm finds a solution after  $\mathcal{O}\left(2^{2n/3}\right)$  operations.

We present a 3XOR algorithm that generalizes an idea of Joux, with complexity  $\mathcal{O}\left(2^{2n/3}/n\right)$ . This algorithm is practical: it is up to  $3\times$  faster than the quadratic algorithm. We also revisit a 3SUM algorithm by Baran-Demaine-Pătraşcu which is asymptotically  $n^2/\log^2 n$  times faster than the quadratic algorithm when adapted to the 3XOR problem, but is otherwise completely impractical.

To gain a deeper understanding of these problems, we embarked on a project to solve actual 3XOR instances for the SHA256 hash function. We believe that this was very beneficial and we present practical remarks, along with a 96-bit 3XOR for SHA256.

Keywords: 3XOR problems, Wagner's algorithm, generalized birthday

## 1 Introduction

The birthday problem is a widely used cryptanalytical tool: given two lists  $\mathcal{L}_1, \mathcal{L}_2$  of bitstrings drawn uniformly at random from  $\{0,1\}^n$ , find  $x_1 \in \mathcal{L}_1$  and  $x_2 \in \mathcal{L}_2$  such that  $x_1 \oplus x_2 = 0$  (the  $\oplus$  notation denotes the bitwise exclusive-or operation). This problem is well understood: a solution exists with constant probability as soon as  $|\mathcal{L}_1| \times |\mathcal{L}_2| \gg 2^n$  holds, and it can be found in  $\mathcal{O}(2^{n/2})$  time by simple algorithms (sorting then scanning  $\mathcal{L}_1$  and  $\mathcal{L}_2$  is a possibility).

Wagner studied in [Wag02], the k-XOR problem and showed that for the 4-XOR problem if we increase the size of the lists from  $2^{n/4}$ , which is the minimal size for having a solution with good probability, to  $2^{n/3}$ , then the number of solutions increases to  $2^{n/3}$  and the problem becomes easier since we do not have to find a needle in a haystack but find one

out of many needles. He showed that it is possible to find one solution in time  $\mathcal{O}(2^{n/3})$ , by searching for a solution that satisfy some property, which leads to the 4-tree algorithm.

In this paper, we are concerned with the 3XOR problem in which there are three lists containing arbitrarily many random bit strings, and where the goal is to find  $(x_1, x_2, x_3) \in \mathcal{L}_1 \times \mathcal{L}_2 \times \mathcal{L}_3$  such that  $x_1 \oplus x_2 \oplus x_3 = 0$ . The specific case where lists are bitstrings and the + is the XOR is also hard [JV13, Vio12]. Nandi [Nan15] exhibited a forgery attack against the COPA mode of operation for authenticated encryption requiring only  $2^{n/3}$  encryption queries and about  $2^{2n/3}$  time. This attack works by reducing the problem of forging a valid ciphertext to that of solving an instance of the 3XOR problem. This attack was later refined by Nikolić and Sasaki [NS14], using an improved 3XOR algorithm, to  $2^{n/2-\varepsilon}$  queries and  $\mathcal{O}\left(2^{n/2-\varepsilon}\right)$  operations, for a small value of  $\varepsilon$ . The 3XOR problem is also interesting for searching parity check relations in fast correlation attack with k=3, while Chose, Joux and Mitton in [CJM02] can only take  $k \geq 4$ .

However, Joux had already proposed a better algorithm five years before Nikolić and Sasaki in [Jou09], which is the best algorithm for the 3XOR problem to this day. The idea is to set  $\mathcal{L}'_i := \{x\mathbf{M}, \ x \in L_i\}$  where  $\mathbf{M}$  is a well-chosen invertible matrix over  $\mathbb{F}_2$ , and then solve the 3XOR problem over  $\mathcal{L}'_1, \mathcal{L}'_2$  and  $\mathcal{L}'_3$ . This results in a solution of the original instance, but the change of coordinates  $\mathbf{M}$  allows a few tricks. Joux's idea was to choose  $\mathbf{M}$  such that n/2 vectors of  $\mathcal{L}'_3$  have their last n/2 bits equal to 0, compute  $\mathcal{L}'_{12} = \mathcal{L}'_1 \bowtie_{n/2} \mathcal{L}'_2$ , the list of the xor of pairs  $\mathcal{L}'_1$  and  $\mathcal{L}'_2$  that matches on the n/2 last bits, and check for collisions between  $\mathcal{L}'_{12}$  and  $\mathcal{L}'_3$ . With correctly chosen list sizes, this yields a  $\sqrt{n/2}$  speedup over Wagner's algorithm. Nikolić and Sasaki [NS14] independently proposed a different variant of Wagner's algorithm: they reduce  $\mathcal{L}_3$  to the elements that have the most frequently-occurring pattern in the  $\ell$  first bits, with  $\ell \simeq n/2$ . Using a probabilistic analysis, notably the well-known maximal bin load when N balls are randomly thrown into N bins [Mit96], they obtain a slightly smaller speedup of  $\sqrt{(n/2)/\ln(n/2)}$ . It seems very hard to combine these two improvements in a new algorithm.

The story is fortunately not over yet. Four years before Joux, an algorithm for the 3SUM problem (over the integers) had been described by Baran, Demaine and Patrascu in [BDP05]. This algorithm, which was the first generic subquadratic algorithm for the 3SUM problem, has a better complexity than all the others, and it can be applied mutatis mutandis to the 3XOR problem with lists of size  $2^{n/3}$  with an expected speedup of  $n^2/\log^2 n$  compared to the folklore quadratic algorithm. On the other hand, the FFT algorithm for 3SUM over the integers given as an exercice in [CLRS01] cannot easily be adapted to the 3XOR problem.

**Our contributions.** We first give a formal definition of the computational model that we consider, which is slightly non-standard... but was apparently adopted by nearly all the previous litterature on generalized birthday algorithms [Wag02, NS14]. We recall several known results about sorting, hashing, etc. which are sometimes implicitly used by previous work.

Armed with these tools, we present a new algorithm, which generalize Joux Algorithm to input lists with arbitrary size. Basically we select a sublist of entries z in the third list, and find an invertible matrix M such that zM start by a fixed number k of zeroes, for all z in this set. Then we search for pairwise elements x,y in the two other lists such that  $(x \oplus y)M$  start by k zeroes. For each of these pairs, we search if  $x \oplus y$  matches an element z of the selected sublist. We iterate the procedure, all over again, until all entries of the third list have been considered. This procedure is always  $\mathcal{O}(n)$  times faster than the quadratic algorithm. We also propose constant-factor improvement of this procedure, that reduces the number of iterations that have to be performed.

We implemented this new algorithm and ran it with input vectors of size n=96 on a 64-bit machine. We used the following strategy: First find all solutions to the 3XOR problem on the first 64 bits of each entries, using our algorithm, then for all triplets of partial solutions thus found, check the last 32 bits of their XOR. In the same conditions, our algorithm run about 2 and 3 times faster than the quadratic algorithm.

We also present a adaptation of the BDP algorithm that we tuned for the 3XOR problem. We show that this algorithm is  $\mathcal{O}\left(n^2/\log^2 n\right)$  faster than the quadratic algorithm, when n is asymptotically large. However, it cannot be used in practice on reasonably small values of n and is mostly of theoretical interest.

We raise several practical considerations that are not really discussed in previous work. For instance, the algorithms designed in [Jou09, NS14] need at least two of the input lists to have size  $2^{n/2-\varepsilon}$ . We claim that, in practice, working with input lists whose size is reduced is better. For instance, to solve the 3XOR problem with 96-bit input vectors, Joux algorithm handles about 0.86 Petabyte of data while Nikolić-Sasaki algorithm handles about 2.55 Petabytes. In this case, even sorting the lists is not very easy. This makes these algorithms quite impractical. We describe algorithms that work faster while requiring only about 384 Megabytes of storage.

We illustrate our result with the computation of a 96-bit 3XOR on the SHA256 hash function. The code used to perform this computation is available from the authors.

# 2 Preliminary

**Notations** Given a list  $\mathcal{L}$ , we denote by  $\mathcal{L}_i$  the *i*-th element of the list, by  $\mathcal{L}[i..j]$  the sublist  $\mathcal{L}_i, \ldots, \mathcal{L}_{j-1}$  and by  $\mathcal{L}[i..]$  the sublist  $\mathcal{L}_i, \mathcal{L}_{i+1}, \ldots$ . We denote by  $\mathcal{L}^{[p]}$  the sublist composed by all vectors whose first bits are p (the size of this prefix is usually clear given the context). For a given vector x, x[i..j] denotes the subvector formed by taking the coordinates from i to j-1. This is similar to the "slice" notation in Python. Let  $\mathcal{A}$  and  $\mathcal{B}$  be two lists. For a given parameter k, we denote by:

$$\mathcal{A} \bowtie_k \mathcal{B} = \{(a, b) \in \mathcal{A} \times \mathcal{B}, \text{ s.t. } a \oplus b \text{ start with } k \text{ zeroes.}\}$$

We denote by L (resp. A, B, C) the size of the list  $\mathcal{L}$  (resp.  $\mathcal{A}$ ,  $\mathcal{B}$ ,  $\mathcal{C}$ ). The log function denotes logarithm in base 2.

#### 2.1 Computational Model and Assumption

We consider that all computations take place inside a RAM (Random Access Machine) model with w-bit words, with  $w = \Theta(n)$  — in other words, the machine is "large enough" to accommodate the problem. We assume that the usual arithmetic and bitwise operations on w bits, as well as memory accesses with w-bit addresses, are elementary operations. This assumption is implicit in previous work on the generalized birthday paradox. This model is sometimes called the transdichotomous model [FW93].

We assume that our machine has black-box access to three oracles  $\mathbf{A}$ ,  $\mathbf{B}$  and  $\mathbf{C}$ . When queried with an n-bit integer i, the oracle  $\mathbf{A}$  returns the n-bit value  $\mathcal{A}_i$  (the i-th element of the list  $\mathcal{A}$ ). The same goes for  $\mathbf{B}$  and  $\mathbf{C}$ . It is understood that, in most cryptographic applications of the 3XOR problem, querying the oracles actually corresponds to evaluating a cryptographic primitive. As such, we assume that the oracles implement random functions.

The machine on which the actual algorithms run is allowed to query the oracles, to store anything in its own memory and to perform elementary operations on w-bit words. An

algorithm running this machine solves the 3XOR problem if it produces a triplet (i, j, k) such that  $A_i \oplus B_j \oplus C_k = 0$ .

The relevant performance metrics of these algorithms are the amount of memory they need (M bits), the number of elementary operations they perform (T) and the number of queries they makes to each oracle.

Most algorithms for the 3XOR problem begin by querying the oracle  $\bf A$  on consecutive integers  $0,1,\ldots,A-1$  (the same goes for  $\bf B$  and  $\bf C$  with respective upper-bounds of B and C) and storing the results in memory. We therefore assume that algorithms start their execution with a "local copy" of the lists, obtained by querying the oracles (except when explicitly stated otherwise).

It is usually simpler in practice to implement algorithms that produce the colliding values  $(A_i, \mathcal{B}_j, \mathcal{C}_k)$  instead of the colliding inputs (i, j, k). The reason for that is that most algorithms sort at least one of the lists, so that  $\mathcal{L}_i$  is not longer at index i in the array holding  $\mathcal{L}$ . It would be possible to store pairs  $(A_i, i)$  in memory, sorting on  $A_i$  and retaining the association with i at the expense of an increased memory consumption. In practice it is much simpler to find the colliding values, then re-query the oracles again to find the corresponding inputs, at the expense of doubling the total number of queries.

To avoid degenerate cases, we assume that the number of queries to the oracles are exponentially smaller than  $2^n$ . More precisely, we assume that there is a constant  $\varepsilon > 0$  such that  $\max\{A,B,C\} < 2^{(1-\varepsilon)n}$ .

# 2.2 Algorithmic Tools

Algorithms for the 3XOR problem process exponentially-long lists of random n-bit vectors. In this section, we review the algorithmic techniques needed to perform three reoccurring operations on such lists: sorting, testing membership and right-multiplication by an  $n \times n$  matrix. We recall that these operations can be performed in linear time.

**Membership Testing.** The simplest algorithm for the 3XOR problem works by considering all pairs  $(a,b) \in \mathcal{A} \times \mathcal{B}$  and checking if  $a \oplus b \in \mathcal{C}$ . This is guaranteed to find all solutions.

It has long been known that hashing allows for dictionaries with  $\mathcal{O}(1)$  access-time and constant multiplicative space overhead [FKS84]. This is enough to make the quadratic algorithm run in time  $\mathcal{O}(AB+C)$ . It is thus beneficial to let  $\mathcal{C}$  be the largest of the three lists.

A simple hash table of size 2C with linear probing allows to test membership with 2.5 probes on average [Knu98, §6.4]. Because the hashed elements are uniformly random, taking some lowest-significant bits yields a very good hash.

Linear hashing is already good, but Cuckoo Hashing improves the situation to 2 probes in the worst case [PR01]: two arrays of size C are sufficient, lookup requires at most one probe in each array and insertion requires an expected constant number of operations on words. In principle, Cuckoo hashing needs a family of universal hash functions: when too many collisions occur during insertion, two new hash functions are randomly chosen and everything is re-hashed with the new hash functions. We may use the usual universal hash functions  $h_{a,b}: x \mapsto (ax+b) \mod p$ , where p is a prime.

We found that Cuckoo hashing was about  $2\times$  faster than linear probing. It allows to make the quadratic algorithm quite efficient, with only a small constant hidden in the  $\mathcal{O}$ 

notation: around 10 CPU cycles are enough to process a pair.

**Sorting and Joining.** The evaluation of the "join"  $\mathcal{A} \bowtie_k \mathcal{B}$  of two lists  $\mathcal{A}$  and  $\mathcal{B}$  of size  $2^{\alpha n}$  is ubiquitous in generalized birthday algorithms following Wagner's ideas. All these algorithms rely on a common idea: they target a solution  $a \oplus b \oplus c = 0$  such that c[0..k] = 0 (for some value of k), which implies that a[0..k] = b[0..k]. These special solutions can be found efficiently by performing a linear-time join operation between  $\mathcal{A}$  and  $\mathcal{B}$ . The following algorithm is thus the workhorse of these techniques.

**Algorithm M** (3XOR with matching prefixes). Given three lists  $\mathcal{A}, \mathcal{B}$  and  $\mathcal{C}$ , where all  $c \in \mathcal{C}$  are such that c[0..k] = 0, returns all triplets  $(a, b, c) \in \mathcal{A} \times \mathcal{B} \times \mathcal{C}$  such that  $a \oplus b \oplus c = 0$ .

- **M1.** [Prepare input] Sort  $\mathcal{A}$  and  $\mathcal{B}$  according to their first k bits. Initialize a hash table with the entries of  $\mathcal{C}$ . Set  $i \leftarrow 0, j \leftarrow 0$ .
- **M2.** [Compare prefix] If i = A or j = B, terminate the algorithm. If  $\mathcal{A}_i[0..k] < \mathcal{B}_j[0..k]$ , increment i and repeat this step. If  $\mathcal{A}_i[0..k] > \mathcal{B}_j[0..k]$ , increment j and repeat this step.
- **M3.** [Save prefix.] Set  $p \leftarrow A_i[0..k]$  and  $j_0 \leftarrow j$ .
- **M4.** [Check pair.] If  $A_i \oplus B_j \in \mathcal{C}$ , then emit  $(A_i, B_j)$ .
- **M5.** [Loop on i, j.] Increment j; if j < B and  $\mathcal{B}_j[0..k] = p$ , return to step M4. Otherwise increment i; if i < A and  $\mathcal{A}_i[0..k] = p$ , then set  $j \leftarrow j_0$  and go back to M4. Otherwise return to M2.

It is well-known that an array of N random k-bit vectors can be sorted in linear time [Knu98, §5.2.5]: the randomness of the input allows to beat the  $\Omega(N \log N)$  lower-bound of comparison-based sorting. Here is a way to do it using  $\mathcal{O}(\sqrt{N})$  extra storage: perform two passes of radix sort on  $0.5 \log N$  bits, then finish sorting with insertion sort.

Each pass of radix sort requires  $\sqrt{N}$  words to store the counters. Besides that, it can permute the input array in-place [Knu98, §5.2.1]. The two passes of radix sort guarantee that the array is sorted according to its first  $\log N$  bits. This reduces the expected number of inversions from  $\approx N^2/4$  to  $\approx N/4$ . Thus, the expected running time of insertion sort will be linear.

This tells us that step M1 runs in time linear in A+B. Steps M2–M3 are repeated at most A+B times and requires a constant number of operations on w-bit words. Steps M4–M5 are executed for each pair  $(a,b) \in \mathcal{A} \times \mathcal{B}$  with a[0..k] = b[0..k]. The expected number of such pairs is  $AB/2^k$ . Considering that we also have to initialize a hash table with the element of C, the total expected running time of the algorithm is thus  $\mathcal{O}\left(A+B+C+AB/2^k\right)$ .

The algorithms designed by Wagner [Wag02], Nikolić–Sasaki [NS14] and Joux [Jou09] use this procedure but differ in how they filter or modify their input to ensure that a solution of the correct form exists with high probability.

**Matrix multiplication.** Given a  $n \times n$  matrix  $\mathbf{M}$  over  $\mathbb{F}_2$ , we need to compute the matrix-matrix product  $\mathbf{LM}$ , where  $\mathbf{L}$  is seen as a  $2^{\alpha n} \times n$  matrix. Performing the product naively would require  $\mathcal{O}(n2^{\alpha n})$  operations. This can improved to  $\mathcal{O}(2^{\alpha n})$  operations, using a trick similar to the "method of Four Russians".

The idea is to divide **M** into slices of  $n\frac{\alpha}{1+\varepsilon}$  rows (for any  $\varepsilon > 0$ ); for each slice, we precompute all the  $2^{\frac{\alpha}{1+\varepsilon}n}$  linear combinations of the rows and store these in a table. This would take  $2^{\frac{\alpha}{1+\varepsilon}n}$  word operations.

Then, the vector-matrix product  $x\mathbf{M}$  can then be evaluated by dividing x in slices of size  $n\frac{\alpha}{1+\varepsilon}$ , looking up the precomputed linear combination of the rows of the corresponding slices of  $\mathbf{M}$ , and xoring together the  $\frac{1+\varepsilon}{\alpha}$  resulting vectors. This step would take only a constant number of word operations and we will need to perform it  $2^{\alpha \cdot n}$  times. Hence, the whole complexity of the procedure is  $\mathcal{O}(2^{\alpha n})$  word operations.

# 3 Finding 3XOR by Linear Changes of Variables

In this section, we propose a new algorithm which is asymptotically n times faster than the quadratic algorithm. It is a generalization of an earlier idea of Joux [Jou09]. While it is asymptotically inferior to the BDP algorithm discussed later, it is faster in practice than the quadratic algorithm for relevant parameter sizes. On the other hand, the BDP algorithm is not practical.

## 3.1 Algorithm Description

This algorithm exploits the fact that when  $\mathbf{M}$  is an  $n \times n$  invertible matrix over  $\mathbb{F}_2$ , then  $a \oplus b \oplus c = 0$  if and only if  $a\mathbf{M} \oplus b\mathbf{M} \oplus c\mathbf{M} = 0$ . The sets of solutions in  $\mathcal{A} \times \mathcal{B} \times \mathcal{C}$  and  $\mathcal{A}\mathbf{M} \times \mathcal{B}\mathbf{M} \times \mathcal{C}\mathbf{M}$  are the same.

The idea of our algorithm is to select a slice of n-k vectors from  $\mathcal{C}$  and to chose a matrix  $\mathbf{M}$  such that the first k entries of the slice become zero. Algorithm  $\mathbf{M}$  then finds all the triplets  $a \oplus b \oplus c = 0$  where c belongs to the slice. Repeating this procedure for all slices yields all the possible solutions.

**Algorithm G** (3XOR by linear changes of variables). Given  $\mathcal{A}, \mathcal{B}$  and  $\mathcal{C}$ , return all  $(a, b, c) \in \mathcal{A} \times \mathcal{B} \times \mathcal{C}$  such that  $a \oplus b \oplus c = 0$ . Let  $k \leftarrow \lceil \log_2 \min(A, B) \rceil$ .

- **G1.** [Loop on i.] Perform steps G2–G4 for  $0 \le i < C/(n-k)$ , then terminate the algorithm.
- **G2.** [Compute change of basis] Set  $u \leftarrow i(n-k)$  and  $v \leftarrow \min\{(i+1)(n-k), C\}$ . Compute an  $n \times n$  matrix **M** such that

$$C[u..v]\mathbf{M} = \begin{pmatrix} 0 & \dots & 0 & \star & \dots & \star \\ \vdots & & \vdots & \vdots & & \vdots \\ 0 & \dots & 0 & \star & \dots & \star \end{pmatrix} \uparrow n-k$$

- **G3.** [Matrix multiplication.] Set  $\mathcal{A}' \leftarrow \mathcal{A}\mathbf{M}$ ,  $\mathcal{B}' \leftarrow \mathcal{B}\mathbf{M}$  and  $\mathcal{C}' \leftarrow \mathcal{C}[u..v]\mathbf{M}$ .
- **G4.** [Join] Run algorithm M to find all pairs (a',b') in  $\mathcal{A}' \times \mathcal{B}'$  such that a'[0..k] = b'[0..k] and  $a' \oplus b' \in \mathcal{C}'$ . For each such triplet, emit  $(a'\mathbf{M}^{-1}, b'\mathbf{M}^{-1})$ .

**Theorem 1.** Algorithm G finds all 3XORs in  $A \times B \times C$  in expected time

$$T = \mathcal{O}\left((A+B)C/n\right)$$

and space linear in A + B + C.

*Proof.* We first observe that the algorithm only needs to store two of copies of the input lists (this could be brought down to a single copy, updated in-place). This establishes the linear space claim. Step G2 can be done in time  $\mathcal{O}(n^3)$ , for instance using the PLUQ factorization  $\mathcal{C}[u..v]^T = \mathbf{PLUQ}$ , where  $\mathbf{P}$  and  $\mathbf{Q}$  are permutation matrices,  $\mathbf{L}$  is lower-triangular with unit diagonal and  $\mathbf{U}$  is upper-trapezoidal. Only the first r rows of  $\mathbf{U}$  are non-zero, where r denotes the rank of  $\mathcal{C}[u..v]$ . It follows that  $\mathbf{M} = \mathbf{P}(\mathbf{L}^T)^{-1}$  is a suitable choice.

Let us assume, without loss of generality, that  $A \leq B$ . Using the techniques described in section 2.2, we know that step G3 costs  $\mathcal{O}(A+B+(n-k))$ . The expected costs of step G4 is  $\mathcal{O}(A+B+AB/2^k)$ . The choice of k at the beginning of the algorithm ensures that this is in fact  $\mathcal{O}(A+B)$ .

Steps G2–G4 are repeated  $C/(n-\log_2 A)$  times, therefore the total complexity of the algorithm is then

$$\mathcal{O}\left(\frac{(A+B)C}{n-\log_2 A}\right).$$

The hypothesis that  $\log_2 A < (1-\varepsilon)n$  guarantees that the denominator is  $\Omega(n)$ , and yields the claimed complexity.

#### 3.2 Constant-Factor Improvements

In step G2, an invertible matrix  $\mathbf{M}$  is found that sends the first k bits of n-k random vectors to zero. In this section, we discuss means to find matrices that have the same effect on *more* than n-k vectors. The result would be that each iteration of steps G2–G4 would process larger slices of  $\mathcal{C}$ . As such, less iterations would be needed.

#### 3.2.1 Finding the Coordinate Change

Given a large collection  $\mathcal{C}$  of independent and uniformly random n-bit vectors x, we are facing the problem of finding a matrix  $\mathbf{M}$  that maximize (or at least increases) the number of  $x \in \mathcal{C}$  such that  $x\mathbf{M}[0..k] = 0$ . In more algebraic terms, let  $V_0$  be the (n-k)-dimensional subspace of  $\{0,1\}^n$  containing all vectors whose first k coordinates are zeroes. The matrix  $\mathbf{M}$  should be chosen such that it sends the largest number of input vectors from  $\mathcal{C}$  to  $V_0$ . Alternatively, let V be the pre-image of  $V_0$  through  $\mathbf{M}$ :  $x\mathbf{M}$  belongs to  $V_0$  if and only if x belongs to V.

Finding a subspace V having a large intersection with the input list is the actual difficult task. Indeed, once a basis  $(b_1, \ldots, b_{n-k})$  of V has been found, building an invertible matrix M that sends V to  $V_0$  is easy. We therefore face the following computational problem:

**Problem 1.** Given a list C of uniformly random vectors in  $\{0,1\}^n$ , find a (n-k) dimensional subspace V of  $\{0,1\}^n$  such that  $V \cap C$  is as large as possible.

In other terms, we are looking for multiple simulatenous linear approximations of the random vectors in  $\mathcal{C}$ . Note that we have to find not only one, but a large number such subspaces (one per iteration of algorithm G).

In the rest of this section, we discuss two way to tackle this problem. The first approach consists in finding all these subspaces at once, during a pre-computation step, then permuting  $\mathcal{C}$  so that for known parameters  $N_0, N_1, \ldots$  the first  $N_0$  entries of  $\mathcal{C}$  span the first subspace, the following  $N_1$  span the second subspace, and so on.

The second approach is iterative: at each iteration of the procedure, we consider the list  $C_{left}$  of the elements of C that we have not treated yet, we search a solution to problem 1 with  $C_{left}$  as input list.

In both cases, we want this extra-computation to be much faster than the actual 3XOR algorithm that it is supposed to speed-up. As such, we give up on finding optimal solutions and instead look for heuristic that produce quick results. We propose two approaches to obtain better results.

#### 3.2.2 All-at-Once Approach Using Wagner's Algorithm

The first method we propose pre-compute sets of linearly dependant vectors in  $\mathcal{C}$ , and then permute the list, so these linearly dependant vectors are stacked together. In the interesting case where  $C=2^{n/2}$ , this enable us to reduce the number of iterations in algorithm G by 25% in exchange for a one-time pre-computation of negligible complexity. We discuss a situation where this happens naturally in section 5.3.

Wagner's celebrated 4-tree algorithm can be used to find quadruplets of distinct indices (r, s, t, u) such that  $C_r \oplus C_s \oplus C_t \oplus C_u = 0$ . These four entries of C are then linearly dependent.

The main idea of this procedure is to find all pairs r < s such that  $C_r \neq C_s$  and  $C_r \oplus C_s$  starts with k zeroes. The expected number of such pairs is  $\mathcal{O}\left(2^k\right)$ . Then, the expected number of distinct couples (r,s) < (t,u), such that  $C_r \oplus C_s \oplus C_t \oplus C_v = 0$  is  $\mathcal{O}\left(2^{3k-n}\right)$ . For instance, if  $C = 2^{n/2}$ , we are in the interesting case where we will find about  $\mathcal{O}\left(2^{n/2}\right)$  such quadruplets.

We then need to isolate a (large) subset of pairwise disjoint quadruplets. Finding the biggest possible subset is the 4D MATCHING problem, which is NP-complete. However, any maximal 4D matching is a a reasonably good approximation, and it can be found efficiently by a greedy algorithm. Let  $\mathcal{M}$  denote a 4D matching, and r denote its cardinality.

We build a permuted list  $\mathcal{D}$  as follows: start from an empty list; for each  $\{r, s, t, u\} \in \mathcal{M}$ , append  $\mathcal{C}_r, \mathcal{C}_s, \mathcal{C}_t$  and  $\mathcal{C}_u$  to  $\mathcal{D}$ . Finally, append  $\mathcal{C} - \mathcal{D}$  to  $\mathcal{D}$ , in any order. Algorithm G can be updated to exploit  $\mathcal{D}$ , with a running time reduced by 25%.

**Algorithm G'** (3XOR by linear change of variables with 4SUM precomputation). Given  $\mathcal{A}, \mathcal{B}$  and  $\mathcal{D}$  (as computed above), return all  $(a, b, d) \in \mathcal{A} \times \mathcal{B} \times \mathcal{D}$  such that  $a \oplus b \oplus d = 0$ . Let  $k \leftarrow \lceil \log_2 \min(A, B) \rceil$ .

- **G'1.** [Loop on i.] Perform steps G'2–G'4 for  $0 \le i < 3r/(n-k)$ , then go to G'5.
- **G'2.** [Compute change of basis] Set  $u \leftarrow 4i(n-k)/3$  and  $v \leftarrow 4(i+1)(n-k)/3$  (note that  $\mathcal{D}[u..v]$  has rank n-k even though it has 4(n-k)/3 rows). Compute an  $n \times n$  matrix  $\mathbf{M}$  such that

$$\mathcal{D}[u..v]\mathbf{M} = \begin{pmatrix} 0 & \dots & 0 & \star & \dots & \star \\ \vdots & & \vdots & \vdots & & \vdots \\ 0 & \dots & 0 & \star & \dots & \star \end{pmatrix} \underbrace{\frac{4}{3}(n-k)}_{k}$$

- **G'3.** [Matrix multiplication.] Set  $\mathcal{A}' \leftarrow \mathcal{A}\mathbf{M}$ ,  $\mathcal{B}' \leftarrow \mathcal{B}\mathbf{M}$  and  $\mathcal{D}' \leftarrow \mathcal{D}[u..v]\mathbf{M}$ .
- **G'4.** [Join] Run algorithm M to find all pairs (a',b') in  $\mathcal{A}' \times \mathcal{B}'$  such that a'[0..k] = b'[0..k] and  $a' \oplus b' \in \mathcal{D}'$ . For each such triplet, emit  $(a'\mathbf{M}^{-1}, b'\mathbf{M}^{-1})$ .

**G'4.** [Finish] Run algorithm G to find all remaining 3-sums in  $\mathcal{A} \times \mathcal{B} \times \mathcal{D}[4r..]$ .

#### 3.2.3 One-at-a-Time Approach Using Decoding Algorithms

The second method we present is an iterative way of finding these vectorial subspace. We discuss how to find *one* hopefully large sublist of the initial list C, that span a (n - k) dimensional subspace, knowing that this procedure will have to be repeated many times with decreasing list sizes.

Recall that an (n-k)-dimensional subspace of  $\{0,1\}^n$  is defined by a system of k linear equations in n variables. Finding a vector space V having a large intersection with C amounts to finding k linear equations that are simultaneously biased over C (i.e. more often simultaneously true than the contrary).

We propose a greedy approach to find these equations: first initialize a list  $\overline{\mathcal{C}}$  to  $\mathcal{C}$ , find a biased equation  $E_1$  over  $\overline{\mathcal{C}}$ , then remove from  $\overline{\mathcal{C}}$  all vectors that do not satisfy  $E_1$ , and re-iterate the method k-1 times. This reduce the problem to that of finding a single biased equation over  $\mathcal{C}$ .

Finding the most biased equation over  $\overline{\mathcal{C}}$  is tempting but too expensive: an exhaustive search would require  $\mathcal{O}\left(2^{n+k}\right)$  operations and FFT-like methods such as the Walsh transform still have a workload of order  $\mathcal{O}\left(k\cdot 2^n\right)$ . We have to settle for a "good", if not optimal bias.

Finding a biased linear equation over C is a decoding problem. Consider the binary linear code spanned by the columns of C, and let y be a low-weight codeword: there is a vector x such that y = Cx and y has a low hamming weight (amongst all possible vectors y). This means that x describes the coefficient of one of the most biased linear equation over C.

Information Set Decoding. A linear binary (error-correcting) code of length n and dimension d is a subspace of  $\mathbb{F}_2^n$  of dimension d. It contains  $2^d$  codewords of n bits. A linear code can be seen as the linear span of the rows of an  $d \times n$  matrix G called the generator matrix of the code. The weight of a codeword is its hamming weight. The minimum-distance of the code is the minimum weight of non-zero codewords. Determining the minimum-distance of a linear code is NP-complete in general. The first algorithm for the minimum-weight codeword problem has been first introduced by McEliece [McE78] in the security analysis of his cryptosystem, and has been widely studied since. One of the simplest way to recover a minimum-weight codeword is to use the Las Vegas randomized Information Set Decoding (ISD) procedure. Although this algorithm has been improved several times [BJMM12, MMT11, MO15], we found out that in our particular case, the Lee-Brickell algorithm [LB88] is nearly as efficient as its later optimizations, while being easier to implement.

Let C be a random binary linear code of dimension k and length N. We denote by d the minimum distance of C and by  $\delta := d/N$  its relative distance. Let R := k/N be the rate of C, and let  $\mathcal{H}$  denote the binary entropy function. The time complexity of the Lee-Brickell algorithm has been analyzed in [CG90] and is about  $\mathcal{O}\left(2^{N\cdot F(R)}\right)$  with:

$$F(R) := (1 - R) \cdot (1 - \mathcal{H}(\delta/(1 - R))). \tag{1}$$

As C is a random code, its minimum distance d is close to the Gilbert-Varshamov bound [Gil52, Var57], and thus we assume that  $\mathcal{H}(\delta) = 1 - R$ .

The function F reaches a maximum of 0.1207 when  $R \simeq 0.454$ . The further improvements of the original Lee-Brickell algorithm mostly aim at reducing its worst case complexity.

However, in our case, we will have to find minimum-weight words in codes of very low rate. As such the values of F(R) we encounter are much smaller.

Using ISD, one can find codewords c of Hamming weight wt(c) = d, where d is the minimum distance of the code. However, the complexity of such algorithms is exponential in the length of the code (the number of vectors in the input list), and therefore, considering our time budget of  $\mathcal{O}(C)$ , we have to set an upper bound  $N_{max}$  on the size of the list that we consider during the procedure. The value of  $N_{max}$  is determined below.

Hence, before using decoding algorithms, we have to reduce the size of the input list below  $N_{max}$ . To this end, we focus the search on the vectors of  $\mathcal{C}$  whose  $\lceil \log C/N_{max} \rceil$  first bits are zero: their expected number is approximaltely  $N_{max}$ , and they already satisfy  $\lceil \log C/N_{max} \rceil$  linear equations...

**Procedure description.** Let  $N_{max}$  and  $\varepsilon$  be fixed parameters such that  $N_{max}$  represent an upper bound on the length of the initial code  $C_0$ , and threshold parameter  $\varepsilon$ . The following procedure finds a set of k biased equations:

**Algorithm F** (*Greedily find simultaneous linear approximations*). Given a list C (sorted in ascending order), two parameters  $N_{max}$  and  $\varepsilon$ , return a (hopefully large) sublist  $\overline{C}$  such that all entries of  $\overline{C}$  satisfy k linear equations.

- **F1.** [Initialization.] Set  $\ell \leftarrow \lceil \log C/N_{max} \rceil$ . Let j be the biggest integer such that  $C_j[0..\ell] = 0$ . Set  $\overline{C} \leftarrow C[0..j]$  and  $t \leftarrow \ell$  (t is the number of equations currently found).
- **F2.** [Loop.] Repeat steps F3–F4 while t < k, then return  $\overline{\mathcal{C}}$  and terminate the algorithm.
- **F3.** [Find equation.] Estimate d, the minimum distance of the code spanned by  $\overline{\mathcal{C}}^T$  (i.e. the matrix whose columns are the entries of  $\overline{\mathcal{C}}$ ), then use the Lee-Brickell ISD algorithm to find a non-zero codeword c, such that  $wt(c) \leq d + \varepsilon$ . Increment t.
- **F4.** [Filter.] For all i such that  $c_i = 1$ , remove  $\overline{C}_i$  from  $\overline{C}$ .

**Theorem 2.** Algorithm F returns a subset of C that spans a (n-k)-dimensional subspace of  $\{0,1\}^n$ .

*Proof.* We prove the following invariant: each time step F3 begins,  $\overline{\mathcal{C}}$  is contained inside a subspace  $E_t$  of dimension n-t.

At the end of step F1,  $\overline{\mathcal{C}}$  belongs to the subspace  $E_{\ell}$  composed by all the vectors whose first  $\ell$  coordinates are zero, so the invariant holds the first time step F3 is performed.

Because c is a codeword that belongs to the code spanned by the columns of  $\overline{\mathcal{C}}$ , there exists x such that  $c = \overline{\mathcal{C}}x$ . Step F4 is to remove from  $\overline{\mathcal{C}}$  the vectors that are not orthogonal to x. It follows that after step F4,  $\overline{\mathcal{C}}$  is contained into the subspace  $E_{t+1} = E_t \cap \{0, x\}^{\perp}$ . Because c is non-zero, at least one vector is removed from the list. This vector belonged to  $E_t$ , and does not belong to  $E_{t+1}$ . As such, the dimension of  $E_{t+1}$  is (at least) one less than that of  $E_t$ .

**Estimation of**  $N_{max}$ . Knowing C the size of the initial list C it is possible to estimate the upper bound  $N_{max}$  over the length of the code that the ISD algorithm can process within our time budget.

Let us denote by f the function  $f(x,y) := x \cdot F(y/x)$  (for 0 < y < x), where F is the function given by equation 1. A minimum-weight codeword in a code C of length  $N_{max}$  and dimension  $n-\ell$  can be found in time  $\mathcal{O}\left(2^{f(N_{max},n-\ell)}\right)$ .

n	k	expected $N_0$	expected $\overline{C}$
64	28	382	81
128	60	1996	154
256	123	9894	285
512	252	60271	545

Table 1: Parameter estimation of the ISD-based algorithm for some value of n and k

If we denote by  $T_{ISD}$ , the total time spent in the  $k-\ell$  iterations of step F3, we claim that

$$T_{ISD} \le (k - \ell) \cdot 2^{f(N_{max}, n - \ell)}. \tag{2}$$

The function f(x, y) grows with x, and also grows with y when  $y < 0.1207 \cdot x$ . This is enough to prove inequality 2, assuming that  $N_{max}$  is much larger than  $n - \ell$ .

To be sure that the complexity of algorithm F does not exceed  $\mathcal{O}(C)$ , we want to find the maximum value of  $N_{max}$  so that  $T_{ISD} < C$ . This can be done numerically (for instance using the bisection algorithm). Table 1 gives an estimation of the parameter for some values of n and k. With k a little lower than n/2, we can hope to find at least n vectors of C in the first subspace (this is to be compared with n/2 without this technique).

# 4 Adaptation of BDP Algorithm

In this section, we revisit an algorithm that was introduced by Baran, Demaine and Pătraşcu (BPD) in [BDP05]. Initially this algorithm was designed for the 3SUM problem over  $(\mathbb{Z}, +)$ , where the size of the three lists is bounded by some parameter N. Their goal was to determine whether a 3SUM exists in the input lists or not and to return it, when appropriate. Their algorithm is subquadratic in N.

The adaptation described here is asymptotically  $n/\log^2 n$  times faster than the algorithm we present in section 3, but is hardly practical. We consider three lists  $\mathcal{A}$ ,  $\mathcal{B}$  and  $\mathcal{C}$  of size  $2^{n/3}$  (this can be generalized to different size of input lists). The high level idea is the following:

- 1. Dispatch  $\mathcal{A}$ ,  $\mathcal{B}$  and  $\mathcal{C}$  in buckets according to their first k bits, and let m denote the expected number of element in each bucket.
- 2. For each triplets of buckets  $(\mathcal{A}^{[i]}, \mathcal{B}^{[j]}, \mathcal{C}^{[i \oplus j]})$ , perform a preliminary constant time test. The test returns false when it is certain that no solution will be found in this triplet.
- For each triplet that has not been rejected, use the quadratic algorithm on this reduced instance.

A full description of the procedure is given in appendix A.

### 4.1 Preliminary Test

Let  $w = \Theta(n)$  denote the size of a word. Let s be such that  $s = \kappa_1 \cdot \log(w)$ , with  $\kappa_1$  a constant to be determined. Let k be a parameter such that k < n/3, and  $P \ge 1$  a parameter to be determined, such that  $3 \cdot P \cdot s \le w$ .

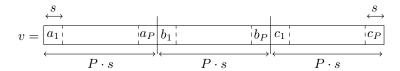


Figure 1: Representation of an index v of the table T

Let T be a table of  $2^{3 \cdot P \cdot s}$  elements. For each index v, that can be represented by figure 1,

$$T[v] = \begin{cases} 1 & \text{if } \exists i, j, k \text{ s.t. } a_i \oplus b_j \oplus c_k = 0 \\ 0 & \text{otherwise.} \end{cases}$$
 (3)

Then, for all three sets of s-bit vectors, that contains each at most P elements, one can check if a solution to the 3XOR problem exists by a constant look-up in T.

By construction,  $|T| = 2^{3 \cdot P \cdot s}$ . We do not want the additional space to exceed the space required to store the lists. We set:

$$3 \cdot P \cdot s = \min(1/3 \cdot n, w). \tag{4}$$

For all  $x \in \mathbb{F}_2^n$ , let h be the function that returns s consecutive bits of x starting from bit k or formally  $h: x \to x[k..k+s]$ . For all  $0 \le \ell < 2^k$  and we denote by  $h(\mathcal{A}^{[\ell]})$  the list

$$h(\mathcal{A}^{[\ell]}) = \{ a = h(x), \text{s.t.} x \in \mathcal{A}^{[\ell]} \}.$$

We define  $h(\mathcal{B}^{[\ell]})$  and  $h(\mathcal{C}^{[\ell]})$  accordingly for all values of  $\ell$ .

Note that a triplet of buckets  $(\mathcal{A}^{[\ell]}, \mathcal{B}^{[t]}, \mathcal{C}^{[r]})$  may contain a solution of the 3XOR problem, only if  $r = \ell \oplus t$ . For all triplet of buckets  $(\mathcal{A}^{[\ell]}, \mathcal{B}^{[t]}, \mathcal{C}^{[\ell \oplus t]})$ , the idea of the algorithm is to check if a solution may exist or not by first checking if there is a solution to the instance  $(h(\mathcal{A}^{[\ell]}), h(\mathcal{B}^{[t]}), h(\mathcal{C}^{[\ell \oplus t]}))$ .

If this test fail, then we know for sure, that there is no solution in  $(\mathcal{A}^{[\ell]}, \mathcal{B}^{[t]}, \mathcal{C}^{[\ell \oplus t]})$ , and we can go to the next instance, without having to perform any other operation. On the other hand, if the test pass, we only know that there is a solution to the 3XOR problem with probability  $(1/2)^{n-k-s}$ .

Solving the small instances  $(h(\mathcal{A}^{[\ell]}), h(\mathcal{B}^{[t]}), h(\mathcal{C}^{[\ell \oplus t]}))$  can be done in constant time, using table T.

Remark 1. One can notice that if a bucket contains more than P vector then these additional vectors are not considered during the preliminary test. Thus, we have to treat them separately afterward. However, if we choose P to be equal to  $\kappa_2 \cdot m$ , with a well chosen  $\kappa_2$ , we can ensure that the number of buckets that will contain more than P element is very small.

Using Chernoff bounds, we figured out that, if we choose  $\kappa_2$  to be around 4.162, the probability that a bucket will contain more than P element will be  $2^{-4 \cdot m}$ .

Remark 2. Taking this in consideration, we have:

$$m = 2^{n-k} = \Theta(n/\log n). \tag{5}$$

This is a direct consequence of the definition of P and s and of equation 4.

## 4.2 Complexity Analysis

**Theorem 3.** With input list of size  $2^{n/3}$ , the time complexity of the BDP algorithm in our model is

$$\mathcal{O}\left(\frac{2^{2n/3} \cdot \log^2(n)}{n^2} + N_{test} \cdot \frac{n^2}{\log^2 n}\right),\,$$

where  $N_{test}$  is the number of triplets that pass the test. The expected value of  $N_{test}$  is:

$$N_{test} \simeq \left(\frac{2^{2 \cdot n/3}}{m} - 1\right) \cdot \left(1 - \left(1 - \frac{1}{w^{\kappa_1}} \cdot \left(1 - \frac{1}{w^{\kappa_1}}\right)\right)^{\Theta(n^3/\log^3(w))}\right) + 1.$$

When n grows to infinity,  $N_{test}$  is equivalent to 1. Thus, the asymptotic complexity of this algorithm is

$$\mathcal{O}\left(\frac{2^{2n/3} \cdot \log^2 n}{n^2}\right).$$

Before we prove this theorem, we need first to introduce some intermediate results.

**Lemma 1.** Let us denote by  $N_{test}$  the number of triplet  $(\mathcal{A}^{[t]}, \mathcal{B}^{[t \oplus \ell]}, \mathcal{C}^{[\ell]})$  The time complexity of the algorithm 1 is:

$$T_{BDP} = \mathcal{O}\left(\frac{2^{2n/3} \cdot \log^2 n}{n^2} + N_{test} \cdot \frac{n^2}{\log^2 n}\right).$$

*Proof.* The time complexity of the full procedure is:

$$T_{BDP} = T_{dispatch} + 2^{2 \cdot k} \cdot T_{test} + N_{test} \cdot T_{solve} + T_{additional},$$

where:  $T_{dispatch}$  is the time it takes to dispatch the elements of the three lists according to their k first bits, As discussed in section 2.2,

$$T_{dispatch} = \mathcal{O}(A + B + C) = \mathcal{O}\left(2^{n/3}\right).$$

 $T_{test}$  is the time complexity of testing one sub-instance  $(\mathcal{A}^{[t]}, \mathcal{B}^{[\ell \oplus t]}, \mathcal{C}^{[\ell]})$ . This is constant time, by a lookup in a precomputed table. We will have to perform this test for all triplets  $(A^{[t]}, B^{[\ell \oplus t]}, C^{[\ell]})$ , that means  $2^{2 \cdot k}$  times.

 $T_{solve}$  is the time required to solve one small instance  $(\mathcal{A}^{[t]}, \mathcal{B}^{[\ell \oplus t]}, \mathcal{C}^{[\ell]})$ .

$$T_{solve} = \mathcal{O}\left(\left(C^{[\ell]} + A^{[t]} \cdot B^{[\ell \oplus b]}\right)\right).$$

Furthermore, as  $C^{[\ell]}$ ,  $A^{[t]}$  and  $B[\ell \oplus b]$  are all  $\mathcal{O}(m)$ .

$$T_{solve} = \mathcal{O}\left(m^2\right)$$
.

 $T_{additional}$  is the time required for additional search, when a bucket contains more than P element. As we have chosen P so that this event is very unlikely, this can be neglected.

All in all we obtain:

$$T_{BDP} = \mathcal{O}\left(2^{n/3} + 2^{2 \cdot k} + N_{test} \cdot m^2\right).$$

In addition, we have  $2^k = 2^{n/3}/m$  and  $m = \Theta(n/\log(n))$ . This is enough to conclude the proof of this lemma.

Lemma 2.

$$N_{test} \simeq \left(2^{2 \cdot k} - 1\right) \cdot \left(1 - \left(1 - \frac{1}{w^{\kappa_1}} \cdot \left(1 - \frac{1}{w^{n-\kappa_1}}\right)\right)^{\Theta(n^3/\log^3(w))}\right) + 1.$$

*Proof.* Let us consider a sub-instance  $(\mathcal{A}^{[t]}, \mathcal{B}^{[\ell \oplus t]}, \mathcal{C}^{[\ell]})$ . Let  $(a, b, c) \in \mathcal{A}^{[t]} \times \mathcal{B}^{[\ell \oplus t]} \times \mathcal{C}^{[\ell]}$ . Recalling that h(a) (resp. h(b), h(c)) represent s arbitrary chosen bits of a (resp. b, c), that are uniformly random, the probability p that  $h(a) \oplus h(b) = h(c)$ , knowing that  $a \oplus b \neq c$  is:

$$p = \frac{1}{2^s} \cdot (1 - \frac{1}{2^{n-s}}).$$

Using Chernoff bounds, we know that the size of the small lists  $h(\mathcal{A}^{[t]})$ ,  $h(\mathcal{B}^{[\ell \oplus t]})$  and  $h(\mathcal{C}^{[\ell]})$  are  $\Theta(m)$  with high probability. Then, we can estimate the probability that a given instance is a false positive to the preliminary test is:

$$\mathbb{P}[\text{pass while should not}] = \left(1 - (1 - p)^{\Theta(m^3)}\right),$$

with  $m = \Theta(n/\log(w))$ .

Furthermore, given the size of the list, we expect only to find one solution. Thus, we expect only one true positive in the preliminary test. Then, there should be only *one* triplet among  $2^{2 \cdot k}$  that contain a solution. The  $2^{2 \cdot k} - 1$  other are expected to contain none.

From here, we can estimate that  $N_{test}$  is:

$$N_{test} = (2^{2 \cdot k} - 1) \cdot \left(1 - (1 - p)^{\Theta(n^3/\log^3(w))}\right) + 1.$$

**Lemma 3.** If  $w(n) \in \Theta(n)$ , and  $\beta$  is a constant, then Choosing  $\kappa_1 = 3$ ,  $N_{test} \underset{0^+}{\sim} 1$ .

Proof. We set:

$$F(n) = \left(1 - \frac{1}{w(n)^{\kappa_1}} \left(1 - \frac{1}{w(n)^{n-\kappa_1}}\right)\right)^{\beta \cdot n^3 / \log^3(w(n))}.$$

Then

$$\ln(F(n)) = \beta \cdot \frac{n^3}{\log^3(w(n))} \cdot \ln\left(1 - \frac{1}{w(n)^{\kappa_1}} \left(1 - \frac{1}{w(n)^{n-\kappa_1}}\right)\right).$$

As  $w(n) \in \Theta(n)$ , with  $\kappa_1 > 0$ :

$$\lim_{n \to +\infty} \frac{1}{w(n)^{\kappa_1}} - \frac{1}{w(n)^n} = 0.$$

Thus,

$$\ln(1 - (\frac{1}{w(n)^{\kappa_1}} - \frac{1}{w(n)^n})) \underset{+\infty}{=} - \left(\frac{1}{w(n)^{\kappa_1}} - \frac{1}{w(n)^n}\right) + \mathcal{O}\left(\frac{1}{n^{2\kappa_1}} + \frac{1}{n^n}\right) \underset{+\infty}{=} -\frac{1}{w(n)^{\kappa_1}} + \mathcal{O}\left(\frac{1}{n^{2\kappa_1}}\right).$$

Then,

$$\ln(F(n)) \underset{+\infty}{=} -\beta \cdot \frac{n^3}{\log^3(w(n))} \cdot \frac{1}{w(n)^{\kappa_1}} + \mathcal{O}\left(\frac{1}{n^3 \log^3(n)}\right).$$

As  $w(n) \in \Theta(n)$ , there are two constant  $\beta_1, \beta_2$  strictly greater than 0, and a certain  $n_0$  such that for all  $n \ge n_0$ ,

$$-\beta_1 \cdot \frac{n^{3-\kappa_1}}{\log^3(w(n))} + \mathcal{O}\left(\frac{n^{3-2\kappa_1}}{\log^3(n)}\right) \le \ln(F(n)) \le -\beta_2 \cdot \frac{1}{\log^3(w(n))} + \mathcal{O}\left(\frac{n^{3-2\kappa_1}}{\log^3(n)}\right).$$

Choosing  $\kappa_1 \geq 3$  will ensure  $\ln(F(n)) \to 0$  when n grows up to infinity. We choose then  $\kappa_1 = 3$ .

From here we can deduce the following equation:

$$\lim_{n \to +\infty} \left( 1 - \frac{1}{w(n)^3} \left( 1 - \frac{1}{w(n)^{n-3}} \right) \right)^{\beta \cdot n^3 / \log^3(w(n))} = 1, \tag{6}$$

Or more simply:

$$(1 - 1/p)^{\Theta(n^3/\log^3(w))} = 1$$

The proof lemma 3, is trivial from here.

Theorem 3 is a direct consequence of all these results.

# 5 Practical Considerations and Experimental Results

#### 5.1 An Academic Exercise

During the preparation of this work, we implemented and compared the relative performances of several algorithms to solve the 3XOR problem. We tried to answer the question: "if someone actually wanted to solve a concrete instance of the 3XOR problem, what would she do?". This is in fact ill-formulated; in all known applications of generalized birthday algorithms, we not only have to solve the instance, but also to create it in the first place. Often, some tradeoffs can be made. This is the case in the aforementioned application to the COPA mode of operation for authenticated encryption: we may either assemble three lists of size  $2^{n/3}$  (and solve the 3XOR instance in time  $\approx 2^{2n/3}$ ) or assemble three lists of size  $2^{n/2}$  (and solve the 3XOR instance in time  $\approx 2^{n/2}$ ).

To better understand these tradeoffs, we chose to tackle an academic "practical" problem: computing a 3XOR on the SHA256 hash function reduced to n bits, for the largest possible value of n.

Most of the literature devoted to the generalized birthday problem is mostly theoretical, in particular because the exponential space requirement of these algorithms makes them quite impractical. One notable exception is  $[BLN^+09]$ , which provides the source code of a high-quality implementation of Wagner's k-list algorithm. Studying this code was enlightening for us.

**Space/Data Constraints are the Hardest.** The algorithms of Joux and Nikolić-Sasaki compute the join of two lists of size  $2^{n/2-\epsilon}$ . Storing and/or moving around such a massive amount of data is the main limiting factor. For instance, with n=96, each list of  $2^{48}$  12-byte entries requires 3 Petabyte of storage. Even by exploiting the fact that  $\mathcal{L}_i$  can be re-computed on-the-fly and does not strictly need to be stored, we came short of a way to compute the join in reasonable time.

**Data/Memory Tradeoffs.** This obstacle can be sidesteped using a well-known technique called "clamping" in [Ber07]. Let us write  $A=2^{\alpha n}, B=2^{\beta n}$  and  $C=2^{\gamma n}$ . Let  $\kappa=(\alpha+\beta+\gamma-1)/2$  (by hypothesis,  $\kappa\geq 0$ ). The idea is that when querying the oracles, we may reject values that are not zero on the first  $\kappa n$  bits – naturally, this only works when  $\kappa\geq \min(\alpha,\beta,\gamma)$ .

This reduces the amount of data that has to be stored in memory by a factor  $2^{\kappa n}$ , and it preserves the existence of a single solution: the product of the sizes of the three "filtered" lists is exactly  $2^{(1-\kappa)n}$ , and entries are  $(1-\kappa)n$  bits wide.

If  $2^{\alpha n}$  queries are allowed to each oracle, then a solution will be found by algorithm G in time  $\mathcal{O}\left(2^{(1-\alpha)n}/n\right)$  using  $\mathcal{O}\left(2^{\frac{1-\alpha}{2}n}/n\right)$  words of memory. The critical case is  $\alpha=\frac{1}{2}$ , where the running time is  $\mathcal{O}\left(2^{n/2}/n\right)$  and storing the inputs lists requires  $\mathcal{O}\left(2^{n/4}\right)$  words. This is the same asymptotic running time as Joux's algorithm, but with a less stringent space pressure. This shows that algorithm G, combined with clamping, is always at least as efficient as Joux's algorithm, and always more efficient that Nikolić-Sasaki's algorithm.

With n=96, this strategy requires  $2^{49.6}$  evaluations of SHA256 to create the lists, about  $2^{48}$  operations to compute the 3XOR using 576 Megabyte of storage. It was practically executed.

The BDP Algorithm is Completely Impractical. While it is asymptotically more efficient, the BDP algorithms fails in practice for reasonable values of n (e.g. n = 96). In fact, for n = 96, and w = 64, from equation 4, we have:

$$3 \cdot P \cdot n = 3 \cdot \kappa_1 \cdot \kappa_2 = 32.$$

Choosing  $\kappa_1 = 3$  as in lemma 3, and  $\kappa_2 = 4.162$ , as in remark 1, we obtain:

$$m = \frac{32}{3 \cdot 4.162 \cdot \log 64} = \frac{32}{74.916} \simeq 0.427.$$

The best we could hope, in that case, is to process "batches" composed of... a half-entry!

Creating the Lists Cannot be Neglected. A single core of a modern CPU is capable of evaluating SHA256 on single-block inputs about 20 million times per second, so that creating the lists takes 11,851 CPU hours sequentially. Solving the 3XOR instance using our implementation of the quadratic algorithm takes 340 CPU hours sequentially, while algorithm G takes 105 hours. Generating the input lists is  $100 \times$  slower than processing them!

It would have been smarter to do clamping on 22 bits instead of 24. This would have reduced the total running time by  $2.5\times$ , at the expense of making the lists 4 times bigger. This is ultimately dependant on the speed at which the "oracles" can be evaluated.

## 5.2 Clusters, Cache, Registers and Other Gory Details

All our implementations are written in plain C, and are relatively concise, totaling 1000 lines of code (not including one-shot tools to generate the hashes, check for potential collisions, etc. which add 1600 additional lines). This section details some of our choices.

n=96 on 64-bit Machines. The transdichotomous RAM model is practical to think about the complexities of algorithms, but it hits limitations when one try to actually implement things. We used the following two-step strategy to cope with values of n larger than 64: first, find and store all triplets (i,j,k) such that  $\mathcal{A}_i[0..64] \oplus \mathcal{B}_j[0..64] \oplus \mathcal{C}_k[0..64] = 0$ . This mean running the quadratic algorithm or algorithm G, and only dealing with 64-bit quantities. This allow for simple and efficient implementations.

In a second step, check if these partial solutions extend to full solutions. This necessitates to deal with the full n bits, but is comparatively much simpler than the first step. In addition, the expected number of partial solutions is small.

**The Quadratic Algorithm.** The summary description given in section 2.2 suggests to allocate a hash table that holds the whole  $\mathcal C$  list. Most accesses to this hash table are likely to incur the penalty of a cache miss, and this can be avoided at no cost. We dispatch the entries of  $\mathcal A, \mathcal B$  and  $\mathcal C$  into buckets of small expected size using their most-significant bits. An eventual 3XOR must lie in  $A^{[i]} \times B^{[j]} \times C^{[i \oplus j]}$  for some i, j. The idea is to process buckets of C one-by-one, storing them in a small hash table. Then, we consider all the pairs from the corresponding bins of A and B. This yields the following algorithm

**Algorithm Q** (*More practical quadratic algorithm*). Find all triplets (a, b, c) from  $\mathcal{A} \times \mathcal{B} \times \mathcal{C}$  such that  $a \oplus b \oplus c = 0$ . The algorithm takes two input parameters k and  $\ell$ .

- **Q1.** [Dispatch.] Dispatch  $\mathcal{A}$ ,  $\mathcal{B}$  and  $\mathcal{C}$  in buckets according to their first k bits (this essentially requires sorting the lists on their first k bits).
- **Q2.** [Loop on u, v.] Perform steps Q3–Q6 for  $0 \le u, v < 2^{\ell}$ , then terminate the algorithm.
- **Q3.** [Loop on i.] Perform steps Q4–Q6 for  $u2^{k-\ell} \le i < (u+1)2^{k-\ell}$ .
- **Q4.** [Hash  $C^{[i]}$ .] Initialize a hash table with the entries of  $C^{[i]}$ .
- **Q5.** [Loop on j.] Perform step Q6 for  $v2^{k-\ell} \le j < (v+1)2^{k-\ell}$ .
- **Q6.** [Check  $\mathcal{A}^{[j]} \times \mathcal{B}^{[i \oplus j]}$ .] For all  $x \in \mathcal{A}^{[j]}$  and all  $y \in \mathcal{B}^{[i \oplus j]}$ , check if  $x \oplus y \in \mathcal{C}^{[i]}$ ; if so, report (x, y) as a solution.

k must be chosen such that the hash table created in step Q4 fits inside L1 cache. Parallelization is easy: all iterations of the loops on u,v can be done concurrently. The two-level loop structure guarantees that one iteration of steps Q3–Q6 only need to read  $2^{n-\ell}$  entries of each list.  $\ell$  is ideally chosen so that the corresponding portions of both  $\mathcal A$  and  $\mathcal B$  fit in L3 cache.  $2^{18}$  entries of  $\mathcal A$  and  $\mathcal B$  fit in 2 megabyte of L3 cache, and a few minutes will be necessary to process the corresponding  $2^{36}$  pairs. Memory bandwidth is not a problem, and the quadratic algorithm scales well on multi-core machines. The algorithm can also be run on machines with limited memory. Algorithm Q runs at 10–11 cycles per pair processed on a "Haswell" Core i5 CPU.

**Algorithm G.** The longest operation of each iteration is the sorting step of Algorithm M, which accounts for 50% of the total running-time. We use three passes of radix-256 sort for the case of n=96 (where the value of the k parameter is 24). The out-of-place version is  $2-3\times$  faster than the in-place version (but requires twice more memory); the technique of "multi-histogramming" helped a little. We use the M4RI library [AB12] to compute the PLUQ factorization. To solve our n=96 problem, a single iteration of algorithm G runs at 75 CPU cycles per list item processed on the same "Haswell" Core i5 CPU.

It is easy to parallelize the loop on i (each iteration takes slightly less than 1s for n=96). The problem is that both the full  $\mathcal{A}$  and  $\mathcal{B}$  must fit in RAM, as they are entirely read in each iteration. When the lists only have  $2^{24}$  entries (as it is the case for n=96), they only require 256 Megabyte. On a multi-core machine, one iterations can be run concurrently per core. One potential problem is that this may saturate the memory bandwidth: each iteration reads 1.25 Gigabytes from memory in about 1s, so on a large chip with 18 cores/36 threads, up to 22.5 Gigabyte/s of memory bandwidth would be required.

A further problem is that, for larger values of n, it becomes impossible to holds many independant copies of the lists in memory. In that case, the sorting and merging operations themselves have to be parallelized, and this is a bit less obvious. If a single copy of the lists does not fit in memory, a distributed sort/merge will be required, and it is likely that the communication overhead will make this less efficient than the quadratic algorithm.

**Distributed Computation.** It is easy to run these algorithms on clusters of lossely-connected machines. We used a simple master-slave approach, in which the iterations to be parallelized are numbered. When a slave node becomes ready, it request an iteration number to the master node. When it has finished processing it, it sends the partial solutions found in the iteration back to the master. The master stores a log of all accomplished iterations. Communications were handled by the  $\varnothing MQ$  library [Hin13]. We actually ran the algorithms on several hundred cores concurrently.

## 5.3 Results and Further Work

Consider the three ASCII strings:

```
\begin{split} x &= \texttt{F00-0x0000B70947f064A1} \\ y &= \texttt{BAR-0x000013f9e450df0b} \\ z &= \texttt{F00BAR-0x0000e9b2cf21d70a} \end{split}
```

The reader can readily check that

After completing this 96-bit 3XOR, we embarked on a project to compute a 128-bit 3XOR on SHA256. To this end, we found a way to use off-the-shelf bitcoin miners, which evaluate SHA256 about one million times faster than a CPU core. Bitcoin miners naturally produce bitstrings whose SHA256 is zero on (at least) the 32 most significant bits. We plan to accumulate three lists of  $2^{32}$  entries, and then to use algorithm G to find  $2^{32}$  partial solutions on 64 bits, amongst which one should lead to a 128-bit 3XOR (thanks to the extra 32 bits of clamping). Note that this setting naturally enables the use of the technique described in section 3.2.2.

We presently accumulated 5% of the input lists, and we keep on mining.

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# A Baran Demaine and Pătrașcu Algorithm

Let k be a parameter such that  $\mathcal{A}$ ,  $\mathcal{B}$  and  $\mathcal{C}$  are dispatched according to their k first bit.

For each small set  $\mathcal{A}^{[\ell]}$  (resp.  $\mathcal{B}^{[t]}$ ,  $\mathcal{C}^{[\ell \oplus t]}$ ), we keep a  $P \cdot s$ -bit vector  $v_A[\ell]$  (resp.  $v_B[t]$ ,  $v_C[\ell \oplus t]$ ), in which the elements of  $h(\mathcal{A}^{[\ell]})$ , (resp.  $h(\mathcal{B}^{[t]})$ ,  $h(\mathcal{C}^{[\ell \oplus t]})$ ) are stored. With this construction, assuming that the number of elements of each buckets does not exceed P, a solution to the instance  $(h(\mathcal{A}^{[\ell]}), h(\mathcal{B}^{[t]}), h(\mathcal{C}^{\ell \oplus t}))$  exists if and only if  $T[v_A[\ell]|v_B[t]|v_C[\ell \oplus t]] = 1$ .

The following procedure is used to initialize the tables  $v_A$ ,  $v_B$  and  $v_C$ :

**Algorithm V** (*Initialise vectors for preliminary test*). Given a list  $\mathcal{A}$  dispatched into  $N_{buck}$  buckets, create the table  $v_A$ .

- V1. [Loop on  $\ell$ .] For  $0 \le \ell < N_{buck}$  perform steps V1–V4.
- **V2.** [Set up.]  $k \leftarrow \max(P, |A^{[\ell]}|)$ .  $v_A[\ell] \leftarrow \bot$
- **V3.** [Loop on i.] For  $0 \le i < k$ , perform step **V4**.
- **V4.** [Update.]  $v_A[\ell] \leftarrow v_A[\ell] | h(A^{[\ell]}[i])$ .

All in all, this leads to algorithm 1:

If there are exceeding elements in the buckets they are to be treated independently.

```
Data: Three lists \mathcal{A}, \mathcal{B} and \mathcal{C} and the precomputed table T
Result: All couples (A_i, \mathcal{B}_i) such that A_i \oplus \mathcal{B}_i is an element of \mathcal{C}
Dispatch \mathcal{A}, \mathcal{B} and \mathcal{C}, according to their k first bits.
Use Algorithm V to create the tables v_A, v_B and v_C;
for 0 \le \ell, t < 2^{\gamma \cdot n} do
    v \leftarrow v_A[t]|v_B[\ell \oplus t]|v_C[\ell];
    if T[v] = 1 ;
                                  // There may be a solution in this sub-instance
     then
        for all couples (A_i, B_i) in A^{[t]} \times B^{[\ell \oplus t]} do
            if A_i \oplus B_j is in C[id_c[\ell]..id_C[\ell+1]] then
              report (\mathcal{A}_i, \mathcal{B}_i);
             end
        end
    \mathbf{end}
    else if A^{[t]} > P; // There is more than P elements of \mathcal{A} that starts
     with the prefix t
     then
        for all couples (A_i, B_i) in A^{[t]}[P..] \times B^{[\ell \oplus t]} do
             if A_i \oplus B_i is in C[id_c[\ell]...id_C[\ell+1]] then
              report (A_i, B_i);
             end
        \mathbf{end}
    end
    else if B^{[\ell \oplus t]} > P;
                                // There is more than P elements of B that
     starts with the prefix t \oplus \ell
     then
        for all couples (A_i, B_i) in A^{[t]} \times B^{[\ell \oplus t]}[P...] do
             if A_i \oplus B_j is in C[id_c[\ell]..id_C[\ell+1]] then
              report (A_i, B_i);
             end
        end
    end
    else if C^{[\ell]} > P; // There is more than P elements of C that starts
     with the prefix \ell
     then
        for all couples (A_i, B_j) in A^{[t]} \times B^{[\ell \oplus t]} do
             if A_i \oplus B_i is in C[id_c[\ell] + P..id_C[\ell+1]] then
              report (A_i, B_j);
             end
        end
    end
end
```

Algorithm 1: Adaptation of BDP algorithm to our problem