Cryptanalytic Applications of the Polynomial Method for Solving Multivariate Equation Systems over GF(2)

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Abstract. At SODA 2017 Lokshtanov et al. presented the first worstcase algorithms with exponential speedup over exhaustive search for solving polynomial equation systems of degree d in n variables over finite fields. These algorithms were based on the polynomial method in circuit complexity which is a technique for proving circuit lower bounds that has recently been applied in algorithm design. Subsequent works further improved the asymptotic complexity of polynomial method-based algorithms for solving equations over the field \mathbb{F}_2 . However, the asymptotic complexity formulas of these algorithms hide significant low-order terms, and hence they outperform exhaustive search only for very large values of n.

In this paper, we devise a concretely efficient polynomial method-based algorithm for solving multivariate equation systems over \mathbb{F}_2 . We analyze our algorithm's performance for solving random equation systems, and bound its complexity by about $n^2 \cdot 2^{0.815n}$ bit operations for d = 2 and $n^2 \cdot 2^{(1-1/2.7d)n}$ for any d > 2.

We apply our algorithm in cryptanalysis of recently proposed instances of the Picnic signature scheme (an alternate third-round candidate in NIST's post-quantum standardization project) that are based on the security of the LowMC block cipher. Consequently, we show that 2 out of 3 new instances do not achieve their claimed security level. As a secondary application, we also improve the best-known preimage attacks on several round-reduced variants of the Keccak hash function.

Our algorithm combines various techniques used in previous polynomial method-based algorithms with new optimizations, some of which exploit randomness assumptions about the system of equations. In its cryptanalytic application to Picnic, we demonstrate how to further optimize the algorithm for solving structured equation systems that are constructed from specific cryptosystems.

1 Introduction

The security of many cryptographic schemes is based on the conjectured hardness of solving systems of polynomial equations over a finite field. This problem is known to be NP-hard even for systems of quadratic equations over \mathbb{F}_2 .

The input to the problem consists of m polynomials in n variables over a finite field \mathbb{F} , denoted by $E = \{P_j(x_1, \ldots, x_n)\}_{j=1}^m$, where each polynomial is given

as a sum of monomials. The algebraic degree of each polynomial is bounded by a small constant d. The goal is to find a solution to the system, namely $\hat{x} = (\hat{x}_1, \ldots, \hat{x}_n) \in \mathbb{F}^n$ such that $P_j(\hat{x}) = 0$ for every $j \in \{1, \ldots, m\}$, or to determine that a solution does not exist.¹

In this paper, we will be interested in the concrete (rather than asymptotic) complexity of algorithms for solving polynomial systems over the field \mathbb{F}_2 and in their applications to cryptanalysis.

1.1 Previous Algorithms for Solving Polynomial Equation Systems

The problem of solving polynomial equation systems over finite fields is widely studied. We give a brief overview of the main algorithms that were applied in cryptanalysis.

Classical techniques developed to solve polynomial systems attempt to find a reduced representation of the ideal generated by the polynomials in the form of a Gröbner basis (e.g., the F4 [17] and F5 [18] algorithms). These methods have had success in solving some very structured polynomial systems that arise from certain cryptosystems (e.g., see [19]), but it is difficult to estimate their complexity in solving arbitrary systems. Related methods such as XL [11] and its variants typically work well only for largely over-defined systems in which $m \gg n$.

In [3] Bardet et al. analyzed the problem of solving quadratic equations over \mathbb{F}_2 and devised an algorithm that combines exhaustive search and sparse linear algebra. The authors estimated the asymptotic complexity of their randomized algorithm for m = n by $O(2^{0.792n})$, under some algebraic assumptions that were empirically found to hold for random systems. However, due to a large overhead hidden in the asymptotic formula, the authors expect their algorithm to beat exhaustive search only when the number of variables is at least 200. Another algorithm based on a different hybrid approach was published by Joux and Vitse [23], who gave experimental evidence that it outperforms in practice previous algorithms for a wide range of parameters. Analyzing the complexity of the algorithm is non-trivial, but according to the recent work of Duarte [16], the algorithm for solving quadratic systems over \mathbb{F}_2 with m = n + 1 does not beat existing algorithms (such as the one of [3]) asymptotically.

Finally, we mention the work of Bouillaguet et al. [8], which devised an optimized exhaustive search algorithm for solving polynomial systems of degree d over \mathbb{F}_2 whose complexity is $2d \log n \cdot 2^n$ bit operations.

1.2 The Polynomial Method

In [27] Lokshtanov et al. presented the first worst-case algorithms for solving polynomial equations over finite fields that have exponential speedup over exhaustive search. These algorithms were based on a technique known as the *polynomial method*. It was borrowed from circuit complexity [4] and recently applied

¹ We denote an assignment to the formal variable vector x in the polynomial $P_j(x)$ by \hat{x} and the value of $P_j(x)$ on this assignment by $P_j(\hat{x})$.

in algorithm design (see [36] for a survey). The randomized algorithm of Lokshtanov et al. for solving equations over \mathbb{F}_2 has runtime of $O(2^{0.8765n})$ for quadratic systems and $O(2^{(1-1/(5d))n})$ in general. Following [27], Björklund, Kaski and Williams [7] reduced the complexity of this algorithm to $O(2^{0.804n})$ for d = 2and $O(2^{(1-1/2.7d)n})$ in general. More recently, these complexities were further improved in [12] by the author to $O(2^{0.6943n})$ for d = 2 and $O(2^{(1-1/2d)n})$ for d > 2.

Although these recent algorithms have better asymptotic complexity than exhaustive search, a close examination reveals that their concrete (non-asymptotic) complexity is above 2^n for parameter ranges that are relevant to cryptography.

1.3 Our Results

We introduce the polynomial method for solving multivariate equation systems over \mathbb{F}_2 as a tool in cryptanalysis. For this purpose, we devise a concretely efficient algorithm for solving such systems.

Our algorithm is relatively simple and its analysis assumes the degree d polynomials are selected uniformly at random. Up to small constants, we bound the complexity of our algorithm by $n^2 \cdot 2^{0.815n}$ bit operations for d = 2, and $n^2 \cdot 2^{(1-1/2.7d)n}$ for $d \ge 2$.

In a straightforward implementation of our algorithm, its memory complexity is significant and only about n times lower than its time complexity. In fact, this is the case for all previous polynomial method-based algorithms. We address this issue by presenting a memory-optimized variant of the algorithm which maintains roughly the same time complexity, but whose memory complexity is reduced to about $n^2 \cdot 2^{0.63n}$ bits for d = 2 and $n^2 \cdot 2^{(1-1/1.35d)n}$ in general.²

Potential fast implementation. Even after the reduction in memory complexity, it remains high and would present a major challenge for obtaining a fast practical implementation of the algorithm. On the other hand, the memory access patterns of the algorithm are fixed and independent of the data, which is an advantage. Moreover, future works may be able to further reduce the memory complexity or utilize time-memory tradeoffs. Taking this optimistic viewpoint, our work may be viewed as a step towards a practically efficient implementation of a polynomial method-based algorithm for solving multivariate equation systems over \mathbb{F}_2 .

We stress, however, that the main goal of the paper is to give a good *analytical* concrete estimate of the complexity of polynomial method algorithms for problem sizes that are too large to be solved in practice. Consequently, they can be used in the security analysis of cryptosystems and serve as a starting point for additional optimizations.

² Asymptotically, the polynomial factor in the memory complexity formula is between n^2 and n^3 , but it is close to n^2 for relevant parameters.

Asymptotic and concrete complexity. Our algorithm can actually be viewed as a concretely efficient variant of the algorithm of [12]. The only reason that [12] seems to have better asymptotic complexity is that it uses a self-reduction to a smaller multivariate system. Essentially, each recursive call reduces the exponent in the complexity formula, where the gain diminishes with the number of recursive calls. On the other hand, each such call increases the lower order terms.

An estimated calculation suggests that a self-reduction is profitable for d = 2starting from about n = 100. Beyond n = 200 for d = 2 the advantage in concrete complexity is made substantial using several recursive calls. On the other hand, for d = 4 the reduction seems profitable only beyond n = 200. We chose not to augment our algorithm with any self-reduction and simply replace it with exhaustive search for two reasons. First, as described below, the applications we present in this paper require solving multivariate systems with d > 2, for which the benefit of the self-reduction seems marginal for relevant parameters. Second, this self-reduction significantly complicates the concrete analysis, whereas we aim for simplicity. Yet, this estimation suggests that the full potential of the algorithm is still to be discovered.

Cryptanalytic applications. We estimate the concrete complexities of our algorithm for solving quadratic systems in 80, 128 and 256 variables by 2^{77} , 2^{117} and 2^{223} bit operations, respectively. In terms of cryptanalysis, the main targets of our algorithm for d = 2 are multivariate public-key cryptosystems (e.g., HFE by Patarin [30] and UOV by Kipnis, Patarin and Goubin [25]), whose security is directly based on the hardness of solving quadratic systems. However, recent multivariate cryptosystems such as GeMSS [9] (an alternate third-round candidate signature scheme in NIST's post-quantum standardization project [29]) were designed with a sufficiently large security margin and resist our attack. Nevertheless, the security margin for some of these cryptosystems seems to be reduced by our algorithm.

Interestingly, the main application of our algorithm is for solving multivariate systems of degree d > 2 which have generally received less attention in the literature compared to quadratic systems. In particular, we apply it to cryptanalyze recently proposed instances [24] of the Picnic signature scheme [10] (an alternate third-round candidate in NIST's post-quantum standardization project) that is based on the security of the LowMC block cipher [1].

We focus on the three Picnic instances where the LowMC block cipher has a full Sbox layer and 4 internal rounds. These instances have claimed security levels of $S \in \{128, 192, 255\}$ bits. The best-known attacks on these instances were recently published by Banik et al. [2], but they are only applicable to weakened variants where the number of LowMC rounds is reduced from 4 to 2. On the other hand, our attacks on the full 4-round instances with $S \in \{128, 192, 255\}$ have complexities of $2^{130}, 2^{188}$ and 2^{245} bit operations, respectively. These attacks are far from being practical, but show that 2 out of the 3 instances do not achieve their claimed security level, while the security of the instance with S = 128 is somewhat marginal. When optimized for time complexity, the attacks for $S \in \{128, 192, 255\}$ require about 2^{112} , 2^{164} and 2^{219} bits of memory, respectively. However, there is no consensus among researchers on a model that takes memory complexity into account and the formal security claims of the Picnic (and LowMC) designers only involve time complexity.³

The authors of [24] also proposed conservative instantiations where the number of LowMC rounds is increased by 1. The attack complexities for these instances with $S \in \{128, 192, 255\}$ bits become $2^{133}, 2^{192}$ and 2^{251} , respectively. Hence the instance with S = 255 still does not achieve the desired security level, while security remains very marginal for the strengthened instance with S = 192.

We also analyze round-reduced variants of the Keccak hash function [5], which was selected by NIST in 2015 as the SHA-3 standard. In particular, we describe the best-known preimage attacks in terms of time complexity on 4 rounds of Keccak-384 and Keccack-512. These have complexities of 2^{374} and 2^{502} bit operations, respectively. We further describe the first collision attack on 4-round Keccak-512 that is (slightly) faster than the birthday bound. We consider the cryptanalysis of round-reduced Keccak as a secondary application since our attacks do not substantially improve upon previous attacks and they are very far from threatening Keccak's security. On the other hand, the attacks only exploit basic properties of Keccak's internal structure and it is likely that they can be improved by using more dedicated techniques.

Complexity evaluation. It is important to emphasize that the complexities of our attacks are measured in *bit operations*. On the other hand, the complexity of exhaustive search for the cryptanalytic problems we consider on a space of size 2^n is larger than 2^n bit operations. Hence the improvement we obtain over exhaustive search is more significant than it may first appear.

In particular, the encryption algorithms of the LowMC instances we cryptanalyze (that are used in Picnic) have complexities of at least 2^{17} bit operations. However, evaluating an attack in terms of the complexity of the LowMC encryption algorithm is misleading, as naive exhaustive search is not the most efficient generic attack on the 4-round LowMC instances. Indeed, breaking these instances is easily reduced to solving a multivariate system in (about) n variables with d = 4, for which the best-known generic attack is the optimized exhaustive search algorithm of Bouillaguet et al. [8], whose complexity is about $2d \log n \cdot 2^n = 8 \log n \cdot 2^n$ bit operations (also see [15] for an alternative algorithm). Overall, in terms of bit operations, our algorithm is more efficient than the one of [8] by a factor which is between 32 and $2^{16} = 65536$ (depending on the LowMC instance considered).

1.4 Comparison to Previous Works

The analysis of our algorithm for random equation systems over \mathbb{F}_2 is simple. In contrast, the analysis of previous cryptanalytic algorithms that beat brute force

³ The Picnic designers have confirmed our findings and plan to update the parameter sets accordingly.

for random equation systems over \mathbb{F}_2 (such as the one by Bardet et al. [3]) is based on heuristic algebraic assumptions that are difficult to analyze.

The algorithm of [3] (applied to systems with m = n) may seem to have a slightly better asymptotic complexity than ours, but this is a misleading comparison, since (as noted above) our algorithm can be extended to have better asymptotic complexity. In terms of concrete complexity for relevant parameters, [3] only beats exhaustive search for d = 2 beyond n = 200, whereas the complexity of our algorithm for d = 2, n = 200 is about 2^{177} .

Unlike our algorithm, the algorithm of Joux and Vitse [23] was shown to perform well in practice. However, its concrete complexity has not been established analytically and it is not clear how to use it in the security analysis of cryptosystems. On the other hand, the complexity of hybrid algorithms (such as [3,23]) for over-defined systems (in which $m \gg n$) is reduced as a function of m, whereas achieving such optimizations for our algorithm is left as an open problem.

Finally, previous polynomial method-based algorithms were only analyzed asymptotically. While it is difficult to calculate their exact concrete complexity, our optimizations reduce complexity by many orders of magnitude for relevant parameters.

1.5 Technical Contribution

Let $E = \{P_j(x_1, \ldots, x_n)\}_{j=1}^m$ be a polynomial system of degree d over \mathbb{F}_2 . Algorithms based on the polynomial method consider the polynomial

$$F(x) = (1 + P_1(x))(1 + P_2(x))\dots(1 + P_m(x))$$

(operations are over \mathbb{F}_2). Note that $F(\hat{x}) = 1$ if and only if \hat{x} is a solution to E. However, the degree of F(x) can be as high as $d \cdot m$ and it generally contains too many monomials to manipulate efficiently. It is thus replaced by a *probabilistic polynomial* $\tilde{F}(x)$ with a lower degree that agrees with F(x) on most assignments. Taking advantage of the low degree of $\tilde{F}(x)$ by using fast polynomial interpolation and evaluation algorithms allows to solve E faster than brute force.

Our main algorithm includes various concrete optimizations and simplifications to previous polynomial method-based algorithms. These are described in detail in Section 3. For example, we reduce the number of polynomials which need to be interpolated and evaluated and show how to jointly interpolate several polynomials with improved amortized complexity.

Then, we show how to reduce the memory complexity of the algorithm by an exponential factor with essentially no penalty in time complexity. The optimization is based on a memory-reduced variant of the Möbius transform over \mathbb{F}_2 which is a fast polynomial interpolation and evaluation algorithm. This variant allows to evaluate a low-degree polynomial on its entire domain with memory complexity proportional to the memory required to store the input polynomial itself (and time complexity proportional to the domain size). Although the Möbius transform is widely used and the variant we describe is simple, it seems not to be well-known. The way this variant is used in our algorithm is, however, slightly more involved.

As an additional technical contribution, we show how to optimize our algorithm for solving *structured* equation systems that are constructed from specific cryptosystems. In particular, we observe that in cryptographic settings, a probabilistic polynomial can be replaced with a deterministic construction of a polynomial that preserves the structure of the polynomials of E. We show that in some cases (e.g., in the analysis of Picnic in Section 5.2) this alternative polynomial has reduced degree, optimizing the attack. We view this optimization as one of the main contributions of this paper, as it may serve as a starting point for future works in cryptanalysis.

Paper structure. Next, we describe some preliminaries. We overview our algorithm in Section 3 and give its details in Section 4. Applications are described in Section 5.

2 Preliminaries

2.1 Boolean Algebra

Given a finite ordered set S, denote by |S| its size and by S[i] its *i*'th element. For a positive integer n, let $[n] = \{1, 2, ..., n\}$.

It is well-known that any Boolean function $F : \mathbb{F}_2^n \to \mathbb{F}_2$ can be uniquely described as a multilinear polynomial, whose algebraic normal form (ANF) is given by $F(x_1, \ldots, x_n) = \sum_{u \in \{0,1\}^n} \alpha_u(F) M_u(x)$, where $\alpha_u(F) \in \{0,1\}$ is the coefficient of the monomial $M_u(x) = \prod_{i=1}^n x_i^{u_i}$ (operations are over \mathbb{F}_2).

We denote by HW(x) the Hamming weight of a vector $x \in \{0,1\}^n$. The algebraic degree of a function F is defined as $\max\{\text{HW}(u) \mid \alpha_u(F) \neq 0\}$. Let W_w^n be the set $\{x \in \{0,1\}^n \mid \text{HW}(x) \leq w\}$. Thus, a function F of degree $d \leq n$ can be described using $|W_d^n| = \sum_{i=0}^d {n \choose i}$ coefficients. We simplify notation by denoting ${n \choose \downarrow w} = \sum_{i=0}^w {n \choose i}$.

For $i \in \{1, \ldots, n\}$ and $b \in \{0, 1\}$, define the function $F_{x_i \leftarrow b} : \mathbb{F}_2^{n-1} \to \mathbb{F}_2$ by

$$F_{x_i \leftarrow b}(x_1, \dots, x_{i-1}, x_{i+1}, \dots, x_n) = F(x_1, \dots, x_{i-1}, b, x_{i+1}, \dots, x_n).$$

Interpolation. Any ANF coefficient $\alpha_u(F)$ can be interpolated by summing (over \mathbb{F}_2) over $2^{\mathrm{HW}(u)}$ evaluations of F: for $u \in \{0,1\}^n$, define $I_u = \{i \in \{1,\ldots,n\} \mid u_i = 1\}$ and let $S_u = \{x \in \{0,1\}^n \mid I_x \subseteq I_u\}$. Then,

$$\alpha_u(F) = \sum_{\hat{x} \in S_u} F(\hat{x}). \tag{1}$$

Indeed, among all monomials only $M_u(\hat{x})$ attains a value of 1 an odd number of times in the expression $\sum_{\hat{x}\in S_u} F(\hat{x}) = \sum_{\hat{x}\in S_u} \sum_{v\in\{0,1\}^n} \alpha_v(F) M_v(\hat{x}).$

Proposition 2.1. Let $F : \mathbb{F}_2^n \to \mathbb{F}_2$ be a Boolean function. For some $1 \leq n_1 \leq n$, partition its n variables into two sets $y_1, \ldots, y_{n-n_1}, z_1, \ldots, z_{n_1}$. Given the ANF of F, write it as $F(y, z) = (z_1 \ldots z_{n_1})F_1(y) + F_2(y, z)$ by factoring out all the monomials that are multiplied with $z_1 \ldots z_{n_1}$. Then, $F_1(y) = \sum_{z \in \{0,1\}^{n_1}} F(y, z)$.

The proposition follows from (1) by considering the polynomial $F_1(y)$ as the symbolic coefficient of the monomial $z_1 \ldots z_{n_1}$. Observe that if F(y, z) is of degree d then $F_1(y)$ is of degree at most $\max(d - n_1, 0)$.

Remark 2.1. Proposition 2.1 is also at the basis of cube attacks [14]. However, in cube attacks, the z variables are public bits controlled by the attacker (e.g., plaintext bits) while the y variables are secret key bits. In our case, we will apply Proposition 2.1 in a setting where all variables are secret key bits.

2.2 Model of Computation

We estimate the complexity of a straight-line implementation of our algorithm by counting the number of bit operations (e.g., AND, OR, XOR) on pairs of bits. This ignores bookkeeping operations such as moving a bit from one position to another (which merely requires renaming of variables in straight-line programs).

2.3 Basic Algorithms

We describe the basic algorithms that our main algorithm uses as sub-procedures.

Möbius transform. Given the truth table of an arbitrary function F (as a bit vector of 2^n entries), the ANF of F can be represented as a bit vector of 2^n entries, corresponding to its 2^n coefficients. This ANF representation can be computed from the truth table of F via the *Möbius transform* over \mathbb{F}_2^n .

A fast algorithm for computing this transform is based on the decomposition

$$F(x_1, \dots, x_n) = x_1 \cdot F_1(x_2, \dots, x_n) + F_2(x_2, \dots, x_n).$$
(2)

Thus, one recursively computes the ANF of $F_1(x_2, \ldots, x_n)$ and $F_2(x_2, \ldots, x_n)$. Given the evaluations of F, for every $(\hat{x}_2, \ldots, \hat{x}_n) \in \{0, 1\}^{n-1}$,

$$F_2(\hat{x}_2,\ldots,\hat{x}_n)=F(0,\hat{x}_2,\ldots,\hat{x}_n)$$

and

$$F_1(\hat{x}_2,\ldots,\hat{x}_n) = F(0,\hat{x}_2,\ldots,\hat{x}_n) + F(1,\hat{x}_2,\ldots,\hat{x}_n).$$

Therefore, computing the evaluations of F_1 requires 2^{n-1} bit operations. Denoting the time complexity by T(n), we have $T(n) = 2T(n-1) + 2^{n-1}$, and hence $T(n) \leq n \cdot 2^{n-1} < n \cdot 2^n$.

By (1), a function F of degree bounded by $d \leq n$ can be interpolated from its evaluations on the set W_d^n . Adapting the Möbius transform for such a function F using the decomposition above gives an algorithm with complexity $T(n, d) \leq$

 $T(n-1,d) + T(n-1,d-1) + \binom{n-1}{\downarrow d}$ and $T(n,n) \leq n \cdot 2^n$. It can be shown by induction that $T(n,d) \leq n \cdot \binom{n}{\downarrow d}$ bit operations.

The Möbius transform over \mathbb{F}_2^n coincides with its inverse which corresponds to evaluating the ANF representation of F (i.e., computing its truth table). More details about the Möbius transform over \mathbb{F}_2^n and its applications in crypography can be found in [22, p.285].

Memory complexity. A standard in-place implementation of the Möbius transform performs n iterations on its input vector, where in each iteration, half of the array entries are XORed to the other half. This requires 2^n bits of memory.

Fast exhaustive search for polynomial systems over \mathbb{F}_2 [8]. At CHES 2010 Bouillaguet et al. presented an optimized exhaustive search algorithm for enumerating over all solutions to a polynomial system over \mathbb{F}_2 . For a polynomial system of degree d with n variables, the complexity of their algorithm is $2d \cdot \log n \cdot 2^n$. The algorithm also requires a preprocessing phase that has complexity of n^{2d} , which is negligible when d is much smaller than n. We note that the analysis of the algorithm makes some randomness assumptions about the polynomial system, and requires that the expected number of solutions to a system with m equations over n variables is about 2^{n-m} .

In this paper we will use this algorithm to find solutions inside sets of the special form $W_w^{n-n_1} \times \{0,1\}^{n_1}$ for some values of w and n_1 in time

$$2d \cdot \log n \cdot |W_w^{n-n_1} \times \{0,1\}^{n_1}| = 2d \cdot \log n \cdot 2^{n_1} \cdot \binom{n-n_1}{\downarrow w}.$$

For an arbitrary value of n_1 , obtaining this complexity may not be trivial because the algorithm of [8] iterates over the search space using a Gray code, and hence it cannot be implemented on the set of low Hamming weight vectors $W_w^{n-n_1}$ (for $w < n - n_1$) in a straightforward manner. We note that there are known monotone Gray codes that traverse the binary vectors (almost) in increasing Hamming weight order [32]. It would be interesting to adapt the algorithm of [8] to use them.

On the other hand, the least significant bits of the vectors in the set $W_w^{n-n_1} \times \{0,1\}^{n_1}$ can be traversed using a standard Gray code, paying a penalty only every 2^{n_1} iterations. Since we can naively enumerate any set of *n*-bit vectors by flipping at most *n* bits at a time, we can conservatively estimate the multiplicative penalty by about *n*. Thus, the amortized penalty over [8] is about $2^{-n_1} \cdot n$. In our setting, $2^{n_1} \gg n$, and the overhead is negligible.

Remark 2.2. Asymptotically, we will set $n_1 = \Theta(n)$, hence $2^{n_1} = \omega(n)$. Concretely, the advantage of the algorithm over exhaustive search will be roughly $\frac{2^{n_1}}{n^2}$. Thus, we may assume that $2^{n_1} \gg n$ holds without loss of generality, as otherwise, the algorithm would not obtain any advantage over exhaustive search.

2.4 Probabilistic Polynomials

Previous algorithms for solving multivariate equation systems based on the polynomial method (starting with [27]) make use of probabilistic polynomials. In particular, these works use the following construction (credited to Razborov [31] and Smolensky [33]). Given m polynomial equations of degree d in the n Boolean variables $x_1, \ldots, x_n, E = \{P_j(x)\}_{j=1}^m$, consider the polynomial

$$F(x) = (1 + P_1(x))(1 + P_2(x))\dots(1 + P_m(x)).$$

Note that \hat{x} is a solution to E if and only if $F(\hat{x}) = 1$. Thus, we call the polynomial F the *identifying polynomial* of E.

The degree of F(x) is generally too high and we work with a probabilistic polynomial with a lower degree, defined as follows. Let $\ell < m$ be a parameter. Pick a uniformly random matrix of full rank ℓ , $A \in \mathbb{F}_2^{\ell \times m}$ and define ℓ degree dpolynomials as

$$R_{i}(x) = \sum_{j=1}^{m} A_{i,j} \cdot P_{j}(x).$$
(3)

We note that previous works [7,12,27] did not restrict the rank of A. In our case, this restriction will slightly simplify the analysis of the algorithm. Let

$$\tilde{F}(x) = (1 + R_1(x))(1 + R_2(x))\dots(1 + R_\ell(x))$$
(4)

be the identifying polynomial of the system $\tilde{E} = \{R_i(x)\}_{i=1}^{\ell}$. Note that the degree of $\tilde{F}(x)$ (denoted by $d_{\tilde{F}}$) is at most $d \cdot \ell$.

Proposition 2.2. For any $\hat{x} \in \{0,1\}^n$, if $F(\hat{x}) = 1$, then $\tilde{F}(\hat{x}) = 1$. Otherwise, $F(\hat{x}) = 0$ and then $\Pr[\tilde{F}(\hat{x}) = 0] \ge 1 - 2^{-\ell}$.

Proof. If $F(\hat{x}) = 1$, then $P_j(\hat{x}) = 0$ for all $j \in [m]$ and therefore $R_i(\hat{x}) = 0$ for all $i \in [\ell]$. Hence, $\tilde{F}(\hat{x}) = 1$.

Otherwise, $F(\hat{x}) = 0$. Let $v \in \mathbb{F}_2^m$ be a vector such that $v_j = P_j(\hat{x})$ and $u \in \mathbb{F}_2^{\ell}$, a vector such that $u_i = R_i(\hat{x})$. Note that $u = A \cdot v$. Since $F(\hat{x}) = 0$, there exists $j \in [m]$ such that $P_j(\hat{x}) = 1$ and thus $v \neq 0$. On the other hand, if $\tilde{F}(\hat{x}) = 1$, then $R_i(\hat{x}) = 0$ for all $i \in [\ell]$, implying that u = 0. Therefore, v is a non-zero vector in the kernel of A. Since A is a uniform matrix of full rank ℓ , any fixed non-zero vector (including v) belongs to its kernel with probability $2^{-\ell} - 2^{-m} < 2^{-\ell}$.

2.5 Previous Polynomial Method Algorithms for Solving Equation Systems over \mathbb{F}_2

In this section we give a short description of the previous polynomial methodbased algorithms of [7,12]. We focus on the parts which are most relevant to this work. The Björklund et al. algorithm [7]. In the algorithm of [7], the search problem of finding a solution to the system E is reduced to the parity-counting problem, where the goal is to compute the parity of the number of solutions.

The first step reduces the search problem to the problem of deciding whether a solution exists. The reduction iteratively fixes one variable of the solution at a time using $\Theta(n)$ calls to the decision algorithm. Then, the decision problem is reduced to the parity-counting problem. This reduction uses the Valiant-Vazirani affine hashing [35], adding random linear equations to the system with the goal of *isolating* some solution to E (assuming a solution exists), such that it is the only solution to the new system. In this case, the output of the parity-counting algorithm is 1. The number of linear equations to add that ensures isolation with high probability depends on the logarithm of the number of solutions to E, which is unknown. Hence, the algorithm exhausts all its possible n values.

We now overview the Björklund et al. parity-counting algorithm. Algebraically, the parity of solutions to $E = \{P_j(x)\}_{j=1}^m$ is computed by $\sum_{\hat{x} \in \{0,1\}^n} F(\hat{x})$, where $F(x) = (1 + P_1(x)) \dots (1 + P_m(x))$. This sum (parity) is computed in parts by partitioning the *n* variables into 2 sets $y = y_1, \dots, y_{n-n_1}$ and $z = z_1, \dots, z_{n_1}$, where $n_1 < n$ is a parameter. Thus,

$$\sum_{\hat{x} \in \{0,1\}^n} F(\hat{x}) = \sum_{\hat{y} \in \{0,1\}^{n-n_1}} \sum_{\hat{z} \in \{0,1\}^{n_1}} F(\hat{y}, \hat{z}).$$

Replace the polynomial F(y, z) with the probabilistic polynomial $\tilde{F}(y, z) = (1+R_1(y,z))\dots(1+R_\ell(y,z))$ for $\ell = n_1+2$, similarly to (4). By Proposition 2.2 and a union bound over all $\hat{z} \in \{0,1\}^{n_1}$, for each $\hat{y} \in \{0,1\}^{n-n_1}$,

$$\Pr\left[\sum_{\hat{z}\in\{0,1\}^{n_1}}\tilde{F}(\hat{y},\hat{z}) = \sum_{\hat{z}\in\{0,1\}^{n_1}}F(\hat{y},\hat{z})\right] \ge 1 - 2^{n_1 - \ell} = \frac{3}{4}$$

In order to compute the sum for each $\hat{y} \in \{0,1\}^{n-n_1}$ efficiently, let

$$G(y) = \sum_{\hat{z} \in \{0,1\}^{n_1}} \tilde{F}(y, \hat{z}).$$

It follows from Proposition 2.1 that the degree of G(y) is at most $d_G = d \cdot \ell - n_1$. Proceed by interpolating G(y) by first computing its values in the set $W_{d_G}^{n-n_1}$. For this purpose, find all solutions to the equation system $\{R_i(y,z)\}_{i=1}^{\ell}$ in $W_{d_G}^{n-n_1} \times \{0,1\}^{n_1}$ via brute force, giving $\tilde{F}(\hat{y}, \hat{z})$ for $(\hat{y}, \hat{z}) \in W_{d_G}^{n-n_1} \times \{0,1\}^{n_1}$. Then, compute $G(\hat{y}) = \sum_{\hat{z} \in \{0,1\}^{n_1}} \tilde{F}(\hat{y}, \hat{z})$ for all $\hat{y} \in W_{d_G}^{n-n_1}$. Given these values of G(y), interpolate it using the Möbius transform.

Next, evaluate G(y) on all $\hat{y} \in \{0,1\}^{n-n_1}$ (using the Möbius transform) to obtain the partial parity of each part $\sum_{\hat{z} \in \{0,1\}^{n_1}} \tilde{F}(\hat{y}, \hat{z})$. However, each partial parity is correct only with probability $\frac{3}{4}$. Thus, repeat the above steps with t = 48n + 1 independent probabilistic polynomials. Obtain t suggestions for each part and take their majority to obtain the true partial parity, except with exponentially small probability. Assuming that all partial parities are computed correctly, return $\sum_{\hat{x} \in \{0,1\}^n} F(\hat{x}) = \sum_{\hat{y} \in \{0,1\}^{n-n_1}} \sum_{\hat{z} \in \{0,1\}^{n_1}} F(\hat{y}, \hat{z})$. Complexity evaluation and optimization by self-reduction. Ignoring low-order (but concretely substantial) terms, the complexity of the above algorithm is dominated by brute force searches on $W_{d_G}^{n-n_1} \times \{0,1\}^{n_1}$ and polynomial evaluations on $\{0,1\}^{n-n_1}$. The complexity is thus

$$O^*(\binom{n-n_1}{\downarrow d_G} \cdot 2^{n_1} + 2^{n-n_1}) = O^*(\binom{n-n_1}{\downarrow d \cdot (n_1+2)-n_1} \cdot 2^{n_1} + 2^{n-n_1}).$$
(5)

 $(O^* \text{ hides polynomial factors in } n)$. The parameter n_1 can be set to balance these terms and optimize complexity.

In [7], the asymptotic complexity is optimized. Instead of brute force, for each $\hat{y} \in W^{n-n_1}_{d_G}$, compute $G(\hat{y})$ by a self-reduction to a parity-counting problem with input $\{R_i(\hat{y}, z)\}_{i=1}^{\ell}$, which is a polynomial system of degree d with n_1 variables z_1, \ldots, z_{n_1} .

The author's algorithm for enumerating solutions [12]. The main result of the followup paper [12] is an asymptotically more efficient algorithm. Essentially, it defines a multiple parity-counting problem (which computes many parities at once), and shows that all the $\binom{n-n_1}{\downarrow d_G}$ parities returned by recursive calls in [7] can be computed more efficiently by a single recursive call to a multiple parity-counting problem. As noted in Section 1.3, we do not use this reduction.

Instead, we focus on the secondary result of [12], which is an algorithm for enumerating *all* solutions to a polynomial system. In our context, we will use a related technique to eliminate the initial reduction of [7] from solving E to parity-counting (which has a high concrete overhead) and replace it with a more direct way of recovering solutions from parity computations.

Isolating solutions. The first observation is that we can isolate many solutions to E at once using a variable partition $x = (y, z) = (y_1, \ldots, y_{n-n_1}, z_1, \ldots, z_{n_1})$.

Definition 2.1 (Isolated solutions). A solution $\hat{x} = (\hat{y}, \hat{z})$ to $E = \{P_j(y, z)\}_{j=1}^m$ is called isolated (with respect to the variable partition (y, z)), if for all $\hat{z}' \neq \hat{z}$, (\hat{y}, \hat{z}') is not a solution to E.

The goal is to "evenly spread" solutions across the different \hat{y} values. Thus, for many solutions (\hat{y}, \hat{z}) , there is no additional solution that shares the same \hat{y} value (namely, (\hat{y}, \hat{z}) is isolated). This is made possible by a random linear change of variables applied to the polynomials of E and a careful choice of n_1 .

Enumerating isolated solutions. The second observation is that all solutions isolated by the variable partition (y, z) can be recovered bit-by-bit by computing $n_1 + 1$ sums (parities) for each $\hat{y} \in \{0, 1\}^{n-n_1}$. Let

$$V_0(y) = \sum_{\hat{z} \in \{0,1\}^{n_1}} F(y, \hat{z}), \text{and let } V_i(y) = \sum_{\hat{z} \in \{0,1\}^{n_1-1}} F_{z_i \leftarrow 0}(y, \hat{z})$$

for $i \in \{1, \ldots, n_1\}$, where $F(y, z) = (1 + P_1(y, z)) \dots (1 + P_m(y, z))$.

Proposition 2.3. Assume that (\hat{y}, \hat{z}) is an isolated solution to E with respect to (y, z). Then, $V_0(\hat{y}) = 1$ and $V_i(\hat{y}) = \hat{z}_i + 1$ for all $i \in \{1, \ldots, n_1\}$.

As a result, given \hat{y} , in order to recover the n_1 least significant bits \hat{z} of the isolated solution (\hat{y}, \hat{z}) (for which $V_0(\hat{y}) = 1$), it is sufficient to compute $V_i(\hat{y})$ for each $i \in \{1, \ldots, n_1\}$. The formal polynomials $V_i(y)$ cannot be directly interpolated since they are derived from F and hence of high-degree. However, for each $i \in \{0, \ldots, n_1\}$, the sums $V_i(\hat{y})$ for $\hat{y} \in \{0, 1\}^{n-n_1}$ can be computed (with negligible error probability) using probabilistic polynomials, similarly to the way that the partial sums $\sum_{\hat{z} \in \{0,1\}^{n_1}} F(\hat{y}, \hat{z})$ are computed in [7] (yet with a more efficient recursive multiple parity-counting algorithm). These sums allow to recover all isolated solutions.

Proof. First, since F is the identifying polynomial of E, then $V_0(\hat{y})$ counts the parity of solutions to the system $\{P_j(\hat{y}, z)\}_{i=1}^m$ (in which y is fixed to \hat{y}). Therefore, if (\hat{y}, \hat{z}) is an isolated solution to E with respect to (y, z), then $V_0(\hat{y}) = 1$.

Next, if $\hat{z}_i = 0$, then the assignment (\hat{y}, \hat{z}) continues to be an isolated solution to E with respect to (y, z) after setting $z_i = 0$ and hence $V_i(\hat{y}) = 1$. Otherwise, $\hat{z}_i = 1$, and E has no solutions after setting $y = \hat{y}$ and $z_i = 0$, implying that $V_i(\hat{y}) = 0$. In both cases, $V_i(\hat{y}) = \hat{z}_i + 1$ for all $i \in \{1, \ldots, n_1\}$ as required.

Testing solutions. For each $\hat{y} \in \{0,1\}^{n-n_1}$, the algorithm computes $n_1 + 1$ sums. Those with $V_0(\hat{y}) = 1$ suggest some solution (\hat{y}, \hat{z}) . The suggestion is correct if (\hat{y}, \hat{z}) is an isolated solution. Otherwise, the suggestion may be a "false alarm". Consequently, all suggested solutions are tested on E.

3 Overview of the New Algorithm

The starting point of our algorithm is the solution enumeration algorithm of [12]. We first notice that in order to find a solution to E, it is, in fact, sufficient to enumerate isolated solutions to \tilde{E} (as they form a superset of the solution set of E), and test each one on E. This has a significant advantage in terms of concrete complexity, as detailed below.

We now describe how we isolate solutions to \tilde{E} with respect to a variable partition x = (y, z) according to a parameter n_1 and then overview our algorithm that outputs all isolated solutions to \tilde{E} and tests them.

Isolating solutions. We use the following proposition.

Proposition 3.1. Let E and \tilde{E} be polynomials systems with identifying polynomials F(x) and $\tilde{F}(x)$, respectively. For $n_1 = \ell - 1$, define a variable partition $x = (y, z) = (y_1, \ldots, y_{n-n_1}, z_1, \ldots, z_{n_1})$. Assume that (\hat{y}, \hat{z}) is an isolated solution to E. Then, $\Pr[(\hat{y}, \hat{z})$ is an isolated solution to $\tilde{E}] \geq 1 - 2^{n_1 - \ell} = \frac{1}{2}$.

Proof. The proposition follows from Proposition 2.2 by a union bound over the set $\{(\hat{y}, \hat{z}') \mid \hat{z}' \neq \hat{z}\}$, whose size is $2^{n_1} - 1$.

Hence, assuming that E has an isolated solution, setting $\ell = n_1 + 1$ ensures that this solution is also isolated in \tilde{E} with probability at least $\frac{1}{2}$. Consequently, the algorithm has to be repeated a few times (with independent probabilistic polynomials) until it is output.

We also need to argue that E has an isolated solution with high probability. The (y, z) variable partition groups a solution to E together with $2^{n_1} - 1$ different assignments. In a cryptographic setting, we may assume that each such assignment satisfies E (containing m equations) with probability 2^{-m} . Thus, a solution to E is isolated with probability at least $1 - 2^{n_1 - m}$, which is typically very close to 1, as in our case we set $n_1 < n/5$ (to optimize complexity) and we usually have $m \gg n/5$.

Enumerating isolated solutions. Similarly to [12], isolated solutions are recovered bit-by-bit by computing n_1 sums, but we enumerate isolated solutions of \tilde{E} rather than E. Define the polynomials

$$U_0(y) = \sum_{\hat{z} \in \{0,1\}^{n_1}} \tilde{F}(y, \hat{z}), \text{and } U_i(y) = \sum_{\hat{z} \in \{0,1\}^{n_1-1}} \tilde{F}_{z_i \leftarrow 0}(y, \hat{z})$$

for $i \in \{1, \ldots, n_1\}$.

Proposition 3.2. Assume that (\hat{y}, \hat{z}) is an isolated solution to \tilde{E} with respect to (y, z). Then, $U_0(\hat{y}) = 1$ and $U_i(\hat{y}) = \hat{z}_i + 1$ for all $i \in \{1, \ldots, n_1\}$.

Proof. The proposition follows from Proposition 2.3, applied with \tilde{E} and \tilde{F} , instead of E and F.

Exploiting the low degree of \tilde{F} , our algorithm interpolates all $n_1 + 1$ polynomials $U_i(y)$ for $i \in \{0, \ldots, n_1\}$ and then evaluates each one on all $\hat{y} \in \{0, 1\}^{n-n_1}$ to recover isolated solutions.⁴ We optimize the interpolation of the polynomials $U_i(y)$, exploiting the following proposition.

Proposition 3.3. Let $\tilde{E} = \{R_i(y, z)\}_{i=1}^{\ell}$ and let \tilde{F} by its identifying polynomial. Denote by $d_{\tilde{F}}$ the degree of \tilde{F} and let $w = d_{\tilde{F}} - n_1$. The polynomial $U_0(y)$ can be interpolated from the solutions to \tilde{E} in the set $W_w^{n-n_1} \times \{0,1\}^{n_1}$, while $U_i(y)$ for $i \in \{1, \ldots, n_1\}$ can be interpolated from the solutions to \tilde{E} in the set $W_{w+1}^{n-n_1} \times \{0,1\}^{i-1} \times \{0\} \times \{0,1\}^{n_1-i}$. Hence, all the $n_1 + 1$ polynomials can be interpolated from the solutions to \tilde{E} in the set $W_{w+1}^{n-n_1} \times \{0,1\}^{i-1} \times \{0\} \times \{0,1\}^{n_1-i}$. Hence, all the $n_1 + 1$ polynomials can be interpolated from the solutions to \tilde{E} in the set $W_{w+1}^{n-n_1} \times \{0,1\}^{n_1}$.

Proposition 3.3 shows that the domains of the system E solved for interpolating $U_i(y)$ for $i \in \{1, \ldots, n_1\}$ overlap. Instead of naively solving \tilde{E} on $n_1 + 1$ overlapping domains, we solve \tilde{E} on one (slightly bigger) domain. Specifically, we use the exhaustive search algorithm of [8] for this purpose. We note that the analysis of [8] requires randomness assumptions about the input system, yet the other optimizations described here for enumerating isolation solutions do not.

⁴ As in the previous algorithms [7,12], we never explicitly interpolate the probabilistic polynomial \tilde{F} itself, but only the polynomials $U_i(y)$ derived from it.

Proof. By Proposition 2.1, the algebraic degree of $U_0(y) = \sum_{\hat{z} \in \{0,1\}^{n_1}} \tilde{F}(y,\hat{z})$ is at most $d_{\tilde{F}} - n_1 = w$, while the algebraic degree of each $U_i(y)$ for $i \in \{1, \ldots, n_1\}$ is at most $d_{\tilde{F}} - n_1 + 1 = w + 1$.

Therefore, $U_0(y)$ can be interpolated from its values in the set $W_w^{n-n_1}$, where the computation of each such value requires 2^{n_1} evaluations of $\tilde{F}(y, z)$. Thus, $U_0(y)$ can be interpolated from the values of $\tilde{F}(y, z)$ in the set $W_w^{n-n_1} \times \{0, 1\}^{n_1}$. Similarly, $U_i(y)$ for $i \in \{1, \ldots, n_1\}$ can be interpolated given the values of $\tilde{F}(y, z)$ in the set $W_{w+1}^{n-n_1} \times \{0, 1\}^{i-1} \times \{0\} \times \{0, 1\}^{n_1-i}$. The proposition follows since \tilde{F} is the identifying polynomial of \tilde{E} , and hence $\tilde{F}(\hat{y}, \hat{z}) = 1$ if and only if (\hat{y}, \hat{z}) is a solution to \tilde{E} .

Testing solutions. Similarly to [12], for each $\hat{y} \in \{0,1\}^{n-n_1}$, the algorithm computes $n_1 + 1$ sums. In our case, those with $U_0(\hat{y}) = 1$ suggest some solution (\hat{y}, \hat{z}) to \tilde{E} (and hence to E), and we need to test each one. However, these tests make expensive evaluations of polynomials, which generally require about $\binom{n}{\downarrow d}$ bit operations. This may lead to a large overhead, particularly for d > 2. In order to reduce this overhead, we repeat the algorithm a small number of times (using independent probabilistic polynomials) and test only candidate solutions that are output more than once. This is an additional concrete optimization over the second algorithm of [12], and makes use of randomness assumptions about the input system to argue that it is unlikely for an incorrect candidate solution to be suggested more than once.

Comparison to the previous works [7,12]. Our algorithm differs from previous works in each of the three elements mentioned above.

First, unlike the worst-case setting of [7,12], isolating solutions in a cryptographic setting is essentially trivial. In particular, there is no need for a random change of variables, and the parameter n_1 will simply be chosen to optimize the complexity. In addition, the procedure of testing solutions is more efficient than the one of [12] as explained above.

Technically, a more interesting modification is that we enumerate isolated solutions to \tilde{E} instead of E as in [12]. As a result, we only need to compute sums of the form $\sum_{\hat{z} \in \{0,1\}^{n_1}} \tilde{F}(\hat{y}, \hat{z})$. This is a significant concrete optimization, as accurate sums of the form $\sum_{\hat{z} \in \{0,1\}^{n_1}} F(\hat{y}, \hat{z})$ are too expensive to compute directly due to the high degree of F. In previous algorithms of [7,12], such sums are computed by majority voting across $48 \cdot n + 1$ evaluations of different polynomials derived from E, which had to be interpolated and evaluated. Consequently, the complexity of our algorithm is reduced by a factor of $\Omega(n)$ while additional savings are obtained via Proposition 3.3. Thus, our algorithm eliminates majority voting altogether and uses probabilistic polynomials in a different and a more direct way to solve E.

The asymptotic complexity of the algorithm is determined by two terms, similarly to (5). It could be improved using the techniques of [12], essentially by recursively solving the multiple parity-counting problem on \tilde{E} for sets of the

form $W_w^{n-n_1} \times \{0,1\}^{n_1}$ (rather than applying brute force). Yet, these recursive calls have a significant concrete overhead (they require working with many more probabilistic polynomials) and we do not use them, as noted in Section 1.3.

4 Details and Analysis of the New Algorithm

The pseudo-code of our main algorithm in given in Algorithm 1. It uses procedures 1 and 2. We now describe it in detail and then analyze its complexity.

Details of Algorithm 1. The main loop of Algorithm 1 runs until we find a solution to E. In each iteration, we define a new probabilistic set of equations \tilde{E} from E and call Procedure 1 to output all candidate solutions to \tilde{E} . The output of Procedure 1 is a 2-dimensional $2^{n-n_1} \times (n_1 + 1)$ array that contains for each $\hat{y} \in \{0, 1\}^{n-n_1}$, the evaluations $U_0(\hat{y})$ and $U_i(\hat{y}) + 1$ for $i \in \{1, \ldots, n_1\}$. Hence, by Proposition 3.2, assuming that (\hat{y}, \hat{z}) is an isolated solution to \tilde{E} , then $U_0(\hat{y}) = 1$ and $U_i(\hat{y}) + 1 = \hat{z}_i$ for $i \in \{1, \ldots, n_1\}$.

We store candidate solutions to \tilde{E} in an array and check whether a potential solution has been output before (for a previous probabilistic set of equations). Such a potential solution is tested against the full system E.

Details of procedures 1 and 2. Procedures 1 and 2 output the potential solutions to a given system \tilde{E} by interpolating the polynomials $U_i(y)$ for $i \in \{0, \ldots, n_1\}$ and evaluating them on all $\hat{y} \in \{0, 1\}^{n-n_1}$.

Recall from Proposition 3.3 that $U_i(y)$ for $i \in \{0, \ldots, n_1\}$ can be interpolated by solving \tilde{E} in the set $W_{w+1}^{n-n_1} \times \{0,1\}^{n_1}$, where $w = d_{\tilde{F}} - n_1$ and $d_{\tilde{F}}$ is the degree of \tilde{F} . In Procedure 2, these solutions are output by the fast exhaustive search algorithm of [8].

We denote by L the number of solutions and store them in memory. Next, we need to compute the values of each $U_i(y)$ in the sets described in Proposition 3.3 (with the aim of interpolating it).⁵ These values are computed by summing the evaluations of $\tilde{F}(y, z)$ on the corresponding subset of $\hat{z} \in \{0, 1\}^{n_1}$. The values of all the polynomials $U_i(y)$ for $i \in \{0, \ldots, n_1\}$ are computed in parallel by iterating over the solutions to the system. For each solution, we calculate its contribution to each of the relevant polynomials. Having calculated the required values of each of the polynomials, we return them.

Then, in Procedure 1, we interpolate $U_i(y)$ for $i \in \{0, \ldots, n_1\}$ using the Möbius transform. Finally, we evaluate all the n_1+1 polynomials on the full range $\hat{y} \in \{0, 1\}^{n-n_1}$ using the Möbius transform and output the potential solutions.

⁵ In practice, we do not need to store all the L solutions in memory at once, but we can interleave the exhaustive search with the computation of the $U_i(y)$ values.

Р	arameters: $n_1, d_{\tilde{F}}$							
In	Initialization: $\ell \leftarrow n_1 + 1, w \leftarrow d_{\tilde{F}} - n_1$							
1:	$PotentialSolutionsList[0] \leftarrow NewList()$							
2:	2: for all $k = 0, 1,$ do							
3:	Pick a uniformly random matrix of full rank ℓ , $A^{(k)} \in \mathbb{F}_2^{\ell \times m}$. Compute $\tilde{E}^{(k)} =$							
	$\{R_i^{(k)}(x)\}_{i=1}^{\ell} = \{\sum_{j=1}^m A_{i,j}^{(k)} \cdot P_j(x)\}_{i=1}^{\ell}$							
4:	$CurrPotentialSolutions \leftarrow OutputPotentialSolutions(\{R_i^{(k)}(x)\}_{i=1}^{\ell}, n_1, w)$							
5:	$PotentialSolutionsList[k] \leftarrow CurrPotentialSolutions$							
6:	for all $\hat{y} \in \{0,1\}^{n-n_1}$ do							
7:	if $CurrPotentialSolutions[\hat{y}][0] = 1$							
	\triangleright test whether $U_0(\hat{y}) = 1$, i.e., entry is valid							
	then							
8:	for all $k_1 \in \{0, \dots, k-1\}$ do							
9:	if $CurrPotentialSolutions[\hat{y}] = PotentialSolutionsList[k_1][\hat{y}]$							
	⊳check whether solution appears twice							
	then							
10:	$sol \leftarrow \hat{y} \ CurrPotentialSolutions[\hat{y}][1 \dots n_1]$							
	\triangleright concatenate n_1 least significant bits							
11:	if TestSolution $(\{P_j(x)\}_{j=1}^m, sol) = \text{TRUE then}$							
12:	return sol							
13:	break \triangleright continue with next \hat{y}							

Algorithm 1: Solve $(\{P_j(x)\}_{j=1}^m)$

4.1 Time Complexity Analysis

We now analyze the expected time complexity of the algorithm, denoted by $T = T_{n_1,d_{\bar{F}}}(n,m,d)$, in terms of bit operations. For this purpose we define the following notation:

- $-N_k$ is the expected number main loop iterations of Algorithm 1.
- $-T_1$ is the expected complexity of an iteration of Algorithm 1, not including the complexity of testing solutions.
- $-N_s$ is the expected total number of solutions tested by Algorithm 1.
- $-T_s$ is the average complexity of testing a solution.

We consider all of these variables (except T_s) as random variables that depend on the randomness of the algorithm and the random choice of E, whose distribution will be defined next. We do not consider T_s as a random variable and will analyze the overall complexity of testing solutions separately.

Throughout the analysis we use the symbol \leq that suppressed factors which are negligible both asymptotically and concretely for relevant concrete parameter choices. Specifically, several terms will be neglected based on assumptions that n_1 is sufficiently large (such as $2^{n_1} \gg n$). We have already used and justified such assumptions (see Remark 2.2).

```
1: (V, ZV[1...n_1]) \leftarrow \text{ComputeUValues}(\{R_i(y, z)\}_{i=1}^{\ell}, n_1, w)
                                                       \trianglerightobtain values for interpolating each U_i(y)
 2: Interpolate U_0(y): apply Möbius transform to V[1 \dots |W_w^{n-n_1}|]
 3: for all i \in \{1, ..., n_1\} do
        Interpolate U_i(y): apply Möbius transform to ZV[i][1 \dots |W_{w+1}^{n-n_1}|]
 4:
 5: Evals[0...n_1][0...2^{n-n_1}-1] \leftarrow \vec{0}
                                                                                   ⊳init evaluation array
 6: for all i \in \{0, 1, \dots, n_1\} do
        Evaluate U_i(y) on \{0,1\}^{n-n_1}
                                                      by Möbius transform. Store result in
 7:
Evals[i][0...2^{n-n_1} - 1]
8: Out[0...2^{n-n_1} - 1][0...n_1] \leftarrow \vec{0}
                                                                                               ⊳init output
 9: for all \hat{y} \in \{0,1\}^{n-n_1} do
        if Evals[0][\hat{y}] = 1 then
10:
11:
           Out[\hat{y}][0] \leftarrow 1
                                                             \trianglerightindicate U_0(\hat{y}) = 1, i.e., entry is valid
12:
           for all i \in \{1, ..., n_1\} do
               Out[\hat{y}][i] \leftarrow Evals[i][\hat{y}] + 1
13:
                                              ⊳copy potential solution by flipping evaluation bit
14: return Out
```

Procedure 1: OutputPotentialSolutions($\{R_i(x)\}_{i=1}^{\ell}, n_1, w$)

Probabilistic setting. We assume that the polynomials of E are chosen independently and uniformly at random from all degree d polynomials, conditioned on having a pre-fixed solution (e.g., a cryptographic key) chosen initially independently of E. The goal is to find the pre-fixed solution, rather than an arbitrary solution, which could be much easier to find (e.g., if E is under-determined, namely $m \ll n$). Thus, the analysis is also relevant in case there are additional (application dependent) constraints on solutions that are not modeled by the equation system, but are enforced when testing solutions.

A formal analysis of Algorithm 1 is given in Appendix A. Below, we give a simple heuristical analysis, assuming each assignment \hat{x} that is not the pre-fixed solution satisfies any polynomial equation in E with probability 1/2 independently of the other assignments and equations.

Theorem 4.1 (Heuristic). For a random equation system, the success probability of Algorithm 1 is at least $1 - 2^{n_1-m}$. Given that $m \ge 2 \cdot (n_1 + 1) + 2$ and $T_s \ll n_1 \cdot n \cdot 2^{n_1}$, its expected running time satisfies

$$T \lesssim 4 \left(2d \cdot \log n \cdot 2^{n_1} \cdot {\binom{n-n_1}{\downarrow d_{\tilde{F}} - n_1 + 1}} + n_1 \cdot n \cdot 2^{n-n_1} \right) \le$$
(6)

$$4\left(2d \cdot \log n \cdot 2^{n_1} \cdot \binom{n-n_1}{\downarrow n_1 \cdot (d-1) + d+1} + n_1 \cdot n \cdot 2^{n-n_1}\right).$$
(7)

In Appendix A (Theorem A.1) we prove a similar result formally for a slightly modified variant of Algorithm 1. The main difference is that we lower bound its

1: $Sols[1...L] \leftarrow BruteForceSystem(\{R_i(y, z)\}_{i=1}^{\ell}, n - n_1, w + 1)$ \triangleright brute force on space $W_{w+1}^{n-n_1} \times \{0,1\}^{n_1}$ 2: $V[1 \dots |W_w^{n-n_1}|] \leftarrow \vec{0}, ZV[1 \dots n_1][1 \dots |W_{w+1}^{n-n_1}|] \leftarrow \vec{0} \quad \rhd \text{init values for each } U_i(y)$ 3: for all $(\hat{y}, \hat{z}) \in Sols[1 \dots L]$ do if $\operatorname{HW}(\hat{y}) \leq w$ 4: \triangleright values of HW more than w do not contribute to $U_0(y)$ then \triangleright get index of \hat{y} in $W_w^{n-n_1}$ 5: $index \leftarrow \text{IndexOf}(\hat{y}, n - n_1, w)$ V[index] = V[index] + 1 \triangleright sum is over \mathbb{F}_2 6: 7:for all $i \in \{1, ..., n_1\}$ do **if** $\hat{z}_i = 0$ 8: \triangleright only values with $z_i = 0$ contribute to $U_i(y)$ for i > 1then $index \leftarrow \text{IndexOf}(\hat{y}, n - n_1, w + 1)$ 9: ZV[i][index] = ZV[i][index] + 110:11: return $V, ZV[1...n_1]$

Procedure 2: ComputeUValues $(\{R_i(y,z)\}_{i=1}^{\ell}, n_1, w)$

success probability by $\frac{5}{8}$ for the sake of simplifying the proof. Regardless, the difference between the heuristic and formal results is small.

The total complexity of Algorithm 1. The complexity formula (7) establishes a tradeoff between two terms. First, the term

$$2d \cdot \log n \cdot 2^{n_1} \cdot \binom{n-n_1}{\downarrow w+1} = 2d \cdot \log n \cdot 2^{n_1} \cdot \binom{n-n_1}{\downarrow d_{\bar{F}} - n_1 + 1}$$
(8)

accounts for the brute force on the space $W_{w+1}^{n-n_1} \times \{0,1\}^{n_1}$ in Procedure 2 (based on the analysis of Section 2.3 for random systems). The second term accounts for the evaluation of the polynomials $U_i(y)$ on $\{0,1\}^{n-n_1}$ in Procedure 1 and is

$$(n_1 + 1) \cdot (n - n_1) \cdot 2^{n - n_1} \le n_1 \cdot n \cdot 2^{n - n_1} \tag{9}$$

(given that $n_1^2 + n_1 \ge n$). The free parameter n_1 is set to balance these terms and optimize the complexity. Assuming the terms are equal, based on the second term, the gain in complexity over 2^n bit operations is roughly $\frac{2^{n_1}}{8 \cdot n_1 \cdot n}$. In Section 4.2 we estimate the total complexity of Algorithm 1 by about $n^2 \cdot 2^{(1-1/2.7d)n}$ bit operations (setting $n_1 \approx \frac{n}{2.7d}$).

Next, we establish Theorem 4.1 and argue that the condition $T_s \ll n_1 \cdot n \cdot 2^{n_1}$ holds both asymptotically and for relevant concrete parameter choices.

Success probability analysis. The analysis is based the assumption that the pre-fixed solution is isolated in E. Note that it is placed in a group with $2^{n_1} - 1$

additional potential solutions that share the same y value. We thus estimate the probability that another solution exists in this group by 2^{n_1-m} , and the pre-fixed solution is isolated with probability at least $1 - 2^{n_1-m}$.

Time complexity analysis. We assume for simplicity that N_k and T_1 are independent. This gives

$$T \le N_k \cdot T_1 + N_s \cdot T_s. \tag{10}$$

The algorithm succeeds once the pre-fixed solution to E is isolated in two tribalistic equation systems. Given that the pre-fixed solution is isolated in E, by Proposition 3.1, every iteration isolates it with probability at least $\frac{1}{2}$. Hence the expected number of iterations is at most $N_k \leq 2 \cdot 2 = 4$.

We analyze the most expensive operations of procedures 2 and 1, showing that the terms (8) and (9) indeed dominate. We then analyze the cost of testing solutions.

Procedure 2. As noted above, using the analysis of Section 2.3, the brute force complexity is given by (8). In addition, we estimate

$$\mathbf{E}[L] = 2^{-\ell} \cdot |W_{w+1}^{n-n_1} \times \{0,1\}^{n_1}| = \frac{1}{2} \cdot \binom{n-n_1}{\downarrow w+1},\tag{11}$$

as the equation systems we solve by brute force have $\ell = n_1 + 1$ equations.

The complexity of computing the values of the arrays V and ZV is slightly more⁶ than $(n_1 + 1) \cdot E[L]$, which is negligible compared to (8), given that $2d \cdot \log n \cdot 2^{n_1} \gg \frac{1}{2}(n+1)$.

Procedure 1. The complexity of interpolating $(U_0(y), U_1(y), \ldots, U_{n_1}(y))$ from their evaluations is $n \cdot \binom{n-n_1}{\downarrow w} + n_1 \cdot n \cdot \binom{n-n_1}{\downarrow w+1} < (n_1+1) \cdot n \cdot \binom{n-n_1}{\downarrow w+1}$, which is negligible compared to (8) given that $2d \cdot \log n \cdot 2^{n_1} \gg (n_1+1) \cdot n$.

The complexity of evaluating these polynomials on $\{0,1\}^{n-n_1}$ is given in (9).

Estimating N_s . Fix an iteration pair $0 \leq i < j < N_k$. Given that we have at least $m \geq 2 \cdot (n_1 + 1) + 2 = 2 \cdot \ell + 2$ equations, the 2ℓ rows of $A^{(i)}$ and $A^{(j)}$ are linearly independent vectors over $\{0,1\}^m$ with high probability⁷ in which case $\tilde{E}^{(i)}$ and $\tilde{E}^{(j)}$ are independent equation systems with ℓ equations. Hence, restricting these systems to a specific $\hat{y} \in \{0,1\}^{n-n_1}$, the pair suggests the same n_1 -bit solution suffix \hat{z} with probability 2^{-n_1} . Considering all $\hat{y} \in \{0,1\}^{n-n_1}$, the expected number of suggested solutions is about 2^{n-2n_1} . As the number of iteration pairs is small and many systems restricted to \hat{y} do no suggest any solution since $U_0(\hat{y}) = 0$, we estimate

$$N_s = 2^{n-2n_1}.$$
 (12)

⁶ We note that the operations of the IndexOf functions can be implemented with small overhead because solutions are output by the brute force algorithm in fixed order.

⁷ Even if the rows of $A^{(i)}$ and $A^{(j)}$ have a few linear dependencies, it does not substantially affect the analysis.

Since (8) and (9) dominate T_1 , up to now we estimate

$$T \le N_k \cdot T_1 + N_s \cdot T_s \lessapprox 4 \cdot (2d \cdot \log n \cdot 2^{n_1} \cdot \binom{n-n_1}{\downarrow w+1} + n_1 \cdot n \cdot 2^{n-n_1}) + 2^{n-2n_1} \cdot T_s.$$

In order to establish Theorem 4.1, we show that

$$T_s \ll n_1 \cdot n \cdot 2^{n_1},\tag{13}$$

so the final term that corresponds to the total complexity of testing candidate solutions may be neglected compared to the term $4 \cdot n_1 \cdot n \cdot 2^{n-n_1}$.

Testing solutions. Naively testing a candidate solution requires evaluating two polynomials in E on average, which has complexity of about $2 \cdot \binom{n}{\downarrow d}$ bit operations. Asymptotically, for constant d, this complexity is negligible compared to $2^{n_1} = 2^{\Omega(n)}$, hence (13) holds.

Concretely, for d = 2, since $n \ll 2^{n_1}$, then $n^2 \ll n_1 \cdot n \cdot 2^{n_1}$ and (13) holds. However, when d > 2 is relatively large compared to n, then (13) may no longer hold for relevant parameter choices (e.g., for d = 4 and n = 128).

On the other hand, we can reduce this complexity such that it becomes negligible for relevant parameter choices by tweaking Algorithm 1. The main idea is to test the potential solutions in batches, reducing the amortized complexity, as described in Appendix B.

In practice, E is constructed from a cryptosystem, and for relatively large d one may simply test candidate solutions by directly evaluating the cryptosystem. Alternatively, given sufficiently many equations, it is possible to run Algorithm 1 for a few more iterations and test only candidate solutions suggested at least 3 times. This mildly increases the time complexity, but reduces the number of candidate solutions to test by another significant factor of about 2^{n_1} .

Experimental validation. The most important probabilistic quantities analyzed are L (11) and N_s (12) (the bound on N_k is rigourous). We have experimentally estimated N_s and L on small random systems with m = n polynomial equations of degree d. Each system was generated uniformly at random, conditioned on having a single solution. For example, we ran 20 independent experiments on systems generated with parameters n = 20, d = 2 and set $n_1 = 4$. The expected value and standard deviation of L were about 13192 and 58, respectively (where we estimated $L \approx 1/2 \cdot \binom{16}{\downarrow 7} = 13166$). The expected value and standard deviation of the number of candidate solutions tested were 1322 and 1078, respectively. We estimated this number in (12) to be about $N_s = 2^{20-8} = 4096$, but the analysis is conservative and based on union bounds. Changing some of the parameters (e.g., to d = 4) did not have a substantial effect on the results.

4.2 Complexity Estimation of Algorithm 1

We bound the complexity of Algorithm 1 up to *small* constants. Although we optimize the more precise formulas (6) and (7) for the applications, the goal of this section is to give a closed-form expression that is a good estimation.

Ignoring (small) constant terms, we write (7) as

$$O\left(\log n \cdot 2^{n_1} \cdot \binom{n-n_1}{\downarrow n_1 \cdot (d-1)+d+1}\right) + O\left(n^2 \cdot 2^{n-n_1}\right).$$

Since $d \ge 2$, we have $n_1 \cdot (d-1) + d + 1 \le (n_1+2) \cdot (d-1) + 1$. It will be convenient to rewrite the expression above with $t = n_1 + 2$ and obtain

$$O\left(\log n \cdot 2^t \cdot \binom{n-t+2}{\downarrow t \cdot (d-1)+1}\right) + O\left(n^2 \cdot 2^{n-t}\right).$$

Focusing on the first term, we have $\binom{n-t+2}{\downarrow t \cdot (d-1)+1} = O\left(n\sqrt{n} \cdot \binom{n-t}{t \cdot (d-1)}\right)$, given that $2(t \cdot (d-1)) \leq n-t+2$. We use the fact that $\binom{u}{v} \leq 2^{uH\left(\frac{v}{u}\right)}$ (where $H(p) = -p\log p - (1-p)\log(1-p)$ is the binary entropy function) and bound the first term by

$$O\left(n\log n \cdot \sqrt{n} \cdot 2^t \cdot 2^{(n-t)H\left(\frac{t \cdot (d-1)}{n-t}\right)}\right).$$

Writing $t = \gamma n$ for some $0 < \gamma < 1$, we bound the total complexity of Algorithm 1 by

$$O\left(n\log n \cdot \sqrt{n} \cdot 2^{\left(\gamma + (1-\gamma)H\left(\frac{\gamma \cdot (d-1)}{1-\gamma}\right)\right)n}\right) + O\left(n^2 \cdot 2^{(1-\gamma)n}\right) = O\left(n^2 \left(2^{\left(\gamma + (1-\gamma)H\left(\frac{\gamma \cdot (d-1)}{1-\gamma}\right)\right)n} + 2^{(1-\gamma)n}\right)\right).$$
(14)

If we choose γ such that $H\left(\frac{\gamma \cdot (d-1)}{1-\gamma}\right) \leq 1 - \frac{\gamma}{1-\gamma}$, then the complexity will be $O(n^2 \cdot 2^{(1-\gamma)n})$. Analysis reveals that a choice of $\gamma = 1/2.7d$ is sufficient for this purpose. Therefore, the complexity is $O\left(n^2 \cdot 2^{0.815n}\right)$ bit operations for d = 2 and $O\left(n^2 \cdot 2^{\left(1-\frac{1}{2.7d}\right)n}\right)$ in general. Comparing against Table 1 (at the end of this section) shows that the formula $n^2 \cdot 2^{(1-1/2.7d)n}$ slightly overestimates the complexity for relevant parameters.

4.3 Optimizing Memory Complexity

The expected memory complexity of the algorithm is about $4 \cdot (n_1+1) \cdot 2^{n-n_1}$ bits, dominated by storing the potential solutions output by the different executions of Procedure 1. A simple way to obtain a time-memory tradeoff is to guess several bits of x and repeat the algorithm for each guess. However, we can improve the memory complexity with essentially no penalty by making use of a memoryefficient implementation of the Möbius transform. **Memory-efficient Möbius transform.** We deal with the problem of evaluating a polynomial $F(x_1, \ldots, x_n)$ of degree d on the space $\{0, 1\}^n$ using the Möbius transform. Assume that d is not too large and the polynomial is represented by a bit array of size $\binom{n}{\downarrow d} \ll 2^n$. Moreover, assume that the application does not need to store the evaluation of the polynomial on the full space, but can work even if the space is partitioned into smaller spaces on which the polynomial is evaluated on the fly. Then, the memory complexity can be reduced as follows. Instead of allocating an array of size 2^n , we work directly with the recursive formula (2), $F(x_1, \ldots, x_n) = x_1 \cdot F_1(x_2, \ldots, x_n) + F_2(x_2, \ldots, x_n)$, by first evaluating $F_2(x_2, \ldots, x_n)$ (i.e., F(x) for $x_1 = 0$), and then calculating and evaluating $F_1(x_2, \ldots, x_n) + F_2(x_2, \ldots, x_n)$ (i.e., F(x) for $x_1 = 1$).

This algorithm does not work in-place, but only keeps in memory the recursion stack. The memory complexity is bounded by the formulas $M(n,d) = M(n-1,d) + \binom{n}{\downarrow d}$ and $M(n,n) = 2^n$. Thus, the total memory complexity is less than $n \cdot \binom{n}{\downarrow d}$. The time complexity in bit operations is bounded by $\binom{n}{\downarrow d} + 2 \cdot \binom{n-1}{\downarrow d} + \ldots + 2^{n-d-1} \cdot \binom{d+1}{\downarrow d} + 2^{n-d} \cdot d \cdot 2^d$.

Remark 4.1. A more precise evaluation reveals that the total number of bit operations is about $d \cdot 2^n$. The exhaustive search algorithm of [8] for enumerating all zeroes of a polynomial of degree d also requires $d \cdot 2^n$ bit operations. It would be interesting to investigate whether the recursive Möbius transform can compete with [8] in practice. We note that it was already observed in [15] that the Möbius transform on degree d polynomials requires $d \cdot 2^n$ bit operations, but the algorithm used a standard implementation with memory complexity of 2^n .

In our context, we will exploit the lower complexity of the top level recursive calls to further reduce the memory complexity (by a small factor), while keeping the time complexity below $n \cdot 2^n$. Specifically, for a parameter $k \approx n - \log \binom{n}{\lfloor d}$, we perform the top k levels of the recursion independently without saving the recursion stack (i.e., we recursively evaluate the input polynomial on all 2^k values of x_1, \ldots, x_k independently). At the bottom levels, we switch to the in-place implementation of the Möbius transform to evaluate the polynomial on all values of x_{k+1}, \ldots, x_n . In order to perform the independent evaluations, we only need to allocate two additional arrays, one for the input to the recursive call and one for its output. The roles of these arrays are interchanged on every recursive call. The memory required for each such additional array is bounded by $\binom{n}{\downarrow d}$. The memory required for the in-place Möbius transform is $2^{n-k} \approx \binom{n}{\downarrow d}$. Therefore, the in-place transform does not require additional memory and the total memory complexity is bounded by $3 \cdot \binom{n}{\downarrow d}$. The time complexity of the procedure is bounded by $2^k \cdot \left(\binom{n}{\downarrow d} + \ldots + \binom{n-k}{\downarrow d}\right) + (n-k) \cdot 2^n < n \cdot 2^n$.

We note that the algorithm of [8] could also be used for the same purpose. However, it requires a preprocessing phase of complexity n^{2d} . On the other hand, we will use the Möbius transform variant with relatively large d (e.g., d = n/3), for which such preprocessing is too expensive. Improving the memory complexity of Algorithm 1. In order to reduce the memory complexity we first interpolate the polynomials $(U_0(y), \ldots, U_{n_1}(y))$ for several executions (e.g., 4 or a bit more) of Procedure 1 in advance. Using the recursive version of the Möbius transform (as described in Section 2.3), the additional memory required is negligible.

The main idea that allows to save memory is to interleave the tasks of evaluating all the polynomials (in parallel) with testing solutions that are suggested at least twice. The parallel evaluation is performed using the memory-optimized implementation of the Möbius transform. This reduces the memory complexity to about 3 times the memory required to store all the polynomials. In fact, for the purpose of testing solutions, we only need to keep the evaluations of each such polynomial on a space proportional to its size, used by the in-place transform. Thus, when sequentially calculating several transforms, we reuse one of the two additional allocated arrays. In total, we require 2 times the memory used for storing all the polynomials, namely

$$8 \cdot (n_1 + 1) \cdot \binom{n - n_1}{\downarrow d_{\bar{\nu}} - n_1 + 1}.$$
(15)

Choosing n_1 that minimizes time complexity (balancing the two terms of (6)), gives $\binom{n-n_1}{\downarrow d_{\tilde{F}}-n_1+1} \approx \frac{n_1 \cdot n}{2d \cdot \log n} \cdot 2^{n-2n_1}$ and total memory complexity of about $\frac{4 \cdot n_1 \cdot (n_1+1) \cdot n}{d \cdot \log n} \cdot 2^{n-2n_1}$ bits. Compared to Algorithm 1, this saves a multiplicative factor of about $\frac{d \cdot \log n}{n_1 \cdot n} \cdot 2^{n_1}$. Since $n_1 < \frac{n}{5}$, $\frac{8 \cdot n_1 \cdot (n_1+1) \cdot n}{2d \cdot \log n} \approx n^2$ for relevant concrete parameters.

Asymptotically, the memory complexity is $O(n^3 \cdot 2^{n-2n_1})$ bits. A choice of n_1 that minimizes time complexity gives $O(n^3 \cdot 2^{0.63n})$ bits for d = 2 and $O(n^3 \cdot 2^{(1-1/1.35d)n})$ in general.

Concrete parameters. In Table 1 we give concrete complexity estimates for interesting parameter sets after optimizing the free parameter n_1 of (7).

Variables	Degree	Internal parameter	Complexity	Memory	Exhaustive search [8]
n	d	n_1	(bit operations)	(bits)	$(2d\log n \cdot 2^n)$
80	2	16	277	2^{60}	2^{84}
100	2	25	2^{117}	2^{91}	2^{133}
120	4	12	2^{129}	2^{112}	2^{134}
100	2	37	2^{170}	2^{132}	2^{197}
192	4	18	2^{188}	2^{164}	2^{198}
256	2	49	2^{223}	2^{173}	2^{261}
200	4	25	2^{246}	2^{219}	2^{262}

Table 1. Concrete complexity of (the memory-optimized variant of) Algorithm 1

5 Cryptanalytic Applications

In this section we describe cryptanalytic applications of Algorithm 1. Our main application is in cryptanalysis of Picnic (and LowMC) variants and our secondary application is in cryptanalysis of round-reduced Keccak. We begin by describing the general optimization method we use in cryptanalysis of Picnic.

5.1 Deterministic Replacement of Probabilistic Polynomials

In cryptanalytic applications, E may have some properties that depend on the underlying cryptosystem and could be ruined in \tilde{E} that mixes the equations of E. We observe that in cryptographic settings, we may replace the randomized construction of \tilde{E} (and \tilde{F}) by a deterministic construction which simply takes subsets of equations from E, thus preserving the properties of E.

Essentially, the randomness of the probabilistic constructions is already "embedded" in E itself. For example, we still (heuristically) expect a variant of Proposition 3.1 to hold: if \hat{x} is a solution to E, it is also a solution to \tilde{E} . On the other hand, since \tilde{E} is a system with $\ell = n_1 + 1$ equations in n_1 variables, we expect it not to have an additional solution with high probability.

In order for the number of tested candidate solutions to remain small, we require the different equations systems \tilde{E} analyzed to be roughly independent. Specifically, the intersections of the equations subsets taken for different "probabilistic" equations systems \tilde{E} should be empty (or small).

5.2 Picnic and LowMC

The Picnic signature scheme [10] is an alternate third-round candidate in NIST's post-quantum standardization project [29]. It uses a zero-knowledge protocol in order to non-interactively prove knowledge of a preimage x to a public value y under a one-way function f, where y is part of the public key and x is the secret signing key. The one-way function is implemented using a block cipher, where the secret signing key is the block cipher's key, while the public key consists of a randomly chosen plaintext and the corresponding ciphertext (the encryption of the plaintext with the secret key). Thus a key-recovery attack on Picnic reduces to finding the block cipher's secret key from one plaintext-ciphertext pair.

Picnic uses the LowMC block cipher family, proposed at EUROCRYPT 2015 by Albrecht et al. [1]. It is optimized for practical instantiations of multiparty computation, fully homomorphic encryption, and zero-knowledge proofs, in which non-linear operations are typically much more expensive than linear ones. Consequently, LowMC uses a relatively small number of multiplications.

LowMC is an SP-network built using several rounds, where in each round, a round-key is added to the state, followed by an application of a linear layer and a non-linear layer (operations are over \mathbb{F}_2). Finally, an additional round-key is added to the state. Importantly, the key schedule of LowMC is linear.

Each non-linear layer of LowMC consists of identical Sboxes $S : \{0, 1\}^3 \rightarrow \{0, 1\}^3$ of algebraic degree 2. The algebraic normal form of an Sbox is

$$\mathcal{S}(a_1, a_2, a_3) = (a_1 + a_2 a_3, a_1 + a_2 + a_1 a_3, a_1 + a_2 + a_3 + a_1 a_2). \tag{16}$$

We note that the inverse Sbox also has algebraic degree of 2.

In this paper, we focus on LowMC instances that were recently integrated into Picnic variants [34]. These instances have internal state and key lengths of 129, 192 and 255 bits, claiming security levels of 128, 192 and 255 bits, respectively and have a full non-linear layer (as opposed to other instances of LowMC that have a partial non-linear layer). All of these instances have 4 rounds, although in a recent publication by some of the designers [24] additional instances with 5 rounds were proposed in order to provide a larger security margin.

Attacks on LowMC instances. As noted above, we analyze LowMC instances with a full Sbox layer given only a single plaintext-ciphertext pair. The best-known attacks on such instances were recently published by Banik et al. [2] where the authors analyzed instances reduced to 2 rounds. Their techniques are based on linearization and it is not clear how to extend them beyond 2 rounds without exceeding the complexity of (optimized) exhaustive search [8].

We focus on the full 4 and 5-round instances. For the sake of completeness, we will also give results on round-reduced LowMC instances.

Fix an arbitrary plaintext-ciphertext pair to a LowMC instance with key and internal state size of n bits. We denote the unknown key by $x = (x_1, \ldots, x_n)$.

Even round number. We begin by considering LowMC instances with an even number of rounds r. We focus on an arbitrary state bit b_i (for some $i \in \{1, \ldots, n\}$) after r/2 rounds. Starting from the plaintext, we symbolically evaluate the encryption process and express b_i as a polynomial $p_i(x)$. Similarly, we symbolically evaluate the decryption process starting from the ciphertext and express b_i as a polynomial $q_i(x)$. This standard meet-in-the-middle approach gives rise to the equation $p_i(x) + q_i(x) = 0$. As the algebraic degree of the LowMC round and its inverse is 2 and the key schedule is linear, the algebraic degree of both $p_i(x)$ and $q_i(x)$ is at most $2^{r/2}$. Repeating this process n times for all intermediate state bits, we obtain an equation systems with m = n equations. For the small values of r we consider, the complexity of calculating the equation system is negligible. We can now apply Algorithm 1 and solve for the secret key.

Odd round number. For an odd number of rounds r, the approach above gives rise to equations of degree at least $2^{(r+1)/2}$, as in general, any intermediate state bit has algebraic degree of at least $2^{(r+1)/2}$ from either the encryption or the decryption side. If we apply Algorithm 1 in a straightforward manner, its complexity compared to an attack on an even number of r-1 rounds would increase substantially. We now show that by a better choice of the equation system and careful analysis we can reduce the algorithm's complexity. Consider the first 3 intermediate state bits b_1, b_2, b_3 that are outputs of the first Sbox in round (r+1)/2. From the decryption side, we can express them as polynomials of degree $2^{(r-1)/2}$, denoted by $q_1(x), q_2(x), q_3(x)$, respectively. From the encryption side, based on (16), we can express these bits as functions of the bits a_1, a_2, a_3 that are inputs the first Sbox in round (r+1)/2,

$$(b_1, b_2, b_3) = \mathcal{S}(a_1, a_2, a_3) = (a_1 + a_2 a_3, a_1 + a_2 + a_1 a_3, a_1 + a_2 + a_3 + a_1 a_2).$$

From the encryption side, we can express each of a_1, a_2, a_3 as a polynomial of degree $2^{(r-1)/2}$ in the key. Equating each bit to its evaluation from the decryption side, we obtain the 3 equations

$$q_1(x) + a_1(x) + a_2(x)a_3(x) = 0,$$

$$q_2(x) + a_1(x) + a_2(x) + a_1(x)a_3(x) = 0,$$

$$q_3(x) + a_1(x) + a_2(x) + a_3(x) + a_1(x)a_2(x) = 0.$$

Each polynomial appearing in these equations is of algebraic degree $2^{(r-1)/2}$.

Recall that the main complexity formula (6) heavily depends on the value of $d_{\tilde{F}}$. This value is upper bounded by $d \cdot \ell$, but it could be lower if we choose \tilde{F} more carefully. Indeed, we will construct the "probabilistic polynomials" deterministically using (non-overlapping) subsets of the original equation system E. The probabilistic analysis is essentially unchanged, as suggested in Section 5.1.

Specifically, in this case, each equation is of degree $2^{(r-1)/2} + 2^{(r-1)/2} = 2^{(r+1)/2}$ due to the multiplication of the a_i 's. However, if we multiply the 3 polynomials for the purpose of calculating \tilde{F} (as in (4), but with R_i 's replaced by the equations above), the term $a_1(x)a_2(x)a_3(x)$ can only be multiplied with at most one of the $q_i(x)$'s and therefore the degree of multiplication of all the equations is at most $4 \cdot 2^{(r-1)/2} = 2^{(r+3)/2}$, rather than the trivial upper bound of $6 \cdot 2^{(r-1)/2}$, reducing the degree by a factor of $\frac{1}{3}$. For example, for r = 3, we get a bound of 8 instead of the general upper bound of 12, whereas for r = 5, we get a bound of 16 instead of the general upper bound of 24.

We proceed to collect n equations as before. However, in Algorithm 1, for some integer ℓ' we analyze equations that are computed as above using ℓ' triplets that are outputs of the Sbox layer of round (r + 1)/2. We now have $\ell = 3\ell'$ equations. The total degree of \tilde{F} in (4) is upper bounded by $d_{\tilde{F}} \leq \ell' \cdot 2^{(r+3)/2} = \ell/3 \cdot 2^{(r+3)/2}$. We choose $n_1 = \ell - 1$ as before. Revisiting the complexity analysis formula of (6), we obtain

$$4\left(2\cdot 2^{(r+1)/2}\cdot \log n\cdot 2^{n_1}\cdot \binom{n-n_1}{\downarrow\frac{\ell}{3}\cdot 2^{(r+3)/2}-n_1+1}+n_1\cdot n\cdot 2^{n-n_1}\right).$$
 (17)

If we take ℓ which is not a multiple of 3, then $d_{\tilde{F}}$ increases more sharply due to the last Sbox. Specifically, if $\ell \mod 3 = 1$, then $d_{\tilde{F}} \leq \frac{\ell-1}{3} \cdot 2^{(r+3)/2} + 2^{(r+1)/2}$, whereas if $\ell \mod 3 = 2$, then $d_{\tilde{F}} \leq \frac{\ell-2}{3} \cdot 2^{(r+3)/2} + 3 \cdot 2^{(r-1)/2}$. We note that standard approaches to deal with the middle Sbox layer (e.g., linearization) result in higher complexity.

Results. Table 2 summarizes our attacks on instances of LowMC. The most important ones have 4 rounds and are used in recent Picnic variants [24]. Attacks on 2-round instances are inferior to those of [2] and are only given for the sake of completeness. On the other hand, attacks on 3-round instances are of interest, as they best demonstrate our advantage over previous works (there are no previous attacks on these instances that are better than exhaustive search [8]).

Remark 5.1. Solutions can be tested simply by evaluating the LowMC encryption process, whose most expensive procedures are the evaluations of the linear layers, each consisting of a multiplication of an *n*-bit state with an $n \times n$ matrix. Naively, this has complexity of $2n^2$ bit operations. It can be checked that (13) holds for the parameters of Table 2.

Security level	Key length	Rounds	Internal parameters	Attack complexity	Memory
S	n	r	$(n_1, d, d_{ ilde F})$	(bit operations)	(bits)
	129	2	(25, 2, 52)	2 ¹¹⁸	2^{92}
128		3	(18, 4, 52)	2^{125}	2^{104}
		4	(12, 4, 52)	2^{130}	2^{113}
	192	2	(37, 2, 72)	2^{170}	2^{126}
100		3	(27, 4, 76)	2^{180}	2^{150}
192		4	(18, 4, 76)	2^{188}	2^{164}
		5	(14, 8, 80)	2^{192}	2^{173}
		2	(49, 2, 100)	2222	2^{173}
055	255	3	(36, 4, 100)	2^{235}	2^{197}
200		4	(25, 4, 104)	2^{245}	2^{218}
		5	(18, 8, 104)	2^{251}	2^{228}

 Table 2. Attacks on LowMC instances

5.3 Keccak

Keccak is a family of cryptographic functions, designed by Bertoni et al. in 2008 [5]. We focus on the Keccak hash function family, selected by NIST in 2015 as the SHA-3 standard. It is built using the sponge construction (cf. [6]) using a permutation that operates on a 1600-bit state. The permutation consists of 24 rounds, where each round consists of an application of a non-linear layer, followed by linear operations over \mathbb{F}_2 . Importantly, the non-linear layer is of algebraic degree 2. We analyze the 4 basic Keccak variants which are parameterized by the output size of k bits and denoted by Keccak-k for $k \in \{224, 256, 384, 512\}$.

A short description of Keccak is given in Appendix C.

Preimage attacks of round-reduced Keccak. We consider messages of length that is smaller than the *rate* of the hash function (so that the output is produced after a single invocation of the permutation). We start by representing the message (preimage) bits as symbolic variables. We then linearize the first round of Keccak by setting some linear constraints on the variables, such that each state bit in the second round of the permutation is a linear polynomial in these variables. Using the linearization technique of [13,20] for Keccak-k (that selects variables that keep the *column parities* constant), this leaves more than 224 and 256 free variables for Keccak-224 and Keccak-256, respectively, and 256 and 128 free variables for Keccak-384 and Keccak-512, respectively (for the SHA-3 versions, the number is slightly smaller).

For Keccak-384 and Keccak-512, we can further partially invert the final nonlinear layer (applied to the first 5×64 Sboxes) on the target image to obtain its input values. For Keccak-224 and Keccak-256, not all these 5×64 output bits are fixed by the image. However, we obtain 192 and 256 linear relations among the input bits of the final Sbox layer, for Keccak-224 and Keccak-256 respectively (e.g., see [20]). Having peeled off 2 out of the 4 non-linear layers, we obtain equations of degree $2^2 = 4$ and solve for the preimage using Algorithm 1.

Results. We begin by considering preimage attacks on 4 rounds of Keccak-224 and Keccak-256. For these instances, we have sufficiently many free variables to obtain systems with 224 and 256 variables (respectively), which we assume to have a solution. Consequently, our preimage attacks have complexities of 2^{217} and 2^{246} bit operations by choosing $n_1 = 22$ and $n_1 = 25$, respectively. On the other hand, the recent independent work [21] obtains complexities of 2^{192} and 2^{218} for Keccak-224 and Keccak-256. The analysis of these attacks ignores the complexity of solving (numerous) linear equation systems over \mathbb{F}_2 with hundreds of variables, but they outperform our attacks nevertheless.

For Keccak-384, the number of free variables is only 256, so we need to solve systems of degree 4 with 256 variables an expected number of $2^{384-256} = 2^{128}$ times (with different initial linear constraints on the variables) to obtain a solution. This requires time $2^{246+128} = 2^{374}$ (by choosing $n_1 = 25$). In terms of bit operations, this improves upon the recent result of [26], which estimated the attack complexity by 2^{375} evaluations of the 4-round Keccak-384 function.

For Keccak-512, we do not linearize the first round, but directly solve a system of degree 8 in 512 variables. This requires 2^{502} bit operations (by choosing $n_1 = 26$) and improves the previous attack [28] that requires 2^{506} Keccak calls.

Collision attacks on round-reduced Keccak. We show how to exploit our algorithm to mount collision attacks on (low-degree) hash functions with an output length of k bits. We then outline a (marginal) collision attack on 4-round Keccak-512. Yet, the main purpose of this part is to introduce the general attack framework which uses our algorithm in a different way compared to its previous applications to key-recovery attacks and to preimage attacks on hash functions.

The problem of finding a collision can be formulated as a non-linear equation system (where the variables are the bits of two colliding messages). However, the complexity of solving such a system is unlikely to be more efficient than a generic birthday attack on the k-bit hash function which takes time $2^{k/2}$. A better idea is to directly speed up the generic birthday attack. In order to do so, for a parameter ℓ , we fix ℓ bits of the output to an arbitrary value v and try to find $2^{(k-\ell)/2}$ messages whose output value on the ℓ bits is equal to v. With high probability, these messages contain a pair whose outputs also agree of the remaining $k - \ell$ bits that we have not fixed, and therefore constitute a colliding pair. In order to find $2^{(k-\ell)/2}$ such messages, we apply Procedure 1.

Suppose we run Procedure 1 with n variables after fixing $\ell = n_1 + 1$ output bits. We expect the output to contain about $\frac{1}{2} \cdot 2^{n-n_1}$ isolated solutions (messages) that satisfy the n_1 constraints. We evaluate the hash function on each output message, and test whether it is indeed a solution. We store all the true solutions and sort them, trying to find a collision among them. If $\frac{1}{2} \cdot 2^{n-n_1} < 2^{(k-\ell)/2}$ (e.g., we lack degrees of freedom and are forced to choose a small value of n), we repeat the procedure several times (with a different set of message variables), while storing all the produced messages until a collision is found.

Since we test all outputs of Procedure 1, the complexity of this attack also directly depends on the number of bit operations required to evaluate the hash function on some message. We denote this number by τ .

Application to Keccak-512. We apply the framework to 4-round Keccak-512, as there are no published attacks better than the birthday bound on this variant. Assuming $\tau = 2^{13}$ (hence the complexity of the birthday attack is $2^{256+13} = 2^{269}$ bit operations), we set d = 4 and n = 128 (unlike the preimage attack, we linearize the first round) and choose $n_1 = \ell - 1 = 12$. Calculation reveals that the complexity is about 2^{263} bit operations, which is roughly 64 times faster than the birthday attack.⁸ If we assume a larger value of τ , the complexity of the attack will increase, but also its relative advantage compared to the birthday attack.

Further optimizations. We leave it to future work to further optimize these attacks by better exploiting the structure of Keccak. For example, some improvement can be obtained by selecting the variables z_1, \ldots, z_{n_1} in a more careful manner in order to reduce the degree of \tilde{F} in Algorithm 1. Additionally, exhaustive search over $W_{w+1}^{n-n_1} \times \{0,1\}^{n_1}$ may be replaced by a more efficient procedure that takes advantage of the structure of the equation system.

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⁸ The complexity of sorting the $2^{(512-13)/2} = 2^{249.5}$ images is estimated to be smaller than 2^{262} bit operations.

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A Formal Analysis of Algorithm 1 with Random Input

In this section we formally analyze Algorithm 1. We begin with worst-case analysis that holds for any input polynomial system E. We then define a distribution over equation systems and analyze a slightly modified variant of Algorithm 1 for this input distribution. Throughout the analysis, we use the notation defined in Section 4.1.

A.1 Worst-Case Analysis

Proposition A.1.

 $T_1 \lesssim T_b + n_1 \cdot n \cdot 2^{n-n_1} + (n_1+1) \cdot (\mathbb{E}[L] + n \cdot \binom{n-n_1}{\downarrow w+1} + N_k \cdot 2^{n-n_1}),$

where T_b is the expected brute force complexity.

Proof. We analyze procedures 2 and 1, and the additional operations in Algorithm 1.

Procedure 2. The expected complexity of brute force is T_b . The expected complexity of computing the values of the arrays V and ZV is slightly more than $(n_1 + 1) \cdot E[L]$.

Procedure 1. The complexity of interpolating $(U_0(y), U_1(y), \ldots, U_{n_1}(y))$ from their evaluations is

$$n \cdot \binom{n-n_1}{\downarrow w} + n_1 \cdot n \cdot \binom{n-n_1}{\downarrow w+1} < (n_1+1) \cdot n \cdot \binom{n-n_1}{\downarrow w+1}.$$

The complexity of evaluating these polynomials on $\{0,1\}^{n-n_1}$ is bounded by $n_1 \cdot n \cdot 2^{n-n_1}$ (as in (9)). The complexity of the final loop that outputs potential solutions is about $(n_1 + 1) \cdot 2^{n-n_1}$, which is negligible compared to (9).

Algorithm 1. The computation time of the probabilistic polynomials is negligible.

The work performed inside the loop in iteration k involves at most k comparison operations, each requires at most $n_1 + 1$ bit operations. The total number of bit operation spent on comparisons (in all iterations) is upper bounded by $\binom{N_k}{2} \cdot (n_1+1) \cdot 2^{n-n_1}$, while the average per iteration is at most $N_k \cdot (n_1+1) \cdot 2^{n-n_1}$.

A.2 Random Equation Systems

We define a random process for choosing a polynomial system $E = \{P_j(x_1, \ldots, x_n)\}_{j=1}^m$ of degree d, conditioned on a pre-fixed solution \hat{x}^* , chosen arbitrarily, but independently of E. This random process is similar to the one used in [3], and is a reasonable model for cryptographic settings.

We first describe how to generate a uniform polynomial of degree d with a pre-fixed solution of \hat{x}^* (in n variables): choose all the $\binom{n}{\downarrow d} - 1$ coefficients of the polynomial independently and uniformly at random, except for the free coefficient (that is not multiplied with any variable). Denote this intermediate polynomial by \bar{P} . Define P by setting all of its $\binom{n}{\downarrow d} - 1$ non-free coefficients equal to those of \bar{P} , and set its free coefficient to $\bar{P}(\hat{x}^*)$ (such that $P(\hat{x}^*) = 0$). Note that P(x) is a random variable that is uniformly distributed in the space of all degree d polynomials that evaluate to 0 at \hat{x}^* . Indeed, there are $2^{\binom{n}{\downarrow d}-1}$ such polynomials and each one is selected with the same probability.

Given \hat{x}^* , we generate E (with parameters n, m, d) by choosing each of its m polynomials as an independent polynomial with the pre-fixed solution \hat{x}^* . We call (the random variable) E generated by the above process a *uniform system* with a pre-fixed solution of \hat{x}^* (with parameters n, m, d).

In the following, we will analyze a modified variant of Algorithm 1, assuming that its input is chosen from such a distribution.

A.3 Analysis of Modified Algorithm 1

We analyze a modified variant of Algorithm 1, where we bound the number of iterations by 4 and return a solution it is found. Otherwise, the algorithm returns failure. Moreover, we define an event \mathcal{U} which holds if for each pair of iterations $0 \leq i < j < 4$, the 2ℓ rows of the matrices $A^{(i)}$ and $A^{(j)}$ are linearly independent. If \mathcal{U} does not hold, we abort and return failure.⁹ Obviously, the distribution of \mathcal{U} only depends on the parameters ℓ, m . We note that these modifications simplify the analysis, but are not required in practice. Our goal will be to bound the expected time complexity T. We have

$$T \le 4 \cdot T_1 + N_s \cdot T_s. \tag{18}$$

Analysis of uniform systems with a pre-fix solution. The following results analyze the probabilistic setting and will be used in the analysis of the algorithm.

Proposition A.2. Let $\hat{x}' \neq \hat{x}^*$ be any n-bit vectors. Let P'(x) be a uniform polynomial of degree $d \ge 1$ with a pre-fixed solution \hat{x}' . Define $P(x) = P'(x + \hat{x}' + \hat{x}^*)$. Then, P(x) is a uniform polynomial of degree d with a pre-fixed solution \hat{x}^* .

Proof. The polynomial P' is distributed uniformly over the space of all degree d polynomials that evaluate to 0 at \hat{x}' . Note that $P(\hat{x}^*) = P'(\hat{x}') = 0$ and the change of variables $P(x) = P'(x + \hat{x}' + \hat{x}^*)$ in an invertible transformation that does not increase the degree of P. Thus, the change of variable shifts the distribution to put zero weight on degree d polynomials that satisfy $P(\hat{x}^*) = 1$. Consequently, P(x) is distributed uniformly in the space of all degree d polynomials that evaluate to 0 at \hat{x}^* , proving the claim.

The following proposition shows that despite the pre-fixed solution, each polynomial of E in this setting essentially behaves as a uniform polynomial on almost all assignments.

Proposition A.3. Let P(x) be a uniform polynomial of degree $d \ge 1$ with a pre-fixed solution \hat{x}^* . Let $\hat{x} \neq \hat{x}^*$. Then, $\Pr[P(\hat{x}) = 0] = \frac{1}{2}$.

⁹ It is also possible to select the matrices $A^{(i)}$ in advance such that the event holds, assuming m is sufficiently larger than 2ℓ .

Proof. The value of $P(\hat{x})$ is computed as a sum $P(\hat{x}) = \sum_{u \in S_{\hat{x}}} \alpha_u(P)$, where $S_{\hat{x}} \subseteq \{0,1\}^n$ is the set of coefficients of monomials of that evaluate to 1 on the assignment \hat{x} .

If $\hat{x}^* = 0$, then the restriction $P(\hat{x}^*) = P(0) = 0$ only sets to 0 the free coefficient of $P(\alpha_0(P))$, while the other $\binom{n}{\downarrow d} - 1$ coefficients are chosen independently and uniformly at random. For any other assignment $\hat{x} \neq \hat{x}^* = 0$, $S_{\hat{x}} \subseteq \{0, 1\}^n$ contains at least one $u \neq 0$ such that the coefficient $\alpha_u(P)$ is uniformly and independently distributed. Hence, $\Pr[P(\hat{x}) = 0] = \frac{1}{2}$.

Otherwise, $\hat{x}^* \neq 0$, but the same argument holds by a change of variables. In particular, pick P(x) by picking P'(x) as a uniform polynomial with a pre-fixed solution of 0 and define $P(x) = P'(x + \hat{x}^*)$. By Proposition A.2 applied with $\hat{x}' = 0$, P is distributed as desired.

Therefore, $\Pr[P(\hat{x}) = 0] = \Pr[P'(\hat{x} + \hat{x}^*) = 0] = \frac{1}{2}$ by the first argument applied to P'(x) evaluated at the non-zero assignment $\hat{x} + \hat{x}^*$.

Proposition A.4. Let $S \subset \{0,1\}^n$ be a subset. Assume that E is a system of polynomials with parameters (n, m, d), such that each $\hat{x} \in S$ satisfies the following condition: for every $P(x) \in E$, $\Pr[P(\hat{x}) = 0] = \frac{1}{2}$ holds independently. Then: (1) the probability that E has a solution in S is at most $|S| \cdot 2^{-m}$, and (2) the expected number of solutions to E in S is $|S| \cdot 2^{-m}$.

Proof. By the above condition, every $\hat{x} \in S$ is a solution to E with probability 2^{-m} . Statement (1) follows by a union bound over all assignments in S and statement (2) follows by linearity of expectation.

The analysis of modified Algorithm 1. Let \mathcal{E} be the event that \hat{x}^* is an isolated solution to E. In the following, we bound $\Pr[\mathcal{E}]$ and complete the analysis of T_1 . We then bound $\Pr[\mathcal{U}]$. These will be used to finish the analysis of the algorithm.

Proposition A.5. $\Pr[\mathcal{E}] \ge 1 - 2^{n_1 - m}$.

Proof. Let $\hat{x}^* = (\hat{y}^*, \hat{z}^*)$ be the pre-fixed solution and define the set $S = \{(\hat{y}^*, \hat{z}) \mid \hat{z} \neq \hat{z}^*\}$. The event \mathcal{E} occurs if E does not have a solution in S. The claim follows by Proposition A.4 applied to S, whose size is $|S| = 2^{n_1} - 1$. The precondition for this proposition holds by Proposition A.3.

Proposition A.6.

$$T_1 \lesssim 2d \cdot \log n \cdot 2^{n_1} \cdot \binom{n-n_1}{\downarrow w+1} + n_1 \cdot n \cdot 2^{n-n_1}.$$

Proof. By Proposition A.1,

$$T_1 \lessapprox T_b + n_1 \cdot n \cdot 2^{n-n_1} + (n_1+1) \cdot (\mathbb{E}[L] + n \cdot \binom{n-n_1}{\downarrow w+1} + N_k \cdot 2^{n-n_1}).$$

We will estimate $T_b = 2d \cdot \log n \cdot 2^{n_1} \cdot {\binom{n-n_1}{\downarrow w+1}}$ and show that the first two terms dominate the complexity. The result will follow by linearity of expectation.

First, since $N_k \leq 4$, then $N_k \cdot (n_1+1) \cdot 2^{n-n_1} \ll n_1 \cdot n \cdot 2^{n-n_1}$, as $n_1 \cdot n \gg 4 \cdot (n_1+1)$. Moreover, $(n_1+1) \cdot n \cdot \binom{n-n_1}{\downarrow w+1} \ll T_b$ given that $2d \cdot \log n \cdot 2^{n_1} \gg (n_1+1) \cdot n$.

Next, observe that in each iteration k, $\tilde{E}^{(k)}$ is itself a uniform system with a pre-fixed solution (with ℓ equations). This follows as E is such a system and since $A^{(k)}$ has full rank ℓ . Therefore, by Proposition A.3 the condition in Proposition A.4 holds (with $S = W_{w+1}^{n-n_1} \times \{0,1\}^{n_1}$, excluding the pre-fixed solution). We can now use the complexity estimate of Section 2.3 for the algorithm of [8]: $T_b = 2d \cdot \log n \cdot 2^{n_1} \cdot {\binom{n-n_1}{lw+1}}$.

Finally, applying Proposition A.4 again with $S = W_{w+1}^{n-n_1} \times \{0,1\}^{n_1}$, we have

$$\mathbf{E}[L] = 2^{-\ell} \cdot |W_{w+1}^{n-n_1} \times \{0,1\}^{n_1}| = \frac{1}{2} \cdot (n_1+1) \cdot \binom{n-n_1}{\downarrow w+1},$$

as $\ell = n_1 + 1$. This is negligible compared to T_b , given that $2d \cdot \log n \cdot 2^{n_1} \gg \frac{1}{2}(n+1)$.

Our next goal is to bound N_s .

Proposition A.7. $N_s \leq \frac{3}{2} \cdot 2^{n-2n_1}$.

Proof. We will show $E[N_s | \mathcal{U}] \leq \frac{3}{2} \cdot 2^{n-2n_1}$. If \mathcal{U} does not hold, then we abort the algorithm and it is clear that the proposition holds.

Fix a pair of iterations $0 \leq i < j < 4$. Conditioned on \mathcal{U} , $\tilde{E}^{(i)}$ and $\tilde{E}^{(j)}$ are independent uniform equation systems with a pre-fixed solution $\hat{x}^* = (\hat{y}^*, \hat{z}^*)$.

Fix $\hat{y} \neq \hat{y}^*$. Denote the restricted systems $\tilde{E}^{(i)}$ and $\tilde{E}^{(j)}$, where $y = \hat{y}$ for all polynomials by $\tilde{E}^i_{\hat{y}}$ and $\tilde{E}^j_{\hat{y}}$, respectively.

Claim (1). Conditioned on \mathcal{U} , the systems $\tilde{E}^i_{\hat{y}}$ and $\tilde{E}^j_{\hat{y}}$, are completely uniform and independent equations system with parameters (n_1, ℓ, d) .

Proof. Conditioned on \mathcal{U} , the 2ℓ polynomials in $\tilde{E}^{(i)} \cup \tilde{E}^{(j)}$ are uniform (and independent) degree d polynomials with a pre-fixed solution $\hat{x}^* = (\hat{y}^*, \hat{z}^*)$. It remains to show that each of these polynomials restricted to $\hat{y} \neq \hat{y}^*$ is a uniform polynomial of degree d.

Let $P(y,z) \in \tilde{E}^{(i)} \cup \tilde{E}^{(j)}$ be a polynomial of degree d and assume first that $\hat{y} = 0$. Suppose we fix all the $\binom{n_1}{\downarrow d}$ coefficients of the polynomial P(0,z). This leaves $P(\hat{y}^*, \hat{z}^*)$ undefined, as it computed by a sum mod 2 of uniform coefficients that are not fixed by P(0,z) (e.g., if $\hat{y}_1^* = 1$, then the computation of $P(\hat{y}^*, \hat{z}^*)$ includes the coefficient of the singleton monomial y_1). In other words, the coefficients of P(0,z) are independent of $P(\hat{y}^*, \hat{z}^*)$ (and the other polynomials in $\tilde{E}^{(i)} \cup \tilde{E}^{(j)}$). Consequently, if P(y,z) is a uniform degree d polynomial with a pre-fixed solution (\hat{y}^*, \hat{z}^*) such that $\hat{y}^* \neq 0$, then P(0,z) is a uniform degree d polynomial.

If $\hat{y} \neq 0$, then suppose we select a polynomial P'(x) as a uniform degree d polynomial with pre-fixed solution $(\hat{y}^* + \hat{y}, \hat{z}^*)$. As claimed above, P'(0, z) is a uniform degree d polynomial. Then, define $P(y, z) = P'(y + \hat{y}, z)$. By Proposition A.2, P(y, z) follows the desired distribution as $P(\hat{y}^*, \hat{z}^*) = P'(\hat{y}^* + \hat{y}, \hat{z}^*) = 0$. Moreover, $P(\hat{y}, z) = P'(0, z)$ is a uniform degree d polynomial. This concludes the proof.

We denote the event that $\tilde{E}^{(i)}$ and $\tilde{E}^{(j)}$ suggest the same solution (\hat{y}, \hat{z}) for some $\hat{z} \in \{0, 1\}^{n_1}$ by \mathcal{E}_1 , and calculate $\Pr[\mathcal{E}_1 \mid \mathcal{U}]$.

Note that Algorithm 1 actually tests a candidate solution (\hat{y}, \hat{z}) if the sum mod 2 of all of the solutions suggested by two iterations i,j with $y = \hat{y}$ is equal to \hat{z} , and the number of solutions to each of $\tilde{E}^{(i)}$ and $\tilde{E}^{(j)}$ is odd.

We define two additional events: \mathcal{E}_2 is the event that $\tilde{E}^{(i)}$ and $\tilde{E}^{(j)}$ have a solution in the space $S = \hat{y} \times \{0, 1\}^{n_1}$. \mathcal{E}_3 is the event the sum mod 2 of all solution to $\tilde{E}^{(i)}$ and of $\tilde{E}^{(j)}$ in S is equal to some $\hat{z} \in \{0, 1\}^{n_1}$ (and thus it may be tested by the algorithm).

Claim (2). $\Pr[\mathcal{E}_2 \mid \mathcal{U}] \leq \frac{1}{4}$.

Proof. Fix some $\hat{y} \in \{0,1\}^{n-n_1}$ such that $\hat{y} \neq \hat{y}^*$ is non-zero and fix any iteration *i*. We bound the probability that the system $\tilde{E}_{\hat{y}}^{(i)}$ has a solution. By Proposition A.4 applied to $S = \hat{y} \times \{0,1\}^{n_1}$, the probability is upper bounded by $2^{n_1-\ell} = \frac{1}{2}$. Similarly, fixing a pair of iterations $0 \leq i < j < 4$ (assuming \mathcal{U}), by Claim (1) about the independence of the systems $\tilde{E}_{\hat{y}}^{(i)}$ and $\tilde{E}_{\hat{y}}^{(j)}$, we use Proposition A.3 (which actually holds under the more restricted setting of uniform systems with a pre-fixed solutions) for each system independently and conclude $\Pr[\mathcal{E}_2 \mid \mathcal{U}] \leq \frac{1}{4}$ as required.

Claim (3). $\Pr[\mathcal{E}_3 \mid \mathcal{E}_2, \mathcal{U}] \leq 2^{-n_1}$.

Proof. By Claim (1), it is sufficient to show that for any uniform system \tilde{E} with parameters (n_1, ℓ, d) that has an odd number of solutions, the sum mod 2 of solutions is uniformly distributed in $\{0, 1\}^{n_1}$.

We do this by re-randomizing the system \hat{E} , to obtain another system \hat{E}' with parameters (n_1, ℓ, d) that has the same probability of being generated by the distribution, but where the sum mod 2 of all of the solutions is set to any fixed $b \in \{0, 1\}^{n_1}$.

Assume that in \tilde{E} the mod 2 sum of solutions is $a \in \{0, 1\}^{n_1}$. Let $b \in \{0, 1\}^{n_1}$ be arbitrary and perform a change of variables: for each $i \in \{1, \ldots, n_1\}$, replace each occurrence of z_i in all polynomials of \tilde{E} by $z_i + a_i + b_i$. This results in a new system \tilde{E}' with the same parameters in which every solution \hat{z} to \tilde{E} is transformed to a solution $\hat{z} + a + b$ to \tilde{E}' . Since the number of solutions to \tilde{E} is odd and their sum mod 2 is a, then the sum mod 2 of solutions to \tilde{E}' is a + a + b = b. Moreover, \tilde{E} and \tilde{E}' have the same probability of being generated. \Box

Finally, we have $\Pr[\mathcal{E}_1 \mid \mathcal{U}] = \Pr[\mathcal{E}_2 \mid \mathcal{U}] \cdot \Pr[\mathcal{E}_3 \mid \mathcal{E}_2, \mathcal{U}] \leq \frac{1}{4} \cdot 2^{-n_1}$. Recall that these events are defined for a fixed $\hat{y} \neq \hat{y}^*$ and iteration pair (i, j). We define an indicator random variable for each such pair (of which we have at most 6) and for each $\hat{y} \in \{0, 1\}^{n-n_1}$ such that $\hat{y} \neq \hat{y}^*$, and sum their expectations (i.e., probabilities). We conclude that $\mathbb{E}[N_s \mid \mathcal{U}] \leq \frac{3}{2} \cdot 2^{n-2n_1}$.

Proposition A.8. $\Pr[\mathcal{U}] \ge 1 - 12 \cdot 2^{-m+2\ell}$

Proof. Fix an iteration pair (i, j). The probability that the 2ℓ rows of the matrices $A^{(i)}$ and $A^{(j)}$ are linearly independent is lower bounded by the probability that they are linearly independent if they are selected uniformly at random (without the rank conditions on $A^{(i)}$ and $A^{(j)}$). The latter probability is

$$(1-2^{-m})(1-2^{-m+1})\dots(1-2^{-m+2\ell-1}) > e^{-2^{-m+1}-2^{-m+2\ell}-\dots-2^{-m+2\ell}} > e^{-2^{-m+2\ell+1}} > 1 - 2^{-m+2\ell+1}$$

where we have used the inequalities $e^{-t} < 1 - \frac{t}{2}$ and $1 - t < e^{-t}$, which hold for all 0 < t < 1.

The result follows by a union bound over all $\binom{4}{2} = 6$ iteration pairs.

Theorem A.1. Assume that the input to (modified) Algorithm 1 is a uniform equation system with pre-fixed solution. Then, it correctly returns this solution with probability at least

$$\frac{11}{16} - 12 \cdot 2^{2\ell - m} - 2^{n_1 - m} = \frac{11}{16} - 48 \cdot 2^{2 \cdot n_1 - m} - 2^{n_1 - m}.$$

In particular, if $m \ge 2 \cdot n_1 + 10$, its success probability is at least $\frac{5}{8}$. Moreover, its expected time complexity satisfies

$$T \lesssim 4 \cdot (2d \cdot \log n \cdot 2^{n_1} \cdot \binom{n-n_1}{\downarrow w+1} + n_1 \cdot n \cdot 2^{n-n_1}) + \frac{3}{2} \cdot 2^{n-2n_1} \cdot T_s$$

(where T_s is the average time to test a solution).

Proof. We assume \mathcal{E} (analyzed in Proposition A.5) and \mathcal{U} and (analyzed in Proposition A.8) occur. These two conditions hold, except with probability at most $2^{n_1-m} + 12 \cdot 2^{-m+2\ell}$.

Then (conditioned on \mathcal{E}), by Proposition 3.1, every iteration isolates the prefixed solution with probability at least $\frac{1}{2}$. The algorithm returns a solution once it is isolated twice. Thus, the probability it succeeds after at most 4 iterations is the sum of probabilities is succeeds after exactly 2, 3, and 4 iterations:

$$\frac{1}{4} + 2 \cdot \frac{1}{8} + 3 \cdot \frac{1}{16} - 2^{n_1 - m} - 12 \cdot 2^{-m + 2\ell} = \frac{11}{16} - 12 \cdot 2^{2\ell - m} - 2^{n_1 - m}.$$

By (18) and based on Proposition A.6 and Proposition A.7,

$$T \leq 4 \cdot T_1 + N_s \cdot T_s \lessapprox$$
$$4 \cdot (2d \cdot \log n \cdot 2^{n_1} \cdot {\binom{n-n_1}{\downarrow w+1}} + n_1 \cdot n \cdot 2^{n-n_1}) + \frac{3}{2} \cdot 2^{n-2n_1} \cdot T_s.$$

Remark A.1. In case $m \ge 4\ell = 4 \cdot (n_1 + 1)$, we can modify the algorithm to select an arbitrary subset of equations of E in each iteration $0 \le i < 4$ to form $\tilde{E}^{(i)}$ such that the 4 selected subsets are non-intersecting (and the analyzed equation systems are completely independent). In this case, Theorem A.1 continues to hold with a slightly better success probability, as $\Pr[\mathcal{U}] = 1$.

B Reducing the Complexity of Testing Candidate Solutions

We describe how to optimize the complexity of testing candidate solutions in Algorithm 1. The expected complexity of naively testing a solution by evaluating it on a polynomial equation is about $\binom{n}{\downarrow d}$ (up to a small constant).

We can optimize the amortized complexity by testing together batches of candidate solutions that share the same value of the most significant bits (MSBs). The evaluation of the equations in E is done via a Horner-like method for multivariate polynomials (by recursively writing polynomials in E as $P_j(x_1, \ldots, x_n) = x_1 \cdot P_j^{(1)}(x_2, \ldots, x_n) + P_j^{(2)}(x_2, \ldots, x_n))$). As an example, for d = 4 and n = 128, a choice of $n_1 = 12$ optimizes the time complexity (as shown in Table 1). Recalling from (12) that we test a total of about $N_s = 2^{n-2n_1} = 2^{104}$ candidate solutions, if we test batches that share the 90 MSBs of y on 16 equations of E (after which a very small fraction of candidates remain), the total testing complexity becomes about $2^{90} \cdot 16 \cdot {\binom{128}{\downarrow 4}} + {\binom{128-90}{\downarrow 4}} \cdot 2^{104} \approx 2^{122}$. This is negligible compared to the total complexity which is about 2^{129} .

C Description of Keccak

We briefly describe the sponge construction and the Keccak hash function. More details can be found in the Keccak specification [5]. The sponge construction [6] works on a state of b bits, split into two parts: the first part contains the first r bits of the state (called the outer part) and the second part contains the last c = b - r bits of the state (called the inner part).

Given a message, it is first padded and cut into r-bit blocks, and the b state bits are initialized to zero. The sponge construction then processes the message in two phases: In the absorbing phase, the message blocks are processed iteratively by XORing each block into the first r bits of the current state, and then applying a fixed permutation on the value of the b-bit state. After processing all the blocks, the sponge construction switches to the squeezing phase. In this phase, k output bits are produced iteratively, where in each iteration the first r bits of the state are returned as output and the permutation is applied to the state.

The Keccak hash function uses multi-rate padding: given a message, it first appends a single 1 bit. Then, it appends the minimum number of 0 bits followed by a single 1 bit, such that the length of the result is a multiple of r. Thus, multi-rate padding appends at least 2 bits and at most r + 1 bits.

We focus on Keccak versions with b = 1600 and c = 2k, where $k \in \{224, 256, 384, 512\}$. The 1600-bit state can be viewed as a 3-dimensional array of bits, a[5][5][64], and each state bit is associated with 3 integer coordinates, a[x][y][z], where x and y are taken modulo 5, and z is taken modulo 64.

The Keccak permutation consists of 24 rounds, which operate on the 1600 state bits. Keccak uses the following naming conventions, which are helpful in describing its round function:

- A row is a set of 5 bits with constant y and z coordinates, i.e. a[*][y][z], or r(y, z).
- A column is a set of 5 bits with constant x and z coordinates, i.e. a[x][*][z].
- A lane is a set of 64 bits with constant x and y coordinates, i.e. a[x][y][*].
- A slice is a set of 25 bits with a constant z coordinate, i.e. a[*][*][z].

Each round of the Keccak permutation consists of five mappings $R = \iota \circ \chi \circ \pi \circ \rho \circ \theta$. The five mappings given below are applied for each x,y, and z (where the state addition operations are over \mathbb{F}_2):

1. θ is a linear map, which adds to each bit in a column, the parity of two other columns.

$$\theta: a[x][y][z] \leftarrow a[x][y][z] + \sum_{y'=0}^{4} a[x-1][y'][z] + \sum_{y'=0}^{4} a[x+1][y'][z-1]$$

- ρ rotates the bits within each lane by T(x,y), which is a predefined constant for each lane.
 ρ: a[x][y][z] ← a[x][y][z + T(x,y)]
- 3. π reorders the lanes.

$$\pi: a[x][y][z] \leftarrow a[x'][y'][z], \text{ where } \begin{pmatrix} x \\ y \end{pmatrix} = \begin{pmatrix} 0 & 1 \\ 2 & 3 \end{pmatrix} \cdot \begin{pmatrix} x' \\ y' \end{pmatrix}$$

4. χ is the only non-linear mapping of Keccak, working on each of the 320 rows independently.

 $\chi: a[x][y][z] \leftarrow a[x][y][z] + ((\neg a[x+1][y][z]) \land a[x+2][y][z])$

Since χ works on each row independently, it can be viewed as an Sbox layer which simultaneously applies the same 5 bits to 5 bits Sbox to the 320 rows of the state. The Sbox function is an invertible mapping in with the algebraic degree of each output bit as a polynomial in the five input bits is 2.

5. ι adds a 64-bit round constant to the first lane of the state. ι : a[0][0][*] \leftarrow a[0][0][*]+ $RC[i_r]$