A SAT-based approach for index calculus on binary elliptic curves

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Abstract. Logical cryptanalysis, first introduced by Massacci in 2000, is a viable alternative to common algebraic cryptanalysis techniques over boolean fields. With XOR operations being at the core of many cryptographic problems, recent research in this area has focused on handling XOR clauses efficiently. In this paper, we investigate solving the point decomposition step of the index calculus method for prime degree extension fields \mathbb{F}_{2^n} , using SAT solving methods. We propose an original XOR-reasoning SAT solver, named WDSAT, dedicated to this specific problem. While asymptotically solving the point decomposition problem with our method has exponential worst time complexity in the dimension l of the vector space defining the factor base, experimental running times show that our solver is significantly faster than current algebraic methods based on Gröbner basis computation. For the values l and n considered in the experiments, WDSAT was up to 300 times faster then MAGMA's F4 implementation, and this factor grows with l and n. Our solver outperforms as well current best state-of-the-art SAT solvers for this specific problem.

Keywords: discrete logarithm, index calculus, elliptic curves, point decomposition, symmetry, satisfiability, DPLL algorithm

1 Introduction

The index calculus algorithm originally denoted a technique to compute discrete logarithms modulo a prime number, but it now refers to a whole family of algorithms adapted to other finite fields and to some algebraic curves. It includes the Number Field Sieve (NFS) [24], dedicated to logarithms in \mathbb{Z}_p and the algorithms of Gaudry [15] and Diem [7] for algebraic curves in \mathbb{F}_{p^n} . Index calculus algorithms proceed in two main steps. The *sieving* (or *point decomposition*) step concentrates most of the number theory and algebraic geometry needed overall. By splitting random elements over a well-chosen factor base, it produces a large sparse matrix, the rows of which are "relations". In a second phase, the *matrix step* produces "good" combinations of the relations by finding a non-trivial vector in the kernel of this matrix. This in turn enables the efficient computation of any discrete logarithm on the input domain. A crucial step of the index calculus on elliptic curves is to solve the *point decomposition problem* (PDP), by generating sufficiently many relations among suitable points on the curve. Using the

so-called summation polynomials attached to the curve, this boils down to solving a system of polynomial equations whose solutions are the coordinates of points. The resulting algorithm has complexity $O(p^{2-2/n})$, but this hides an exponential factor in n which comes from the hardness of solving the point decomposition problem.

Consequently, when p is large and n is small, the Gaudry-Diem algorithm has a better asymptotic complexity then generic methods for solving the discrete logarithm problem and Gröbner basis algorithms have become a well-established technique [19] to solve these systems. Since a large number of instances of PDP needs to be solved, most of the research in the area has focused on improving the complexity of this step. It was noticed that there are two degrees of freedom when we model the problem the number of points in a relation and the size of the factor base. The hardness of the problem for the Gröbner approach is determined by the choice of these parameters, and also by the choice of an equation for the elliptic curve. Several simplifications such as symmetries and polynomials with lower degree obtained from the algebraic structure of the curve have been proposed [9].

When we consider elliptic curves defined over \mathbb{F}_{2^n} with n large, solving the PDP system via Gröbner basis quickly becomes a bottleneck, and index calculus algorithms are slower than generic attacks, from a theoretical and a practical point of view. Moreover, when n is prime, it is not known how to define the factor base in order to exploit all the symmetries coming from the algebraic structure of the curve, without increasing the number of variables when solving PDP [35]. Finally, note that for random systems, pure Gröbner basis algorithms are both theoretically and practically slower than simpler methods, typically exhaustive search [5,25], hybrid methods [2] and SAT solvers. It is thus natural that we turn our attention towards combinatorics tools to solve the PDP in characteristic 2.

Until recent years, SAT solvers have been proven to be a powerful tool in the cryptanalysis of symmetric schemes. They were successfully used for attacking secret key cryptosystems such as Bivium, Trivium, Grain, AES etc. [16,22,17,33,32]. Their use in public key cryptosystems, although little explored, suggests that they may yield better performance than algebraic attacks in some cases. Recently, Galbraith and Gebregiyorgis [14] observed experimentally that generic SAT solvers (such as MINISAT) are more efficient than available Gröbner basis implementations, on CNF instances derived from the polynomial system for the PDP over binary curves.

In this paper, we take important steps towards fully replacing Gröbner basis techniques for solving PDP with constraint programming ones. First, we model the point decomposition problem as a logical formula, with a reduced number of clauses, when compared to the model used in [14]. Secondly, we propose a dedicated SAT solver for our model. Specifically, our solver is adapted for XOR-reasoning, which is one of the core operations in polynomial systems. We implement a DPLL-based backtracking procedure with three modules - a state-of-the-art unit propagation module and two XOR-reasoning modules. We experimented with the index calculus attack on the discrete logarithm for elliptic curves over prime degree binary extension fields. We obtain an important speedup in comparison with the best currently available implementation of Gröbner basis (F4 [10] in MAGMA [4]) and generic solvers [34,33]). In addition, our experiments show that Gröbner basis cannot compete with SAT solvers techniques in

terms of memory requirements. To illustrate, a system, which is solved with a SAT solver using between 60MB and 200MB of memory, requires 38GB when using the Gröbner basis method.

We estimate that our solver can be used for any attack that can be reduced to the problem of solving a multivariate boolean polynomial system. Notable examples are the HFE cryptosystem [26] and HFE-inspired schemes which are promising candidates for the NIST Post-Quantum Cryptography Standardization [27,13,1].

This paper is organized as follows. Section 2 gives an overview of the index calculus algorithm on elliptic curves, introduces the PDP problem and briefly recalls algebraic and combinatorial techniques used in the literature to solve this problem. Sections 3 and 4 detail the reasoning models used in our experiments and our SAT solver. In Section 5 we give worst time complexity estimates for solving a PDP instance and derive the complexity of our SAT-based index calculus algorithm. Finally, Section 6 presents benchmarks obtained with our implementation. We compare this against results obtained using MAGMA's F4 implementation and several available best generic SAT-solvers, such as MINISAT, CRYPTOMINISAT and GLUCOSE.

2 An Overview of Index Calculus

Given a finite group (G, +) and two elements $g, h \in G$, the discrete logarithm problem (DPL) consists in finding $x \in \mathbb{Z}$ such that $h = x \cdot g$. The security of a great number of asymmetric cryptosystems relies on this problem.

The index calculus method yields subexponential complexity for solving the DLP in the multiplicative group of a finite field. We briefly recall the three phases of this method, using the additive notation for the group law.

- 1. Choice of an appropriate factor base $\mathcal{B} = \{g_1, ..., g_n\}$, such that $\mathcal{B} \subseteq G$.
- 2. Decomposition phase: compute random integers a_i , b_i and try to decompose $[a_i]g + [b_i]h$ into the factor base. This is also called the *relation search phase*, since every successful decomposition of the form $[a_i]g + [b_i]h = \sum_{j=1}^n [c_{ij}]g_j$ is called a *relation*.
- 3. Linear algebra: when more than $k=\#\mathcal{B}$ linearly independent relations are found and the matrices $A=(a_ib_i)$ and $M=(c_{ij})$ are stored, use linear algebra to find a kernel vector $v=(v_1...v_k)$ of the matrix M. The discrete log of h can be solved by computing $x=-(\sum_i a_iv_i)/(\sum_i b_iv_i)$.

Performing this attack on elliptic curve groups is technically more difficult, notably because it is hard to find a factor base over which we can perform the decomposition phase efficiently. In 2008 and 2009, Gaudry [15] and Diem [7] independently proposed a technique to perform point decomposition for elliptic curves over extension fields, using Semaev's summation polynomials [29]. Since this paper focuses on binary elliptic curves, we introduce Semaev's summation polynomials here directly for these curves.

Let \mathbb{F}_{2^n} be a finite field and E be an elliptic curve defined by the equation

$$E: y^2 + xy = x^3 + ax^2 + b, (1)$$

with $a,b\in\mathbb{F}_{2^n}$. Using standard notation, we take $\bar{\mathbb{F}}_{2^n}$ to be the algebraic closure of \mathbb{F}_{2^n} and $E(\mathbb{F}_{2^n})$ (resp. $E(\bar{\mathbb{F}}_{2^n})$) to be the set of points on the elliptic curve defined over \mathbb{F}_{2^n} (resp. $\bar{\mathbb{F}}_{2^n}$). Let \mathcal{O} to be the point at infinity on the elliptic curve. For $m\in\mathbb{N}$, the m^{th} -summation polynomial is a multivariate polynomial in $\mathbb{F}_{2^n}[X_1,\ldots,X_m]$ with the property that, given $P_1,\ldots,P_m\in E(\bar{\mathbb{F}}_{2^n})$ points, then $P_1+\ldots+P_m=\mathcal{O}$ if and only if $S_m(\mathbf{x}_{P_1},\ldots,\mathbf{x}_{P_m})=0$. We have that

$$S_2(X_1,X_2) = X_1 + X_2, S_3(X_1,X_2,X_3) = X_1^2 X_2^2 + X_1^2 X_3^2 + X_1 X_2 X_3 + X_2^2 X_3^2 + b (2)$$

and for $m \ge 4$ we have the following recursive formula:

$$S_m(X_1, \dots, X_m) = Res_X(S_{m-k}(X_1, \dots, X_{m-k-1}, X), S_{k+2}(X_{m-k}, \dots, X_r, X))$$

The polynomial S_m is symmetric and has degree 2^{m-2} in each of the variables. Let V be a vector subspace of $\mathbb{F}_{2^n}/\mathbb{F}_2$, whose dimension l will be defined later. We define the factor basis \mathcal{B} to be :

$$\mathcal{B} = \{ (\mathbf{x}, \mathbf{y}) \in E(\mathbb{F}_{2^n}) | \mathbf{x} \in V \}.$$

Heuristically, we can easily see that the factor base has $O(2^l)$ elements.

Definition 1. Let V be a vector subspace of $\mathbb{F}_{2^n}/\mathbb{F}_2$. Given a point $R \in E(\mathbb{F}_{2^n})$, the point decomposition problem is to find m points $P_1, \ldots, P_m \in \mathcal{B}$ such that $R = P_1 + \ldots + P_m$.

Using Semaev's polynomials, this problem is reduced to the one of solving a multivariate polynomial system.

Definition 2. Given $s \ge 1$ and a l-dimensional vector subspace V of $\mathbb{F}_{2^n}/\mathbb{F}_2$ and $f \in \mathbb{F}_{2^n}[X_1,\ldots,X_m]$ any multivariate polynomial of degree bounded by s, find $(\mathbf{x}_1,\ldots,\mathbf{x}_m) \in V^m$ such that $f(\mathbf{x}_1,\ldots,\mathbf{x}_m)=0$.

Using the fact that \mathbb{F}_{2^n} is a n-dimensional vector space over \mathbb{F}_2 , the equation $f(\mathbf{x}_1,\ldots,\mathbf{x}_m)=0$ can be rewritten as a system of n equations over \mathbb{F}_2 , with ml variables. In the literature, this is called a *Weil restriction* [15] or *Weil descent* [28]. The probability of having a solution to this system depends on the ratio between n and l. Roughly, when $n/l \sim m$ the system has a reasonable chance to have a solution.

Recent work on solving the decomposition problem has focused on using advanced methods for Gröbner basis computation such as Faugère's F_4 and F_5 algorithms [10,11]. This is a natural approach, given that similar techniques for small degree extension fields in characteristic > 2 yielded index calculus algorithms which are faster than the generic attacks on the DLP.

2.1 Solving the decomposition problem using Gröbner basis

In this section, we give a brief account on the complexity of the index calculus algorithm obtained using the Gröbner approach. It is well known that the complexity of Gröbner basis algorithms based on Lazard's Gaussian elimination on Macaulay matrices [23]

strongly depends on the maximal degree D reached during the computation, called the degree of regularity of the system.

In 2012, Faugère et~al.[12] proposed a dedicated Gröbner basis algorithm based on techniques "à la Lazard" which solves PDP in time $O(2^{\omega\tau})$ and memory $O(2^{2\tau})$, where ω is the linear algebra constant and $\tau\approx n/2$. This running time is obtained under the unproven yet plausible heuristic assumption that a certain set of equations arising in the algorithm is linearly independent and also by establishing a certain bound on the degree of regularity. As a consequence, they derive an index calculus algorithm solving the DLP over an elliptic curve defined over a binary field \mathbb{F}_{2^n} having a worst time complexity $O(2^{\omega\tau+l})$.

Later on, experiments carried by Petit and Quisquater [28] using MAGMA's F_4 [4] implementation suggested that the degree of regularity of the system derived from a PDP instance via the Weil restriction is much smaller than the bound obtained in [12]. Going further, the authors of [28] introduce new heuristic assumptions on the bound on the degree of regularity of such a system and claim the existence of a subexponential index calculus algorithm under these assumptions. If correct, the result of Petit and Quisquater would be remarkable. Unfortunately, the experimental data available in the literature concerns all small values of l and n and several papers in this area [18,30] debate on the validity of the assumptions made in [28]. For this reason, we only compare ourselves to the results obtained in Faugère $et\ al$.

A common technique when working with Semaev's polynomials is to use a symmetrization process in order to further reduce the degree of the polynomials appearing in the PDP system. In short, since S_m is symmetric, we can rewrite it in terms of the elementary symmetric polynomials

$$e_{1} = \sum_{1 \leq i_{1} \leq m} X_{i_{1}},$$

$$e_{2} = \sum_{1 \leq i_{1}, i_{2} \leq m} X_{i_{1}} X_{i_{2}},$$

$$\cdots$$

$$e_{m} = \prod_{1 \leq i \leq m} X_{i}.$$
(4)

We denote S'_{m+1} the polynomial obtained after symmetrizing $S_{m+1}(X_1, X_2, \dots, X_m, X_{m+1})$ in the first m variables, i.e. we have

$$S'_{m+1} \in \mathbb{F}_{2^n}[e_1, \dots, e_m, X_{m+1}].$$

In [35], the authors report on experiments lead on systems obtained using a careful choice of the vector space V and application of the symmetrization process. Using Magma's F_4 available implementation, we experimented with both the symmetric and the non-symmetric version for PDP systems and found, as in [35], that the symmetric version yields better results. Therefore, in order to set the notation, we detail this approach here.

¹ In [12] the authors state the complexity is $O(2^{\omega \tau})$. In the proof of their Theorem 3, the exact time is $2^{m \log m + l + \omega \tau}$. Since l is large, we cannot ignore it here.

Let t be the root of the defining polynomial of \mathbb{F}_{2^n} over \mathbb{F}_2 . Following [35], we choose the vector space V to be the l-dimension subspace generated by $1, t, t^2, \ldots, t^{l-1}$. Therefore we can write:

$$e_{1} = d_{1,0} + \dots + d_{1,l-1}t^{l-1}$$

$$e_{2} = d_{2,0} + \dots + d_{2,2l-2}t^{2l-2}$$

$$\dots$$

$$e_{m} = d_{m,0} + \dots + d_{m,m(l-1)}t^{m(l-1)},$$
(5)

where the $d_{i,j}$ with $1 \le i \le m, 0 \le j \le i(l-1)$ are binary variables. After choosing $\mathbf{x}_{m+1} \in \mathbb{F}_{2^n}$ and substituting e_1, \ldots, e_m as in Equation (5), we get:

$$S'_{m+1}(e_1,\ldots,e_m,\mathbf{x}_{m+1})=f_0+\ldots+f_{n-1}t^{n-1},$$

where f_i , $0 \le i \le n-1$ are polynomials in the binary variables $d_{i,j}$, $1 \le i \le m$, $0 \le j \le i(l-1)$. After a Weil descent, we obtain the following polynomial system

$$f_0 = f_1 = \dots = f_{n-1} = 0.$$
 (6)

One can see that with this approach, the number of variables is increased by a factor m, but the degrees of the polynomials in the system are seriously reduced. The system in Equation (6) is also the starting point for our SAT model of the point decomposition problem.

2.2 Solving the decomposition problem using SAT solvers

Before presenting our approach for finding solutions of the PDP using SAT solvers, we give preliminaries on the Satisfiability problem, its terminology and solving techniques.

Propositional variables can take two possible truth values: TRUE and FALSE. We denote a propositional variable by x.

- A *literal* is a signed propositional variable. Therefore, it can be positive (denoted by x) or negative (denoted by $\neg x$). A literal x (resp. $\neg x$) is *satisfied* if it is assigned to TRUE (resp. FALSE). A literal x (resp. $\neg x$) is *falsified* if it is assigned to FALSE (resp. TRUE);
- An OR-clause is a non-exclusive disjunction (\vee) of literals (e.g. $x_1 \vee \neg x_2 \vee x_3$). An OR-clause is said to be falsified if all of its literals are falsified and it is set to be satisfied if at least one of its literals is satisfied;
- A XOR-clause is an exclusive disjunction (⊕) of literals. (e.g. x₁ ⊕ ¬x₂ ⊕ x₃). A XOR-clause is said to be satisfied (resp. falsified) if an even (resp. odd) number of its literals is satisfied;
- An *interpretation* of a given propositional formula consists in assigning a truth value to a set of its variables;
- A CNF formula is a conjunction (∧) of one or more OR-clauses. A CNF formula is said to be *satisfiable* if there exists at least one interpretation which satisfies all of its OR-clauses, and it is said to be *unsatisfiable* when the opposite is true. Every formula in propositional logic has a closed-CNF form.

In the remainder of this paper, we will refer to an OR-clause simply by a clause, since CNF is the standard form used in SAT solvers. A clause where the operation between literals is an exclusive OR, will be referred to as a XOR-clause.

Propositional satisfiability or SAT is the problem of determining whether a (usually CNF) formula is satisfiable. A SAT solver is a special purpose program to solve the SAT problem. The most straightforward method for solving the SAT problem is to complete the truth table associated to the formula in question. This is equivalent to an exhaustive search method and thus impractical. Luckily, in some cases a *partial* assignment on the set of variables can determine whether a clause is satisfiable. Assigning l, a literal from the partial assignment, to TRUE will lead to:

- 1. Every clause containing l is removed (since the clause is satisfied).
- 2. In every clause that contains $\neg l$ this literal is deleted (since it can not contribute to the clause being satisfied).

The second rule above can lead to obtaining a clause composed of a single literal, called a *unit* clause. Since this is the only literal left which can satisfy the clause, it must be set it to TRUE. A new truth value is *propagated* by the unit clause. Hence, the described method is called *unit propagation*. The reader can refer to [3] for more details.

In the case where the unit clause literal is already assigned to the opposite truth value, we consider that we have encountered a *conflict*. The last partial assignment that lead to this conflict has to be undone. This is commonly known are *backtracking*.

Example 1. For instance, these two atomic operations can be illustrated thanks to the following sample built on a set of 10 clauses numbered C_1 to C_{10} .

$$\begin{array}{ccccc} C_1: x_1 \vee x_2 \vee x_3 & C_2: \neg x_3 \vee x_4 \vee \neg x_5 & C_3: \neg x_3 \vee x_1 \vee \neg x_4 \\ C_4: \neg x_2 \vee x_5 \vee x_6 & C_5: \neg x_6 \vee \neg x_4 \vee x_5 & C_6: x_1 \vee \neg x_4 \vee x_6 \\ C_7: x_3 \vee x_4 \vee \neg x_5 & C_8: x_1 \vee \neg x_2 \vee \neg x_4 & C_9: x_4 \vee \neg x_5 \vee \neg x_6 \\ & C_{10}: \neg x_2 \vee x_3 \vee x_4 & \end{array}$$

Assigning the variable x_5 to TRUE leads the clauses C_4 and C_5 to be satisfied by the literal x_5 . As well and as a consequence, clauses C_2 , C_7 and C_9 cannot be satisfied thanks to the literal x_5 . Hence, $\neg x_5$ can be deleted from these clauses. At this step, our formula is:

$$\begin{array}{ccccccc} C_1: x_1 \vee x_2 \vee x_3 & C_2: \neg x_3 \vee x_4 & C_3: \neg x_3 \vee x_1 \vee \neg x_4 \\ C_4: & C_5: & C_6: x_1 \vee \neg x_4 \vee x_6 \\ C_7: x_3 \vee x_4 & C_8: x_1 \vee \neg x_2 \vee \neg x_4 & C_9: x_4 \vee \neg x_6 \\ & C_{10}: \neg x_2 \vee x_3 \vee x_4 & \end{array}$$

Using the same reasoning, assigning the variable x_3 to TRUE leads the clauses C_1 , C_7 and C_{10} to be satisfied and the literal $\neg x_3$ to be deleted from the clauses C_2 and C_3 . At this step, our formula is :

Then, C_2 is a unit clause composed of the literal x_4 and as a consequence, x_4 has to be assigned to TRUE. We say that the truth value of x_4 is inferred through unit propagation.

The basic backtracking search with unit propagation that we described composes the Davis-Putnam-Logemann-Loveland (DPLL) algorithm [6], which is a state-of-the-art complete SAT solving technique. DPLL works by trying to assign a truth value to each variable in the CNF formula, recursively building a binary search tree of height equivalent (at worst) to the number of elementary variables. After each variable assignment, the formula is simplified by unit propagation. If a *conflict* is met, a backtracking procedure is launched and the opposite truth value is assigned to the last assigned literal. If the opposite truth value results in conflict as well, we backtrack to an earlier assumption or conclude that the formula is *unsatisfiable* - when there are no earlier assumptions left. The number of conflicts is a good measure for the time complexity of a SAT problem solved using a DPLL-based solver. If the complete search tree is built, the worst case complexity is $O(c2^v)$, where c is the number of clauses in the formula and v is the number of variables.

A common variation of the DPLL is the conflict-driven clause learning (CDCL) algorithm [31] which consists in deriving a new clause each time a conflict is met. Both methods are *complete* methods, which means they are guaranteed to give an answer. State-of-the-art CDCL solvers, such as MINISAT and GLUCOSE, have been shown to be a powerful tool for solving CNF formulas. However, they are not equipped to handle XOR-clauses and thus parity constraints have to be translated into CNF. Since handing CNF-clauses derived from XOR constraints is not necessarily efficient, recent works have concentrated on coupling CDCL solvers with a XOR-reasoning module. Furthermore, these techniques can be enhanced by Gaussian elimination, as in the works of Soos *et al.* (resulting in the CRYPTOMINISAT solver) [33,32], Han and Jiang [17], Laitinen *et al.*[22,21]. The CRYPTOMINISAT solver, specifically designed for exploiting the XOR operation in cryptographic problems, offers the possibility either to perform only top-level Gaussian elimination or to perform this operation dynamically during the CDCL process. One has to choose the best strategy depending on the problem at hand.

3 Model description

This section gives in full detail the three models we used in our experiments: the algebraic one used by Yun-Ju *et al* [35], the CNF model used by Galbraith and Gebregiyorgis [14] and the model we propose.

3.1 The algebraic model

Since the logical models are constructed starting from the algebraic one, we present first the model used when solving the PDP problem using Gröbner basis. The elementary symmetric polynomials e_i are written in terms of the $d_{i,j}$ binary variables, as in Equation (5). Similarly, since we look for a set of solutions $(\mathbf{x}_1,\ldots,\mathbf{x}_m)\in V^m$, the X_i variables are written formally as follows:

$$X_{1} = c_{1,0} + \dots + c_{1,l-1}t^{l-1}$$

$$X_{2} = c_{2,0} + \dots + c_{2,l-1}t^{l-1}$$

$$\dots$$

$$X_{m} = c_{m,0} + \dots + c_{m,l-1}t^{l-1}$$

where $c_{i,j}$, with $1 \le i \le m$, $0 \le j \le l-1$, are binary variables. Using Equation (4), we derive the following equations:

$$d_{1,0} = c_{1,0} + \dots + c_{m,0}$$

$$d_{1,1} = c_{1,1} + \dots + c_{m,1}$$

$$\dots$$

$$d_{m,m(l-1)} = c_{1,l} \cdot \dots \cdot c_{m,l}.$$
(7)

The remaining equations correspond to polynomials f_i , $0 \le i \le n-1$, obtained in Equation (5) via the Weil descent on S'_{m+1} . Recall that these are polynomials in the binary variables $d_{i,j}$. We now describe how we derive logical formulas from this system.

3.2 The XOR model

When creating constraints from a boolean polynomial system, the multiplication of variables becomes a conjunction of literals and the sum of multiple terms becomes a XOR-clause. From the two sets of equations in the algebraic model, we obtain two sets of XOR-clauses, where the terms are single literals or conjunctions. To illustrate, the logical formula derived from Equation (7) is as follows:

$$\neg d_{1,0} \oplus c_{1,0} \oplus \ldots \oplus c_{m,0}
\neg d_{1,1} \oplus c_{1,1} \oplus \ldots \oplus c_{m,1}
\dots
\neg d_{m,m(l-1)} \oplus (c_{1,l} \wedge \ldots \wedge c_{m,l}).$$
(8)

SAT solvers adapted for XOR reasoning in the literature contain XOR clauses formed by xoring single literals, and not conjunctions of several ones. Since our solver follows this paradigm, we have to transform the system above further. We substitute all conjunctions in a XOR clause by a newly added variable. Let c^\prime be the variable substituting

a conjunction $(c_{i_1,j_1} \wedge c_{i_2,j_2} \wedge ... \wedge c_{i_k,j_k})$. We have $c' \Leftrightarrow (c_{i_1,j_1} \wedge c_{i_2,j_2} \wedge ... \wedge c_{i_k,j_k})$, which rewrites as

$$(c' \vee \neg c_{i_1,j_1} \vee \neg c_{i_2,j_2} \vee \dots \vee \neg c_{i_k,j_k}) \wedge (\neg c' \vee c_{i_1,j_1}) \wedge (\neg c' \vee c_{i_2,j_2}) \wedge \dots (\neg c' \vee c_{i_k,j_k}) \wedge$$

$$(9)$$

For clarity, variables introduced by substitution of monomials containing exclusively the variables $c_{i,j}$ will be denoted c' and clauses derived from these substitutions are said to be in the X-substitutions set of clauses. Similarly, substitutions of the monomials containing only the $d_{i,j}$ variables are denoted by d' and the resulting set is referred to as the E-substitutions set of clauses.

Note from Equation (9) that the number of clauses obtained by substitution of a k-degree monomial is k + 1. This will be further discussed in our complexity analysis.

After substituting conjunctions, we will refer to the set of clauses obtained from Equation (8) as the E-X-relation set of clauses. Finally, the equations corresponding to polynomials f_i , $0 \le i \le n-1$, are derived in the same manner and the resulting clauses will be referred to as the F set of clauses.

That concludes the four sets of clauses in our SAT model. This model does not represent a CNF formula, since the E-X-relation set and the F set are made up of XOR-clauses. Hence, it will be referred to as the XOR model. The WDSAT solver is adapted for XOR reasoning and takes a XOR model as input.

Proposition 1. Assigning all $c_{i,j}$ variables, for $1 \le i \le m$ and $1 \le j \le l$, leads to the assignment of all variables in the XOR model through unit propagation.

Proof. Let us examine the unit propagation process for each set of clauses separately.

- 1. Clauses in the X-substitutions set are obtained by transforming $c' \Leftrightarrow (c_{i_1,j_1} \land c_{i_2,j_2} \land ... \land c_{i_k,j_k})$. We note that on the right of these equivalences there are only $c_{i,j}$ variables and on the left there is one single c' variable. The assignment of all of the $c_{i,j}$ variables will yield the assignment of all variables on the left of the equivalences, i.e. all c' variables.
- 2. Clauses in the E-X-relations set are obtained by transforming the algebraic system in (7). We observe that on the right of the equations there are only $c_{i,j}$ and c' variables and on the left there is one single $d_{i,j}$ variable. When all $c_{i,j}$ and all c' variables are assigned, all $d_{i,j}$ variables will have their truth value assigned through unit propagation on the E-X-relation set.
- 3. Clauses in the E-substitutions set are obtained by transforming $d' \Leftrightarrow (d_{i_1,j_1} \land d_{i_2,j_2} \land ... \land d_{i_k,j_k})$. Similarly as with the X-substitutions set, we have only $d_{i,j}$ variables on the right of these equivalences and one single d' variable on the left. The assignment of all of the $d_{i,j}$ variables will thus yield the assignment of all d' variables.
- Finally, parity constraints in set F decide whether the obtained interpretation satisfies the formula.

Transformation of the polynomial S_{m+1}' to a reasoning model is certainly not the bottleneck of the PDP. It is however done once for each point decomposition, hence at least 2^l times in a full ECDLP computation. To optimize the transformation, we split the process in two phases. The first phase is done after choosing a factor base, and before starting point decomposition. At this point, we can do the whole Weil descent process except for the last step - multiplying all monomials with the coefficients of vector X_{m+1} . Three sets of clauses (X-substitutions, E-substitutions and E-X-relations) are created during this phase.

When a random point to decompose is chosen, we can do the last step and deduce the F set. Without giving too much detail, we implement the transformation in a way that the second phase is computationally faster at the expense of the first one.

3.3 The CNF model

Since most of the modern SAT solvers can read and process CNF formulas, we explain the classical technique for transforming a XOR model to a CNF model. In fact, this is also the technique used in MAGMA's available implementation for deriving a CNF model from a boolean polynomial system.

A XOR-clause is said to be satisfied when it evaluates to TRUE, i.e. when there are an odd number of literals set to TRUE. The CNF-encoding of a ternary XOR-clause $(x_1 \oplus x_2 \oplus x_3)$ is

$$(x_1 \lor \neg x_2 \lor \neg x_3) \land (\neg x_1 \lor x_2 \lor \neg x_3) \land (\neg x_1 \lor \neg x_2 \lor x_3) \land (x_1 \lor x_2 \lor x_3)$$

$$(10)$$

П

Similarly, a XOR-clause of size k can be transformed to a conjunction of 2^{k-1} OR-clauses of size k. Since the number of introduced clauses grows exponentially with the size of the XOR-clause, it is a good practice to cut up the XOR-clause into manageable size clauses before proceeding with the transformation. To cut a XOR-clause $(x_1 \oplus \ldots \oplus x_k)$ of size k in two, we introduce a new variable x' and we obtain the following two XOR-clauses:

$$(x_1 \oplus \ldots \oplus x_i \oplus \boldsymbol{x'}) \land (x_{i+1} \oplus \ldots \oplus x_k \oplus \neg \boldsymbol{x'}).$$

In our our experiments with MINISAT in Section 6, we used a CNF model obtained after cutting into ternary XOR-clauses, since any XORSAT problem reduces in polynomial time to a 3-XORSAT problem [3]. To the best of our knowledge, MAGMA's implementation adopts a size 5 for XOR clauses. Determining the optimal size at which to cut the XOR-clauses was out of scope for our work and we did not experiment with this parameter.

Table 1 serves as a comparison on the number of variables, equations and clauses between the three models described in this section. Values for the algebraic and XOR model are exact, whereas those for the CNF model are averages obtained from experiments presented in Section 6.

		Gröb	ner model	C	NF model	XOR model		
1	n	#Vars	#Equations	#Vars	#CNF-clauses	#Vars	#CNF-clauses	#XOR-clauses
	13	42	40	2474	9552	502	1505	40
5	15	42	42	2687	10409	502	1505	42
	17	42	44	2941	11433	502	1505	44
	17	51	50	4686	18237	767	2364	50
6	18	51	51	4752	18505	767	2364	51
	19	51	52	5019	19577	767	2364	52
	19	60	58	6981	27216	1101	3466	58
7	21	60	60	7566	29565	1101	3466	60
L	23	60	62	8223	32201	1101	3466	62
	23	69	68	11036	43210	1510	4835	68
8	24	69	69	11318	44339	1510	4835	69
	26	69	71	12074	47374	1510	4835	71
	27	78	78	16130	63325	2000	6495	78
	37	78	88	20969	82721	2000	6495	88
9	47	78	98	25456	100709	2000	6495	98
	59	78	110	31942	126702	2000	6495	110
	67	78	118	35917	142632	2000	6495	118
	30	87	87	22472	88397	2577	8470	87
	47	87	104	32866	130040	2577	8470	104
10	59	87	116	40203	159437	2577	8470	116
	67	87	124	45394	180232	2577	8470	124
	79	87	136	52510	208743	2577	8470	136
	33	96	96	29700	116976	3247	10784	96
	59	96	122	49538	196434	3247	10784	122
11	67	96	130	55310	219553	3247	10784	130
	79	96	142	63531	252485	3247	10784	142
	89	96	152	71556	284626	3247	10784	152

Table 1: Number of variables and equations/clauses in three models.

In 2014, Galbraith and Gebregiyorgis [14] used MAGMA's implementation to compute the equivalent CNF logical formulas of the polynomial system resulting from the Weil descent of a PDP system and ran experiments using the general-purpose MINISAT solver to get solutions for these formulas. One can see from Table 1 that the model they used has a significantly larger number of clauses and variables, when compared to the XOR model. This motivated our choice of the XOR model for this work.

4 A dedicated SAT solver for solving PDP

In this Section, we present a built-from-scratch SAT solver, named WDSAT, dedicated to solving the model described in Section 3. We recall from Proposition 1 that assigning all $c_{i,j}$ variables in the XOR model leads to the assignment of all variables through unit propagation. Following the DPLL algorithm explained in Section 2, we construct a binary search tree by trying to recursively assign a truth value to each $c_{i,j}$ variable. The solver implements three reasoning modules - one module for the CNF-part and two for XOR-reasoning. They are briefly detailed below.

- CNF module This module, dedicated to reasoning from the CNF-part of the model, contains structures designed for fast unit propagation on CNF-clauses.
- XORSET module This module performs unit propagation on the parity constraints.
 When all except one literal in a XOR clause is assigned, we infer the truth value of the last literal according to parity reasoning.
- XORGAUSS module The second XOR-reasoning module performs Gaussian elimination on the XOR system dynamically once before starting the solving process and then on each level of the binary search tree. During the initialization process, XOR-clauses are normalized and represented as equivalence classes. A clause in normal form contains only positive literals and does not contain more than one occurrence of the same variable. Presenting the XOR-clauses simply as equivalence classes allows us to avoid performing matrix operations, that can be computationally intensive.

All modules implement a SET_IN function which takes as input a list of literals and a propositional formula F. It consists in setting all literals from the list to TRUE and in consequence, simplifying F and inferring truth values for other literals through the appropriate unit propagation technique. This function sends a conflict signal when a conflict is reached. In addition, all modules have a LAST_ASSIGNED function which returns the list of literals that were assigned during the last call to the respective SET_IN function.

When an assumption of a truth value for a literal x is made, the solver initially tries to set it in the CNF module. If the literal x is set successfully (it does not result in conflict), it will infer truth values for other literals. All of the inferred, as well as the initial literal are set in the XORSET module. The new list of XOR-implied literals is then set in the CNF module. We do this process back and forth until there are no more new implications.

When the process stops, the list of all inferred literals since the beginning of the process, as well as the initial x, are transferred to the XORGAUSS module. If the XORGAUSS module finds new XOR-implied literals, the list is sent to the CNF module and the process is restarted. When all modules are stabilized and there are no more inferred literals to set, we go one level further in the search tree and a new assumption is made.

If a conflict is reached in any of the reasoning modules, the process is stopped and a backtracking procedure is launched.

This concludes the entire process of assigning a truth value which is presented as the ASSIGN function detailed in Algorithm 1. In the algorithm, *to_set* is a list containing

literals to be set in the CNF and XORSET modules, whereas $to_set_in_XORGAUSS$ is a list containing literals to be set in the XORGAUSS module.

Algorithm 1 Function ASSIGN(F, x): Assigning a truth value to a literal x in a formula F, simplifying F and inferring truth values for other literals.

Input: Propositional formula F, Literal x

 ${f Output}$: FALSE if a conflict is reached, TRUE and F simplified thanks to unit propagation within the three modules otherwise.

```
1: to\_set \leftarrow \{x\}.
2: to\_set\_in\_XORGAUSS \leftarrow \{x\}.
3: while to\_set \neq \emptyset do
4:
         while to\_set \neq \emptyset do
5:
             if SET_IN_CNF (to\_set, F) \rightarrow \text{conflict then}
6:
                 return (FALSE, -).
7:
             end if
8:
             to\_set \leftarrow LAST\_ASSIGNED\_IN\_CNF().
9:
             to\_set\_in\_XORGAUSS \leftarrow to\_set.
10:
             if SET_IN_XORSET(to\_set, F)\rightarrowconflict then
11:
                  return (FALSE, -).
12:
             end if
13:
             to\_set \leftarrow \texttt{LAST\_ASSIGNED\_IN\_XORSET}().
14.
             to\_set\_in\_XORGAUSS \leftarrow to\_set \cup to\_set\_in\_XORGAUSS.
15:
         end while
16:
         if SET_IN_XORGAUSS(to\_set\_in\_XORGAUSS, F)\rightarrowconflict then
17:
             return (FALSE, -).
18:
19:
         to\_set \leftarrow LAST\_ASSIGNED\_XORGAUSS().
20: end while
21: return (TRUE, F).
```

The ASSIGN function is called by a recursive SOLVE function which is at the core of the WDSAT solver and is detailed in Algorithm 2. On line 4 of Algorithm 1, we arbitrarily choose one of the $c_{i,j}$ variables. At first, we try to assign a truth value of FALSE to the chosen $c_{i,j}$ variable. If the formula can not be satisfied after this assumption, we apply the classic backtracking technique (function BACKTRACK) and the opposite truth value is set. If the formula can not be satisfied with a value of TRUE for $c_{i,j}$ either, we conclude that the formula is unsatisfiable.

Remark 1. The conflict-driven clause learning (CDCL) variation [31] of the DPLL algorithm has been shown to yield a significant performance improvement for a number of SAT problems. Indeed, it is at the core of many modern general-purpose SAT solvers. However, because of the nature of SAT models coming from a boolean polynomial system, we estimate that the cost of CDCL would outweigh its eventual benefit for this particular problem. Indeed, the only information that the CNF-part holds is the equivalence between a literal and a conjunction. A set of clauses of the form (9) is independent from other such sets. Consequently, WDSAT does not implement CDCL techniques. Compar-

Algorithm 2 Function SOLVE(F): Recursive function for solving a SAT formula derived from PDP.

```
Input: Propositional formula F
Output: TRUE if formula is satisfiable, FALSE otherwise.
 1: if all clauses and all XOR-clauses are satisfied then
2:
        return TRUE.
3: end if
4: choose one c_{i,j}.
 5: (contradiction, F') \leftarrow ASSIGN(F, \neg c_{i,j}).
 6: if contradiction then
        BACKTRACK().
8: else
        if SOLVE(F') returns FALSE then
9:
10:
            BACKTRACK().
11:
        else
12:
           return TRUE.
13:
        end if
14: end if
15: (contradiction, F') \leftarrow ASSIGN(F, c_{i,j}).
16: if contradiction then
17:
        BACKTRACK().
        return FALSE.
18:
19: end if
20: return SOLVE(F').
```

isons of running times obtained with WDSAT to running times for CDCL based solvers, in Section 6, confirm this assumption.

4.1 Breaking symmetry

From the symmetry of Semaev's summation polynomials we have that when $\{\mathbf{x}_1,\ldots,\mathbf{x}_m\}$ is a solution, all permutations of this set are a solution as well. These solutions are equivalent and finding more that one is of no use for the PDP. We observe redundancy in the binary search tree. Indeed, for m=3 when a potential solution $\{\mathbf{x}_1,\mathbf{x}_2,\mathbf{x}_3\}$ has been eliminated, $\{\mathbf{x}_2,\mathbf{x}_1,\mathbf{x}_3\}$ does not need to be tried out. To avoid this redundancy, we establish the following constraint $\mathbf{x}_1 \leq \mathbf{x}_2 \leq \ldots \leq \mathbf{x}_m$.

It would be tedious to add this constraint in the model itself. Any solution implies adding clauses and weighing the SAT model. Instead, we decided to implement this constraint in the solver using a tree-pruning-like technique. We apply this technique on top of the recursive function SOLVE in Algorithm 2. In the function SOLVE we were trying out both FALSE and TRUE for the truth value of a chosen variable. In the breaking symmetry variation of SOLVE, denoted SOLVE_BR_SYM, in some cases the truth value of FALSE will not be tried out as all potential solutions after this assignment would not satisfy the constraint $\mathbf{x}_1 \leq \mathbf{x}_2 \leq \ldots \leq \mathbf{x}_m$. The new algorithm for the SOLVE_BR_SYM function is detailed in Algorithm 3 and the line numbers that distinguish it from Algorithm 2 are in bold. Note that one crucial difference between the two algorithms is the

choice of a variable on line 4. While this choice was arbitrary in Algorithm 2, in Algorithm 3 variables need to be chosen in the order from the leading bit of \mathbf{x}_1 to the trailing bit of \mathbf{x}_m . If this is not respected, SOLVE_BR_SYM does not yield a correct answer.

Continuing with the notation from Section 3, $c_{i,j}$ corresponds to the j^{th} bit of the i^{th} **x**-vector, where $2 \leq i \leq m$ and $1 \leq j \leq l$. In Algorithm 3, we decide whether to try out the truth value of FALSE for $c_{i,j}$ or not by comparing two **x**-vectors bit for bit, in the same way that we would compare binary numbers. When we are deciding on the truth value of $c_{i,j}$ we have the following reasoning:

- If $c_{i-1,j}$ is FALSE, we try to set $c_{i,j}$ both to FALSE and TRUE (if FALSE fails). When $c_{i,j}$ is set to FALSE, all of the potential \mathbf{x}_i solutions are greater than or equal to \mathbf{x}_{i-1} , thus we continue with the same bit comparison on the next level. However, when $c_{i,j}$ is set to TRUE, all of the potential \mathbf{x}_i solutions are strictly greater than \mathbf{x}_{i-1} and we no longer do bit comparison on further levels.
- If $c_{i-1,j}$ is TRUE, we only try out the truth value of FALSE and we continue to do bit comparison since the potential \mathbf{x}_i solutions are greater than or equal to \mathbf{x}_{i-1} at this point.

Lastly, we give further information which explain in full detail Algorithm 3. We use a flag denoted compare to instruct whether to do bit comparison at the current search tree level or not. On line 6 we reset the compare flag to TRUE since $c_{i,j}$, when j=0, corresponds to a leading bit of the next **x**-vector. Lastly, if-conditions on line 8 have to checked in the specified order.

Algorithm 3 hides a depth-first transversal of a binary search tree with a symmetry breaking technique. We specifically designed it for the PDP, but it can be applied to similar problems that deal with symmetry.

5 Time complexity analysis

As we explained in Section 2, the time complexity of a SAT problem in a DPLL context is measured by the number of conflicts. This essentially corresponds to the number of leaves created in the binary search tree. The worst case complexity of the algorithm is thus 2^h , where h is the height of the tree.

As per Proposition 1, we only reason on $c_{i,j}$ variables from the XOR model. Therefore, h=ml and the worst-case complexity for the PDP is 2^{ml} .

Furthermore, with the symmetry breaking technique explained in Section 4.1, we optimize this complexity by a factor of m!. Indeed, out of the m! permutations of the solution set $\{\mathbf{x}_1, \ldots, \mathbf{x}_m\}$, only one satisfies $\mathbf{x}_1 \leq \mathbf{x}_2 \leq \ldots \leq \mathbf{x}_m$ (neglecting the equality).

This concludes that the worst-case number of conflicts reached for one PDP computation is

$$\frac{2^{ml}}{m!}. (11)$$

Going further in the time complexity analysis, we observe that to find one conflict we go through (in the worst case) all clauses in the model during unit propagation. Hence, the running time per conflict grows linearly with the number of clauses. First, **Algorithm 3** Function SOLVE_BR_SYM(F, compare): Recursive function for solving a SAT formula derived from PDP.

```
Input: Propositional formula F and a flag compare
Output: TRUE if formula is satisfiable, FALSE otherwise.
 1: if all clauses and all XOR-clauses are satisfied then
2:
        return TRUE.
3: end if
4: choose one c_{i,j}.
5: if j=0 then
6:
        compare \leftarrow \texttt{TRUE}.
7: end if
8: if (compare is FALSE) or (i = 1) or (c_{i-1,j}) is set to FALSE) then
        (contradiction, F') \leftarrow ASSIGN(F, \neg c_{i,j}).
10:
        if contradiction then
11:
            BACKTRACK().
12:
            compare \leftarrow \texttt{FALSE}.
13:
        else
14:
            if SOLVE\_BR\_SYM(F', compare) returns FALSE then
15:
                BACKTRACK().
16:
                compare \leftarrow \texttt{FALSE}.
            else
17:
18:
                return TRUE.
19:
            end if
20:
        end if
21: end if
22: (contradiction, F') \leftarrow ASSIGN(F, c_{i,j}).
23: if contradiction then
24:
        BACKTRACK().
25:
        return FALSE.
26: end if
27: return SOLVE_BR_SYM(F', compare).
```

let us count the number of clauses in the X-substitution set. For every $2 \leq d \leq m$ there exist $\binom{m}{d} \cdot l^d$ monomials of degree d given by products of variables $c_{i,j}$, and they each yield d+1 clauses (see Equation (9)). In total, the number of clauses in the X-substitutions set is

$$(\sum_{d=2}^{m} {m \choose d} \cdot l^d)(d+1).$$

Recall that degree one monomials are not substituted and thus do not produce new clauses. We can adapt this reasoning for the E-substitutions set as well.

The number of XOR-clauses in the XOR model is equivalent to the number of equations in the algebraic model. We have $\frac{m(m+1)}{2}(l-1)+m$ in the E-X-relation set and n in the F set.

Remark 2. Using this analysis, we approximate the number of clauses for m=3, as all experiments presented in this paper are performed using the fourth summation

polynomial.

$$#C \approx {3 \choose 2} \cdot 3l^2 + {3 \choose 3} \cdot 4l^3 + {3 \choose 2} \cdot 3(3l-2)^2 + (6l-3) + n \approx (12)$$
$$\approx 4l^3 + 171l^2 - 210l + n + 69.$$

In practice, many monomials have no occurrence in the system after the Weil descent process. In fact, the value in (12) is a huge overestimate and exact values for $l \in \{6, \ldots, 11\}$ are shown in Table 1.

Assuming that we take m small, we conclude that the number of clauses in our model, denoted by C, is polynomial in l. Let t be a constant representing the time to process one clause. The running time of the PDP is bounded by

$$t \cdot C \cdot 2^{ml}/m!$$
.

This allows us to establish the following result on the complexity of our SAT-based index calculus algorith.

Theorem 1. The complexity of the index calculus algorithm for solving ECDLP on a curve defined over \mathbb{F}_{2^n} , using a factor base given by a vector space of dimension $l \approx \frac{n}{m}$, is $\tilde{O}(2^{n+l})$, where the \tilde{O} hides a polynomial factor in l.

Proof. In order to perform a whole ECDLP computation, one has to find 2^l relations. Following [8], when $l \approx \frac{n}{m}$ the probability that a random point can be written as a sum of m factor basis elements is heuristically approximated by $\frac{2^{ml}}{m!2^n}$. The time complexity for the full decomposition phase, using our dedicated WDSAT solver is:

$$\#Ct2^{n+l}$$
.

This worst case complexity is to be compared to the $O(2^{\omega \frac{n}{2} + l})$ complexity of Faugère *et al* [12]. Moreover, we underline here that Faugère *et al*'s proof of this result is based on heuristic assumption on the Gröbner basis computation for PDP, while our analysis for the SAT-based approach simply relies on the rigourously proved worst case for the DPLL search tree.

6 Experimental Results

We conducted experiments using S'_4 on binary Koblitz elliptic curves [20] defined over \mathbb{F}_{2^n} . We experimented with Gröbner basis and SAT approaches. All tests were performed on a 2.40GHz Intel Xeon E5-2640 processor.

The Gröbner basis approach takes as input an algebraic model. We used the *grevlex* ordering, as this is considered to be optimal in the literature. MINISAT and GLUCOSE solvers process a CNF model input, whereas CRYPTOMINISAT and WDSAT use the

XOR model. Using the XOR model is a huge advantage, as it has far less clauses and variables than the CNF model.

Gaussian elimination can be beneficial for SAT instances derived from cryptographic problems. However, it has been reported to yield slower runtimes for some instances as performing the operation is very costly. For this reason, CryptoMiniSat does not include Gaussian elimination by default, but the feature can be turned on explicitly. We experimented with both variants, as we did for WDSAT.

In Table 2 CRYPTOMINISAT $_{\rm XG}$ and CRYPTOMINISAT denote CryptoMiniSat with and without Gaussian elimination respectively. Same notation is adopted for the WD-SAT solver. These experiments were performed using WDSAT before applying the breaking symmetry technique as Table 2 serves as comparison between the different approaches and breaking symmetry is a feature that none of the other solving tools possess. The performance of the complete WDSAT solver is shown in Table 3 where we increase l up to 11. With other approaches, we could not handle these larger values of l in reasonable time.

We experimented with different values of n for each l and we performed tests on 20 instances for each parameter size. Half of the instances have a solution and the other half do not. We show averages on satisfiable and unsatisfiable instances separately, since running times differ between the two cases. SAT solvers stop as soon as they find a solution and if this is not the case they need to respond with certainty that a solution does not exist. Hence, running times of SAT solvers are significantly slower when there is no solution. On the other hand, [35] indicates that the computational complexity of Gröbner basis is lower when a solution does not exist.

We set a timeout of 10 hours for each run and $\#Runs_{XG}$ corresponds to the number of instances that were successfully solved within this time frame. Note however, that Gröbner basis computations for l=8 did not halt because they reached a timeout. They were stopped because of the memory limit of 200GB.

Other information in Table 2 are the average runtime in seconds and the average number of conflicts.

		,	SATisfiable		UNSATisfiable			
Approach	1	n	Runtime	#Conflicts	#Runs _{succ}	Runtime	#Conflicts	#Runs _{succ}
		17	207.220	NA	10	142.119	NA	10
	6	18	208.859	NA	10	155.765 NA 1	NA	10
		19	215.187	NA	10		NA	10
		19	3854.708	NA	10		10	
Gröbner	7	21	3485.273	NA	10	2444.487	NA	10
		23	3128.844	NA	10	2286.136 NA	NA	10
		23			0			0
	8	24			0			0
		26			0			0

			,	SATisfiable		UI	le	
Approach	1	n	Runtime	#Conflicts	#Runs _{succ}	Runtime	#Conflicts	#Runs _{succ}
	6	17	62.702	408189	10	270.261	1463309	10
		18	45.220	297995	10	238.731	1211297	5
		19	229.055	1778377	10	388.719	2439933	10
		19	406.918	1919565	10	6777.431	25180492	10
MINISAT	7	21	1324.932	6809454	10	9933.959		
		23	12945.613		10	13260.586	59289671	10
		23	8027.974	63384411	8			0
	8	24			0			0
		26			0			0
		17	81.898	711918	10	119.694	815185	10
	6	18	50.981	387141	10	104.625	655006	6
		19	299.175	2332066	10	269.212	2077689	10
	7	19	908.091	5357976	10	1356.990	5884897	9
GLUCOSE		21	1626.805	9467330	8	3138.480	14451643	8
		23	2585.200	12528231	7	3760.138	16898505	
	8	23	6755.026	20886673	7			0
		24			0			0
		26			0			0
	6	17	133.983	775948	10	363.513	1709971	10
		18	107.568	629571	10	3097.879	13549468	6
		19	560.080	3396192	10	1172.740	5726372	10
		19	1210.612	5713259	10		26079224	
CRYPTOMINISAT	7	21	5298.106	22233588	10	23009.998		
		23	3637.032	12159752	10	19857.454	47086152	10
		23	9846.554	18509058	10			0
	8	24	7902.745	12121156	10			0
		26	6905.477	13269631	10			0
		17	119.866	677336	10	436.811	1877699	10
	6	18	168.428	956679	10	1469.945	6304762	10
		19	224.484	1219840	10	615.952	2763754	10
	7	19	893.425	3722805	10	3587.929	8642108	10
CRYPTOMINISAT _{XG}		21	737.768	2366264	10	4272.053	12340513	10
		23	580.007	1753040	10	3253.786	8183887	10
	8		11265.010	19604250	10			0
		24		11582517	10			0
		26	3933.637	7920920	9			0

			,	SATisfiable		U	UNSATisfiable		
Approach	1	n	Runtime	#Conflicts	#Runs _{succ}	Runtime	#Conflicts	#Runs _{succ}	
	6	17	.601	49117	10	3.851	254686	10	
		18	1.295	97283	10	3.856	254602	10	
		19	.470	38137	10	3.913	255491	10	
		19	9.643	534867	10	44.107	2073089	10	
WDSAT	7	21	7.636	403434	10	45.780	2072053	10	
		23	9.303	477632	10	47.347	2067168	10	
		23	68.929	2646071	10	525.057	16666331	10	
	8	24	114.945	4291229	10	524.715	16691522	10	
		26	185.480	6261107	10	533.607	16684378	10	
		17	9.193	48178	10	56.718	253123	10	
	6	18	20.043	95283	10	58.001	252985	10	
		19	7.041	36835	10	58.876	252799	10	
		19	169.629	528383	10	736.863	2062232	10	
$WDSAT_{XG}$	7	21	131.272	397758	10	763.570	2061408	10	
		23	159.101	473223	10	779.432	2060501	10	
		23	1290.702	2630567	10	9124.361	16639322	10	
	8	24	2145.313	4256384	10	9421.994	16638399	10	
		26	3404.765	6231289	10	9623.677	16636122	10	

Table 2: Comparing different approaches for solving the PDP.

As expected, the Gröbner basis approach was outperformed by state-of-the-art SAT solvers when a solution exists. When there is no solution, it yields faster running times than solvers, but it quickly becomes impractical because of the memory requirements. For comparison, the memory used by SAT solvers for l=7 is between 60MB and 200MB, whereas Gröbner basis require 38GB.

Our dedicated WDSAT solver yields significantly faster running times than any of the state-of-the-art tools. As we explained in Section 5, the number of conflicts found by WDSAT is bounded by 2^{3l} . We observe however, that running times are slower for the WDSAT $_{XG}$ variant. This is explained by observing that the number of conflicts is only slightly better when Gaussian elimination is used. The cost of performing a Gaussian elimination at every level of the binary search tree outweighs the benefit of having reached less conflicts.

Choosing the WDSAT variant without Gaussian elimination as optimal, we continued experiments for bigger size parameters using this variant coupled with the breaking symmetry technique. Table 3 shows results for $l=\overline{6,11}$ and n sizes up to 89. As before, all values are an average of 10 runs. If we compare the number of conflicts for the first three l sizes of the complete WDSAT solver with its symmetrical variant in Table 2, we observe a speedup factor that rapidly approaches $6.^2$ This confirms our claims in Section 5.

Comparing results for l=6 and l=7 in Table 3 with the equivalent results for the Gröbner basis method in Table 2, we observe that WDSAT is up to 300 times faster

² We compare the cases where there is no solution, as these have more stable averages.

than Gröbner basis for the cases where there is no solution and up to 1300 times faster for instances allowing a solution. This is a rough comparison, as the factor grows with parameters l and n and running times for satisfiable instances can vary remarkably.

6.1 Whole Point Decomposition Phase Computation

Previously shown experiments for solving PDP are done with arbitrary choices of parameters n and l. However, when performing a whole ECDLP attack, choosing the factor base is a crucial step in the index calculus method. The number of relations that needs to be found is exponential in l, as is the running time for one point decomposition (see 11). However, taking a smaller l decreases the probability of successfully decomposing a randomly chosen point, and thus increases the number of times we solve the PDP.

To understand better the optimal ratio n/l, we computed the whole point decomposition phase for n=24 using different l sizes. The experiment consists essentially in computing the PDP on instances for randomly chosen X_{m+1} until we find 2^l valid decompositions. Instances that turn out to not have a solution are tossed.

Results from these experiments are in Table 4. We present average running times in hours (Runtime), the number of generated satisfiable (#Generated SAT) and unsatisfiable (#Generated UNSAT) instances and the probability that a random point can be decomposed (P). This probability is, in fact, the ratio between the number of generated instances that have a solution and the total number of generated instances. We compare this to the heuristically approximated probability $\frac{2^{ml}}{m!2^n}$, denoted by (P_{approx}). #Runs denotes the number of times we ran the experiment for each n/l ratio.

Further research is needed to fully understand the best choice of l, but current experimental results suggest that it is not $l \approx n/m$. With this ratio, even though there is a good chance of finding a composition for a randomly chosen point, the runtime of PDP and the number of relations needed are too high. The cost of the final phase linear algebra increases with l as well.

At the time of submission of this paper, we do not have results for bigger values of n and l. For this reason, we refrain from drawing a conclusion from Table 4, as results can be ambiguous for such small parameters. As an example, the time to compute the Weil descent and derive the XOR model, is not negligible in the cases of l=4 and l=5 where the time to solve the instance is less than 60 milliseconds long. This might be the only reason we observe a decrease of the runtime between l=4 and l=5, whereas from the remaining results in Table 4 we observe that the time increases with l.

7 Conclusions and Future Work

Gröbner basis methods have been shown powerful in solving the PDP in the index calculus attack for elliptic curves defined over small degree extension fields in characteristic > 2. In this paper we argue that for finite fields in characteristic 2 a SAT-based approach yields more practical results. We started by explaining that general-purpose SAT solvers cannot yield considerably faster running times because the number of variables in a SAT model is significantly larger than the number of variables in the algebraic model.

		SAT	risfiable	UNSA	ATisfiable
1	n	Runtime	#Conflicts	Runtime	#Conflicts
	17	.233	18517	.645	43867
6	18	.395	29115	.642	43798
	19	.200	15715	.641	43995
	19	3.138	176931	7.107	351329
7	21	2.576	138494	7.337	351133
	23	3.280	169294	7.622	350068
	23	24.642	967945	83.797	2800507
8	24	41.603	1563509	83.673	2805893
	26	52.181	1800948	86.122	2803872
	27	371	10204950	915	22412545
	37	511	11890506	1047	22395287
9	47	736	15300468	1170	22378061
	59	753	13462450	1333	22387826
	67	652	11058755	1420	22392359
	30	4137	82804327	9760	179133743
	47	7244	116867105	12129	179024622
10	59	8829	124714894	13951	179065509
	67	6437	85315820	14381	179054617
	79	6387	77450682	15979	179033814
	33	51271	780737411	101832	1432516919
	59	92322	1011596679	136314	1432242942
11	67	52514	577957022	143946	1432211311
	79	48268	491455095	157801	1432104989
	89	53292	485794746	171274	1432071999

Table 3: Running times and number of conflicts using the complete WDSAT solver.

n	1	Runtime	#Generated SAT	#Generated UNSAT	P	P_{approx}	#Runs
	4	2.59	16	149475	.00010	.00004	10
	5	1.83	32	49484	.00064	.00032	10
24	6	2.49	64	11478	.00557	.00260	10
	7	6.45	128	2994	.04275	.02083	10
	8	17.99	256	688	.37209	.16666	8

Table 4: Whole Point decomposition phase computation.

Our first contribution is to propose a PDP XOR model with only ml core variables, whose assignment propagate all remaining variables in the model. Consequently, with appropriate solving methods the time complexity of the PDP is in fact 2^{ml} . To this end, we conceived a SAT solver dedicated to solving systems derived from a Weil descent, named WDSAT. We further optimized the time complexity of this solver by a factor of m! using a symmetry breaking technique.

We presented experiments for the PDP on prime degree extension fields using parameter sizes of up to l=11 and n=89. In addition, our experiments with a full ECDLP computation for small size parameters suggest that further research is needed to find the optimal n/l ratio for the index calculus attack on these curves.

Another perspective would be to find and set an early abort threshold for our solver, since experiments suggest that PDP instances that have no solution have slower running times than instances allowing a solution.

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