

Accelerating ID-based Encryption based on Trapdoor DL using Pre-computation

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Abstract. The existing identity-based encryption (IBE) schemes based on pairings require pairing computations in encryption or decryption algorithm and it is a burden to each entity which has restricted computing resources in mobile computing environments. An IBE scheme (MY-IBE) based on a trapdoor DL group for RSA setting is one of good alternatives for applying to mobile computing environments. However, it has a drawback for practical use, that the key generation algorithm spends a long time for generating a user's private key since the key generation center has to solve a discrete logarithm problem.

In this paper, we suggest a method to reduce the key generation time of the MY-IBE scheme, applying modified Pollard rho algorithm using significant pre-computation (**mPAP**). We also provide a rigorous analysis of the **mPAP** for more precise estimation of the key generation time and consider the parallelization and applying the tag tracing technique to reduce the wall-clock running time of the key generation algorithm.

Finally, we give a parameter setup method for an efficient key generation algorithm and estimate key generation time for practical parameters from our theoretical analysis and experimental results on small parameters. Our estimation shows that it takes about two minutes using pre-computation for about 50 days with 27 GB storage to generate one user's private key using the parallelized **mPAP** enhanced by the tag tracing technique with 100 processors.

Keywords: Identity-based Encryption, Trapdoor DL Groups, Discrete Logarithm, Pre-computation.

1 Introduction

In [26], Shamir suggested the concept of identity-based cryptosystem and proposed identity-based signature schemes. Since then, there have been several proposals [9, 16, 28, 31] to construct an identity-based encryption (IBE) scheme. However, these proposals are not fully satisfactory. Some schemes require the condition that users cannot collude for the security and other schemes spend a long time to generate a user's private key. Later, the first sufficiently secure and efficient IBE scheme was proposed using pairings by Boneh and Franklin [4]. Since Boneh and Franklin's construction, many IBE schemes [3, 10, 33, 34] based on pairings have been proposed to enhance the security or efficiency. Also, some IBE schemes [2, 11] based on lattices have been also proposed under the necessary of lattice-based cryptosystems.

However, the existing IBE schemes based on pairings or lattices are not proper on some systems in which each entity has restricted computing resources, such as mobile computing environments. In case of pairing-based IBE schemes, an encryption algorithm or a decryption algorithm requires one or more pairing computations and it spends much computation cost than modular exponentiations over finite fields. Also, although the encryption and decryption

algorithm of lattice-based IBE schemes are more efficient than those of pairing-based IBE schemes, the public key size and private key size of lattice-based IBE schemes are quite large to utilize in mobile entities.

Among previous IBE schemes, Maurer and Yacobi's suggestion is one of good alternatives for applying to mobile computing environments. In [16], the authors proposed a non-interactive key distribution (NIKD) algorithm based on a trapdoor DL group which is a maximal cyclic subgroup G of \mathbf{Z}_N^* where an integer N is hard to factor. Also the authors provided the IBE (MY-IBE) scheme from their NIKD algorithm. The most hard computation in encryption and decryption algorithms of the MY-IBE scheme is a modular exponentiation in \mathbf{Z}_N^* and the user's private key consists of one element less than N , hence, they are no burden for a mobile entity.

However, the MY-IBE scheme has some problems for practical use. First, the security proof of the MY-IBE scheme is insufficient to formal security notion of IBE schemes, provided in [4]. Second, there are no secure full domain hash functions into a maximal cyclic subgroup G of \mathbf{Z}_N^* . Moreover, the most serious obstacle for practical use of the MY-IBE scheme is that the key generation time is quite long since the key generation center (KGC) has to solve a discrete logarithm problem (DLP) in G for generating a private key of a user.

Later, Paterson and Srinivasan resolved the above three problems. In [20], they proposed a full domain hash function into G and proved the security of the proposed hash function. Then, they refined the MY-IBE scheme using their hash function and provided the security proof of the refined MY-IBE scheme based on the formal security definition of IBE schemes. Finally, they suggested that the use of the index calculus algorithm with significant pre-computation to reduce the key generation time. However, although KGC utilizes the index calculus algorithm with significant pre-computed data, there are no known concrete methods reducing DLP solving time and we cannot estimate the expected key generation time and required resources such as the amount of memory and the number of processors.

1.1 Our Contribution

The authors in [20] noted that there were no proper variants of the Pollard rho algorithm [23] using pre-computation. However there have been some proposals [12, 14] to reduce DLP solving time, modifying the parallelized Pollard rho algorithm [32] using pre-computation. In this paper, we suggest the method to reduce the key generation time of the MY-IBE scheme applying a modified Pollard rho algorithm using pre-computation (**mPAP**).

Moreover, although there have been some complexity analyses [12, 14] related to the **mPAP**, they did not give precise relations between the required memory size, the pre-computation time and the online time. Especially, all previous analyses did not provide concrete required memory size in the **mPAP**. Hence, one cannot precisely estimate the required memory size of the MY-IBE scheme with respect to DLP solving time in the practical system. In this paper, we provide more rigorous complexity analysis of the **mPAP** including relations between the amount of pre-computation, online computation and the memory size and give experimental results of the **mPAP** on small parameters.

Also, we discuss about two extensions for reducing the wall-clock running time of the online phase in the mPAP. First, we consider the possibility of the parallelization of the mPAP and provide the complexity analysis of the parallelized mPAP. According to our analysis, while the pre-computation phase of the mPAP can trivially be parallelized with speedup linear in the number of processors, the online phase of the mPAP can be parallelized with linear increments of storage for speedup linear in the number of processors. Second, we consider applying the tag tracing technique [7] to the mPAP for solving DLPs over a finite field. From experimental results, we confirm that the tag tracing technique works well with the mPAP.

Lastly, we suggest a parameter setup method for reducing key generation time with maintaining security level and estimate the key generation time and the memory size of the MY-IBE scheme in the practical system from our theoretical analysis and implementation results on small parameters. According to our estimation on parameters for 2^{80} security, the parallelized mPAP enhanced by the tag tracing technique [7] requires about one minutes and 46 seconds with 100 processors when 100 processors are used for pre-computation for about 50 days with 27 GB storage. Since the key generation process is performed by the key generation center, 100 processors are quite practical.

1.2 Organization

We introduce the MY-IBE scheme and give some related works on DL algorithm using pre-computation in Section 2. Section 3 describes the mPAP and provides more rigorous complexity analysis of the algorithm. We also give experimental results on small parameters and consider the parallelization and applying the tag tracing technique. In Section 4, we suggest a parameter setup method for reducing key generation time and estimate the key generation time of the MY-IBE scheme from our theoretical analysis and experimental results on small parameters.

2 Preliminaries

In this section, we introduce an identity-based encryption scheme (MY-IBE) based on trapdoor DL groups for RSA setting, proposed by Maurer and Yacobi [16] and refined by Paterson and Srinivasan [20].

The key generation center has to solve one DLP in the key generation algorithm of the MY-IBE scheme and it takes a long time for practical use. To reduce the key generation time the authors of [20] suggested the re-use of significant pre-computed elements for solving previous DLPs. However, their suggestion was not enough to make the MY-IBE scheme practical for real systems. At the end, we briefly introduce related works on our suggestion to solve DLPs using significant pre-computation, which is a modification of the parallelized Pollard rho algorithm [32].

2.1 Identity-based Encryption based on Trapdoor DL

First, we define trapdoor DL groups and trapdoor DL group generators.

Definition 1 ([8]). We define the two following algorithms:

- **TDLGen** : a polynomial-time algorithm which takes as input a security parameter λ and outputs a finite cyclic group G , a generator g of G and its trapdoor information τ .
- **SolveTDL** : a polynomial-time algorithm which takes as input a finite cyclic group G , a generator g of G , a trapdoor information τ and a target element h of a DLP and outputs the DL of h based on g .

We also define a polynomial-time algorithm **Gen** which takes as input a security parameter λ , runs **TDLGen**(λ) algorithm and then outputs (G, g) . If (G, g, τ) is an output of **TDLGen** and a DLP over G is hard for the output of **Gen**, we call G a trapdoor DL group.

As examples of trapdoor DL groups, one considers a maximal cyclic subgroup G of \mathbf{Z}_N^* where N is a product of primes and its trapdoor information is the factorization of N [16]. Then one who knows a factorization of N can efficiently solve a DLP over G using Index Calculus algorithm or Pohllig-Hellman algorithm with Pollard rho algorithm. However, if N is a product of more than three primes, there are no secure full domain hash function into G . Also, one may consider trapdoor DL groups over an elliptic curve $E(\mathbf{F}_{2^{161}})$, whose trapdoor information are isogenies [30]. But an elliptic curve $E(\mathbf{F}_{2^{161}})$ is the only currently known possible parameter and it is not yet known how to generalize construction to higher security level.

In this paper, among these trapdoor DL groups, we will only deal with a maximal cyclic subgroup G of \mathbf{Z}_N^* where N is a product of two primes p, q that are roughly the same size and satisfy $p \equiv 3 \pmod{4}$, $q \equiv 1 \pmod{4}$, and $p - 1, q - 1$ are B -smooth integers.

Note that **TDLGen** and **SolveTDL** are both polynomial-time algorithms in the above definition. To our knowledge, there are no polynomial-time algorithms to solve a DLP in a maximal cyclic subgroup G of \mathbf{Z}_N^* although the trapdoor information is given. Hence, the above group G does not satisfy the definition of trapdoor DL groups. However, there are some algorithms to solve DLP over G , whose complexity is sub-exponential or exponential in the security parameter λ but it is quite practical. Hence, we will apply the definition of **SolveTDL** relaxedly.

Identity-based Encryption based on trapdoor DL groups In [16], Maurer and Yacobi proposed an IBE scheme based on a trapdoor DL group G which is a maximal cyclic subgroup of \mathbf{Z}_N^* where N is a product of primes and whose trapdoor information is a factorization of N . However, when N is a product of more than three primes, the authors could not give an efficient full domain hash function from a set $\{0, 1\}^*$ to G . Also, although they presented two efficient full domain hash functions from $\{0, 1\}^*$ to G when N is a product of two primes, it was proved that their suggestions were not secure [15, 17–19]. In [19], the authors presented another full domain hash function into G , however they did not give a security proof of the presented hash function.

Later, Paterson and Srinivasan [20] provided a security proof of a full domain hash function presented in [19]. Based on this full domain hash function, they provided the IND-ID-CPA se-

cure IBE scheme based on the security notion in [4]¹. Here, we present the MY-IBE scheme modified by Paterson and Srinivasan. The scheme consists of four algorithms:

- **Setup(λ)**: this algorithm runs **TDLGen** algorithm to obtain (G, g, τ) where G is a maximal cyclic subgroup of \mathbf{Z}_N^* , g is a generator of G and trapdoor information τ which is a factorization of N . (We assume that N is a product of two primes that are roughly the same size and satisfy $p \equiv 3 \pmod{4}$, $q \equiv 1 \pmod{4}$ and $\gcd(p-1, q-1) = 2$.) Let H be a hash function from $\{0, 1\}^*$ to \mathbf{Z}_N and define $H_1 : \{0, 1\}^* \rightarrow G$ by

$$H_1(\text{ID}) = \left(\frac{H(\text{ID})}{N} \right) H(\text{ID})$$

where $\left(\frac{x}{N} \right)$ denotes the Jacobi symbol. Let $H_2 : G \rightarrow \{0, 1\}^\ell$ be a hash function where ℓ is the bit size of messages. Then this algorithm outputs

$$\begin{aligned} \text{params} &= (\lambda, G, N, g, H, H_1, H_2, \ell) \\ \text{msk} &= \tau = (p, q). \end{aligned}$$

- **KeyGen(params, msk, ID)**: this algorithm computes $H_1(\text{ID})$. Then it runs **SolveTDL** $(G, g, \tau, H_1(\text{ID}))$ and obtains the private key s_{ID} such that $g^{s_{\text{ID}}} = H_1(\text{ID})$. It outputs s_{ID} .
- **Encrypt(params, ID, M)**: this algorithm computes $H_1(\text{ID})$ and chooses $r \in \mathbf{Z}_N$ uniformly at random. Then it outputs $C = (U, V)$ where $U = g^r$ and $V = M \oplus H_2(H_1(\text{ID})^r)$.
- **Decrypt(params, s_{ID}, C)**: for a ciphertext $C = (U, V)$, this algorithm outputs $M' = V \oplus H_2(U^{s_{\text{ID}}})$.

Note that the above scheme is IND-ID-CPA secure under the CDH assumption and one-wayness of the hash function H_1 . Also the CDH problem in the above group G is at least hard to factor N . Therefore, the above scheme is IND-ID-CPA secure if factoring N is hard.

Although the encryption and decryption algorithms are efficient since those require two and one modular exponentiation, respectively, KGC has to solve one DLP over G to generate the private key for one user. Hence, the weak point of the above scheme is to spend a long time in the key generation algorithm. To overcome this obstacle, the authors in [20] suggested the use of the index calculus algorithm with a large amount of pre-computation.

2.2 Discrete Logarithm Algorithm using Pre-computation

To reduce key generation time in KeyGen algorithm of the MY-IBE scheme, the authors in [20] suggested the use of the index calculus algorithm [1] with significant pre-computation over \mathbf{Z}_p and \mathbf{Z}_q for solving a DLP in a maximal cyclic subgroup G of \mathbf{Z}_N^* where $N = pq$ and p, q are primes. The index calculus algorithm for solving DLPs divides into two parts, the pre-computation phase and the online phase. In the pre-computation phase, when the cyclic group G and a generator g are given, one computes DLs of all elements in the factor base \mathcal{B} . Then in the online phase, one finds some relations between the target element of the

¹ Throughout this paper, we follow the security notion of IBE schemes in [4].

DLP and elements in the factor base \mathcal{B} and then solves the DLP using DLs of elements in the factor base \mathcal{B} , which are pre-computed in the pre-computation phase.

However, the complexity of the online phase is the same with the complexity of the pre-computation phase in the original index calculus algorithm. Hence, it also requires a large amount of computations for the online phase. For achieving 80-bit security, we assume that the bit size of p and q is 512. Then, according to the analysis of the number field sieve method [6] which is a variant of the index calculus algorithm, it is required about $O(2^{78.04})$ multiplications² in \mathbf{Z}_p and \mathbf{Z}_q in the online phase and hence it is impractical to realize in the system.

To make the MY-IBE scheme more practical, we will propose the use of the Pohlig-Hellman algorithm and the modified Pollard rho algorithm with a large amount of pre-computation. In [20], the authors noted that there were no proper algorithms to solve DLPs using the Pollard rho algorithm with pre-computation. However, there have been some modified Pollard rho algorithms [12, 14] to solve a DLP efficiently using pre-computation. In [14], the authors provided the algorithm for solving multiple DLPs using elements which were computed for solving previous DLPs, modifying the parallelized Pollard rho algorithm [32]. Then the authors in [12] modified the above multiple DLPs solving algorithm to the DLP solving algorithm with pre-computation, by solving randomly generated DLPs before the target element of the original DLP is given.

In the rest of this section, we introduce some basic concepts, r -adding walk iterating function and distinguished point (DP) collision detection method, which are composing variants of Pollard rho algorithm.

When G is a cyclic group of order q generated by a generator g , if we know the factorization of $q = \prod_{i=1}^{\ell} p_i^{e_i}$ where p_i 's are primes, one can reduce the DLP over G to DLPs over a group of order p_i 's using the Pohlig-Hellman algorithm. Hence, it is assumed that the group order q is prime from here in this subsection. The group element h denotes the target element of the DLP, in other words, we are looking for the value $\log_g h$.

r-adding walk iteration function We briefly look into r -adding walk iterating functions. Partition G into r roughly same sized subsets G_1, \dots, G_r so that $G = G_1 \cup \dots \cup G_r$. The index function $s : G \rightarrow \{1, 2, \dots, r\}$ is defined to be almost pre-image uniform and efficiently computable. Then choose r pairs $(u_i, v_i) \in \mathbf{Z}_q \times \mathbf{Z}_q$ and set r multipliers M_i to $g^{u_i} h^{v_i}$. (In [12], the authors suggested the use of multipliers which are independent of the target element. Hence we will set multipliers of r -adding walk iterating function in the mPAP to the form $g^{u_i} h^0$ where u_i is a randomly chosen integers in \mathbf{Z}_q .) Define r -adding walk iterating function $F_r : G \times \mathbf{Z}_q \times \mathbf{Z}_q \rightarrow G \times \mathbf{Z}_q \times \mathbf{Z}_q$ by

$$F_r(y, a, b) = (y \cdot M_{s(y)}, a + u_{s(y)}, b + v_{s(y)})$$

where $y = g^a h^b$. Throughout this paper, F_r will denote an r -adding walk iterating function.

² The heuristic complexity of the number field sieve for solving one DLP over \mathbf{Z}_p^* is $\exp((c + o(1))(\log p)^{1/3}(\log \log p)^{2/3})$ bit operations where $c = (64/9)^{1/3} \approx 1.9233$. And it takes $O(\log^2 p)$ bit operations for one modular multiplication between $\log p$ -bit integers. Hence we obtain that the computational complexity of the number field sieve for solving one DLP over 512-bit prime field is $O(2^{78.04})$ multiplication.

It was shown [25] that the expected number of iterations for finding a collision in a walk generated by r -adding walk for $r \geq 8$ is $O(\sqrt{q})$. Experimental results [29] over elliptic curve groups show that the expected number of iterations required to find a collision with 20-adding walk is very close to $\sqrt{\frac{q}{2}\pi}$, which is that with a random function.

DP collision detection method Let us introduce the DP technique [24], which was originally used in time memory tradeoff techniques. One sets the distinguishing property that is easy to check and define a DP by a point satisfying the distinguishing property in G . For example, one may define the distinguishing property to be that a certain number of the most significant are all zeros under a fixed encoding of G . One starts with an empty table, and the walk is computed iteratively until the walk encounters a DP. Then one searches for the same point with the occurred DP in the table. If it is not in the table, one stores it and generates another walk. The DP collision detection method is required about t additional iterations for noticing a collision after a collision occurs in a walk when t^{-1} is the proportion of DPs in G . It is straightforward to apply the DP method to multiple walks, so the DP method has an advantage that admits n -times speedup with an n -processor parallelization [32].

3 Analysis of Discrete Logarithm Algorithm using Pre-computation

In this section, we describe the mPAP, which is a modification of the parallelized Pollard rho algorithm [32]. Then, we provide more rigorous complexity analysis of the algorithm and give experimental results on small parameters. Finally, we consider two extensions to reduce wall-clock running time of the online phase in the mPAP, the parallelization of the mPAP and applying the tag tracing technique.

3.1 Algorithm Description

Now, we are ready to describe the mPAP, which is a modification of the parallelized Pollard rho algorithm [32]. Our description is modified in three points compared with the algorithm in [12]. First, in the pre-computation phase one generates chains which are started from a random starting point and are ended at a DP, not randomly generating DL instances and solving these DLPs. Second, multipliers of an r -adding walk iterating function are given by a special form. Third, one does not store created DPs in the online phase because its advantage is almost negligible.

Let us describe the mPAP. Choose positive integers m, t such that $mt^2 = \alpha q$ where m is the number of generated chains in the pre-computation phase, t^{-1} is the proportion of DPs in the group, and α is not a large constant. The parameters are assumed not to be extreme, in the sense that $1 \ll m, t \ll q$ and a typical value is $\alpha \approx 1$. We shall later determine the proper size of parameters m, t, α . Fix an r -adding walk iterating function F_r so that the multipliers have the form $g^u h^0$ for some random $u \in \mathbf{Z}_q$. Since the pre-computation phase starts before a target element is given, the multipliers of the r -adding walk iterating function F_r must have the above form. Determine the distinguishing property so that a proportion of distinguished points in G is t^{-1} .

In the pre-computation phase, one chooses m random starting points $\mathbf{g}_{i,0} = g^{a_{i,0}}h^0 \in G$ for $1 \leq i \leq m$ and iteratively computes $\mathbf{g}_{i,j+1} = F_r(\mathbf{g}_{i,j})$. Each chain is terminated at its first encounter with a DP. We call by a *DP chain* a chain ended at a DP. Then one stores the occurred DPs and the exponents corresponding to each DP in a DP table DT.

The average number of iterations to generate a DP chain is t , but some of chains may fall into a loop that never reaches a DP. In order to detect a chain falling into an infinite loop, we set a chain length bound \hat{t} . Any chain longer than this bound is discarded and one can choose to regenerate a chain from a different starting point. Note that the probability for a chain not to reach a DP until its \hat{t} -th iteration is $(1 - \frac{1}{t})^{\hat{t}+1} \approx \exp(-\frac{\hat{t}+1}{t})$. This shows that setting \hat{t} to a reasonable multiple of t will suffice in removing the effect of any such discarded chains for any practical purpose.

After the pre-computation phase, one obtains the following matrix.

$$\begin{array}{ccccccc} \mathbf{g}_{1,0} & \xrightarrow{F_r} & \mathbf{g}_{1,1} & \xrightarrow{F_r} & \cdots & \xrightarrow{F_r} & \mathbf{g}_{1,t_1} \\ \mathbf{g}_{2,0} & \xrightarrow{F_r} & \mathbf{g}_{2,1} & \xrightarrow{F_r} & \cdots & \xrightarrow{F_r} & \mathbf{g}_{2,t_2} \\ & & \vdots & & & & \\ \mathbf{g}_{m,0} & \xrightarrow{F_r} & \mathbf{g}_{m,1} & \xrightarrow{F_r} & \cdots & \xrightarrow{F_r} & \mathbf{g}_{m,t_m} \end{array}$$

Such matrix consisting of m DP chains is called the *DP matrix*.

In the online phase, when a target element h of DLP is given, one starts to generate a DP chain from an element h^r for a random $1 \leq r < q$. When a DP occurs in a chain, one compares it with the stored DPs in the table DT. If a collision is not found, one generates another DP chain from a distinct starting point $h^{r'}$ for a random $1 \leq r' \neq r < q$. Otherwise, one can get the DL of h from the relation $a_i \equiv xb_j + a_j \pmod{q}$ where x is the DL of h , a pair (g_i, a_i, b_i) and (g_j, a_j, b_j) is the collision, (g_i, a_i, b_i) is the stored point in the table DT, and (g_j, a_j, b_j) is the created DP in the online phase.

In the modification of the mPAP described in [12], created DPs in the online phase are also added to the DP table DT. However, unless α is not extremely small, the number of newly created DPs in the online phase is just one or two. It does not give a big help of accelerating DL computation of a present target element and hence we do not consider saving created DPs in the online phase.

3.2 Complexity Analysis

The analysis of the mPAP can be inferred from DL algorithm [14] for multiple instances. According to the analysis, the expected number of group operations to solve k DLPs sequentially with their algorithm is about $\sqrt{2kq}$. From these results, we simply guess that the expected number of group operations to solve the $(k+1)$ -th DLP is about $\sqrt{2q}/(\sqrt{k+1} + \sqrt{k}) = \sqrt{2(k+1)q} - \sqrt{2kq}$ when $\sqrt{2kq}$ group operations was done in the pre-computation phase since the expected number of group operations to solve $k+1$ DLPs with their algorithm is about $\sqrt{2(k+1)q}$.

Moreover, Hitchcock et al. [12] provided the expected online time complexity of the mPAP associated with the amount of pre-computation. However, their analysis has two missing points:

- Although collisions may occur in the pre-computation phase, it does not consider this fact in their analysis. Hence they regard that the number of distinct elements in a DP matrix is the same with the number of iterations in the pre-computation phase.
- It does not consider the additional iterations to reach a DP for collision detection after a collision occurs in the online phase.

Now, let us correct the complexity of the mPAP. We shall exploit the analysis technique of time-memory tradeoff. Recently, in [13], the authors gave the expected number of distinct entries in a DP matrix for various \hat{t} . However, it does not give an error bound of the approximation value. In the following lemma, we present the precise limit value of the expected number of distinct entries in a DP matrix when \hat{t} is sufficiently large and provide the difference between the limit value and the expected number when we use a specific \hat{t} . This lemma will make up for the first missing point of the analysis provided by Hitchcock et al.

Lemma 1. *Consider a DP matrix created with parameters satisfying $mt^2 = \alpha q$. When the iterating function is taken to be the random function and \hat{t} is sufficiently large, we can expect the DP matrix to contain*

$$\frac{\sqrt{1 + 2\alpha + O(\frac{1}{\hat{t}})} - 1}{\alpha} m(t - 1). \quad (1)$$

distinct entries. Moreover, the difference between the expected number of distinct entries contained in the DP matrix and the above limit value is bounded by $m(\hat{t} + t) \exp(-\hat{t}/t)$.

Proof. We follow the proof of Proposition 10 provided in [13]. Consider a DP matrix generated in the pre-computation phase. Let m_j be the number of elements which first appear at j -th column in a DP matrix. Now we assume that chains not reaching a DP until \hat{t} steps remain on a DP matrix for an analysis and however we will give the error bound considering that these chains are discarded. Then the recurrence relation

$$\frac{m_j}{q} = \left(1 - \exp\left(\frac{-m_{j-1}}{q}\right)\right) \left(1 - \frac{1}{t}\right) \left(1 - \frac{\sum_{i=0}^{j-1} m_j}{q(1 - 1/t)}\right)$$

with $m_0 = m$ is satisfied from Lemma 6 in [13]. Differently from the analysis of [13], the initial value of the sequence $(m_i)_{i=0}^\infty$ is m since our DP table are generated from the exact m starting points without making up for the discarded chains to store the exact m DPs in a table.

Let $\mu_i = \frac{m_i}{q(1 - 1/t)}$ and $\sigma_j = \sum_{i=0}^{j-1} \mu_i$. Then the recurrence relation

$$\sigma_{j+1} - \sigma_j = \frac{m_0}{q} - \frac{1}{t} \sigma_j - \frac{1}{2} \sigma_j^2 + O\left(\frac{1}{t^3}\right) \quad \text{with } \sigma_0 = 0$$

is also satisfied from Lemma 7 in [13]. The sequence $(\sigma_j)_{j=0}^{\infty}$ is monotone increasing from the definition of the sequence $(\sigma_j)_{j=0}^{\infty}$ and all σ_j 's are bounded since the sum $\sum_{j=0}^{\infty} m_j$ does not exceed the group order q . Hence the sequence $(\sigma_j)_{j=0}^{\infty}$ converges and the limit S of this sequence is

$$-\frac{1}{t} + \sqrt{\left(\frac{1}{t}\right)^2 + 2\left(\frac{m}{q} + O\left(\frac{1}{t^3}\right)\right)}$$

and all σ_j 's are less than S . Therefore, the expected number of distinct entries in a DP table is

$$\sum_{i=1}^{\infty} m_i = q(1 - 1/t) \left(-\frac{1}{t} + \sqrt{\left(\frac{1}{t}\right)^2 + 2\left(\frac{m}{q} + O\left(\frac{1}{t^3}\right)\right)} \right) \quad (2)$$

$$= \frac{\sqrt{1+2\alpha+O(\frac{1}{t})}-1}{\alpha} m(t-1). \quad (3)$$

Let E be the expected number of distinct elements in a DP table that chains whose length exceeds \hat{t} are discarded. Then,

$$\begin{aligned} E &\leq \sum_{i=1}^{\hat{t}} m_i \leq q(1 - 1/t) \lim_{j \rightarrow \infty} \sigma_j \\ &= q(1 - 1/t) S \leq \frac{\sqrt{1+2\alpha}-1}{\alpha} m(t-1). \end{aligned}$$

Now, we consider the lower bound of E . Note that the probability that a chain reaches a DP at j steps is $(1 - \frac{1}{t})^{j-1} \frac{1}{t}$. Let E_1 be the expected number of entries belonging to the discarded chains. Then the relation $E \geq q(1 - 1/t) S - E_1$ is satisfied.

$$E_1 = m \sum_{i=\hat{t}}^{\infty} (1 - \frac{1}{t})^{i-1} \frac{1}{t} i = \frac{m}{t} \sum_{i=\hat{t}}^{\infty} (1 - \frac{1}{t})^{i-1} i \quad (4)$$

$$= m \left(\hat{t} \left(1 - \frac{1}{t}\right)^{\hat{t}-1} + \left(1 - \frac{1}{t}\right)^{\hat{t}} t \right) \quad (5)$$

from $\frac{x^{\hat{t}}}{1-x} = \sum_{i=\hat{t}}^{\infty} x^n$ and its derivation and

$$\begin{aligned} (5) &\leq \left(1 - \frac{1}{t}\right)^{\hat{t}} \left(m\hat{t} - m\frac{\hat{t}}{t} + mt\right) \\ &\leq m(\hat{t} + t) \exp(-\hat{t}/t) \end{aligned}$$

since $f(x) = (1 - \frac{1}{x})^x$ is an increasing function and $\lim_{x \rightarrow \infty} f(x) = \exp(-1)$. Hence

$$E_1 \geq \frac{\sqrt{1+2\alpha}-1}{\alpha} m(t-1) - m(\hat{t} + t) \exp(-\hat{t}/t).$$

Therefore, the difference between the limit value and the number of distinct entries on practical parameters is bounded by $m(\hat{t} + t) \exp(-\hat{t}/t)$.

Considering the fact that the authors ignored $(1 - \frac{1}{t})$ in the proof of Proposition 10 in [13], Lemma 1 shows that the exact limit value of the expected number of distinct entries in a DP matrix is the same with the approximation value presented in [13]. Also it shows that the difference between the limit value and the expected number on practical parameters is negligible since $m(\hat{t} + t) \exp(-\hat{t}/t)$ is sufficiently smaller than $\frac{\sqrt{1+2\alpha}-1}{\alpha} m(t-1)$ when \hat{t} is sufficiently larger than t .

We are now ready to discuss about the probability for solving a DLP with generating a single DP chain in the online phase.

Theorem 1. *Fix parameters satisfying $mt^2 = \alpha q$ and generate a DP matrix. Then generate another chain from a random starting point, terminating at its first DP occurrence. When the iterating function F is taken to be the random function, we can expect the ending DP of the new chain to be equal to one of the ending DPs for the pre-generated chains with probability*

$$1 - \frac{1}{\sqrt{1+2\alpha}},$$

and the error term is bounded by $\frac{7\alpha}{t} + \frac{\alpha}{\sqrt{1+2\alpha}}(c+1) \exp(-c)$ when $\hat{t} = ct$ for some constant c .

Proof. Let us write **DP** for the set of all DPs in G , **DM** (DP matrix) for the set of all elements belonging to the pre-generated DP chains, and **DT** (DP table) for the set of ending points of the pre-generated DP chains.

Once the DP matrix **DM** is ready, one is told to create a new chain

$$\mathbf{h}_0 \xrightarrow{F} \mathbf{h}_1 \xrightarrow{F} \cdots \xrightarrow{F} \mathbf{h}_j \xrightarrow{F} \cdots$$

from a random starting point \mathbf{h}_0 . At each iteration, as a new \mathbf{h}_j is created, one of the following three events may occur.

- E1. $\mathbf{h}_j \in \text{DM}$: The new chain has merged with a pre-generated chain. The rest of the chain will automatically follow the pre-generated chain and terminate with a DP belonging to **DT**.
- E2. $\mathbf{h}_j \in \text{DP} \setminus \text{DM} = \text{DP} \setminus \text{DT}$: The new chain has terminated with a DP without merging with a pre-generated chain. The chain cannot reach a point belonging to **DT**.
- E3. $\mathbf{h}_j \notin \text{DM} \cup \text{DP}$: The new chain has neither merged with one of the pre-generated chains nor reached a DP. One needs to continue onto the next iteration.

Hence the new chain terminates with an element from **DT** if and only if each iteration of the chain results in event E3, before finally sinking into event E1.

Let $\delta = \frac{\sqrt{1+2\alpha}-1}{\alpha} m + m(\hat{t} + t) \exp(-\hat{t}/t) + \frac{mO(1)}{\alpha(\sqrt{1+2\alpha}+\sqrt{1+2\alpha+O(\frac{1}{t})})}$. According to Lemma 1, for a random function, we can expect event E1 to happen, at each iteration,

with probability

$$\Pr[E1] = \frac{\sqrt{1+2\alpha} - 1}{\alpha} \frac{mt}{q} + \frac{\delta}{q} = \frac{\sqrt{1+2\alpha} - 1}{t} + \frac{\delta}{q},$$

where the equality follows from a use of $mt^2 = \alpha q$. We can also state

$$\begin{aligned}\Pr[E3] &= 1 - \frac{\#(\text{DM} \cup \text{DP})}{q} = 1 - \frac{\#\text{DM} + \#\text{DP} - \#(\text{DM} \cap \text{DP})}{q} \\ &= 1 - \frac{\frac{q}{t} + \frac{\sqrt{1+2\alpha}-1}{\alpha} mt}{q} + \epsilon \\ &= 1 - \frac{\sqrt{1+2\alpha}}{t} + \epsilon\end{aligned}$$

as the probability for event E3's occurrence when ϵ is $\frac{\#(\text{DM} \cap \text{DP})}{q} - \frac{\delta}{q^2} \leq \frac{m}{q}$.

Finally, gathering the above information and argument, we can compute the probability for the new chain to terminate at an element of DT to be

$$\begin{aligned}&\sum_{k=0}^{\infty} \left(1 - \frac{\sqrt{1+2\alpha}}{t} + \epsilon\right)^k \left(\frac{\sqrt{1+2\alpha} - 1}{t} + \frac{\delta}{q}\right) \\ &= 1 - \frac{1 - t\epsilon}{\sqrt{1+2\alpha} + t\epsilon} + \frac{\alpha\sqrt{1+2\alpha} + \alpha\delta}{mt(\sqrt{1+2\alpha} + \epsilon)} \\ &\approx 1 - \frac{1}{\sqrt{1+2\alpha}}.\end{aligned}$$

When $\hat{t} = ct$ for some constant c , the error term of the last approximation is bounded by $2t\epsilon + \frac{5\alpha}{t} + \frac{\alpha}{\sqrt{1+2\alpha}}(c+1)\exp(-c) < \frac{7\alpha}{t} + \frac{\alpha}{\sqrt{1+2\alpha}}(c+1)\exp(-c)$.

Those familiar with time-memory tradeoff techniques can interpret this theorem as giving the probability of false alarms occurrence³ during the processing of a single non-perfect DP table. This high probability is an annoyance in the time-memory tradeoff. While every collision of the online DP chain with the pre-computed table will always bring about a solution to the DLP since the r -adding walk does not modify the exponent of h . Hence this high probability is a good thing for DLP solving.

Execution Complexity As given by Theorem 1, the mPAP succeeds using a single online DP chain with probability $1 - \frac{1}{\sqrt{1+2\alpha}}$.⁴ Taking the inverse of this value, we can state the expected number of online DP chain creations until successful DL retrieval to be $\frac{\sqrt{1+2\alpha}}{\sqrt{1+2\alpha}-1}$. Since

³ One should consider whether allowing the starting point to be the ending point will produce any difference in the final result.

⁴ Although we consider the online time complexity with the error term of probability in Theorem 1, that increases less than $\frac{7(\sqrt{1+2\alpha}+1)^3}{4\alpha} + \frac{(\sqrt{1+2\alpha}+1)^2}{2}(c+1)\exp(-c)t$, i.e., $T < \frac{\sqrt{1+2\alpha}}{\sqrt{1+2\alpha}-1} t + \frac{7(\sqrt{1+2\alpha}+1)^3}{4\alpha} + \frac{(\sqrt{1+2\alpha}+1)^2}{2}(c+1)\exp(-c)t$. This error comes from setting the chain length bound \hat{t} . However, since t is sufficiently large and $(c+1)\exp(-c)$ is sufficiently smaller than t , we ignore this error on our analysis.

the average length of a DP chain is t , we can state the expected online time complexity T , storage complexity⁵ M , and the pre-computation time complexity P as

$$T \approx \frac{\sqrt{1+2\alpha}}{\sqrt{1+2\alpha}-1} t, \quad M \approx m, \quad \text{and} \quad P = mt. \quad (6)$$

Tradeoff Curve Using the equation $mt^2 = \alpha q$, we obtain

$$PT \approx \frac{\alpha\sqrt{1+2\alpha}}{\sqrt{1+2\alpha}-1} q. \quad (7)$$

This shows that any increase in pre-computation is awarded by a corresponding decrease in online time. Interpretation of the following equations give us more detail.

$$\sqrt{M}T \approx \frac{\sqrt{\alpha}\sqrt{1+2\alpha}}{\sqrt{1+2\alpha}-1} \sqrt{q} \quad \text{and} \quad P \approx \sqrt{\alpha} \sqrt{M} \sqrt{q}. \quad (8)$$

With minimal storage, the expected pre-computation and online time are both $O(\sqrt{q})$. By utilizing a storage of size M , we can reduce online time by a factor of \sqrt{M} at the price of a \sqrt{M} factor increase in pre-computation time.

Note that for any fixed q , the right-side of tradeoff curve between storage and online time is optimal when the constant $\frac{\sqrt{\alpha}\sqrt{1+2\alpha}}{\sqrt{1+2\alpha}-1}$ is at its minimum. Explicitly, one can show that the minimum value of 2.35480 is attained at $\alpha = \frac{1+\sqrt{5}}{4} \approx 0.809017$.

3.3 Experiments

We have tested the complexity analysis of the **mPAP** by running it with small parameters. This was implemented on a dual-core AMD Opteron 2.6 GHz system and the NTL library [27] was used to provide the required finite field arithmetics. Throughout the test, the cyclic group $G = \langle g \rangle$ was taken to be a subgroup of \mathbf{Z}_p^* , where p was taken to be a random 1024-bit prime. The 20-adding walk served as the iterating function. As is customary in time memory tradeoff tables, sequential points g^1, g^2, \dots, g^m were used in place of the m random points to simplify implementation. Likewise, the first online chain started from DLP target h and sequential powers of h were used thereafter. The chain length bound was set to $10t$ to detect random walks that fell into loops without reaching a DP. Chains were not regenerated to replace the discarded chain.

In Table 1, we compare our theory against experiment results under variations of α . The success probability entries test the validity of Theorem 1 and the last two rows of the table test the T value as given by (6).

Table 2 shows a similar comparison under a fixed α and varying m and t parameters. This table gives an indication of how well the relation (8) is observed, i.e., whether the increase in storage results in the predicted reduction of online time complexity.

⁵ We are ignoring the fact that duplicate entries may be removed to reduce storage.

Table 1. Online phase success probability and complexity for various α (q : 42-bit prime, $t = 2^{14}$)

α		0.001	0.01	0.5	0.81	1	1.5	2	10
success probability at 1st chain (%)	experiment	0.10	1.30	28.10	37.50	38.10	48.10	54.10	76.90
	theory	0.10	0.99	29.29	38.22	42.26	50.00	55.28	78.18
success probability at 2nd chain (%)	experiment	0.00	0.70	21.50	24.10	25.80	24.80	23.80	17.40
	theory	0.10	0.98	20.71	23.61	24.40	25.00	24.72	17.06
average iterations until solution (unit: t)	experiment	1085.7	99.37	3.46	2.62	2.51	2.00	1.87	1.33
	theory	1001.5	101.50	3.41	2.62	2.37	2.00	1.81	1.28

Table 2. Online phase success probability and complexity for various m and t (q : 42 bits prime, $\alpha = 0.81$)

$\log t$	8	11	14	17	20	theory
m	46121039	586463	12564	207	2	
storage size (MB)	11806.986	150.135	3.216	0.053	0.001	
success probability (%)	1st chain	36.90	33.80	37.50	37.60	24.50
	2nd chain	23.00	25.50	24.10	22.30	16.40
ave. iterations until solution (unit: t)	2.70	2.65	2.62	2.78	15.30	2.62

Test results given above are average values done over multiple table creations and multiple DLP target solving per created table. Each pre-computation table on experiments was created with a different random multiplier set for the 20-adding walk. All tests used 10 tables and 100 DLPs per table.

The implementation results support our theoretic analysis very well for a wide range of parameters. The only visible exception corresponds to when m is extremely small. In this case the expected DP chain length approaches the expected rho length of a random walk and thus too many chains are being discarded. Unlike the small m case, the results for large m follows the theory closely. This implies that the mPAP will work well even when extremely large storage is employed.

3.4 Parallelization

The pre-computation phase of the mPAP can trivially be parallelized. While pre-computation is certainly the computationally intensive part of the mPAP, there may be situations where the parallelization of the online phase is desired, possibly to reduce the wall-clock running time of the online algorithm further. Let us investigate this matter in this subsection.

Readers familiar with time-memory tradeoff techniques will have read the previous material with $\alpha \approx 1$ in their minds, but one can confirm through a careful re-reading of all proofs that everything we wrote is true for even extremely small α .

With Theorem 1 confirmed for small α , let us consider the parallel use of n processors at the online phase. We take parameters m and t such that $mt^2 = \frac{\alpha}{n}q$, where $\alpha \approx 1$. Then, based on Theorem 1, one can state that the probability of success at the first run of the online phase, i.e., after the n DP chains have been produced, is $1 - (\frac{1}{1+2\alpha/n})^{n/2}$. Inverting this, we

can state

$$T_{1/n} \approx \left\{ 1 + \frac{1}{(1 + \frac{2\alpha}{n})^{\frac{n}{2}} - 1} \right\} t$$

as the average time spent by each processor, until the DLP is solved. As before, we require $M \approx m$ storage and $P = mt$ pre-computation time.

Applying the equation $mt^2 = \frac{\alpha}{n} q$, we can write the analogue to equation (7) as follows.

$$P T_{1/n} \approx \frac{\alpha}{n} \left\{ 1 + \frac{1}{(1 + \frac{2\alpha}{n})^{\frac{n}{2}} - 1} \right\} q. \quad (9)$$

It is easy to check that $1 + \frac{1}{e^\alpha - 1} \leq 1 + \frac{1}{(1+2\alpha/n)^{n/2}-1} \leq 1 + \frac{1}{\sqrt{1+2\alpha}-1}$, so that we can treat the term inclosed between the braces as an insignificant constant. Noting that $n T_{1/n}$ is the total online processing time, one can see that (9) is almost identical to (7). In other words, regardless of parallelization, with the same pre-computation effort, one needs the same total online processor time to solve a given DLP.

Analogue of (8) is as follows.

$$\sqrt{M} T_{1/n} \approx \sqrt{\frac{\alpha}{n}} \left\{ 1 + \frac{1}{(1 + \frac{2\alpha}{n})^{\frac{n}{2}} - 1} \right\} \sqrt{q} \quad (10)$$

and

$$P \approx \sqrt{\frac{\alpha}{n}} \sqrt{M} \sqrt{q}. \quad (11)$$

This says that with n -processor parallelization, one can achieve \sqrt{n} times online (wall-clock) time reduction and also an \sqrt{n} factor reduction in pre-computation. A more practical view is to say that, by deciding to invest n -times more on the online processing power, one reduces the required cost of the pre-computation phase. Since the pre-computation is much more resource consuming than the online phase, in most situations, the less-than-linear speedup of the online phase will be justified by the reduction of pre-computation cost. In addition, if the parameters were such that the storage cost outweighed the processor cost, then one can surely justify the cost of additional processors.

This brings us to the subject of what can be done while maintaining the pre-computation cost. In that case, one can read from (9) that we achieve linear speedup of the online phase, but one must be slightly more careful before jumping to this conclusion. The equation (11) shows that the storage must also be increased by a factor of n in addition to the number of processors. In other words, by increasing the storage and processor used during the online phase by a factor of n , one achieves n -times speedup of the online phase. If the parameters were such that the storage cost is much smaller than the processor cost, we may claim linear speedup through parallelization. If otherwise, this may not be what most would call linear speedup, but n -times investment on what is used during the online phase results in an n -times speedup of online phase.

3.5 Applying Tag Tracing Technique

In order to reduce the time for solving DLP over a finite field more, we consider applying tag tracing technique [7] to the mPAP. Since the mPAP consists of r -adding walk iterating function and DP collision detection method, the tag tracing technique can be applied to the mPAP well. In this subsection, we confirm that the tag tracing technique works well in the mPAP from experimental results.

Test environment and implementation method of the mPAP were the same with Section 3.3. For the tag tracing technique, we stored all possible products of multipliers of 4-adding walk up to 40 steps using pre-computation with 54.9 MB and 36.951 seconds on average of all tests. Other parameters were set equal to experiments of Section 4.5 in [7] for fair comparison.

Table 3. Comparison of 20-adding walk and tag tracing ($\alpha = 0.81$)

q	phase	20-adding walk	tag tracing	20-add./tag.	storage
42bit ($t=2^{14}$)	pre-computation	1032.221 s	77.995 s	13.234	5 MB
	online	0.256 s	0.027 s	9.487	
48bit ($t=2^{16}$)	pre-computation	17005.030 s	1274.148 s	13.346	17 MB
	online	1.038 s	0.105 s	9.886	
54bit ($t=2^{18}$)	pre-computation	210312.400 s	15439.350 s	13.622	68 MB
	online	4.115 s	0.419 s	9.821	
80bit* ($t=2^{30}$)	pre-computation	137 y 104.360 d	10 y 136.282 d		268 MB
	online	4 h 39.620 m	29 m 29.472 s		
80bit† ($t=2^{30}$)	pre-computation	1 y 136.097 d	37 d 20.707 h		27 GB
	online	2 m 47.772 s	17.695 s		

estimated value (*with 1 processor, †parallel processing with 100 processors)

Table 3 gives experimental results of applying 4-adding walk with tag tracing technique and 20 adding walk to the mPAP. All results are average values done over 10 table creations and 100 DLP targets solving per created table. Each table was produced with a different random multipliers for use of different r -adding walks.

We know that the use of 4-adding walk with the tag tracing technique is about 12 times faster than the use of 20-adding walk without the tag tracing technique from the result of [7] in our implementation environment. The last row in Table 3 shows that the tag tracing technique makes the mPAP 9.487 – 13.622 times faster. This shows that applying tag tracing technique to the mPAP works well.

Moreover, we also estimate the time to solve a DLP on a group of 80-bit order from experimental results. In Table 3, we confirm that the pre-computation time and the online time are about $2^{2d/3}$ and $2^{d/3}$ times increased respectively, when d is the difference between bit sizes of group orders. From this fact, we give the estimation of solving a DLP on a 80-bit order group using the mPAP with 20-adding walk iterating function and 4-adding walk with the tag tracing technique in the eighth and ninth rows in Table 3. The last two rows in

Table 3 also provide the estimation time to solve a DLP using the parallelized mPAP with 100 processors.

4 Key Generation Time Estimation

In this section, we propose a parameter setup method of the MY-IBE scheme for achieving 2^{80} security and efficiently generating a user's private key. Then we estimate the key generation time on proposed parameters when applying the mPAP to KeyGen algorithm based on our complexity analysis and experimental results on small parameters.

4.1 Parameter Setup

In the MY-IBE scheme, the base group G is taken to be a maximal cyclic subgroup of \mathbf{Z}_N^* where N is a product of roughly same sized primes p, q such that $p \equiv 3 \pmod{4}$, $q \equiv 1 \pmod{4}$, and $\gcd(p-1, q-1) = 2$. In order to obtain a secure MY-IBE scheme, as mentioned before, it is assumed that the CDH assumption in G holds and hence factoring N is to be hard.

Consider required conditions that N is hard to factor using known integer factorization algorithms within 2^{80} time complexity. First, N is to be at least 1024-bit integer (and hence both p, q are at least 512-bit primes) to endure against the number field sieve factorization algorithm [6]. Second, there are some factoring algorithms whose time complexity depends on the factorization of $p-1$ and $q-1$. In case of the Pollard $p-1$ algorithm [22], it takes $O(B \log N / \log B)$ group operations to factor N where $p-1$ and $q-1$ are B -smooth integers. Also, if factors of $p-1$ or $q-1$ consist of one ($\log B$)-bit prime and other $(\log B)/2$ -bit primes, then one can factor N within $O(\sqrt{B})$ operations using Brent's method [5]. Hence one has to set p, q to be primes such that both $p-1$ and $q-1$ are at least 2^{80} -smooth integers and have at least two prime factors of 80-bit or more.

However, the number of large prime factors of $p-1$ and $q-1$ has a significant effect on the key generation time. Let us look into a process of solving a DLP in KeyGen algorithm of the MY-IBE scheme, which applies Pohlig-Hellman algorithm [21] with the mPAP. When an identity ID is given to KGC, he tries to compute the DL of $H_1(\text{ID})$ to the base g . Since KGC knows the factorization of N as trapdoor information, he can apply the mPAP or the Pollard rho algorithm to obtain $H_1(\text{ID})^{c_i}$'s to the base g^{c_i} for all c_i 's where $c_i = p-1/p_i$ or $q-1/q_i$, $p-1 = \prod_{i=1}^{\ell} p_i$, and $q-1 = \prod_{i=1}^{\ell} q_i$, i.e., he solves DLPs over subgroups of G , whose order are p_i 's and q_i 's. Thereafter, he computes the DL of $H_1(\text{ID})$ to the base g using Chinese remainder theorem. In this process, large prime factors of $p-1$ and $q-1$ cause more key generation time or much memory size for the mPAP to reduce key generation time. Hence, we recommend the use of p, q such that both p, q are 512-bit primes and each $p-1, q-1$ has two 80-bit primes and other less than 40-bit primes.

4.2 Key Generation Time Estimation from Experimental Results

We shall give the estimated key generation time with experimental results applying the mPAP. Test environment was the same with Section 3.3. Throughout the test, G was taken to be

a maximal cyclic subgroup of \mathbf{Z}_N^* where N was an 1024-bit integer which was a product of 512-bit primes p, q and two prime factors of $p - 1$ and $q - 1$ were $(\log B)$ -bit and other prime factors were less than $(\log B)/2$ -bit each. We utilized the mPAP enhanced by the tag tracing technique [7] for solving DLPs on groups of four large prime factors and applied Pollard rho algorithm with 20-adding walk and DP collision detection method for solving DLPs on the other groups. Parameters for tag tracing technique were set equal to Section 3.5 and those for the mPAP were set to $t = 2^{(\log B)/3}$, $\alpha = 0.81$.

Table 4. Key generation time for various B

security level	pre-comp. size	pre-comp. time	online time			storage
			small factors	large factors	total	
42	$2^{29.9}$	3 m 25 s	0.266 s	0.079 s	0.345 s	17 MB
48	$2^{33.9}$	54 m 45 s	0.495 s	0.310 s	0.805 s	68 MB
80*	2^{51}	13 y 231 d	3 m 12.605 s	2 h 52.578 m	2 h 55.788 m	270 MB
80†	2^{51}	49 d 19 h	1.926 s	1 m 43.547 s	1 m 45.473 s	27 GB

estimated value (*with 1 processor, †parallel processing with 100 processors)

In Table 4, the third and fourth rows provide our experimental results of the key generation time corresponding to B , which are average values done over 10 table generations and 100 DLP targets solving per a created table. From these, when the difference between $\log B$'s is d -bit, we are sure that the pre-computation time and the online-time for large prime factors are about $2^{2d/3}$, $2^{d/3}$ times increased respectively and the online time for small prime factors is about $2^{d/4}$ times increased.

We estimate the pre-computation time and the online time for 2^{80} security. Assume that we perform $P = 2^{51}$ pre-computation for four large prime factors and store 2^{18} elements in each table. Then it takes about 13 years and 231 days for pre-computation with 270 MB storage. Since the pre-computation can be parallelized, one can generate the table in 49 days and 19 hours using 100 processors and 270 MB storage, in addition to the factors of n , as the trapdoor information. When an instance of DLP is given, it can be solved in 2 hours and 56 minutes on one processor.

If, as described in Section 3.4, we decide to parallelize the online phase across 100 processors while keeping the pre-computation at 2^{51} , the online phase reduces to one minute and 46 seconds with the use of 27 GB storage. Since this process is to be done by the key generation center to handle just key extractions, the presented pre-computation and online time may both be quite practical.

5 Conclusion

In this paper, we provided the method to reduce the key generation time of the MY-IBE scheme for practical use. To achieve this, we suggested the use of the mPAP with significant pre-computation for solving DLPs over a trapdoor DL group in the key generation algo-

rithm and gave more rigorous complexity analysis of the mPAP. Also we discussed about the parallelization of the mPAP and applying the tag tracing technique.

Finally, we gave the parameter setup method so that KGC efficiently solves a DLP and we estimated the key generation time on proposed parameter from our theoretical analysis and experimental results on small parameters.

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