Secure Poisson Regression

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Abstract

We introduce the first construction for secure two-party computation of Poisson regression, which enables two parties who hold shares of the input samples to learn only the resulting Poisson model while protecting the privacy of the inputs.

Our construction relies on new protocols for secure fixed-point exponentiation and correlated matrix multiplications. Our secure exponentiation construction avoids expensive bit decomposition and achieves orders of magnitude improvement in both online and offline costs over state of the art works. As a result, the dominant cost for our secure Poisson regression are matrix multiplications with one fixed matrix. We introduce a new technique, called correlated Beaver triples, which enables many such multiplications at the cost of roughly one matrix multiplication. This further brings down the cost of secure Poisson regression.

We implement our constructions and show their extreme efficiency. In a LAN setting, our secure exponentiation for 20-bit fractional precision takes less than 0.07ms with a batch-size of 100,000. One iteration of secure Poisson regression on a dataset with 10,000 samples with 1000 binary features needs about 65.82s in the offline phase, 55.14s in the online phase and 17MB total communication. For several real datasets this translates into training that takes seconds and only a couple of MB communication.

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1 Introduction

Privacy preserving computation technologies aspire to enable a wide range of modern computations used to analyze data, while providing strong privacy guarantees for the input data, which is often partitioned across multiple parties. Approaches based on cryptographic techniques for secure multiparty computation (MPC) have maintained the invariant of strong privacy guarantees while progressively supporting more complex functionality. In recent years, such approaches have taken on some of the most powerful available tools for data analysis which come from machine learning (ML). These tools bring functionalities with new levels of complexity to be supported in secure computation.

Existing MPC systems that support ML computations have mostly considered algorithms that aim to solve classification tasks, the most prominent of which are neural networks [4, 21, 26, 30]. In this work, we focus on a different type of computation: modeling Poisson processes. These processes are used to represent counts of rare independent events which happen at random but at a fixed rate. In such a process, the rate of events can be characterized by an underlying Poisson distribution. Poisson distributions are used to describe processes across many life and social sciences. Some examples include the number of bacteria over time in a petri dish, the number of mutations of a strand of DNA of a certain length, the number of losses and claims in insurance policies in a certain period of time, and the number of purchases a user makes after being shown online advertisements.

It is common to model response variables that follow the Poisson distribution by assuming their dependence on a set of explanatory (predictor) variables. Specifically, it is often assumed that the logarithm of the expected response is some linear combination of the explanatory variables. In this setting, the relationship between a response variable and the corresponding explanatory variables can be learned using Poisson regression. When the explanatory variables represent features which are conjectured to affect the counts, the regression model can be interpreted as uncovering the statistical significance of the effect of different features on the response variable. For example, Poisson regression has been used to model the dependence of the mortality rate from lung cancer on the age and smoking habits of people [15], the frequency at which voters engage in political discussion as a function of the method they use for voting (e.g., in person or by mail), their demographics, political affiliations, news exposure and others [27], the effect of age, gender, preexisting conditions such as diabetes and obesity on the mortality rate from COVID-19 [29], the number of payment defaults in credit scoring based on socio-economic characteristics [20], and the number of purchases that users make influenced by online advertisements they have been shown [28].

Traditionally, Poisson regression is performed by collecting all the examples (observed response variable counts together with the observed explanatory variable values), and performing training. However, in many of the above examples the information reflected in the predictor variables comes from different sources that hold health or financial data, which is highly sensitive information that is often subject to privacy regulations. Thus, while the final output model could be a useful tool for drawing insights about the underlying processes and events, providing the input data in the clear for the training is not an option. In this paper, we propose a solution that enables the computation while keeping all the inputs hidden from the parties performing the computation, revealing only the final Poisson regression model.

We introduce a system for secure computation that enables two parties who hold different parts of the training samples for Poisson regression to compute the final Poisson model. We assume the most general setting where the two parties hold cryptographic shares of the input training data, and obtain cryptographic shares of the resulting model. This representation can capture any partition of the input among the parties and also enables computation with the output model that does not reveal the model parameters to either party. Our new two party computation construction for Poisson regression leverages several new constructions for its building block components that offer improved efficiency. These functionalities have numerous uses beyond Poisson regression and thus are of independent interest as tools for secure computation.

Secure Exponentiation. A key component of Poisson regression involves the exponential (e^x) function. This step constitutes the nonlinear portion of Poisson regression and is not part of existing MPC implementations for ML functionalities. Nonlinear computations have traditionally been very challenging for secure computation techniques, since such techniques are generally better suited for evaluating linear functions or low-degree polynomials. Indeed, in existing MPC frameworks for ML functionalities [4, 21], the nonlinear components of the computation (e.g., the logistic function or the RELU function) are the core challenge that these works solve, and they contribute the most significant part of the cost of the final constructions. Adding to the challenge is the fact these nonlinear functions work on real numbers, which are quite difficult to support in MPC. Most approaches replace the nonlinear function with an approximation such as a low-degree polynomial or a piecewise linear function, which is easier to evaluate in MPC. However, such approximations could lead to significant degradation in the quality of the learned model (i.e., higher model error compared to training in the clear), and thus, the evaluation of the resulting constructions needs to consider jointly efficiency and accuracy.

In our work, we present a new construction for secure fixed-point exponentiation. It leverages a close approximation of the exact function with high precision that enables a significant efficiency improvement compared to existing constructions. In particular, all existing secure exponentiation approaches rely either on inaccurate polynomial approximations, or on bit decomposition of the exponent, which comes with a significant computation and communication cost. Our techniques avoid this multi-round computation step by leveraging ideas that enable the parties to obtain approximate multiplicative shares of the output only with local operations. We can control the accuracy and failure probability by appropriate parameter adjustment, only assuming knowledge of bounds on the input range. These bounds arise naturally in the context of Poisson regression. We introduce a new way to split the computation of the exponentiation into computation that depends only on the integer part of the exponent and computation that depends only on the fractional part of the exponent. Furthermore, we provide a novel way to combine the two computations with only local operations to obtain multiplicative shares of the output. Our only communication requirement is to transform the multiplicative shares of the output of the exponentiation into additive shares, which can be used for any further computation. For this, we leverage an existing protocol from Ghodosi et al. [16] that relies on a small amount of offline precomputation and a single round of online computation.

Since there are no prior works that consider (fixed-point) exponentiation in the two-party semihonest model, as comparison points, we consider state-of-the-art works that achieve a similar functionality in the malicious setting [9, 10] or in the semi-honest setting for $n \ge 3$ parties using a floating-point representation [7]. Although the comparison is not direct (see Section 7.1 for details), our protocol achieves orders of magnitude improvement on both the online throughput and the offline cost which indicates the possibility of substantial gains even when comparing in the same setting. In terms of accuracy, we can tune the parameters of our construction so that the output is arbitrarily close to the "true" exponentiation on the values in the clear without significant efficiency penalty (for example we can go from error 0.006% to error 0.0002% with 5 additional bits of precision). Our construction is so efficient that the nonlinear component of our Poisson regression protocol is no longer the cost bottleneck, and no longer degrades the quality of the computation, which stands in stark contrast to other works in the area of secure ML.

Optimized Secure Matrix Multiplication. Poisson regression makes extensive use of matrix multiplications. For secure multiplication on shared values, a well-known work [11] uses precomputed random *Beaver* triples followed by a single online communication round. In similar fashion, state of the art techniques for secure matrix multiplication [21] generalize Beaver triples to matrices and optimize the online communication and amount of preprocessing required; only one (matrix) Beaver triple is required for each matrix multiplication.

We make the observation that the matrix multiplication operations used in the Poisson regression training have a specific structure that can be exploited to further optimized the communication cost of the matrix multiplications: the same matrix \mathbf{X} is used in many multiplications with many different matrices \mathbf{Y}_i . While we can use independently generated Beaver triples for each multiplication, we show a more efficient way to precompute multiplication triples which takes advantage of the structure of the online matrix multiplications. We call these correlated Beaver triples, and they enable multiple online multiplications with the same matrix.

Using correlated Beaver triples results in improvements in the online phase f matrix multiplications: the communication cost is reduced by up to a factor of (n + 1) (where n is the number of training samples; see Section 5). Thanks to our very efficient secure exponentiation, the dominant cost (more than 90% for both computation and communication) in the secure Poisson regression protocol comes from secure matrix multiplication operations. Consequently, the use of correlated Beaver triples translates directly to a significant overall improvement of the cost of the whole secure Poisson regression protocol.

Experimental Results. We implemented all our constructions and provide detailed benchmarking. Our secure exponentiation protocol achieves significant efficiency improvements over existing approaches. Our implementation uses a 127-bit modulus for the computation field, which suffices for our Poisson regression evaluation. For this modulus, in a LAN setting with 1.5GB/s bandwidth, secure exponentiation for shared exponents with 20-bit precision takes less than 0.07ms when 100K evaluations are batched for communication. SCALE-MAMBA [9], which offers malicious security but is our most relevant point of comparison, uses a larger 245-bit modulus and a 40-bit precision, partly motivated by numerical instability for smaller sizes. Our construction does not have such instabilities and achieves online throughput that is 200x more efficient. The improvement in the offline phase is even greater, where our protocol requires 2000x less offline preprocessing and has a 500,000x improvement in offline computation.

We evaluate our secure Poisson regression implementation using three real datasets: Somoza's data on infant and child survival in Colombia, time to Ph.D. data, and data on the three-year survival status of breast-cancer patients [1]. We further evaluate the scalability of our system using larger synthetic datasets. The accuracy of our secure regression is essentially identical to that of plaintext computation of the regression. In our LAN setting, the total training (with 1000 iterations) for each of the three datasets takes less than 8s in the online phase, 120s in the offline phase, and 121MB total communication. The computation and communication overhead for our

construction scales roughly linearly with the size of the training data. For a dataset with 10,000 samples with 1000 binary features, and evaluation with a 127-bit modulus and 20-bit fractional precision, one training iteration requires 65.82s in the offline phase, 23.73s in the online phase and 17MB total communication. We also estimate the efficiency for secure Poisson regression for datasets used to predict COVID-19 case fatality rate, credit default rates and ad campaign conversion rates (see Section 8).

Paper organization. The rest of this paper is organized as follows. We introduce some standard background and preliminaries in Section 2. Our basic computation model and operations are provided in Section 3. Section 4 describes our full secure Poisson regression protocol; more details on our correlated Beaver triples and our secure fixed-point exponentiation protocol are given in Section 5 and Section 6 respectively. Section 7 provides a comprehensive set of experiments and benchmarks for our full regression protocol as well as for the correlated triple and exponentiation sub-components. Finally, in Section 8, we analyze the performance of our protocol for several concrete applications.

2 Preliminaries and Background

We start with some preliminaries and introduce standard background techniques on regression, and secure computation.

Basic notation. \mathbb{Z} denotes the integers and \mathbb{R} denotes the real numbers. \mathbb{Z}_N denotes the ring of integers modulo N. For a prime q, \mathbb{F}_q denotes the field with q elements, and \mathbb{F}_q^{\times} denotes its multiplicative group. We use bold uppercase letters (e.g., \mathbf{M}) to denote matrices and bold lowercase letters (e.g., \mathbf{u}, \mathbf{v}) to denote (row) vectors. Throughout the paper, e denotes Euler's constant. In some places, we abuse function notation slightly, and write $f(\mathbf{u})$ to denote the vector resultant from applying f to each element in \mathbf{u} separately.

2.1 Poisson Regression and Gradient Descent

In this section, we provide a brief overview of the gradient descent technique in the context of Poisson regression.

Poisson regression. Regression is a common statistical technique to learn a function $g(\mathbf{x}_i) \approx y_i$, given *n* training samples \mathbf{x}_i (each with *m* features), and corresponding output labels y_i . Different forms of regression model different classes of functions *g*. For example, machine learning has extensively used linear regression (to model linear outputs) and logistic regression (to model binary outputs).

When the response variable y is count or rate-based (rather than continuous), using Poisson regression makes more sense. For Poisson regression, the expected response is modeled as a Poisson distribution, and therefore, $g(\mathbf{x}_i) = e^{\langle \boldsymbol{\theta}, \mathbf{x}_i \rangle}$, where $\boldsymbol{\theta}$ is the coefficient or weights vector, and $\langle \cdot, \cdot \rangle$ is the dot product. Rate-data can be modeled by an extra multiplicative factor t_i denoting the time "exposure" for each sample over which the response variable was computed.

Gradient descent. Gradient descent is a standard machine learning technique used to train a model iteratively. A model can be defined by a set of parameters $\boldsymbol{\theta} = (\theta_1, \dots, \theta_m)$. To learn the model parameters from data, the algorithm iteratively attempts to minimize a predetermined convex function. At each step, the parameters are updated based on the gradient.

In this paper, we focus specifically on Poisson regression with *exposure*, which allows for modeling of rate-based data. For this, training data is provided as $(\mathbf{X}, \mathbf{Y}, \mathbf{T})$, where $\mathbf{X} \in \mathbb{R}^{n \times m}$ contains data for the explanatory variables, $\mathbf{Y} \in (\mathbb{R}^+)^{n \times 1}$ contains data for the response variable, and $\mathbf{T} \in (\mathbb{R}^+)^{n \times 1}$ is the exposure data. n is the number of training samples and m is the number of features (or explanatory variables). Henceforth, unless specified, we use Poisson regression with exposure. Poisson regression attempts to learn model parameters $\boldsymbol{\theta}$, by minimizing $-L(\boldsymbol{\theta}|\mathbf{X},\mathbf{Y},\mathbf{T})$ where $L(\cdot)$ is the log likelihood function. For this, the gradient will be computed as:

$$\frac{\partial L(\boldsymbol{\theta}|\mathbf{X},\mathbf{Y},\mathbf{T})}{\partial \boldsymbol{\theta}} = \sum_{i=1}^{n} \mathbf{x}_{i} (y_{i} - t_{i} e^{\langle \boldsymbol{\theta}, \mathbf{x}_{i} \rangle})$$

where (\mathbf{x}_i, y_i, t_i) is the *i*th data point. The training will now update the parameters iteratively. For the (k+1)th iteration, $\boldsymbol{\theta}^{(k+1)}$ is computed as follows:

$$\boldsymbol{\theta}^{(k+1)} = (1-\beta)\boldsymbol{\theta}^{(k)} + \alpha \sum_{i=1}^{n} \mathbf{x}_{i}(y_{i} - t_{i} \cdot e^{\langle \boldsymbol{\theta}, \mathbf{x}_{i} \rangle})$$
$$= (1-\beta)\boldsymbol{\theta}^{(k)} + \alpha \mathbf{X}^{T} \left(\mathbf{Y} - \mathbf{T} \circ e^{\mathbf{X}\boldsymbol{\theta}^{(k)}} \right)$$

where the exponential function is applied to each element in $\mathbf{X}\boldsymbol{\theta}^{(k)}$, \circ is the Hadamard (elementwise) product, and the constants α and β denote the learning rate and the regularization parameter respectively. $\boldsymbol{\theta}^{(0)}$ is usually initialized either as the zero vector, or with random weights.

2.2 Secure Computation Functionalities

Secure computation protocols enable functionalities where parties can compute a function on their joint private inputs in a way that is guaranteed only the output of the computation. Our protocol constructions are in a two-party setting and provide semi-honest security [17], i.e., the parties are assumed to follow the prescribed protocol. We denote the two parties by P_0 and P_1 . We use $[\![x]\!]^{\mathbb{Z}_N}$ to denote an (additive) sharing of x over \mathbb{Z}_N . We drop the superscript when it is clear from context. We write $[\![x]\!] = ([\![x]\!]_0, [\![x]\!]_1)$ where P_0 holds $[\![x]\!]_0$ and P_1 holds $[\![x]\!]_1$ such that $[\![x]\!]_0 + [\![x]\!]_1 = x \mod N$. The sharing is chosen randomly, for example by first choosing $[\![x]\!]_0$ uniformly at random in \mathbb{Z}_N and then assigning $[\![x]\!]_1 = x - [\![x]\!]_0 \mod N$.

We use the notation $F(\llbracket x \rrbracket, \llbracket y \rrbracket)$ to denote that the parties P_0 and P_1 engage in a computation of some functionality F, with P_0 contributing $\llbracket x \rrbracket_0$ and $\llbracket y \rrbracket_0$ as input, and P_1 contributing $\llbracket x \rrbracket_1$ and $\llbracket y \rrbracket_1$ as input, with each party receiving its corresponding secret share of the result as output.

Next we overview some secure computation techniques that we use in our protocols.

Multiplication using Beaver triples. Suppose that P_0 and P_1 are given shares $[\![x]\!]$, and $[\![y]\!]$, over \mathbb{Z}_N . To compute $[\![z]\!] = [\![xy]\!]$, a common technique in the preprocessing model is to use Beaver's multiplication trick [11]. For this, a randomly sampled tuple $([\![a]\!], [\![b]\!], [\![c]\!])$, also called a "Beaver triple", such that $c = ab \mod N$ is provided to the two parties. Now, to compute $[\![z]\!]$, P_0

and P₁ start by locally computing $\llbracket u \rrbracket = \llbracket x \rrbracket - \llbracket a \rrbracket$ and $\llbracket v \rrbracket = \llbracket y \rrbracket - \llbracket b \rrbracket$. Next, they reconstruct u and v by communicating their share to the other party. Finally, P_i can compute its share $\llbracket z \rrbracket_i = i \cdot uv + u \llbracket b \rrbracket_i + v \llbracket a \rrbracket_i + \llbracket c \rrbracket_i$. This works since $\llbracket z \rrbracket_0 + \llbracket z \rrbracket_1 = (x-a)(y-b) + (x-a)(b) + (y-b)(a) + c = xy - ab + c = xy \mod N$. Note that the same technique works for multiplying fixed-point numbers. A different Beaver triple is needed for each secure multiplication to be performed in the protocol. Each secure multiplication needs a preprocessing of 3 ring elements per party, and has an online communication of 2 elements per party. We use $\mathcal{F}_{\mathsf{mult}}(\llbracket x \rrbracket, \llbracket y \rrbracket)$ to denote executing the secure multiplication functionality.

Vectorization for Beaver triples. Beaver triples also work for secure matrix multiplication, where the two matrices to be multiplied are secret shared between P_0 and P_1 [21]. Consider a matrix multiplication between an $n \times m$ size matrix **X** and an $m \times k$ size matrix **Y**. Naively, the matrix multiplication **XY** requires nmk multiplications, and therefore nmk Beaver triples would be necessary. However, as [21] notes, this can be optimized by sharing a matrix Beaver triple ([[A]], [[B]]], [[C]]), where **A** and **B** are matrices with the same dimension as **X** and **Y** respectively, and **C** = **AB** mod N. Here, the preprocessing (using Ring-LWE) and online costs per party are nk + mk and nm + mk ring elements respectively. We use $\mathcal{F}_{matMult}([[X]], [[Y]])$ to denote executing the secure matrix multiplication functionality.

We use standard Ring-LWE based techniques to generate the Beaver triples, and compress the real number of bits required for preprocessing and communication. We provide a background on Ring-LWE in Section 2.2.1.

2.2.1 Ring-LWE-based Encryption

Ring-Learning-With-Errors [19] (RLWE) is a hardness assumption based on which efficient homomorphic encryption schemes have been constructed. We use the additive homomorphic properties of the encryption scheme proposed by Brakersky et al. [12], based on RLWE, to generate the Beaver triples in our preprocessing phase. For a positive integer N, the scheme is defined over the ring $R = \mathbb{Z}[X]/\Phi_N(X)$ where $\Phi_N(X)$ is an N^{th} cyclotomic polynomial of degree $\phi(N)$ ($\phi(\cdot)$ is the Euler's totient function). We define the ring $R_t = R/tR$, and use p, q to denote the plaintext and ciphertext modulus respectively. Choosing p and q carefully allows us to pack $\phi(N)$ plaintexts $(m_1, \dots, m_{\phi(N)})$ into a single ring element $m \in R_p$ and enables SIMD operations (addition, multiplication) over the packed plaintexts. Packing is used to generate preprocessing data for our secure exponentiation protocol.

Basic definition. We now describe the operations for Ring-LWE-based encryption in the twoparty setting.

- Key Generation. One party samples a key pair $(\mathsf{sk}, \mathsf{pk})$ such that $\mathsf{sk} = (1, -s)$, where $s \in R$ with coefficients in $\{-1, 0, 1\}$ and s has low Hamming weight (e.g., $\mathrm{H}(s) = 64$) and $\mathsf{pk} = (a, b)$, where $a \leftarrow R_q$ and $b = as + t\epsilon \in R_q$ with ϵ drawn from a small noise distribution χ .
- Encryption. Given a packed plaintext $m \in R_p$, its fresh ciphertext can be given by (c_0, c_1) where $c_0 = m + bv + p\epsilon_0$ and $c_1 = av + p\epsilon_1$ (where $v, \epsilon_0, \epsilon_1$ are drawn from the noise distribution).

- Decryption. The party that holds the secret key can decrypt the ciphertext to recover the underlying plaintext. Given a ciphertext $c \equiv (c_0, c_1) \in R_q^2$, the plaintext can be computed as $\text{Dec}_{sk}(c) = c_0 + c_1 s \mod p$.
- Plaintext addition. Given a ciphertext $c \equiv (c_0, c_1) = \text{Enc}(m) \in R_q^2$ and a plaintext message $m' \in R_p$, one can produce the encryption of (m + m') as $c' = (c_0 + m', c_1) = \text{Enc}(m + m')$.
- Scaling. Given a ciphertext $c \equiv (c_0, c_1) = \operatorname{Enc}(m) \in R_q^2$ and a scalar $a \in \mathbb{Z}_p$, one can produce the encryption of $am = (am_1, \cdots, am_N)$ as $c' = (a \cdot c_0, a \cdot c_1) = \operatorname{Enc}(a \cdot m)$.
- Shifting. Given a ciphertext $c \equiv (c_0, c_1) = \operatorname{Enc}(m) \in R_q^2$ where $m = (m_1, \dots, m_K, 0, \dots, 0)$, we can produce the encryption of $m' = (0, \dots, 0, m_1, \dots, m_K, 0, \dots, 0)$ where m_1 is shifted by a distance t and $t + K \leq N$. Let $v = (0, \dots, 0, 1, 0, \dots, 0)$ where v is zero everywhere except for the t^{th} position. Then $c' = (v \circ c_0, v \circ c_1) = \operatorname{Enc}(m \circ v) = \operatorname{Enc}(m')$ where \circ denotes polynomial multiplication operation in $R_q = \mathbb{Z}_q[X]/\Phi_N(X)$.

Choosing parameters for Ring-LWE. Following the parameters suggested by [6], we use a ciphertext prime q = 160 bits for our RLWE scheme when generating Beaver triples for ring of size p = 63 bits ($\mathbb{Z}_{2^{63}}$) and q = 295 bits when generating triples for ring of size p = 127 bits ($\mathbb{Z}_{2^{127}}$). In the first case, we use a polynomial modulus of degree $2^{12} = 4096$, while for the later case $2^{14} = 16384$. This is sufficient for security of at least 128 bits. To allow efficient encryption and decryption via the use of number theoretic transform, we choose a ciphertext modulus q such that $q \equiv 1 \mod 2N$.

3 Secure Computation over Fixed-Point Rings

Poisson regression operates over the real numbers. When the computation is done in the clear, one can leverage floating point representation to achieve high precision. For secure computation, while there are techniques that emulate floating point representation [7], they are often expensive. A more efficient approach that is commonly used is to adapt the actual computation to work with a fixed-point representation while preserving accuracy. We adopt this approach in our work as well and similarly to other works [21], we will compute over fixed-point numbers mapped onto an integer ring.

We start by defining a fixed-point ring that will be used to represent our fixed-point numbers.

Fixed-point ring. A fixed-point ring \mathcal{R} is a tuple $(\mathbb{Z}_{2^l}, l_x, l_f)$ where \mathbb{Z}_{2^l} is the ring of integers modulo 2^l , and l_x, l_f are positive integers with $l_f \leq l_x \leq l-1$. \mathcal{R} will be used to represent fixed-point numbers with at most l_f (binary) fractional bits, and whose absolute value is less than $2^{l_x-l_f}$. Non-negative numbers will be in the range $\mathcal{R}^+_* = [0, 2^{l_x})$ and negative numbers will be in the range $\mathcal{R}^-_*(2^l - 2^{l_x}, 2^l)$ in their two's complement representation. $\mathcal{R}_* = \mathcal{R}^+_* \cup \mathcal{R}^-_*$ is the total part of \mathcal{R} wherein the fixed-point numbers are represented.

For a real number r, with $|r| < 2^{l_x - l_f}$, we will use the hat operator, as in \hat{r} , to denote its representation in the ring \mathcal{R} . Note that $\hat{r} = \lfloor 2^{l_f} \cdot r \rfloor$ when $r \ge 0$ and $\hat{r} = 2^l - \lfloor 2^{l_f} \cdot |r| \rfloor$ when r < 0. For example, in $\mathcal{R} = (\mathbb{Z}_{2^{10}}, 3, 2)$, a real number x = 1.25 will be represented in \mathcal{R} by $\hat{x} = \lfloor 2^2 \cdot 1.25 \rfloor = 5$, and y = -1.25 will be represented by $\hat{y} = 2^{10} - \lfloor 2^2 \cdot 1.25 \rfloor = 1019$. Note that something like z = 1.26 will also be represented by $\hat{z} = 5$ due to truncation.

Similarly, for a ring element $x \in \mathcal{R}_*$, i.e., $x \in [0, 2^{l_x}) \cup (2^l - 2^{l_x}, 2^l)$, we will use the under-tilde operator, as in \underline{x} , to denote its canonical real number representation. By canonical, we mean the real number which gives no truncation error when represented in the ring. For instance, in the previous example, $\underline{5} = 1.25$ and not 1.26.

Secure operations. We define secure arithmetic operations on values that have been secret shared between the two parties, P_0 and P_1 , in our protocol. We distinguish between two types of operations: (1) Basic ring operations are operations over shares in the ring \mathbb{Z}_{2^l} treating elements as integers; (2) Fixed-point or FP operations, on the other hand, are operations that manipulate shares in the ring \mathbb{Z}_{2^l} , treating the underlying elements as fixed-point numbers. For a given $\mathcal{R} = (\mathbb{Z}_{2^l}, l_x, l_f)$, we will use $[\![x]\!]^{\mathcal{R}}$ or $[\![x]\!]^{\mathbb{Z}_{2^l}}$ to denote additive shares of $x \in \mathbb{Z}_{2^l}$. P₀ and P₁ will hold the shares $[\![x]\!]_0$ and $[\![x]\!]_1$ respectively. We will drop the superscript \mathcal{R} when it is clear from context. With this notation, we now define some basic useful secure ring operations.

- 1. Basic operations:
 - (Addition Add). Given shared values [x] and [y], Add([x], [y]) outputs [x + y].
 - (Multiplication Mult). Given shared values [x] and [y], Mult([x], [y]) outputs [xy].

Addition can be done non-interactively by each party locally adding its shares modulo 2^{l} . Multiplication is modulo 2^{l} and can be done in one interactive round using Beaver triples.

- 2. Fixed-Point operations: These operations are for elements in \mathcal{R}_* . Intuitively, the functionality here can be thought of as first retrieving the real numbers corresponding to the ring elements (using the under-tilde operator), then computing the result in real numbers, and finally casting back into the fixed-point ring (using the hat operator).
 - (FP Addition FPAdd). Given shared values $\llbracket x \rrbracket$ and $\llbracket y \rrbracket$, FPAdd($\llbracket x \rrbracket, \llbracket y \rrbracket$) outputs $\llbracket (\widehat{x}) + (\widehat{y}) \rrbracket$.
 - (Public FP Multiplication PubFPMult). Given [x] and a public element $c \in \mathcal{R}_*$, PubFPMult([x], c) outputs $[\widehat{(c)(x)}]$.
 - (FP Multiplication FPMult). Given shared values $\llbracket x \rrbracket$ and $\llbracket y \rrbracket$, FPMult($\llbracket x \rrbracket$, $\llbracket y \rrbracket$) outputs $\llbracket \widehat{(\underline{x})(\underline{y})} \rrbracket$.
 - (Public FP Division PubFPDiv). Given [x] and a public positive integer $c \in \mathbb{Z}^+$, PubFPDiv([x], c) outputs $[\widehat{(x)/c}]$.
 - (FP Exponentiation FPExp). Given a public positive base element $b \in [0, 2^{l_x})$, and a shared exponent $[\![x]\!]$, FPExp $(b, [\![x]\!]$) outputs $[\![(\widehat{b})^{(\underline{x})}]\!]$.

Note that the basic addition and multiplication operations are over \mathbb{Z}_{2^l} but for FP operations, they are over reals. It is easy to see though that Add and FPAdd provide the same functionality when the underlying shares represent valid fixed-point elements. To avoid overflow for FP operations, we will require that the underlying real numbers represented by any FP operation will still be smaller

in absolute value than the $2^{l_x-l_f}$. In practice, this can be done be choosing a large enough ring to handle the range of values necessary for any computation.

Similar to the basic operations, FPAdd can be done non-interactively, and FPMult can be done using Beaver triples. Due to truncation, FPMult can have an error of at most 2^{-l_f} in the underlying computation. Public fixed-point multiplication and division can both be done non-interactively with an error of at most 2^{-l_f} , and we provide protocols to do so in Section 3.1. The non-interactive protocols also have an associated failure probability, which as similar to e.g., [21] can be made arbitrarily small by increasing the gap between l and l_x (see Section 3.1). The exponentiation protocol is a novel contribution of our paper and we provide the full details in Section 6.

We can also use a prime modulus q for our fixed-pointring (instead of 2^l), embed fixed-point numbers into $[0, 2^{l_x}) \cup (q - 2^{l_x}, q)$ in \mathbb{F}_q , and define all of the above operations similarly over \mathbb{F}_q .

Failure probability and approximation errors. The secure computation of FP operations may come inbuilt with some probability of failure as well as errors as a result of truncation. We say that a protocol has failure probability p_{fail} and error ϵ if, except with probability p_{fail} , the error in the underlying fixed-point computation is bounded by ϵ . The failure probability, similar to e.g., [21] can be made arbitrarily small by increasing the gap between l and l_x .

Ring change. A final useful operation we introduce is to switch between rings with different moduli. Given N and N', and a shared value $[\![x]\!]^{\mathbb{Z}_N}$, the operation $\operatorname{RingChange}([\![x]\!]^{\mathbb{Z}_N}, \mathbb{Z}_{N'})$ will output $[\![x]\!]^{\mathbb{Z}_{N'}}$, a sharing of $x \pmod{N'}$ in $\mathbb{Z}_{N'}$. We will only require the operation for N' > N and when x is small $(x < 2^{l_x})$ which allows us to do this without any interaction. We detail a non-interactive protocol for this in Section 3.1.

3.1 Detailed Secure Functionalities

We provide more details on the public fixed-point multiplication, and division functionalities, as well as the RingChange operation, and RLWE encryption.

Public fixed-point division. Consider a sharing $[\![x]\!]$ over \mathcal{R}^+_* with modulus N, and a public positive divisor $c \in \mathbb{Z}^+$. Except with probability $2^{l_x}/N$, the sharing is such that $[\![x]\!]_0 + [\![x]\!]_1 = x + N$. Now, to compute the fixed-point division by c, P_0 computes $[\![z]\!]_0 = N - \left\lfloor \frac{N - [\![x]\!]_0}{c} \right\rfloor$ and P_1 computes $[\![z]\!]_1 = \left\lfloor \frac{[\![x]\!]_1}{c} \right\rfloor$. Notice now that, $\frac{x}{c} + N - 1 \leq [\![z]\!]_0 + [\![z]\!]_1 \leq \frac{x}{c} + N + 1$. Therefore, $([\![z]\!]_0, [\![z]\!]_1)$ is a sharing of the representation of \underline{x}/c in \mathcal{R} , with an error at most 2^{-l_f} . Note that while our regression protocol does not require it, a public negative divisor can easily be handled by first dividing by the absolute value and then subtracting the shares from the modulus N.

Public fixed-point multiplication. Consider a sharing $\llbracket x \rrbracket$ over \mathcal{R}_* with modulus N, and a positive public element $c \in \mathcal{R}_*^+$. Let $\llbracket x \rrbracket = (r, (x - r) \mod N)$. Let $\llbracket z \rrbracket_0 = \left\lfloor \frac{c \llbracket x \rrbracket_0 - cN}{2^{l_f}} \right\rfloor \mod N$ and $\llbracket z \rrbracket_1 = \left\lfloor \frac{c \llbracket x \rrbracket_1}{2^{l_f}} \right\rfloor \mod N$. Let $\frac{c \llbracket x \rrbracket_0 - cN}{2^{l_f}} = w_0 - d_0$, and $\frac{c \llbracket x \rrbracket_1}{2^{l_f}} = w_1 + d_1$, where w_i are the integer parts and $0 \le d_i < 1$ are the fractional parts. Note the negative sign on d_0 since $N > \llbracket x \rrbracket_0$. We show that $(\llbracket z \rrbracket_0, \llbracket z \rrbracket_1)$ form a sharing of $(\widehat{c})(\widehat{x})$. Recall that this is $\frac{cx}{2^{l_f}}$ when $x \in \mathcal{R}_*^+$ and $N - \frac{c(N-x)}{2^{l_f}}$ when $x \in \mathcal{R}_*^-$.

Case 1) $x \in \mathcal{R}^+_*$. Then, when $r \in [2^{l_x}, N)$, the sharing is such that $[\![x]\!]_0 + [\![x]\!]_1 = x + N$. Now, $[\![z]\!]_0 + [\![z]\!]_1 \mod N \equiv w_0 + w_1 \equiv (w_0 - d_0 + w_1 + d_1) + (d_0 - d_1) \equiv \frac{cx}{2^{l_f}} + (d_0 - d_1)$. Therefore, $(cx)/2^{l_f} - 1 \leq [\![z]\!]_0 + [\![z]\!]_1 \mod N \leq (cx)/2^{l_f} + 1$.

Case 2) $x \in \mathcal{R}_*^-$. Then, when $r \in [0, N - 2^{l_x}]$, the sharing is such that $[\![x]\!]_0 + [\![x]\!]_1 = x$ (without the modulo). Now, $[\![z]\!]_0 + [\![z]\!]_1 \mod N \equiv w_0 + w_1 \equiv (w_0 - d_0 + w_1 + d_1) + (d_0 - d_1) \equiv \frac{-c(N-x)}{2^{l_f}} + (d_0 - d_1)$. Therefore, $(N - \frac{c(N-x)}{2^{l_f}}) - 1 \leq [\![z]\!]_0 + [\![z]\!]_1 \mod N \leq (N - \frac{c(N-x)}{2^{l_f}}) + 1$.

Consequently, when $r \in [2^{l_x}, N - 2^{l_x}]$, i.e., except with probability less than 2^{l_x+1} , this results in a sharing of the representation of $(\underline{x})(\underline{c})$, with an error of at most 2^{-l_f} . Note that a negative ccan also be handled analogously to PubFPDiv.

Ring change. We only require the RingChange operation to switch rings between \mathbb{Z}_N and $\mathbb{Z}_{N'}$ where N' > N, and only for positive fixed-point numbers. Consider a random sharing of $x \in [0, 2^{l_x})$ in \mathbb{Z}_N and denote the two shares by $[\![x]\!]_0 = r$ and $[\![x]\!]_1 = x - r \mod N$. Note that when $r \in [0, 2^{l_x})$, $[\![x]\!]_0 + [\![x]\!]_1 = x$ (even without a mod N). For any other r, $[\![x]\!]_0 + [\![x]\!]_1 = x + N$. This means that for a random sharing of x, the addition "wraps around" N with probability $1 - \frac{2^{l_x}}{N}$. Now, if we set $[\![x]\!]_0^{\mathbb{Z}_{N'}} = [\![x]\!]_0^{\mathbb{Z}_N} + N' - N$ and $[\![x]\!]_1^{\mathbb{Z}_{N'}} = [\![x]\!]_1^{\mathbb{Z}_N}$, then $([\![x]\!]_0^{\mathbb{Z}_{N'}}, [\![x]\!]_1^{\mathbb{Z}_{N'}})$ forms a sharing of x in $\mathbb{Z}_{N'}$ and wraps around N'. Consequently, except for a failure probability of at most $2^{l_x}/N$, the above protocol switches the sharing of x from \mathbb{Z}_N to $\mathbb{Z}_{N'}$ with no error.

If necessary, the range for both the shares of both parties can be expanded to all of $\mathbb{Z}_{N'}$ by using a PRG. Specifically, both parties can agree on a PRG G the outputs values in $\mathbb{Z}_{N'}$, and a seed $g_0 = s$. For the j^{th} RingChange, they can compute the next PRG value g_j . Then P₀ adds g_j modulo N' to its share, and P₁ subtracts g_j modulo N' to its share.

4 Secure Poisson Regression Protocol

We detail our full secure Poisson regression protocol in this section.

Protocol input. Recall that for Poisson regression (with exposure), each of the *n* training samples is of the form (\mathbf{x}_i, t_i, y_i) where \mathbf{x}_i contains *m* features, t_i is the exposure value, and y_i is the response output. We use **X** to denote the $n \times m$ matrix of training samples, **T** to denote the $n \times 1$ vector of exposures, and **Y** to denote the $n \times 1$ vector of response values. We assume that all entries are already represented as fixed-point elements and shared between the two protocol parties. We use $[\mathbf{X}], [[\mathbf{Y}]], [[\mathbf{T}]]$ to denote the sharings.

Protocol parameters. Prior to the protocol, we require P_0 and P_1 to agree on the following parameters: (1) A fixed-point ring $\mathcal{R} = (\mathbb{Z}_{2^l}, l_x, l_f)$; (2) An *l* bit prime *q*, and an exponent bound (for the exponentiation protocol); (3) The regression parameters α (learning rate), β (regularization term), and the number of iterations *K*.

4.1 Basic Design

The goal of the regression protocol is to output a sharing of a weights vector $\boldsymbol{\theta}$. For this, we use gradient descent, which updates the weights at every iteration. Three variants are commonly used,

which differ in the way the weights are updated: (1) Standard, where the entire dataset is used for each iteration; (2) Mini-batch, where a small random sample is used for each iteration; and (3) Stochastic, where a single random sample is used for each iteration. For this paper, we use the standard gradient descent for our secure Poisson regression, but note that our protocol can be adapted for any variant. We discuss the alternative mini-batch variant in Section 4.3.

Recall that in the update step of our gradient descent, the weights for the k^{th} iteration are updated as follows:

$$\boldsymbol{\theta}^{(k+1)} = (1-\beta)\boldsymbol{\theta}^{(k)} + \alpha \mathbf{X}^T \left(\mathbf{Y} - \mathbf{T} \circ e^{\mathbf{X}\boldsymbol{\theta}^{(k)}} \right)$$

Let $\llbracket \boldsymbol{\theta}^{(k)} \rrbracket$ denote a sharing of the weights vector after the k^{th} iteration. Parties start with $\llbracket \boldsymbol{\theta}^{(0)} \rrbracket$ initialized randomly or as shares of 0. Now, each iteration of our regression proceeds as follows: (1) First, P₀ and P₁ compute the (fixed-point) matrix multiplication $\llbracket \mathbf{U} \rrbracket = \llbracket \mathbf{X} \boldsymbol{\theta}^{(k)} \rrbracket$. (2) Next, each element in \mathbf{U} is exponentiated (*n* exponentiations in total). Let $\llbracket \mathbf{V} \rrbracket$ be the sharing of the result after each term in $\llbracket \mathbf{U} \rrbracket$ is exponentiated; (3) Then, P₀ and P₁ compute an element-wise product $\llbracket \mathbf{W} \rrbracket = \llbracket \mathbf{T} \circ \mathbf{V} \rrbracket$; (4) Next, P₀ and P₁ compute the (fixed-point) matrix multiplication by α), and addition by $(1 - \beta) \boldsymbol{\theta}^{(k)}$ can be computed locally, to end up with shares of the updated weights $\boldsymbol{\theta}^{(k+1)}$. Our protocol requires 4 rounds, one for each of the first four steps. Figure 1 contains a detailed description of our protocol. The element-wise product and matrix multiplications, can be computed using the functionality $\mathcal{F}_{\mathsf{mult}}$ and $\mathcal{F}_{\mathsf{matMult}}$ respectively, and implemented using matrix Beaver triples as preprocessing. The fixed-point exponentiations are computed using the functionality $\mathcal{F}_{\mathsf{FPexp}}$, which we describe in detail in Section 6.

Basic protocol cost. From the previous description, we note that each gradient descent iteration computes 2 matrix computations (of sizes $(n \times m, m \times 1)$ and $(m \times n, n \times 1)$, 1 element-wise product for n size vectors, and n secure exponentiations. By using the matrix Beaver triples optimization from [21], a total of 2nm+n triples are enough in the preprocessing stage (per iteration). In addition to this, we utilize further optimizations for batched multiplication that substantially improve the performance of our protocol, when amortized over multiple iterations. Our key observation for this optimization is that the matrix multiplications in each iteration have \mathbf{X} , or \mathbf{X}^T as one of the multiplicands. In other words, for K iterations, we have K multiplications of the form (\mathbf{X}, \cdot) and K of the form (\mathbf{X}^T, \cdot) . This allows us to batch together the multiplications in separate iterations using correlated randomness where one of the matrices in the Beaver triple is reused. We detail this optimization in Section 5, and show that it does not leak any extra information about the multiplicands.

The *n* secure exponentiations in each iteration require a total preprocessing of 2n field elements per party, and a communication of *n* field elements per party (see Section 6). Note that all of the exponentiations are independent and can be done in parallel in a single round.

Failure probability. The fixed-point multiplication, and exponentiation operations have a small failure probability, which depends on the chosen parameters. We compute the overall failure probability for our regression protocol, which will be helpful to choose appropriate parameters for a given acceptable failure probability.

Secure Poisson Regression

Setup. P_0 and P_1 agree on a fixed-point ring \mathcal{R} , a prime q, and parameters for the Poisson regression: learning rate α , a regularization term β , number of iterations K. Input. Two parties have shares $([\mathbf{X}]_i, [\mathbf{Y}]_i, [\mathbf{T}]_i)$ over \mathcal{R} . $\mathbf{X} \in \mathcal{R}^{n \times m}$ is the feature matrix where n is the number of samples and m is the number of explanatory variables, $\mathbf{Y} \in \mathcal{R}^n$ is the label vector, $\mathbf{T} \in \mathcal{R}^n$ is the exposure vector.

Protocol.

- 1. Both parties initialize shares $\llbracket \boldsymbol{\theta}^{(0)} \rrbracket$ to 0^m .
- 2. For k = 1 to K do:
 - (a) The parties make a call to $\mathcal{F}_{matMult}$, and set $\llbracket \mathbf{U} \rrbracket \leftarrow \mathcal{F}_{matMult}(\llbracket \mathbf{X} \rrbracket, \llbracket \boldsymbol{\theta}^{(k-1)} \rrbracket)$.
 - (b) The parties make a call to $\mathcal{F}_{\mathsf{FPexp}}$ on each element of $\llbracket \mathbf{U} \rrbracket$. Let $\llbracket \mathbf{V} \rrbracket \leftarrow \mathcal{F}_{\mathsf{FPexp}}(e, \llbracket \mathbf{U} \rrbracket)$.
 - (c) The parties make calls to \mathcal{F}_{mult} on corresponding element of the vectors $[\![\mathbf{T}]\!]$ and $[\![\mathbf{V}]\!]$. Let $[\![\mathbf{W}]\!] \leftarrow \mathcal{F}_{mult}([\![\mathbf{T}]\!], [\![\mathbf{V}]\!])$.
 - (d) The parties compute $\llbracket \mathbf{S} \rrbracket \leftarrow \llbracket \mathbf{Y} \mathbf{W} \rrbracket$ locally.
 - (e) The parties make a call to $\mathcal{F}_{\mathsf{matMult}}$, and set $\llbracket \mathbf{Z} \rrbracket \leftarrow \mathcal{F}_{\mathsf{matMult}}(\llbracket \mathbf{X}^T \rrbracket, \llbracket \mathbf{S} \rrbracket)$.
 - (f) The parties update their share for $\boldsymbol{\theta}$ locally:

$$\llbracket \boldsymbol{\theta}^{(k)} \rrbracket \leftarrow (1 - \beta) \cdot \llbracket \boldsymbol{\theta}^{(k-1)} \rrbracket + \alpha \cdot \llbracket \mathbf{Z} \rrbracket$$

Output. Party P_i outputs its share $[\boldsymbol{\theta}^{(K)}]_i$.

Figure 1: 2PC protocol for Secure Poisson Regression.

Consider $\mathcal{R} = (\mathbb{Z}_{2^l}, l_x, l_f)$, and \mathbb{F}_q as parameters for our regression protocol. Each fixed-point multiplication has a failure probability of at most 2^{l_x+1-l} due to truncation. For matrix multiplication between a $(n \times m)$, and a $(m \times k)$ matrix, the failure probability is at most $nk \cdot 2^{l_x+1-l}$ (see [21]).

For each iteration of the regression, there are a total of 2(n+m) truncations for the multiplication steps (n each from steps 2a and 2c, and m each for steps 2e and 2f), which add up to a failure probability of $(2n + 2m) \cdot 2^{l_x+1-l}$. Additionally, there are n exponentiations in step 2c, each of which has a failure probability of at most $2^{l_x+1}/q$ (see Section 6 for details). Therefore, by the union bound, the total failure probability of our regression protocol for K iterations is at most $K(2(n+m) \cdot 2^{l_x+1-l} + n \cdot 2^{l_x+1}/q)$. This dictates the parameter choices for the fixed-point ring and the prime field required for an acceptable failure probability, say $p_{fail} < 2^{-40}$. Note that the failure probability can be made arbitrarily small by increasing l and q.

Standard Poisson regression. The secure regression protocol we described so far is for the general version of Poisson regression with exposure. Standard Poisson regression does not contain the exposure data (**T**). This means that for standard Poisson regression, the element-wise product between **T** and $e^{\mathbf{X}\boldsymbol{\theta}^{(k)}}$ is no longer necessary. Therefore, we can reduce one communication round, resulting in a 3-round protocol. The other steps of our protocol remain exactly the same.

Secure Inference Protocol

Setup. P₀, P₁ agree on a fixed-point ring \mathcal{R} and a prime q. **Input.** P_i has share $\llbracket \boldsymbol{\theta} \rrbracket_i$ of the weights vector.

Protocol.

- 1. The parties make a call to $\mathcal{F}_{matMult}$ to compute the dot product between $\boldsymbol{\theta}$ and \mathbf{x} . Let $\llbracket u \rrbracket \leftarrow \mathcal{F}_{matMult}(\llbracket \mathbf{x}^T \rrbracket, \llbracket \boldsymbol{\theta} \rrbracket)$.
- 2. The parties make a call to $\mathcal{F}_{\mathsf{FPexp}}$ on $\llbracket u \rrbracket$. Let $v \leftarrow \mathcal{F}_{\mathsf{FPexp}}(e, \llbracket u \rrbracket)$.
- 3. The parties make a call to $\mathcal{F}_{\mathsf{mult}}$ to multiply $\llbracket t \rrbracket$ and $\llbracket u \rrbracket$. Let $\llbracket y \rrbracket \leftarrow \mathcal{F}_{\mathsf{mult}}(\llbracket t \rrbracket, \llbracket u \rrbracket)$.

Output. Party P_i outputs its share $\llbracket y \rrbracket_i$.

Figure 2: Secure inference protocol.

4.2 Secure Inference

A useful functionality, after the regression is complete is to use the learned weights to predict, or infer the value of the response variable for future samples. Formally, suppose that P_0 and P_1 hold a sharing $\llbracket \theta \rrbracket$ of the weights. Now, given a new sample ($\llbracket x \rrbracket, \llbracket t \rrbracket$) that is shared between the two parties, the goal is to use the learned weights to compute a sharing $\llbracket y \rrbracket$ of the response variable. Note that $\llbracket y \rrbracket = \llbracket e^{\mathbf{x}^T \theta} \rrbracket$, which can be computed securely along the same lines as our secure Poisson regression protocol. The inference protocol is detailed in Figure 2.

4.3 Additional Considerations

Learning rate. It is important to choose a good learning rate for the Poisson regression to converge efficiently. A large learning rate may cause the regression to not converge, while a small value can cause it to converge slowly. For regression done in the clear, the learning rate can be adjusted according to the magnitude of the gradient to maximize the efficiency of the training and to avoid divergence. In a secure setting however, this needs to be done carefully in order not to leak information. Testing whether the regression is diverging/converging, or revealing the magnitude of the gradient descent risks exposing sensitive information from the training dataset. All the probes and adjustments for the learning rate must be done securely. For example, we can add a secure function to clip the gradient if its magnitude is larger than certain bound before updating the weights. This is useful as the gradient tends to be large at the beginning of the training and could cause divergence even when the learning rate is relatively small.

Mini-batches. Mini-batched gradient descent, where a small random batch of training samples is used for every iteration, is usually more efficient than batched gradient descent. In this paper, however, we do not run experiments with mini-batch gradient descent as our datasets are all small (each has less than 75 features).

Our correlated Beaver triples also work for mini-batched gradient descent, and allow for similar improvements in both the offline and online phase. For iteration j, the parties generate Beaver triples $([\mathbf{A}]], [\mathbf{B}_j]], [[\mathbf{C}_j]])$ and sample a random permutation π_j together. The rows of $([[\mathbf{A}]]], [[\mathbf{C}_j]])$,

Secure Batched Matrix Multiplication Functionality $\mathcal{F}_{\mathsf{batchMult}}$

Input. P_i has shares $(\llbracket \mathbf{X} \rrbracket_i, \llbracket \mathbf{Y}_1 \rrbracket_i, \cdots, \llbracket \mathbf{Y}_K \rrbracket_i)$. Functionality.

- 1. Wait for shares $(\llbracket \mathbf{X} \rrbracket_i, \llbracket \mathbf{Y}_1 \rrbracket_i, \cdots, \llbracket \mathbf{Y}_K \rrbracket_i)$ from party \mathbf{P}_i , for $i \in \{0, 1\}$.
- 2. Reconstruct $\mathbf{X}, \mathbf{Y}_1, \cdots, \mathbf{Y}_K$ and compute $\mathbf{Z}_i = \mathbf{X} \cdot \mathbf{Y}_i$.
- 3. Secret share $\mathbf{Z}_j = \llbracket \mathbf{Z}_j \rrbracket_0 + \llbracket \mathbf{Z}_j \rrbracket_1$ for $j \in [1, K]$ where $\llbracket \mathbf{Z}_j \rrbracket_0$ is sampled uniformly at random and give $\llbracket \mathbf{Z}_j \rrbracket_i$ to \mathbf{P}_i .

Output. Party P_i outputs shares $[\![\mathbf{Z}_1]\!]_i, \cdots, [\![\mathbf{Z}_K]\!]_i$ where $\mathbf{Z}_j = \mathbf{X} \cdot \mathbf{Y}_j$.

Figure 3: Ideal functionality for batched matrix multiplication

and the training data are then shuffled according to π_j . Now, the triples and the training data can be partitioned into mini batches, with each gradient descent iteration run over a different mini-batch.

5 Optimized Batched Multiplication

We now describe our optimized batched multiplication protocol for efficient computation of many multiplications where one of the multiplicands stays the same. More specifically, we want to compute K multiplications of the form \mathbf{XY}_j for secret shared matrices where **X** has size $n \times m$, and all \mathbf{Y}_i have size $m \times k$. \mathbf{P}_i is provided shares $[\![\mathbf{X}]\!]_i, [\![\mathbf{Y}_1]\!]_i, \cdots, [\![\mathbf{Y}_K]\!]_i$, and the goal now is to compute shares of the multiplications $[\![\mathbf{Z}_j]\!] = [\![\mathbf{XY}_j]\!]$ (for $j \in [1, K]$) more efficiently. We provide the ideal functionality $\mathcal{F}_{\mathsf{batchMult}}$ for this in Figure 3.

5.1 Online Phase

To realize the $\mathcal{F}_{batchMult}$ functionality, instead of using independent beaver triples $(\llbracket \mathbf{A}_j \rrbracket, \llbracket \mathbf{B}_j \rrbracket, \llbracket \mathbf{C}_j \rrbracket)$ (one for each multiplication), we prove (in Theorem 1) that we can use correlated randomness across the multiplications and therefore need only a single matrix sharing $\llbracket \mathbf{A} \rrbracket$ for the **X** multiplicand. Formally, our preprocessing requirement is now the shares $\llbracket \mathbf{A} \rrbracket, \llbracket \mathbf{B}_1 \rrbracket, \cdots, \llbracket \mathbf{B}_K \rrbracket, \llbracket \mathbf{C}_1 \rrbracket, \cdots, \llbracket \mathbf{C}_K \rrbracket$. We call these *correlated Beaver triples*. The batched multiplication protocol that uses these triples is detailed in Figure 4.

If **X** is large compared to the \mathbf{Y}_j (as is the case in Poisson regression), this optimization is significant since we only need one matrix to mask **X** across all multiplications. Note that we can use the same batch multiplication technique to compute the element-wise product in our protocol.

Theorem 1. The protocol $\Pi_{batchMult}$ in Figure 4 securely realizes the ideal functionality $\mathcal{F}_{batchMult}$ in Figure 3 in the $\mathcal{F}_{batchBeaver}$ -hybrid world and in the presence of a semi-honest adversary.

Proof. Let P_i be the corrupted party. The simulator S queries $\mathcal{F}_{\mathsf{batchBeaver}}$ to obtain shares of the Beaver triples and hands them to P_i . This means that the distribution of the Beaver triple shares is identical in both the hybrid and ideal world. Next, the simulator opens $(\mathbf{X} - \mathbf{A})$ to a random matrix, and uses the output from the ideal functionality $\mathcal{F}_{\mathsf{batchMult}}$ to compute the values $(\mathbf{Y}_j - \mathbf{B}_j)$. This allows S to simulate the shares $[[(\mathbf{Y}_j - \mathbf{B}_j)]]_{1-i}$ of the other party in the online

Secure Batched Matrix Multiplication Protocol $\Pi_{batchMult}$

Input. P_i has shares $(\llbracket \mathbf{X} \rrbracket_i, \llbracket \mathbf{Y}_1 \rrbracket_i, \cdots, \llbracket \mathbf{Y}_K \rrbracket_i)$. Offline Phase. P_i retrieves correlated Beaver triples $(\llbracket \mathbf{A} \rrbracket_i, \llbracket \mathbf{B}_1 \rrbracket_i, \cdots, \llbracket \mathbf{B}_K \rrbracket_i, \llbracket \mathbf{C}_1 \rrbracket_i, \cdots, \llbracket \mathbf{C}_K \rrbracket_i)$ such that $\mathbf{C}_j = \mathbf{A} \cdot \mathbf{B}_j$ for $j \in [1, K]$. This is done by making a call to $\mathcal{F}_{\mathsf{batchBeaver}}$.

Online Protocol.

- 1. P_i computes shares $[\![\mathbf{X} \mathbf{A}]\!]_i, [\![\mathbf{Y}_1 \mathbf{B}_1]\!]_i, \cdots, [\![\mathbf{Y}_K \mathbf{B}_K]\!]_i$ locally. Then, P_0 and P_1 exchange the shares to reconstruct $(\mathbf{X} \mathbf{A}), (\mathbf{Y}_1 \mathbf{B}_1), \cdots, (\mathbf{Y}_K \mathbf{B}_K)$.
- 2. For each $j \in [1, K]$, party \mathbf{P}_i computes

$$\left[\!\left[\mathbf{Z}_{j}\right]\!\right]_{i} = i \cdot (\mathbf{X} - \mathbf{A})(\mathbf{Y}_{j} - \mathbf{B}_{j}) + (\mathbf{X} - \mathbf{A})\left[\!\left[\mathbf{B}_{j}\right]\!\right]_{i} + \left[\!\left[\mathbf{A}\right]\!\right]_{i}(\mathbf{Y}_{j} - \mathbf{B}_{j}) + \left[\!\left[\mathbf{C}_{j}\right]\!\right]$$

Output. Party P_i outputs shares $\llbracket \mathbf{Z}_1 \rrbracket_i, \cdots, \llbracket \mathbf{Z}_K \rrbracket_i$

Figure 4: Protocol for batched matrix multiplication

phase. Note that $[\![(\mathbf{Y}_j - \mathbf{B}_j)]\!]_{1-i}$ is a function of the Beaver triple shares, the output $[\![\mathbf{Z}_j]\!]_i$ and $(\mathbf{X} - \mathbf{A})$. In both worlds, these matrices are uniformly random and independent from one another. Therefore, the joint distributions between the hybrid world and the ideal world are identical. \Box

Online cost improvement. Correlated Beaver triples improve the cost of our protocol significantly. In the online phase, since $\mathbf{X} - \mathbf{A}$ only needs to be reconstructed once instead of for each multiplication, the amortized total online communication per multiplication for our technique is $\frac{2nm}{K} + 2mk$ ring elements, compared to 2nm + 2mk using standard matrix Beaver triples from [21]. In the setting of Poisson regression, the most significant factor in these costs is $\Theta(nm)$ as k = 1 typically, which results in a $\frac{n+1}{(n/K)+1}$ factor improvement for the online phase. When K is large, this is very close to (n + 1). For example, if the number of training samples n = 1000 and the model is trained over K = 1000 iterations, the communication cost of the online phase is reduced by 500 times if correlated Beaver triples are used.

5.2 Offline Phase

We now discuss how to generate the correlated Beaver triples. The ideal functionality is given in Figure 5. Without any need for optimization, the triples can be generated in the offline phase using the two approaches from SecureML [21]: OT-based and additive homomorphic encryption (AHE) based. The latter, which uses Paillier encryption, requires 190x less communication than the former, but is more expensive computationally. Experiments from [21] show that the AHE-based approach is better in WAN network, while the OT-based is 20-30x faster in LAN setting (See Table 2 in [21]).

Here, we show how to significantly improve triple generation via the use of Ring-LWE. Our approach works for any ring \mathbb{Z}_N $(N = 2^l \text{ in our case})$ and does not rely on packing techniques for \mathbb{Z}_{2^l} as in [24] (where the number of slots is only $\phi(m)/5$, resulting in 80% space being wasted) or on the embedding of plaintext values in a larger prime field of length $2 \cdot l + \sigma + 2$ as in [25] (which increases the communication and computation cost by at least $(2 + (\sigma + 2)/l)$ times). In more detail, to generate K correlated Beaver triples, our protocol proceeds as follows: First, P₀

Correlated Beaver Triple Functionality $\mathcal{F}_{\mathsf{batchBeaver}}$

Parameters. Let n, m, k, K, N be functionality parameters, where n, m, k are used to define matrix sizes, K is the number of triples to generate, and N defines the ring \mathbb{Z}_N .

Functionality.

- 1. Sample uniformly at random $\llbracket \mathbf{A} \rrbracket_i \in \mathbb{Z}_N^{n \times m}$, $\llbracket \mathbf{B}_j \rrbracket_i \in \mathbb{Z}_N^{m \times k}$, $\llbracket \mathbf{C}_j \rrbracket_0 \in \mathbb{Z}_N^{n \times k}$ for $i \in \{0, 1\}, j \in [1, K]$.
- 2. Compute $\llbracket \mathbf{C}_j \rrbracket_1 = (\llbracket \mathbf{A} \rrbracket_0 + \llbracket \mathbf{A} \rrbracket_1)(\llbracket \mathbf{B}_j \rrbracket_0 + \llbracket \mathbf{B}_j \rrbracket_1) \llbracket \mathbf{B}_j \rrbracket_0.$
- 3. Send $(\llbracket \mathbf{A} \rrbracket_i, \llbracket \mathbf{B}_j \rrbracket_i, \llbracket \mathbf{C}_j \rrbracket_i)$ to \mathbf{P}_i .

Output.

- 1. P₀ outputs $([[A]]_0, [[B_1]]_0, [[C_1]]_0, \cdots, [[B_K]]_0, [[C_K]]_0).$
- 2. P₀ outputs $([[\mathbf{A}]]_1, [[\mathbf{B}_1]]_1, [[\mathbf{C}_1]]_1, \cdots, [[\mathbf{B}_K]]_1, [[\mathbf{C}_K]]_1)$ where $[[\mathbf{C}_j]]_1 = -\mathbf{R}_j$.

Figure 5: Ideal functionality for generating correlated Beaver triples

and P_1 sample random matrices to be shares of \mathbf{A} and $\mathbf{B}_1, \dots, \mathbf{B}_K$. Let the shares held by P_i be $[\![\mathbf{A}]\!]_i, [\![\mathbf{B}_1]\!]_i, \dots, [\![\mathbf{B}_K]\!]_i$. In order to obtain the shares of $\mathbf{C}_j = \mathbf{A}\mathbf{B}_j$, the parties need to compute the shares of $[\![\mathbf{A}]\!]_i [\![\mathbf{B}_j]\!]_{1-i}$ as $[\![\mathbf{A}]\!]_i [\![\mathbf{B}_j]\!]_i$ can then be computed locally by each party P_i (since $[\![\mathbf{C}_j]\!] = [\![\mathbf{A}\mathbf{B}_j]\!] = [\![\![\mathbf{A}]\!]_0 [\![\mathbf{B}_j]\!]_0 + [\![\mathbf{A}]\!]_1 [\![\mathbf{B}_j]\!]_1 + [\![\mathbf{A}]\!]_0 [\![\mathbf{B}_j]\!]_1 + [\![\mathbf{A}]\!]_1 [\![\mathbf{B}_j]\!]_0]\!]$). We propose two different ways to compute the shares of $[\![\mathbf{A}]\!]_i [\![\mathbf{B}_j]\!]_{1-i}$. The first approach works better when the number of training samples n is large while the second approach works better when the number of explanatory variables m is small. We benchmark the cost to generate triples using both approaches in Table 3 (Section 7.2), and also compare to the Paillier encryption based approach used in [21]. Overall, both of our approaches are significantly better than Paillier encryption in terms of both communication and computation.

Approach I: Column-wise encryption. P_0 encrypts each column of the the matrix $[\![\mathbf{A}]\!]_0$ separately using Ring-LWE and sends the encrypted columns to P_1 . Define $[\![\mathbf{B}_j]\!]_1 = (b_{1j}, \cdots, b_{mj})^T$, \mathbf{A}_i as the i^{th} column of $[\![\mathbf{A}]\!]_0$, and \mathbf{E}_i as $\operatorname{Enc}_{sk}(\mathbf{A}_i)$ for $1 \leq i \leq m$. P_1 uses the additive homomorphic properties of Ring-LWE to compute the encryption $\mathbf{D}_j = \sum_{i=1}^m b_{ij} \mathbf{E}_i + \mathbf{R}_j = \operatorname{Enc}([\![\mathbf{A}]\!]_0 [\![\mathbf{B}_j]\!]_1 + \mathbf{R}_j)$. P_1 sends the ciphertexts to P_0 who decrypts them to obtain $[\![\![\mathbf{A}]\!]_0 [\![\mathbf{B}_j]\!]_1]\!]_0 = [\![\mathbf{A}]\!]_0 [\![\mathbf{B}_j]\!]_1 + \mathbf{R}_j$ while P_1 has $[\![\![\mathbf{A}]\!]_0 [\![\mathbf{B}_j]\!]_1]\!]_1 = -\mathbf{R}_j$. If the number of training samples n is much smaller than the length of the ciphertext (say N, the degree of the cyclotomic polynomial used in the Ring-LWE scheme), P_1 can pack multiple \mathbf{D}_j into a single ciphertext to optimize communication. Assuming N = 2n, $\mathbf{D}_0 = \operatorname{Enc}(d_1, \cdots, d_n, 0, \cdots, 0)$, and $\mathbf{D}_1 = \operatorname{Enc}(d'_1, \cdots, d'_n, 0, \cdots, 0)$, we can produce the ciphertext $\mathbf{D}'_1 = \operatorname{Enc}(0, \cdots, 0, d'_1, \cdots, d'_n)$ by multiplying \mathbf{D}_1 with the plaintext message $(0, \cdots, 0, 1, 0, \cdots, 0)$ which is zero everywhere except for the n^{th} position. Now, $\mathbf{D}_{01} = \mathbf{D}_0 + \mathbf{D}'_1$ is the ciphertext containing $(b_1, \cdots, b_n, b'_1, \cdots, b'_n)$. Similarly, if t = N/n, we can pack t ciphertexts \mathbf{D}_j into one ciphertext. The parties now reverse roles to compute shares of $[\![\!\mathbf{A}]\!]_1 [\![\!\mathbf{B}_j]\!]_0$, and finally shares of \mathbf{AB}_j . Figure 6 contains a full description of this protocol.

To analyze the efficiency of this approach, first notice that each party sends m ciphertexts and receives $K \cdot n/N$ ciphertexts. Now, if $\log(q)$ is the bitlength of the Ring-LWE ciphertext

Correlated Beaver Triple Protocol $\Pi^{I}_{batchBeaver}$ (Approach I)

Setup. P_0 and P_1 agree on a ring $\mathbb{Z}_{\widehat{N}}$ and the parameters of an additive homomorphic encryption (Ring-LWE) scheme. They also agree on the number of Beaver triple matrices to generate. **Protocol.**

- 1. P₀ and P₁ sample shares of random matrix **A** and of random vectors $\mathbf{B}_1, ..., \mathbf{B}_K$ (denoted $(\llbracket \mathbf{A} \rrbracket_0, \llbracket \mathbf{B}_1 \rrbracket_0, \cdots, \llbracket \mathbf{B}_K \rrbracket_0)$ and $(\llbracket \mathbf{A} \rrbracket_1, , \llbracket \mathbf{B}_1 \rrbracket_1, \cdots, \llbracket \mathbf{B}_K \rrbracket_1)$ respectively). **A** is a $n \times m$ matrix and \mathbf{B}_i is a column vector of size m.
- 2. Compute $\llbracket \llbracket \mathbf{A} \rrbracket_0 \llbracket \mathbf{B}_j \rrbracket_1 \rrbracket$ for $1 \le j \le K$.
 - (a) Denote $\llbracket \mathbf{B}_j \rrbracket_1 = (b_{1j}, \cdots, b_{mj})^T$ and \mathbf{A}_i the i^{th} column of $\llbracket \mathbf{A} \rrbracket_0$.
 - (b) P₀ samples a secret key sk of the Ring-LWE scheme and computes $\mathbf{E}_j \leftarrow \operatorname{Enc}_{\mathsf{sk}}(\mathbf{A}_j)$ for $1 \leq j \leq m$. It then sends $(\mathbf{E}_1, \dots, \mathbf{E}_m)$ to P₁. (P₀ shares the public key with P₁.)
 - (c) Upon receiving $(\mathbf{E}_1, \dots, \mathbf{E}_m)$ from P₀, P₁ samples random matrix $\mathbf{R} = (\mathbf{R}_1, \dots, \mathbf{R}_K)$ and uses the additive homomorphic property of the encryption scheme to compute $\mathbf{D}_j = \sum_{i=1}^m b_{ij}\mathbf{E}_i + \mathbf{R}_j$ for $1 \le j \le K$. P₁ sends $\mathbf{D}_1, \dots, \mathbf{D}_K$ back to P₀.
 - (d) P₀ decrypts \mathbf{D}_j and obtains $\llbracket \llbracket \mathbf{A} \rrbracket_0 \llbracket \mathbf{B}_j \rrbracket_1 \rrbracket_0 = \llbracket \mathbf{A} \rrbracket_0 \cdot \llbracket \mathbf{B}_j \rrbracket_1 + \mathbf{R}_j$ while P₁ sets $\llbracket \llbracket \mathbf{A} \rrbracket_0 \llbracket \mathbf{B}_j \rrbracket_1 \rrbracket_1 = -\mathbf{R}_j$
- 3. Compute $\llbracket \llbracket \mathbf{A} \rrbracket_1 \llbracket \mathbf{B}_j \rrbracket_0 \rrbracket$ for $1 \le j \le K$. The parties reverse roles in the previous step to compute $\llbracket \llbracket \mathbf{A} \rrbracket_1 \llbracket \mathbf{B}_j \rrbracket_0 \rrbracket$.

4.
$$P_i \text{ sets } [\![\mathbf{C}_j]\!]_i = [\![\mathbf{A}]\!]_i [\![\mathbf{B}_j]\!]_i + [\![\![\mathbf{A}]\!]_0 [\![\mathbf{B}_j]\!]_1]\!]_i + [\![\![\mathbf{A}]\!]_1 [\![\mathbf{B}_j]\!]_0]\!]_i \text{ for } 1 \le j \le K.$$

Output.

- 1. P₀ outputs $([[A]]_0, [[B_1]]_0, [[C_1]]_0, \cdots, [[B_K]]_0, [[C_K]]_0).$
- 2. P₁ outputs $([[\mathbf{A}]]_1, [[\mathbf{B}_1]]_1, [[\mathbf{C}_1]]_1, \cdots, [[\mathbf{B}_K]]_1, [[\mathbf{C}_K]]_1).$



modulus, then the cost to generate K correlated Beaver triples is $2(m + Kn/N)N(2\log(q)) = 4(mN + Kn)\log(q)$ bits. When K is very large ($\gg mN$), the amortized cost per triple is roughly $4n\log(q)$ bits. For 127-bit input, we use a ciphertext modulus with length $\log(q) = 295$. For Paillier encryption, the amortized cost per triple is $2(m + n)\log(q')$ bits where q' = 6144 is the length of the Paillier ciphertext. Since m < n is typical for training data, our protocol uses at least 10x less bandwidth than the AHE approach from [21]. Besides the smaller communication cost, our approach also provides significant gains in the computation time. To multiply a constant with a ciphertext in Ring-LWE we only need to perform multiplications over field of size 295 bits. However, if Paillier encryption is used, an exponentiation in group of size 6144 bits needs to be computed which is much more expensive.

In terms of computational cost, our protocol requires $m \times K$ multiplications between a scalar and a ciphertext and K(1 - n/N) shift operations.

Approach II: Row-wise encryption. While the previous approach is efficient in terms of communication, it results in a lot of wasteful computation if the number of training samples n is much smaller than the degree N of the cyclotomic polynomial used for Ring-LWE. Our second

Correlated Beaver Triple Protocol $\Pi^{II}_{batchBeaver}$ (Approach II)

Setup. P_0 and P_1 agree on a ring $\mathbb{Z}_{\widehat{N}}$ and the parameters of an additive homomorphic encryption (Ring-LWE) scheme. They also agree on the number of Beaver triple matrices to generate, K. **Protocol.**

- 1. P₀ and P₁ sample shares of random matrix **A** and of random vectors $\mathbf{B}_1, ..., \mathbf{B}_K$ (denoted $(\llbracket \mathbf{A} \rrbracket_0, \llbracket \mathbf{B}_1 \rrbracket_0, \cdots, \llbracket \mathbf{B}_K \rrbracket_0)$ and $(\llbracket \mathbf{A} \rrbracket_1, , \llbracket \mathbf{B}_1 \rrbracket_1, \cdots, \llbracket \mathbf{B}_K \rrbracket_1)$ respectively). **A** is a $n \times m$ matrix and \mathbf{B}_i is a column vector of size m.
- 2. Compute shares $\llbracket \llbracket \mathbf{A} \rrbracket_1 \llbracket \mathbf{B}_j \rrbracket_0 \rrbracket$ for $1 \le j \le K$.
 - (a) Denote $\llbracket \mathbf{A} \rrbracket_1 = (a_{ij})$ and $\mathbf{B} = (\llbracket \mathbf{B}_1 \rrbracket_0, \cdots, \llbracket \mathbf{B}_K \rrbracket_0)$ and \mathbf{B}^i the i^{th} row of \mathbf{B} .
 - (b) P_0 samples a secret key sk of the Ring-LWE scheme and computes $\mathbf{E}^j \leftarrow \operatorname{Enc}_{\mathsf{sk}}(\mathbf{B}^j)$ for $1 \leq j \leq m$. It then sends $(\mathbf{E}^1, \dots, \mathbf{E}^m)$ to P_1 . (P_0 shares the public key with P_1 .)
 - (c) Upon receiving $(\mathbf{E}^1, \dots, \mathbf{E}^m)$ from \mathbf{P}_0 , \mathbf{P}_1 samples random matrix $\mathbf{R} = (\mathbf{R}^1, \dots, \mathbf{R}^n)^T$ and uses the additive homomorphic property of the encryption scheme to compute $\mathbf{D}^i = \sum_{j=1}^m a_{ij} \mathbf{E}^j + \mathbf{R}^i$ for $1 \le i \le n$. \mathbf{P}_1 sends $\mathbf{D}^1, \dots, \mathbf{D}^n$ back to \mathbf{P}_0 .
 - (d) P₀ decrypts \mathbf{D}^i and obtains $\llbracket \llbracket \mathbf{A} \rrbracket_1 \llbracket \mathbf{B}_j \rrbracket_0 \rrbracket_0 = \llbracket \mathbf{A} \rrbracket_1 \cdot \llbracket \mathbf{B}_j \rrbracket_0 + \mathbf{R}_j$ while P₁ sets $\llbracket \llbracket \mathbf{A} \rrbracket_1 \llbracket \mathbf{B}_j \rrbracket_0 \rrbracket_1 = -\mathbf{R}_j$
- 3. Compute $\llbracket \llbracket \mathbf{A} \rrbracket_0 \llbracket \mathbf{B}_j \rrbracket_1 \rrbracket$ for $1 \le j \le K$. The parties reverse roles in the previous step to compute $\llbracket \llbracket \mathbf{A} \rrbracket_0 \llbracket \mathbf{B}_j \rrbracket_1 \rrbracket$.

4.
$$P_i \text{ sets } [\![\mathbf{C}_j]\!]_i = [\![\mathbf{A}]\!]_i [\![\mathbf{B}_j]\!]_i + [\![\![\mathbf{A}]\!]_0 [\![\mathbf{B}_j]\!]_1]\!]_i + [\![\![\mathbf{A}]\!]_1 [\![\mathbf{B}_j]\!]_0]\!]_i \text{ for } 1 \le j \le K.$$

Output.

- 1. P₀ outputs $([[\mathbf{A}]]_0, [[\mathbf{B}_1]]_0, [[\mathbf{C}_1]]_0, \cdots, [[\mathbf{B}_K]]_0, [[\mathbf{C}_K]]_0).$
- 2. P₁ outputs $([[\mathbf{A}]]_1, [[\mathbf{B}_1]]_1, [[\mathbf{C}_1]]_1, \cdots, [[\mathbf{B}_K]]_1, [[\mathbf{C}_K]]_1).$



approach therefore, will be geared towards settings when $n \ll N$.

For this, P_0 first encrypts each row of the matrix $\mathbf{B} = (\llbracket \mathbf{B}_1 \rrbracket_0, \cdots, \llbracket \mathbf{B}_K \rrbracket_0)$ separately and sends the ciphertexts $\mathbf{E}^j \leftarrow \operatorname{Enc}_{sk}(\mathbf{B}^j)$ for $1 \leq j \leq m$ to P_1 . Now, P_1 uses the additive property of Ring-LWE to compute $\mathbf{D}^i = \sum_{j=1}^m a_{ij} \mathbf{E}^j + \mathbf{R}^i$ for $1 \leq i \leq n$, where \mathbf{D}^i is the encryption of the i^{th} row of the matrix $\llbracket \mathbf{A} \rrbracket_1 \mathbf{B} + \mathbf{R}$ and \mathbf{R} is a random matrix sampled by P_1 , and sends the ciphertexts to P_0 . Note that when K is much smaller than N, P_1 can pack multiple ciphertexts into one before sending them back to P_0 to reduce the communication cost. The packing is done by simply shifting the ciphertexts as described in the first approach. P_0 now decrypts the ciphertexts to obtain $\llbracket \llbracket \mathbf{A} \rrbracket_1 \llbracket \mathbf{B}_j \rrbracket_0 \rrbracket_0 = \llbracket \mathbf{A} \rrbracket_1 \llbracket \mathbf{B}_j \rrbracket_0 + \mathbf{R}_j$, while P_1 sets $\llbracket \llbracket \mathbf{A} \rrbracket_1 \llbracket \mathbf{B}_j \rrbracket_0 \rrbracket_1 = -\mathbf{R}_j$. Similar to the first approach, the two parties now reverse roles to compute shares of $\llbracket \mathbf{A} \rrbracket_1 \llbracket \mathbf{B}_j \rrbracket_0$, and finally shares of $\mathbf{A} \mathbf{B}_j$. Figure 7 contains a full description of this protocol.

Assume $K \ll N$ (in our experiments, K = 1000 and $N = 2^{14}$). To analyze the efficiency of this approach, first notice that P₁ sends *m* ciphertexts to P₁ and receives $n \cdot K/N$ ciphertexts. The communication cost to generate *K* correlated Beaver triples is therefore $2(m+n \cdot K/N)N(2\log(q)) = 4(mN + nK)\log(q)$ bits. When mN < nK (for example, $n = 1000, m = 10, N = 2^{14}, K = 1000$), the amortized cost for one triple is less than $8n\log(q)$ bits, which is around 5x cheaper than the

AHE-based approach from [21]. In terms of computation, our protocol requires $n \times m$ multiplications between a scalar and a ciphertext and n(1 - K/N) shift operations. The second approach is faster than the first one when n < K.

In our secure Poisson regression protocol, we also need to generate the correlated Beaver triples for the multiplication between shares of scalars \mathbf{T}_i and \mathbf{V}_i where \mathbf{T}_i is fixed during the training process. This is equivalent to having n = m = 1, so the second approach will be used to generate these correlated Beaver triples.

Theorem 2. The protocols $\Pi^{I}_{batchBeaver}$ (Figure 6) and $\Pi^{II}_{batchBeaver}$ (Figure 7) both securely realize the ideal functionality $\mathcal{F}_{batchBeaver}$ in Figure 5 in the presence of a semi-honest adversary.

Proof. The security proofs for both approaches are identical and directly follow from the security of the additive homomorphic scheme. First, suppose that P_0 is the corrupted party. Then, the simulator S queries the ideal functionality and obtains P_0 's output. In Step 3, the simulator sends P_0 the encryption of a random matrix \mathbf{A}'_1 , receiving $[[\mathbf{A}']_1 \cdot [\mathbf{B}_j]_0]_1$. The simulator can infer $[[\mathbf{A}']_1 \cdot [\mathbf{B}_j]_0]_0$ as it knows P_0 's input. The simulator rewinds to Step 2 and sends P_0 the ciphertexts \mathbf{D}_j which encrypt the messages $[\mathbf{C}_j]_0 - [[\mathbf{A}]_0 \cdot [\mathbf{B}_j]_0 - [[[\mathbf{A}']_1 \cdot [\mathbf{B}_j]_0]_0]_0$. It is easy to see that the joint distributions in both worlds are computationally indistinguishable by a reduction to the security of the additive homomorphic encryption scheme.

Now, suppose that P_1 is the corrupted party. To simulate P_1 , it is enough to send P_1 the encryption of random messages in Step 2. In Step 3, the simulator just ignores what it receives from P_1 and sends back the encryption of $[\![\mathbf{C}_j]\!]_1 - [\![\mathbf{A}]\!]_1 \cdot [\![\mathbf{B}_j]\!]_1 - [\![[\mathbf{A}']\!]_0 \cdot [\![\mathbf{B}_j]\!]_1]\!]_1$. Again, the ability to distinguish the real and ideal execution will imply the ability to break the additive homomorphic encryption scheme.

We conclude that the joint distributions in both worlds are computationally indistinguishable.

6 Secure Fixed-Point Exponentiation

In this section, we detail our novel secure fixed-point exponentiation protocol. To simplify our analysis, our protocol will mirror $\mathcal{F}_{\mathsf{FPexp}}$ functionality (Figure 8) rather than the previously defined FPExp operation. Note that due to truncation errors, the two functionalities are not identical. However, we will show later (in Section 6.4) that the result computed by $\mathcal{F}_{\mathsf{FPexp}}$ is close to the actual fixed-point exponentiation result. Similar to the FPExp operation, the functionality $\mathcal{F}_{\mathsf{FPexp}}$ will take as inputs a public base and a secret shared exponent. Since we are working in a fixed-point ring, we will consider our inputs to be the fixed-point representations rather than the real numbers themselves. Given a fixed-point ring $\mathcal{R} = (\mathbb{Z}_{2^l}, l_x, l_f)$, a public base $b \in \mathcal{R}_*$, and a shared exponent $[\![x]\!], \mathcal{F}_{\mathsf{FPexp}}(b, [\![x]\!])$ will compute a sharing of something "close" to $(\underline{b})^{(\underline{x})}$. We compare our work to existing techniques in Section 6.5 and benchmark our protocol in Section 7.1.

6.1 Protocol Construction

It is straightforward to construct a protocol that realizes the \mathcal{F}_{FPexp} functionality. First, we note that the PubFPMult, FPAdd, RingChange, and PubFPDiv operations used in steps 1, 2, 4, 9, and 10 of \mathcal{F}_{FPexp} can all be computed by locally manipulating the shares. Steps 3, 5 and 6 are also

Secure Fixed-Point Exponentiation Functionality $\mathcal{F}_{\mathsf{FPexp}}$

Public Parameters. P₀ and P₁ agree on a fixed-point ring $\mathcal{R} = (\mathbb{Z}_{2^l}, l_x, l_f)$, an *l*-bit prime $q < 2^l$, and an exponent bound $A \in \mathbb{Z}^+$. **Input.** P₀ and P₁ have shares $[\![x]\!] \in \mathcal{R}_*$, and a public real base $b \in \mathbb{R}^+$, satisfying $\underline{x} \log_2(b) > 1 - A$

Functionality.

1. Let $\llbracket x' \rrbracket \leftarrow PubFPMult(\llbracket x \rrbracket, \widehat{\log_2(b)})$	// Convert to base 2 exponentiation
2. Let $\llbracket z \rrbracket \leftarrow FPAdd(\llbracket x' \rrbracket, \widehat{A})$	// Make exponent > 1
3. Let $(z_i^{\text{int}}, z_i^{\text{frac}}) \leftarrow \left(\left\lfloor \llbracket z \rrbracket_i / 2^{l_f} \right\rfloor, (\llbracket z \rrbracket_i / 2^{l_f}) - z_i^{\text{int}} \right)$	// Split into integer and fractional parts
4. Let $(z_0^{\text{int}}, z_1^{\text{int}}) \leftarrow RingChange((z_0^{\text{int}}, z_1^{\text{int}}), \mathbb{Z}_{q-1})$	// RingChange from $\mathbb{Z}_{2^{l-l_f}}$ to \mathbb{Z}_{q-1}
5. Let $(v_i^{\text{int}}, v_i^{\text{frac}}) \leftarrow (2^{z_i^{\text{int}}} \mod q, 2^{z_i^{\text{frac}}})$	// Exponentiate both parts
6. Let $v_i \leftarrow \left(v_i^{\text{int}} \cdot \lfloor 2^{l_f} v_i^{\text{frac}} \rfloor\right) \mod q$	// Get each party's local share
7. Let $y' \leftarrow v_0 v_1 \mod q$	// Combine shares of both parties
8. Create a random additive sharing $\llbracket y' \rrbracket$ in \mathbb{F}_q	// Convert to additive shares
9. Let $\llbracket y \rrbracket^{\mathbb{F}_q} \leftarrow PubFPDiv(\llbracket y' \rrbracket, 2^{l_f + A})$	// Divide by the remaining factor
10. Let $\llbracket y \rrbracket^{\mathbb{Z}_{2^l}} \leftarrow RingChange(\llbracket y \rrbracket^{\mathbb{F}_q}, \mathbb{Z}_{2^l})$	// RingChange from \mathbb{F}_q to \mathbb{Z}_{2^l}

Figure 8: Functionality $\mathcal{F}_{\mathsf{FPexp}}$

purely local computations. The only point at which communication will be necessary is to retrieve an additive sharing of y' (steps 7, 8). Effectively, here, P_0 and P_1 need to go from a multiplicative sharing of $y' \in \mathbb{F}_q$ to an additive sharing of the same y'.

To accomplish this, we use a 2-party variant of the efficient MTA (multiplicative to additive) protocol from Ghodosi et al. [16] (also given in Figure 9). Suppose that P₀ and P₁ hold multiplicative shares m_0 and m_1 of a secret s in \mathbb{F}_q . The protocol requires a tuple (α_i, β_i) of preprocessed values (in \mathbb{F}_q) such that $\alpha_0\alpha_1 + \beta_0\beta_1 = 1$. Now, the MTA protocol proceeds as follows: First, P₀ and P₁ simultaneously send $v_0 = \beta_0 m_0$ and $v_1 = \alpha_1 m_1$ respectively to the other party. Then, P₀ and P₁ can compute $a_0 = \alpha_0 m_0 v_1$ and $a_1 = \beta_1 m_1 v_0$. Note that a_0 and a_1 are the required additive shares of s since $a_0 + a_1 = \alpha_0 m_0 \alpha_1 m_1 + \beta_1 m_1 \beta_0 m_0 = m_0 m_1 (\alpha_0 \alpha_1 + \beta_0 \beta_1) = s$. [16] also shows that the shares are individually uniformly random.

The source of the preprocessed values is not provided in [16] but they are nevertheless easy to compute even without a trusted dealer. For this, first, P₀ samples u_0, w_0 and P₁ samples α_1, β_1 uniformly at random from \mathbb{F}_q^{\times} . Next, the two parties can securely compute $r = u_0\alpha_1 + w_0\beta_1$, and resample if r = 0. The probability that a resample is necessary is at most 1/(q-1). Finally, P₀ can set $\alpha_0 = u_0r^{-1}$ and $\beta_0 = w_0r^{-1}$, where r^{-1} is the multiplicative inverse of r in \mathbb{F}_q^{\times} . Notice now, that $\alpha_0\alpha_1 + \beta_0\beta_1 = rr^{-1} = 1$, as required. Note that since the resample probability is negligible, the distribution of r is negligibly close to uniformly random.

Since the only communication is through the MTA protocol, the security of our protocol securely realizing the \mathcal{F}_{FPexp} functionality is a direct consequence of the security of the MTA protocol. In total, our protocol requires only one round, and a single field element sent by each party.

Multiplicative to Additive Conversion: MTA

Public Parameters. A finite field \mathbb{F}_q where q is prime. All operations will be in \mathbb{F}_q . Preprocessing. P_0 is given (α_0, β_0) and P_1 is given (α_1, β_1) satisfying $\alpha_0 \alpha_1 + \beta_0 \beta_1 = 1$. Input. P_0 and P_1 have multiplicative shares m_0 and m_1 of s, i.e., $s = m_0 m_1$. Required Output. P_i outputs a_i such that $a_0 + a_1 = s$. Protocol Description. $\begin{array}{c}
\frac{P_0}{v_0 \leftarrow \beta_0 m_0} & P_1 \\
\hline v_1 \leftarrow \alpha_1 m_1 \\
\hline v_1 \\
\hline a_0 \leftarrow \alpha_0 m_0 v_1 \\
\hline \end{array}$

Figure 9: Protocol MTA to convert from multiplicative to additive shares

Output a_1

6.2 Protocol Details

Output a_0

We now describe the main technical components of why our protocol is a useful proxy for computing the fixed-point exponentiation. We defer the concrete error analysis of our protocol to Section 6.4. We begin with a simplified version of our protocol where $\underline{b} = 2$, and the exponent satisfies $\underline{x} > 1$, and handle other exponents and other (positive) bases later.

Our strategy is as follows: (1) First, we split the exponentiation into two parts: an integer part and a fractional part. (2) Next, each part is exponentiated separately (and locally) to get multiplicative shares of the final result (along with an extra factor). (3) We then use a single round of interaction to convert the multiplicative shares to additive shares. (4) Finally, each party can locally remove the extra factor to obtain additive shares of the final result. We detail each of these steps below.

Splitting the exponent. Let $[\![z]\!]$ be a sharing of the fixed-point exponent, where P₀ holds $[\![z]\!]_0$ and P₁ holds $[\![z]\!]_1$. We use z here (instead of x) to follow along with functionality $\mathcal{F}_{\mathsf{FPexp}}$, and standardize the notation for a general base, since the first two steps there reduce the problem to a base 2 exponentiation (of a positive exponent). The party P_i first splits its share $[\![z]\!]_i$ as $(z_i^{\text{int}}, z_i^{\text{frac}})$ where $z_i^{\text{int}} = \lfloor [\![z]\!]_i / 2^{l_f} \rfloor$ and $z_i^{\text{frac}} = [\![z]\!]_i / 2^{l_f} - z_i^{\text{int}} = ([\![z]\!]_i \mod 2^{l_f}) / 2^{l_f}$. Notice now that $z = ([\![z]\!]_0 + [\![z]\!]_1 \mod 2^l) = 2^{l_f} ((z_0^{\text{int}} + z_1^{\text{int}} \mod 2^{l-l_f}) + (z_0^{\text{frac}} + z_1^{\text{frac}}))$. Therefore,

$$2^{\tilde{z}} = \left(2^{(z_0^{\mathrm{int}} + z_1^{\mathrm{int}}) \bmod 2^{l-l_f}}\right) \cdot \left(2^{z_0^{\mathrm{frac}} + z_1^{\mathrm{frac}}}\right)$$

This allows us to exponentiate the integer and fractional parts separately and combine them at a later step. Note that the two integer and fractional exponent shares may not always sum up to the actual integer and fractional parts of z respectively. This is because the two fractional shares could add up to more than 1, leaving the integer shares to sum to $\lfloor z/2^{l_f} \rfloor - 1 \mod 2^{l-l_f}$. Furthermore, our integer exponentiation requires the exponent to be positive. This leads to our requirement of z > 1. We will relax this assumption later.

Integer exponentiation. First, we observe that $\llbracket w \rrbracket_0 = z_0^{\text{int}}$ and $\llbracket w \rrbracket_1 = z_1^{\text{int}}$ form a sharing of $w = (z_0^{\text{int}} + z_1^{\text{int}} \mod 2^{l-l_f})$ over the ring $\Re = \mathbb{Z}_{2^{l-l_f}}$. Denote this sharing by $\llbracket w \rrbracket^{\Re}$. Now, we can use existing integer ring exponentiation techniques (such as [8, 23, 31]) to compute 2^w . These techniques however require a few rounds of communication even for a public base. Instead, here, we will describe an alternative method that can be done locally in a way that will seamlessly combine with the fractional exponentiation part.

For this, we assume that the parties have agreed on an *l*-bit prime q (i.e., $2^{l-1} < q < 2^l$). We will first convert the sharing of w in \mathfrak{R} to a sharing in \mathbb{Z}_{q-1} using the **RingChange** operation. Note that the ring size increases if at least 1 fractional bit is present. Recall that since w is positive (from our exponent assumption), with probability $(1 - 2^{l_x}/q)$, the new sharing $\llbracket w \rrbracket^{\mathbb{Z}_{q-1}}$ will satisfy $w + (q-1) = \llbracket w \rrbracket_0^{\mathbb{Z}_{q-1}} + \llbracket w \rrbracket_1^{\mathbb{Z}_{q-1}}$. Now, the two parties can exponentiate their shares locally (mod q) to directly get a multiplicative sharing of 2^w . This works since,

$$\left(2^{\llbracket w \rrbracket_0^{\mathbb{Z}_{q-1}}} \mod q\right) \cdot \left(2^{\llbracket w \rrbracket_1^{\mathbb{Z}_{q-1}}} \mod q\right) \mod q$$
$$= \left(2^{w+q-1}\right) \mod q = 2^w \mod q$$

where the last step is due to Fermat's little theorem. Let $v_0^{\text{int}} = 2^{\llbracket w \rrbracket_0^{\mathbb{Z}_{q-1}}} \mod q$ and $v_1^{\text{int}} = 2^{\llbracket w \rrbracket_1^{\mathbb{Z}_{q-1}}} \mod q$ be the final multiplicative shares (in \mathbb{F}_q) of 2^w held by P_0 and P_1 .

Fractional exponentiation. Let z_0^{frac} and z_1^{frac} be the fractional exponents held by P_0 and P_1 respectively. Notice that if both parties locally exponentiate (in \mathbb{R}) their shares, they would end up with multiplicative shares (in \mathbb{R}) of the fractional exponentiation result. Specifically, if P_i computes $v_i^{\text{frac}} = 2^{z_i^{\text{frac}}}$, then $v_0^{\text{frac}} \cdot v_1^{\text{frac}} = 2^{z_0^{\text{frac}} + z_1^{\text{frac}}}$. To allow for seamless integration with the integer exponentiation part, we have P_i later compute $\lfloor 2^{l_f} \cdot v_i^{\text{frac}} \rfloor$. A crucial observation here is that since $2^0 \leq v_i^{\text{frac}} < 2^1$, $\lfloor 2^{l_f} \cdot v_i^{\text{frac}} \rfloor$ is small and positive, and therefore it can also be viewed as an element in \mathbb{F}_q . Furthermore, the multiplication (now in \mathbb{F}_q), will not wrap around the modulus q. This will allow v_i^{frac} and v_i^{int} to be combined easily. Note that the product will include an extra factor is necessary when first combining the integer and fractional parts and will be divided out later. This will become evident in our error analysis.

Combining the two parts. At this stage, P_i holds the result of the integer exponentiation v_i^{int} , and the result of the fractional part v_i^{frac} . Let $d_i = \lfloor 2^{l_f} \cdot v_i^{\text{frac}} \rfloor$. Ignoring errors due to truncation for now, we have:

$$(v_0^{\text{int}} \cdot v_1^{\text{int}} \cdot d_0 \cdot d_1) \mod q \approx \left(2^{(z_0^{\text{int}} + z_1^{\text{int}}) \mod 2^{l-l_f}}\right) \left(2^{2l_f}\right) \left(2^{z_0^{\text{frac}} + z_1^{\text{frac}}}\right) \mod q$$
$$= 2^{2l_f} 2^{\mathbb{Z}} \mod q$$
$$= 2^{l_f} \widehat{(2\mathbb{Z})} \mod q$$

This means that barring any truncation errors, if P_i computes $y'_i = v_i^{\text{int}} \cdot d_i \mod q$, then $y'_0 y'_1 \mod q \approx (2^{l_f})(2^{\underline{z}})$. Now, P_0 and P_1 convert the multiplicative shares of $y' = y'_0 y'_1$ to additive ones through the MTA protocol which requires one round of interaction. The leftover 2^{l_f} factor can be divided out

through local computation using PubFPDiv. Finally, both parties can locally use the RingChange protocol to convert their shares back to \mathbb{Z}_{2^l} . Note that this conversion is once again from a smaller to a larger ring since $q < 2^l$. We will bound the error resultant from truncation in Section 6.4.

Working with bases other than 2. Our Poisson regression usecase requires secure base e exponentiation, but so far our protocol only works for base 2. To make it work for any positive base b, we first observe that given a real exponent u, $b^u = 2^{u \log_2(b)}$. Consequently, as the first protocol step, the sharing $[\![x]\!]$ of the (base b) exponent in \mathcal{R} will be converted to a sharing $[\![z]\!]$, of the equivalent base 2 exponent, where $(z) = (x) \log_2(b)$. This can be computed as $[\![z]\!] =$ PubFPMult $([\![x]\!], \log_2(b))$ and requires no interaction.

Working with exponents ≤ 1 . As previously mentioned, we initially required our fixed-point exponent to be greater than 1 since this guarantees correctness for the integer ring exponentiation. To handle other exponents, we will assume that there is an agreed upon exponent bound $A \in \mathbb{Z}^+$, such that for base b and exponent sharing $[\![x]\!]$, it holds that $(\underline{x} \log_2(b)) > 1 - A$, i.e., the most negative exponent for base 2 exponentiation still has an absolute value of less than A - 1. Suppose that $[\![x']\!]$ is the sharing of the exponent after converting to a base 2 exponentiation. We now need to ensure that $\underline{x'} > 1$. This can be done by adding A to the exponent, or equivalently, adding \widehat{A} to the sharing using FPAdd to get a new sharing $[\![z]\!]$. At the end of the protocol, the extra 2^A factor will be divided out. We note that since the 2^A factor will be present in intermediate steps, both \mathcal{R} and \mathbb{F}_q will need to be large enough to accommodate it.

Protocol cost and other considerations. Our exponentiation protocol has a total online cost of 2 \mathbb{F}_q elements (1 per party), and a preprocessing cost of 4 \mathbb{F}_q elements (2 per party). We note that our protocol can easily be adapted to working solely in the field \mathbb{F}_q (with appropriately defined fixed-point representation), rather than switching between our defined fixed-point ring and \mathbb{F}_q . This design is simpler but usually much slower since common operations like multiplication, truncation etc., are much faster over a ring \mathbb{Z}_{2^l} , as compared to a field. Therefore, for our purpose, it is far more cost efficient to work mostly in \mathbb{Z}_{2^l} (and $\mathbb{Z}_{2^{l-l_f}}$), and only switch to \mathbb{F}_q inside of the exponentiation subprotocol.

Assumption on the exponent bound. We emphasize that our assumption of a minimum allowable exponent is not unreasonable in the context of fixed-point exponentiation. Given l_f fractional bits, $2^{(-z)}$ where $z > l_f$ is already not representable in the fixed-point ring. Consequently, this gives us a natural bound of $2^{(-z)}$ on how negative the exponent can be for the computation to even make sense. Of course, a tighter bound A can be chosen if appropriate. This observation allows our protocol to be orders of magnitude faster than prior work, since it does not require an expensive bit decomposition to first detect whether the exponent is negative; we can simply add the bound to all exponents to always work with positive exponents for the main protocol. One caveat is that we lose the ability to detect if our predefined bound has been violated without resorting to a bit decomposition, and our protocol may produce incorrect results when the bound is incorrectly defined or is exceeded during protocol execution. We point out though, that this assumption is not unlike a standard assumption of a large enough ring modulus to hold the fixed-point computations, and similar assumptions appear in [4, 21].

Failure probability. We analyze the total failure probability of our base 2 exponentiation protocol. Note that this is different from our error analysis in Section 6.4. In particular, we say that our exponentiation protocol has failure probability p_{fail} and error ε if, except with probability p_{fail} , the protocol error is bounded by ε . To compute the failure probability, first suppose that the (positive) base 2 exponent z is secret shared as $([[z]]_0, [[z]]_1)$. With probability at least $1 - 2^{l_x - l}$, we have $[[z]]_0 + [[z]]_1 = z + 2^l$, i.e., the two shares wrap around \mathbb{Z}_{2^l} . When this happens, the integer components will also wrap around $\mathbb{Z}_{2^{l-l_f}}$, and after the RingChange to \mathbb{Z}_{q-1} , z_0^{int} and z_1^{int} will wrap around \mathbb{Z}_{q-1} .

Next, after the integer and fractional parts are exponentiated combined, and converted from multiplicative to additive shares, the random additive sharing of y' in \mathbb{F}_q will also wrap around \mathbb{F}_q with probability at least $1 - \frac{2^{l_x}}{q}$. Finally, the later PubFPDiv and RingChange back to \mathbb{Z}_{2^l} steps will work smoothly when the sharing of y' wraps around \mathbb{F}_q .

Therefore, using the union bound, we can bound the total failure probability of the exponentiation protocol as $2^{l_x-l} + 2^{l_x}/q < 2^{l_x+1}/q$, since we use $q < 2^l$. Given the exponent bound A, choosing $l_x = 2A + 2l_f$ is sufficient, and therefore, we can rewrite the bound as $2^{2A+2l_f+1}/q$. Note that the failure probability can easily made as small as necessary by increasing the size of \mathbb{F}_q , and our fixed-point ring. For example, to achieve $p_{\text{fail}} < 2^{-40}$, with $l_f = 15$ bits of precision, and A = 5, roughly an 81-bit modulus will be required.

6.3 Other Considerations

Alternate 2-round protocol. We also describe a alternate 2-round variant of our exponentiation protocol. Here, instead of combining the integer and fractional exponentiation shares locally first, the MTA protocol is used to retrieve additive shares of the integer and fractional result separately. Note that this can be done simultaneously in 1 round. Finally, in the second round, shares of both results can be combined through a single secure multiplication. In total, 8 \mathbb{F}_q elements are transmitted in the online phase, and 14 \mathbb{F}_q elements are required for preprocessing. While the communication cost is larger than the previously described 1-round protocol, one upshot of this construction is that it can tolerate a smaller ring size. Recall that in the 1-round protocol, the full result along with an extra 2^{l_f} factor needs to fit in the ring. This is no longer necessary for the 2-round protocol and depending on the usecase and the number of fractional bits used, the trade-off may be acceptable. For our regression usecase however, there are other constraints that increase the size of the fixed point ring. Furthermore, in practice, the computational gain as a result of a smaller ring size (in the order of microseconds for our construction), will almost certainly be overshadowed by the extra communication round (usually in the order of milliseconds). Therefore, we use the 1-round protocol that optimizes for communication cost.

Malicious security. Although our secure fixed-point exponentiation protocol operates exclusively in the semi-honest setting, we comment briefly on the challenges of extending it to a maliciously secure version. One possible technique is for the protocol parties to operate on authenticated shares [22] and use generic zero-knowledge proofs to prove that each party performs their steps correctly. However, doing so would likely reduce the efficiency gains of our protocol substantially. In particular, a key step in our protocol is separating the exponentiation into integer and fractional parts, following which the fractional part can be exponentiated locally in real numbers (or floating point) and still be seamlessly combined with the integer exponentiation part. In the malicious setting, it is expensive to prove that these steps were performed correctly, and it may be more efficient to use a polynomial approximation instead for the fractional exponentiation, together with cut-and-choose or ZK techniques to prove correctness. We leave these explorations for future work.

6.4 Error Analysis

We will now compute a bound on the error of our exponentiation protocol for base 2. For this, we will compute the difference between the result computed by $\mathcal{F}_{\mathsf{FPexp}}$ and the actual exponentiation (in real numbers).

Let $[\![z]\!]$ be a sharing of the (base 2) exponent in the fixed-point ring $\mathcal{R} = (\mathbb{Z}_{2^l}, l_x, l_f)$, that computes the exponentiation $2^{(z)}$ (in \mathbb{R}). First, we note that the integer exponentiation produces no error; the only error results from the truncation in the fractional part and its subsequent combination with the exponentiation of the integer part. Let z_i^{int} and z_i^{frac} denote the integer and fractional parts of the underlying fixed-point of the share $[\![z]\!]_i$, after $(z_0^{\text{int}}, z_1^{\text{int}})$ has undergone a RingChange to become a sharing in \mathbb{Z}_{q-1} . Note that no error is added by the RingChange. It is easy to see that the true computation 2^{z} can be written as $2^{(z_0^{\text{int}}+z_1^{\text{int}} \mod q-1)}2^{z_0^{\text{frac}}}2^{z_1^{\text{frac}}}$.

easy to see that the true computation 2^{z} can be written as $2^{(z_{0}^{\text{int}}+z_{1}^{\text{int}} \mod q-1)} 2^{z_{0}^{\text{frac}}} 2^{z_{1}^{\text{frac}}}$. Following $\mathcal{F}_{\text{FPexp}}$, we first compute $v_{i}^{\text{int}} = 2^{z_{i}^{\text{int}}} \mod q$, and $v_{i}^{\text{frac}} = 2^{z_{i}^{\text{frac}}}$, and combine them to get $v_{i} = (v_{i}^{\text{int}} \cdot \lfloor 2^{l_{f}} v_{i}^{\text{frac}} \rfloor) \mod q$. Since v_{i}^{frac} is a positive real, suppose that $v_{i}^{\text{frac}} = d_{i} + \varepsilon_{i}$, where $0 \leq \varepsilon_{i} < 2^{-l_{f}}$. In other words, ε_{i} is the part not representable in l_{f} fractional bits. Now, $v_{i} = (v_{i}^{\text{int}} \cdot 2^{l_{f}} \cdot (v_{i}^{\text{frac}} - \varepsilon_{i})) \mod q$. Consequently,

$$\begin{aligned} y' &= 2^{2l_f} \cdot v_0^{\text{int}} \cdot v_1^{\text{int}} \cdot (v_0^{\text{frac}} - \varepsilon_0) \cdot (v_1^{\text{frac}} - \varepsilon_1) \mod q \\ &= 2^{2l_f} \cdot 2^{(z_0^{\text{int}} + z_1^{\text{int}} \mod q - 1)} \cdot (v_0^{\text{frac}} - \varepsilon_0) \cdot (v_1^{\text{frac}} - \varepsilon_1) \\ &= 2^{2l_f} \left[2^{(z_0^{\text{int}} + z_1^{\text{int}} \mod q - 1)} (v_0^{\text{frac}} v_1^{\text{frac}} - \varepsilon_0 v_1^{\text{frac}} - \varepsilon_1 v_0^{\text{frac}} + \varepsilon_0 \varepsilon_1) \right] \\ &= 2^{2l_f} \left[2^{(z)} + 2^{(z_0^{\text{int}} + z_1^{\text{int}} \mod q - 1)} (-\varepsilon_0 v_1^{\text{frac}} - \varepsilon_1 v_0^{\text{frac}} + \varepsilon_0 \varepsilon_1) \right] \end{aligned}$$

where the mod q can be removed from step 2 onwards, since \mathbb{F}_q is large enough to accommodate the entire intermediate result. Now, $2^{(z_0^{\text{int}}+z_1^{\text{int} \mod q-1)}} = 2^{(\underline{z})}/(v_0^{\text{frac}} \cdot v_1^{\text{frac}})$, and $1 \leq v_i^{\text{frac}} < 2$ and therefore,

$$2^{2l_f} \left[2^{(z)} - 2^{(z)} \cdot 2^{-l_f} \frac{(v_0^{\text{frac}} + v_1^{\text{frac}})}{v_0^{\text{frac}} v_1^{\text{frac}}} \right] < y' < 2^{2l_f} \left[2^{(z)} + 2^{(z)} \cdot 2^{-2l_f} \right]$$

This gives,

$$2^{l_f} 2^{(z)} (2^{l_f} - 2) < y' < 2^{l_f} 2^{(z)} (2^{l_f} + 2^{-l_f})$$

Now, $y \leftarrow \mathsf{PubFPDiv}(\llbracket y' \rrbracket, 2^{l_f})$ results in an additional potential error of at most ± 1 . That is,

$$-1 + 2^{(z)}(2^{l_f} - 2) < y < 1 + 2^{(z)}(2^{l_f} + 2^{-l_f})$$

In other words, the computed fixed-point number $y = y/2^{l_f}$ differs from the real value $2^{(z)}$ as,

$$\left| \underbrace{y}{y} - 2^{(\underline{z})} \right| < 2^{-l_f} (2 \cdot 2^{(\underline{z})} + 1)$$

To put this in perspective, a computation of $2^{10.125} \approx 1116.68$ in a fixed point ring with $l_f = 15$, will result in a maximum possible error of 0.068, or at most 0.006%. With $l_f = 20$, the maximum error reduces to 0.0002%. This should be more than reasonable for most practical settings, and indeed fits our regression usecase well, since regression is resistant to small errors. Furthermore, we emphasize that the error can always be made arbitrarily small by increasing the number of fractional bits available for the computation. Also note that this error is achieved for the worst possible sharing of the exponent, and may be much smaller for a random sharing.

Error dependence on actual value. The astute reader might observe that the above computed error (in the fixed-point ring) is bounded by a small multiple of the actual real number result $2^{\tilde{z}}$. We highlight that this is not unlike the error of chaining two truncated secure multiplications. For example, suppose that $[[\hat{a}]], [[\hat{b}]], [[\hat{c}]], [[\hat{d}]]$ are sharings held by P₀ and P₁ of fixed-point numbers a, b, c, d. Recall that secure multiplication can result in an error of at most ± 1 in the fixed-point ring. This means that the secure multiplication of a, b can result in a sharing of $\hat{ab} + 1$, while the secure multiplication of c, d can result in a sharing of $\hat{cd} + 1$. At this point, if the two resultant shares are also multiplied, the complete result can be at most $\hat{abcd} + \hat{ab} + \hat{cd} + 2$. In other words, the error here can also depend on the actual numbers involved in the computation.

6.5 Comparison to Existing Techniques

To highlight the strong improvements of our secure exponentiation protocol, we provide a comparison to existing techniques in literature. Later, in Section 7.1, we also benchmark our protocol and compare the results to existing works.

Modular and integer exponentiation. There is a long line of work [8, 14, 23, 31] on secure integer and modular exponentiation. Here, the goal is to compute $b^a \mod N$ where both a, b are secret shared in \mathbb{Z}_N . Other variants have also been studied; for example, where b is assumed to be public, or when N is prime. Classical approaches, such as the one put forth by Damgård et al. [14], required a full decomposition of the shared values, which is computationally expensive. Later approaches [8, 23, 31] are more practical and did not require decomposition. Still, for semihonest adversaries, even when the base b is public, previous approaches require more than 1 round of communication (e.g., [8] describes a 2-round protocol, and [31] describes a 3-round protocol).

A byproduct of our protocol, is a modular/integer exponentiation protocol that works for a public base, and when the exponent is small. In the online phase, it requires a single (simultaneous) communication round, and an exchange of a single field element in either direction. In other words, for exponentiation modulo a prime q, our protocol has a total communication cost of $2 \log q$ bits, and a total preprocessing cost of $4 \log q$ bits.

Fixed-point exponentiation. There are substantially fewer works tackling fixed-point exponentiation. [13] provides several common fixed-point operations but does not describe exponentiation.

The most relevant comparison for fixed-point exponentiation is the protocol from the SCALE-MAMBA package [10]. The authors mention that they could not find a secure fixed-point exponentiation protocol in academic literature and resorted to constructing their own protocol for base 2 (and by extension any public base) exponentiation. While [10] primarily targets malicious security, we briefly compare the overall technique of their protocol to ours.

Their technique is as follows: First, the protocol compares the exponent to zero and proceeds differently depending on positive and negative exponents. Next, each exponent share is split into its integer and fractional parts and the exponentiation is done separately. For the integer part, earlier known techniques are used. The fractional computation is approximated using a degree-9 polynomial $P_{1045}(x)$, described by Hart [18]. Finally, the two computations are combined using a single multiplication, and a division is performed if the exponent was negative.

The overall structure is similar to our technique of splitting the computation into an integer and a fractional part. However, their approach has some notable drawbacks compared to our protocol. First, as far as we are aware, a comparison to zero requires a full bit decomposition of each parties shares, which as mentioned earlier, is expensive. In contrast, our work does not need to compare to zero to be able to handle negative exponents. Instead, here, we assume that the exponent is bounded from below, which allows us to convert the exponents to a suitable positive value beforehand. The extra multiplicative factor (2^A if A is the exponent bound) is later divided out. This assumption is suitable for our application, and results in a much faster protocol. Second, for the fractional exponentiation, SCALE-MAMBA uses a polynomial approximation, using a degree-9 polynomial, which will require several rounds of communication and/or a large communication cost. Here, for our protocol, we make a crucial observation that the fractional exponentiation (in real numbers) is small enough, such that the multiplicative shares of the product do not wrap around the fixedpoint ring. This allows us to directly compute shares of the fractional approximation without any polynomial approximation. Finally, we also realize that the integer and fractional parts can be combined locally first, giving rise to our 1-round protocol. For this, we do require a slightly larger ring $(l_f \text{ more bits})$ to ensure that our intermediate computations can be appropriately represented. but we think that this tradeoff is appropriate. Furthermore, as we point out in Section 7.1, SCALE-MAMBA chooses a larger ring to reduce numerical instability that was observed, which already results in a larger ring size requirement than our protocol.

7 Experimental Evaluation

Implementation details. We implemented our protocols in C++, and compiled the code using the open-source Bazel [2] build tool. We support moduli up to 127-bit for both the fixed-point ring and the field. For the operations, we use the native C++ uint64_t type for moduli smaller than 64-bits, and uint128 from Google's abseil library [3] for larger moduli. We give users the option to decide the base integer size (64-bit or 128-bit) and provide experimental results for both.

Experimental setup. We ran all of our experiments on two compute-optimized c2-standard-8 Google cloud instances with 3.1 GHz base frequency and 32 GB RAM. Our code is single-threaded and only uses a single core of the instance. For the LAN setting, both instances were deployed in the us-central1 region where the mean network latency was 0.15ms and the bandwidth was about 1.5GB/s. For the WAN setting, one instance was in us-central1 while the other was in us-west2; the mean network latency was 49ms and the bandwidth was about 50MB/s.

7.1 Secure Exponentiation Experiments

We benchmark our secure exponentiation protocol separately and present our results and comparisons here.

(l_f, l)		64-bit BA	seInt		128-bit BaseInt				
	Offino	Onlino	End-to-End		Offino	Onlino	End-to-End		
	Onnie	Omme	LAN	WAN	Omme	Omme	LAN	WAN	
(5, 32)	3.09	0.004	4.21	11.06	7.68	1.01	9.56	16.29	
(10, 63)	3.22	0.99	5.43	12.07	9.24	14.9	25.02	32.17	
(15, 63)	3.22	1.01	5.49	12.11	9.24	16.0	26.29	32.95	
(20, 100)	-	-	-	-	11.15	33.2	44.91	52.35	
(20, 127)	-	-			12.9	54.9	68.63	75.91	

Table 1: Timing benchmarks (in μ s) for the exponentiation protocol, for base 64-bit and 128-bit int sizes. Exponents in the range [-5, 5] were randomly sampled and shared in the fixed-point ring. Offline and online phase computation times (in μ s) are averaged over 1 million runs, and don't include communication. End-to-end times per exponentiation are given in the LAN and WAN settings where 100K exponentiations are batched for communication. End-to-end times include communication and communication costs for both the online and offline phases.

1.	Our ap	proach	Polynomial Approx. [10]				
^t f	μ	σ	μ	σ			
5	0.0286	0.011	0.0932	0.0202			
10	0.0009	0.0003	0.0051	0.0019			
20	9.2×10^{-7}	3.5×10^{-7}	6.6×10^{-6}	5×10^{-6}			
30	9.1×10^{-10}	3.5×10^{-10}	1.3×10^{-8}	2×10^{-8}			
40	8.9×10^{-13}	3.4×10^{-13}	1.8×10^{-9}	7.4×10^{-9}			

Table 2: Mean (μ) and standard deviation (σ) of the fractional exponentiation error as a ratio of the actual result for both our approach and polynomial approximation (as in [10]). Exponents are sampled and shared randomly. The error is averaged over 1 million runs.

Timing experiments. We provide the offline and online computation times as well as end-to-end benchmarks (both LAN and WAN) for several (l, l_f) parameters and for both 64-bit and 128-bit base integer sizes. The results are shown in Table 1. We find that especially when batching the communication for several exponentiations together, the impact of the network is quite minimal, primarily due to the small amount of communication our protocol requires.

Accuracy experiments. As mentioned earlier, our exponentiation has smaller error than standard techniques. Since, the error comes only from the fractional part, we implement both our fractional exponentiation as well as a degree-9 polynomial approximation used in [10] and compare the errors in Table 2. Our errors are smaller by 1 to 2 orders of magnitude and the difference gets wider with more fractional bits.

Apart from the smaller errors, we also note that our technique only requires a single round of communication while a degree-d polynomial approximation usually takes d rounds when implemented using Horner's method (as in done in [10] to reduce total communication). One shortcoming however, is that while it is straightforward to extend polynomial approximation to the malicious setting, it is not obvious how to efficiently do the same for our technique.

Benchmark comparisons. We did not find any prior work on fixed-point exponentiation that targets the same setting we do. The most relevant protocols are the ones in SCALE-MAMBA [10] (benchmarked in [9]) and Aliasgari et al. [7]. The protocol from [10] uses fixed-points but focuses primarily on active security; the one from [7] is in the semi-honest setting but uses floating-points, is described only for $n \ge 3$ parties and only for Shamir shares. Consequently, while we provide some comparison points, our comparison is not direct and comes with significant caveats. We intend the comparison to be primarily directional, and to highlight the difference in broader protocol approaches. Specifically, we believe the comparison shows the simplicity of our design in the 2-party semi-honest setting, and the corresponding performance gains (often μ s vs ms or s).

First, we compare to the 2-party protocol from [9, 10]. As noted earlier, this protocol targets the active-security setting, while we target the semi-honest setting, so the comparison is not direct. [10] requires a full bit decomposition and uses a polynomial approximation for the fractional part, which incurs a larger error than our approach. For fixed-point exponentiation with l = 245 and $l_f = 40$, [9] shows an online runtime of 15 ms, an offline runtime of 18000 ms, and an offline cost of 1337 Beaver triples, 1 square tuple, and 7688 shared bits, which comes out to ~2MB per exponentiation. In contrast, for those parameters, our total offline cost is 980 bits, i.e., a 2000x improvement. Our implementation only supports a maximum of l = 127, and therefore our comparison is not direct, but for $(l_f, l) = (20, 127)$, our online runtime was 0.055 ms, and our offline time was 0.013 ms.

For comparable parameters, the offline time for the arctan operation in [9], which requires fewer preprocessed bits than exponentiation, is still around 7000 ms, which implies a more than 500,000x improvement. [9] notes that the reason large parameters were chosen specifically for exponentiation (as opposed to (20, 128) for other functions like square-root, sine, cosine etc.), was the high numerical instability. This is not observed in our protocol for the parameters (20, 127), which is in part due to our ability to effectively exponentiate the fractional component in \mathbb{R} rather than using polynomial approximation.

Our protocol also has a large throughput advantage. While [9] reports 76 ops/s when 50 invocations are run in parallel, we achieve $\sim 15,000$ ops/s run sequentially for our 127-bit modulus. We also note that the implementation from [9] leverages multiple threads while all our code is single threaded and could potentially be optimized further.

Aliasgari et al. [7] provide a secure exponentiation protocol in the semi honest setting. They consider floating-point exponentiation in the 3-party setting with Shamir shares. This is significantly different from our setting, since we target fixed-point exponentiation in the 2-party setting. The comparison therefore comes with significant caveats, but we provide a brief analytical comparison here to highlight the differences in techniques, and therefore efficiency. In particular, the protocol from [7] requires a full bit decomposition, 4 comparison tests, and l_f floating-point multiplications (where l_f is number of significant bits). For a l_f -bit comparable precision (for their best setting where l_f is more than the number of exponent bits k), it requires at least $16 + 12 \log l_f + \log \log l_f$ rounds and $\mathcal{O}(k) + \mathcal{O}(l_f \log l_f)$ interactive operations (involving exchange of a secret share) taking preprocessing and parallel computation into account. In comparison, our protocol requires a single round and only one interactive operation (i.e., only one secret share is exchanged) regardless of land l_f .

	l = 63										
Deteget		m	Paillier [21]			Corr	elated Trip	les A	Correlated Triples B		
Dataset	n	m	LAN (s)	WAN (s)	Comm.	LAN (s)	WAN (s)	Comm.	LAN (s)	WAN (s)	Comm.
	1	1	0.0638	0.0732	3 KB	0.0043	0.0044	0.48 KB	0.00006	0.00017	$0.48~\mathrm{KB}$
	100	10	4.2724	4.6407	0.17 MB	0.0165	0.0175	9.60 KB	0.0022	0.0033	9.60 KB
		100	13.469	14.046	$0.30 \ \mathrm{MB}$	0.1392	0.1435	24.0 KB	0.0162	0.0178	24.0 KB
		1000	105.48	105.98	1.65 MB	1.3659	1.4024	168 KB	0.1621	0.1652	168 KB
Roplicated	1000	10	40.085	40.462	1.52 MB	0.0210	0.0227	81.6 KB	0.0202	0.0224	81.6 KB
Replicated		100	106.94	107.62	1.65 MB	0.1556	0.1561	96.0 KB	0.1484	0.1509	96.0 KB
		1000	776.68	777.24	3.00 MB	1.5033	1.5065	240 KB	1.4914	1.4928	240 KB
		10	395.42	396.04	15.0 MB	0.0926	0.1091	805 KB	0.1994	0.2183	$805~\mathrm{KB}$
	10000	100	1041.3	1041.7	$15.2 \ \mathrm{MB}$	0.5213	0.5352	848 KB	1.4749	1.4885	848 KB
		1000	7492.0	7498.4	16.5 MB	5.5495	5.5641	1.28 MB	14.740	14.796	1.28 MB
Somoza	21	11	1.1794	1.2350	48 KB	0.0176	0.0182	3.68 KB	0.0007	0.0011	3.68 KB
PhD	73	17	3.7691	3.8559	135 KB	0.0259	0.02659	8.48 KB	0.0025	0.0035	8.8 KB
Cancer	36	14	1.9553	2.0278	75 KB	0.0218	0.0225	$5.12~\mathrm{KB}$	0.0012	0.0018	$5.12~\mathrm{KB}$

l = 127

Deteret				Paillier [21]		Corr	related Trip	les A	Corr	related Trip	les B
Dataset	n		LAN (s)	WAN (s)	Comm.	LAN (s)	WAN (s)	Comm.	LAN (s)	WAN (s)	Comm.
	1	1	0.06563	0.0752	3 KB	0.0048	0.0051	0.89 KB	0.00007	0.00024	0.89 KB
		10	5.0371	5.3857	0.17 MB	0.0180	0.0192	17.8 KB	0.0024	0.0037	17.8 KB
	100	100	20.088	20.676	0.30 MB	0.1488	0.1495	44.4 KB	0.0179	0.0195	44.4 KB
		1000	171.18	171.85	1.65 MB	1.4597	1.4620	311 KB	0.1759	0.1814	311 KB
Doplicated		10	47.739	48.342	1.52 MB	0.0229	0.0263	151 KB	0.0227	0.0265	151 KB
Replicated	1000	100	173.58	174.84	1.65 MB	0.1602	0.1642	178 KB	0.1601	0.1640	178 KB
		1000	1433.1	1434.4	3.00 MB	1.5889	1.6150	444 KB	1.6003	1.6006	444 KB
		10	474.36	476.59	15.0 MB	0.1077	0.1331	1.49 MB	0.2234	0.2506	1.49 MB
	10000	100	1709.5	1710.3	15.2 MB	0.5737	0.6008	$1.57 \ \mathrm{MB}$	1.5840	1.6158	$1.57 \ \mathrm{MB}$
		1000	14053	14056	16.5 MB	5.9126	6.0368	2.36 MB	15.920	15.956	2.36 MB
Somoza	21	11	1.3610	1.4202	48 KB	0.0192	0.0200	6.81 KB	0.0007	0.0012	6.81 KB
PhD	73	17	4.6749	4.7559	135 KB	0.0282	0.0296	15.7 KB	0.0028	0.0042	16.3 KB
Cancer	36	14	2.3283	2.4085	75 KB	0.0233	0.0245	9.47 KB	0.0013	0.0021	9.47 KB

Table 3: Micro benchmarks for generation of correlated Beaver triples $([\mathbf{A}], [\mathbf{B}_i], [\mathbf{C}_i] = [\mathbf{AB}_i])$ in the offline phase for l = 63 and l = 127 bits. A has dimension $n \times m$; the \mathbf{B}_i have dimension $m \times 1$. The plaintext modulus used is 2^{l} . Times (in seconds) and communication cost for correlated triples are amortized for one triple over 1000 iterations. The baseline cost for triple generation via Paillier encryption (with a 3072-bit keysize) is averaged over 5 iterations. All of our code is single threaded and is run in the LAN setting.

7.2**Offline Phase Experiments**

We provide micro benchmarks for the offline phase generation of correlated Beaver triples in both the LAN and WAN settings in Table 3. As a baseline, we also compare to the cost when using Paillier encryption. Both our approaches (see Section 5.2) have 5x-10x less communication cost and are 520x-4200x faster than Paillier AHE-based approaches. Experiments in [21] suggest that an OT-based approach is 20x-30x faster than the Paillier AHE-based one in the LAN setting which highlights that our protocols would also be faster than OT-based triple generation.

7.3 Poisson Regression Experiments

We now measure the performance of our end-to-end secure Poisson regression protocol and compare its accuracy with plaintext regression, where the data is provided without encryption. Our secure regression is implemented with fixed-point numbers, while the C++ double type is used in the plaintext version.

Datasets. We run our regression experiments on three datasets (detailed next) from the Princeton University course on Generalized Linear Models [1].

- 1. Somoza. This dataset contains infant and child survival rates in Colombia, based on the World Fertility Survey. Survival is modeled as a function of sex, cohort, and age range. The dataset tracks 2000 infants over several years, and provides aggregate exposures and counts over 21 distinct feature combinations.
- 2. *Time to PhD.* This dataset predicts PhD graduation as a function of years in graduate school, university, and residence status. We encode the explanatory variables into 17 binary features. Data from 35,000 PhD students is used to calculate the aggregate exposure period and graduation counts for 73 distinct feature combinations.
- 3. Smoking and Cancer. This dataset contains information from a Canadian study of mortality by age and smoking status. There are 14 different binary features, corresponding to different age buckets and smoking statuses. There are 36 distinct feature combinations, containing counts and exposure periods from a total of 92,000 respondents.

All of the datasets are publicly available at [1].

Accuracy evaluation. To quantify accuracy, we benchmark our secure Poisson regression protocol against a plaintext regression baseline for different learning rates and fixed-point precision. See Figures 10 and 11. We observe that our secure protocol performs almost exactly as well as the plaintext regression: the lines plotted for model error versus number of iterations are nearly coincident.

When we take a closer look at the learned parameter θ , we find that the actual weights learned by the secure protocol are also nearly exactly the same as those from plaintext learning. See Table 5: the root mean square error between the secure weights and the plaintext weights is very small regardless of the dataset being tested on.

Performance evaluation. We also benchmark the computation and communication efficiency of our end-to-end protocol regression protocol in Table 4. In addition to the earlier datasets, we also run our experiments on larger synthetic datasets. For this, we replicate the Somoza dataset to obtain a new dataset of the appropriate size $(n \times m)$. We report our timing results for this under the "Replicated" dataset header.

As there is no previous work done on secure Poisson regression, it is not possible for us to compare efficiency of our protocol with other work. Instead, we compare our protocol with a "basic" version that does not use correlated Beaver triples. We still use our exponentiation protocol. For correlated triples, since the gain is only when multiple gradient descent iterations are run, for our timing values, we run 1000 iterations, and report the amortized time for 1 iteration.



(c) Smoking and Cancer

Figure 10: Convergence of the RMSE for plaintext regression versus Secure Poisson regression with 15-bit fixed-point precision.

We find that our protocol performs well, even for larger datasets. For example, in the LAN setting, for a dataset with 10,000 elements and 100 features, it has an amortized cost of 3.116 seconds of offline time, 5.501 seconds of online time, and 14.8 MB of communication. Over 100 iterations, the cost is about 5 minutes of offline time, 9 minutes of online time, and 1.48 GB of communication.

8 Applications

In this section, we give several concrete applications for secure Poisson regression, and discuss performance of our protocol in each of these scenarios.

8.1 COVID-19 Case Fatality Rate

Recent work [29] performs an analysis of COVID-19 case fatality using Poisson Regression. They measure the effect of 9 binary variables on the counts of COVID-19 fatalities, using 2070 cases as training examples. Variables include age-range (≥ 60 years), presence of cardiovascular disease, and presence of neurologic diseases. The regression model is used to compute the incidence rate



(c) Smoking and Cancer

Figure 11: Convergence of the RMSE for plaintext regression versus Secure Poisson regression with 20-bit fixed-point precision.

ratio (IRR) for each variable, that is, the ratio between predicted fatalities when that variable is present versus not.

This case provides a good example for health data, where multiple hospitals may hold slices of the data, and may not want it to be centralized in the clear. To compute over this data privately, hospitals could send shares of the data to two servers who could perform Poisson regression securely, and compute shares of the model parameters. The model could then be sent to each hospital which would individually compute the IRR for each variable, and release the aggregate IRRs.

On a synthetic dataset with similar shape, in the LAN setting, our protocol takes 0.268 seconds in the offline phase and 0.103 seconds in the online phase per iteration of gradient descent, with a total communication cost of 3.52 MB. Assuming 100 iterations of gradient descent are needed in order to converge, this results in 26.8 seconds in the offline phase, 10.3 seconds in the online phase, and a communication of 352 MB.

8.2 Predicting Credit Default Rates

[20] use Poisson Regression to model the rate of default payments by borrowers. They measure the effect of 6 variables, including income, age, monthly credit card expenditure, and home-ownership on the monthly rate of defaulted loan payments using a sample of 1002 individuals. After regression,

					Stan	dard			Correlated Triples					
Detect			Offline			Online			Offline			Online		
Dataset		111	LAN	WAN	Comm.	LAN	WAN	Comm.	LAN	WAN	Comm.	LAN	WAN	Comm.
			(s)	(s)	(MB)	(s)	(s)	(MB)	(s)	(s)	(MB)	(s)	(s)	(MB)
		10	15.110	16.415	0.638	0.001	0.198	0.038	0.025	0.026	0.075	0.0006	0.196	0.006
	100	100	33.317	35.408	0.908	0.006	0.199	0.314	0.039	0.054	0.104	0.002	0.199	0.008
		1000	218.80	220.92	3.61	0.059	0.241	3.075	0.318	0.335	0.392	0.022	0.217	0.049
	1000	10	142.56	152.30	6.11	0.006	0.205	0.382	0.129	0.241	0.721	0.002	0.190	0.062
Repl.		100	276.21	286.77	6.38	0.056	0.240	3.130	0.381	0.502	0.750	0.021	0.207	0.062
		1000	1617.2	1627.6	9.08	0.556	1.002	30.61	3.053	3.172	1.04	0.209	0.405	0.107
		10	1420.1	1514.5	60.8	0.055	0.238	3.815	1.119	2.296	7.19	0.020	0.208	0.614
	10000	100	2734.0	2833.2	61.1	0.542	1.032	31.28	2.844	4.050	7.28	0.202	0.389	0.643
		1000	15896	16000	63.8	5.91	10.902	306.0	61.746	64.371	8.14	2.101	2.469	0.931
Sozoma	21	11	3.6838	4.007	0.161	0.0005	0.188	0.008	0.003	0.006	0.020	0.0004	0.188	0.001
PhD	73	17	12.089	13.053	0.495	0.0009	0.194	0.040	0.011	0.021	0.063	0.0007	0.188	0.002
Cancer	36	14	6.1600	6.690	0.261	0.0006	0.193	0.017	0.005	0.011	0.032	0.0005	0.188	0.001

 $(l_f, l) = (15, 63)$

 $(l_f, l) = (20, 127)$

					Stan	dard					Correlat	ted Triple	es		
Deteret				Offline			Online			Offline			Online		
Dataset	n		LAN	WAN	Comm.	LAN	WAN	Comm.	LAN	WAN	Comm.	LAN	WAN	Comm.	
			(s)	(s)	(MB)	(s)	(s)	(MB)	(s)	(s)	(MB)	(s)	(s)	(MB)	
		10	16.700	18.006	0.644	0.008	0.201	0.077	0.028	0.034	0.139	0.006	0.200	0.013	
	100	100	46.740	48.877	0.914	0.065	0.271	0.629	0.044	0.065	0.192	0.062	0.254	0.016	
		1000	351.32	354.22	3.61	0.653	1.172	6.149	0.344	0.371	0.725	0.618	0.963	0.049	
	1000	10	158.52	168.74	6.17	0.059	0.260	0.763	0.277	0.324	1.33	0.056	0.245	0.123	
Repl.		100	410.40	421.94	6.44	0.593	1.035	6.259	0.414	0.604	1.39	0.558	0.846	0.131	
		1000	2931.8	2944.1	9.14	5.888	8.474	61.22	3.256	3.460	1.92	5.562	6.045	0.214	
		10	1582.2	1680.5	61.5	0.584	1.023	7.630	2.650	3.107	13.3	0.549	1.001	1.227	
	10000	100	4077.7	4181.3	61.7	5.825	8.342	62.56	3.116	4.997	13.5	5.501	5.989	1.285	
		1000	29040	29154	64.4	59.745	72.234	611.9	65.82	70.61	15.1	55.144	55.862	1.862	
Sozoma	21	11	4.067	4.421	0.162	0.002	0.199	0.016	0.003	0.008	0.037	0.002	0.197	0.002	
PhD	73	17	13.960	15.005	0.500	0.009	0.202	0.081	0.012	0.028	0.116	0.008	0.201	0.004	
Cancer	36	14	6.950	7.526	0.263	0.004	0.199	0.034	0.006	0.014	0.060	0.004	0.199	0.003	

Table 4: Benchmarks for the end-to-end Poisson regression protocol for different datasets. n is the number of examples and m is the number of features. For larger values of n and m, the Somoza dataset was replicated. Times (in seconds) are given per iteration of gradient descent over the entire dataset. For the "Standard" column, we use standard Beaver triples generated via Paillier encryption (e.g., as in [21]) along with our fixed-point exponentiation protocol. For correlated triples, the protocol is amortized over 1000 iterations. All of our code is single threaded.

the authors propose using the model inference to data of loan applicants to compute predicted defaults, and thereby characterize risk level.

This case involves training on sensitive financial data, which may be distributed across several institutions. Securely computing regression on these values would then consist of two phases: combining the records from multiple institutions, followed by performing secure regression on the joint data. The former task can be handled using techniques like privacy-preserving record linkage [5]. Our secure protocol is a good fit for the latter part, as well as the subsequent inference.

On a synthetic dataset with similar shape, in the LAN setting, our protocol incurs 5.8 seconds of offline time and 3.4 seconds of online time to perform 100 iterations of secure gradient descent, with a total communication cost of 157 MB. Each iteration would incur 0.058 seconds and 0.034

		RMSE b	etween plaintext	weights and secure weights		
Learning rate	Iterations	Somoza	Time to PhD	Smoking and		
		Somoza	Time to PhD	Lung Cancer		
	100	0.00064	-	0.00016		
0.0001	500 1000	0.00259	-	0.00048		
		0.00456	-	0.00097		
	100	0.00034	0.00031	0.00021		
0.00005	500	0.00160	0.00123	0.00057		
	1000	0.00346	0.00200	0.00150		
	100	0.00029	0.00030	0.00023		
0.00003	500	0.00131	0.00126	0.00060		
	1000	0.00294	0.00228	0.00107		

Table 5: Root mean square errors between the weights obtained from secure regression and those from plaintext regression. This table shows that the learned weights from secure regression are nearly the same as those obtained from plaintext regression.

seconds of offline and online time, with 1.57 MB of communication.

8.3 Modeling Ad Campaign Conversion Rates

Google researchers [28] describe a system for measuring ad campaign conversion rates using Poisson regression. A "conversion" corresponds to an individual buying an item after seeing one or more ads. [28] give several ways to model multiple ad channels having a combined effect on an individual, with the ad effects decaying over time. One is to use a "step" decay: assigning each ad channel 3 binary attributes, corresponding to whether an individual was exposed to the ad in the short term (1 day prior), medium term (2-7 days prior) or long term (7-30 days prior). The conversion rate is then learned via Poisson regression using such attributes for some combination of ad channels. Credit for a conversion is proportionally distributed to each ad channel according to the relative change in predicted conversion rate when that ad channel is switched from exposed to unexposed. The total credit per ad channel is computed as the sum of its proportional credit across all conversions in the dataset.

This problem is an excellent case for the use of secure computation techniques, since it involves sensitive business and user data that may be held by different ad companies and transaction data providers. A secure solution would require privately joining the records, securely performing regression, and then securely computing the aggregate credit for each ad channel. The private join could be achieved using privacy-preserving record linkage techniques [5]. Our work is well-suited for regression as well as the subsequent inference.

On a synthetic dataset with 5 ad channels and 3 binary attributes per channel for a total of 15 binary attributes, and assuming 100,000 training points, our regression takes 6.90 seconds of offline time and 8.197 seconds of online time per iteration of gradient descent, with 156.7 MB of total communication. For 100 iterations of gradient descent, we incur 11.5 minutes of offline time and about 14 minutes of online time, with 15.67 GB of total communication.

9 Conclusion

Poisson regression is a widely used technique for modeling Poisson processes that occur across the life and social sciences. In many settings, the inputs for training Poisson models are sensitive health or financial data held by different parties. The secure Poisson regression protocol introduced in this paper enables computation on private data which reveals only the output Poisson model while protecting the inputs. Our construction achieves this with great efficiency while preserving accuracy comparable to computation in the clear. For several real datasets, this means execution in just a few seconds with a couple MB of communication. At the crux of our protocol is a new construction for secure fixed-point exponentiation and a new technique for correlated matrix multiplication, both of which are of independent interest with applications far beyond Poisson regression.

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