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Have you forgotten? A method to assess if machine learning models have forgotten data

Xiao Liu¹ and Sotirios A. Tsaftaris^{1,2}

¹ School of Engineering, University of Edinburgh, Edinburgh EH9 3FB, UK ² The Alan Turing Institute, London, UK Xiao.Liu@ed.ac.uk, S.Tsaftaris@ed.ac.uk

Abstract. In the era of deep learning, aggregation of data from several sources is a common approach to ensuring data diversity. Let us consider a scenario where several providers contribute data to a consortium for the joint development of a classification model (hereafter the target model), but, now one of the providers decides to leave. This provider requests that their data (hereafter the query dataset) be removed from the databases but also that the model 'forgets' their data. In this paper, for the first time, we want to address the challenging question of whether data have been forgotten by a model. We assume knowledge of the query dataset and the distribution of a model's output. We establish statistical methods that compare the target's outputs with outputs of models trained with different datasets. We evaluate our approach on several benchmark datasets (MNIST, CIFAR-10 and SVHN) and on a cardiac pathology diagnosis task using data from the Automated Cardiac Diagnosis Challenge (ACDC). We hope to encourage studies on what information a model retains and inspire extensions in more complex settings.

Keywords: Privacy · Statistical measure · Kolmogorov-Smirnov

1 Introduction

Deep learning requires lots of, diverse, data and in healthcare likely we will need to enlist several data providers (e.g. hospital authorities, trusts, insurance providers who own hospitals, etc) to ensure such data diversity. To develop models, data from several providers must thus be either centrally aggregated or leveraged within a decentralized federated learning scheme that does not require the central aggregation of data (e.g. [1, 13, 14, 17, 19]). Thus, several providers will contribute data for the development of a deep learning model e.g. to solve a simple classification task of normal vs. disease. Suddenly, one of the providers decides to leave and asks for the data to be deleted but more critically that *the model 'forgets' the data*.

Let us now assume that the model has *indeed* not used the data and has 'forgotten' them (we will not care how herein) but we want to *verify* this. In other words, and as illustrated in Fig. 1, our problem is to assess whether the model has not used the data under question in the training set.



Fig. 1. Our goal. We aim to develop an approach that an 'auditor' can use to ascertain whether a model $f(x; D^*; \theta)$ has used data D_Q during training.

We consider that an *auditor* will have access to the query dataset D_Q and the model $f(\mathbf{x}; D^*, \theta^*)$, (trained on data D^*) to render a decision whether f() retains information about D_Q . We assume that the training dataset D^* is *unknown* but both D^* and D_Q are sampled from a *known* domain \mathbb{D} for a given *known* task.¹ We emphasize that we are *not* proposing a method that forgets data, as for example the one in [8]. We believe that assessing whether data have been forgotten is an important task that should come first (because without means to verify claims, anyone can claim that data have been forgotten).

Related work is limited. In fact, our communities (machine learning or medical image analysis) have not considered this problem yet in this form. The closest problem in machine learning is the *Membership Inference Attack* (MIA) [4, 5, 16, 18, 20]. While these works borrow concepts from MIA, as we detail in the related work section, MIA sets different hypotheses that have considerable implications in returning false positives when data sources overlap. Our major novelty is setting appropriate test hypotheses that calibrate for data overlap.

Our approach builds on several key ideas. First, inspired by [18], we adopt the Kolmogorov-Smirnov (K-S) distance to compare statistically the similarity between the distribution of a model f()'s outputs and several purposely constructed reference distributions. To allow our approach to operate in a blackbox setting (without precise knowledge of f(), we construct 'shadow' models (inspired by [20]) to approximate the target model f(); however, we train the shadow models on D_Q but also on another dataset D_C sampled from domain \mathbb{D} which does not overlap (in element-wise sense) D_Q .

Contributions:

- 1. To introduce a new problem in data privacy and retention to our community.
- 2. To offer a solution that can be used to detect whether a model has forgotten specific data including the challenging aspect when data sources may overlap.
- 3. Experiments in known image classification benchmark datasets to verify the effectiveness of our approach. Experiments on the pathology classification component of the ACDC simulating a healthcare-inspired scenario.

¹ Knowledge of the task (e.g. detect presence of a pathology in cardiac MRI images), implies knowledge of the domain \mathbb{D} (e.g. the space of cardiac MRI images). Without this assumption, D^* can be anything, rendering the problem intractable.

2 Related work

The purpose of *Membership Inference Attack* is to learn which data were used to train a model. The hypothesis is that a model *has* used some data and the goal is to find which part of the query data are in the training set. Here we briefly review key inspirational approaches.²

In [20] they train a model f_{attack} that infers whether some data **D** were used by f. f_{attack} accepts as inputs the decisions of f (the softmax outputs), **D** and the ground-truth class. Their premise is that machine learning models often behave "differently on the data that they were trained on vs. data that they 'see' for the first time". Thus, the task reduces to training this attack model. They rely on being able to generate data that resemble **D** (or not) and train several, different, 'shadow' models that mimic f to obtain outputs that can be used to train f_{attack} . We rely here on the same premise and use the idea of training shadow models of f() to enable gray-box inference on f() using only its outputs.

In [18], in the context of keeping machine learning competitions fair, using the same premise as [20], they propose a statistical approach to compare model outputs, and eschew the need for f_{attack} . They use Kolmogorov-Smirnov (K-S) distance to measure the statistical similarity between emissions of the classifier layer of a network between query data and a reference dataset to see whether models have used validation data to train with. However, their approach assumes that both query (validation set) and reference datasets, which in their context is the testing set, are known *a priori*. We adopt (see Section 3) the K-S distance as a measure of distribution similarity but different from [18] we construct reference distributions to also calibrate for information overlap.

Why our problem is not the same as MIA: An MIA algorithm, by design, assumes that query data may have been used to train a model. Whereas, we care to ascertain if a model has not used the data. These are different hypotheses which appear complementary but due to data source overlap have considerable implications on defining false positives and true negatives.³ When there is no information overlap between datasets, a MIA algorithm will return the right decision/answer. However, when D_Q and D^* statistically overlap (i.e. their manifolds overlap), for a sample both in D_Q or D^* , a MIA algorithm will return a false positive even when the model was trained on D^* alone. In other words, MIA cannot tell apart data overlap. Since overlap between D_Q and D^* is likely and a priori the auditor does not know it, the problem we aim to address has more stringent requirements than MIA. To address these requirements we assume that we can sample another dataset D_C from \mathbb{D} (that we can control not to at least sample-wise overlap with D_Q) to train shadow models on D_C .

 $^{^{2}}$ We do not cover here the different task of making models and data more private, by means e.g. of *differential privacy*. We point readers to surveys such as [9, 10] and a recent (but not the only) example application in healthcare [14].

³ In fact, overlap will frequently occur in the real world. For example, datasets collected by different vendors can overlap if they collaborate with a same hospital or if a patient has visited several hospitals. Our method has been designed to address this challenging aspect of overlap.

Algorithm 1 The proposed method.

Input: the target model $f(\mathbf{x}; D^*, \theta^*)$, query dataset D_Q and dataset domain \mathbb{D} . Output: answer to hypothesis if $f(\mathbf{x}; D^*, \theta^*)$ is trained with D_Q . Step 1. Train $\tilde{f}(\mathbf{x}; D_Q, \theta_Q)$ with the same design of $f(\mathbf{x}; D^*, \theta^*)$. Step 2a. Sample the calibration dataset D_C from domain \mathbb{D} but without overlapping (in sample-wise sense) D_Q . Step 2b. Train $\tilde{f}(\mathbf{x}; D_C, \theta_C)$ with the same design of $f(\mathbf{x}; D^*, \theta^*)$. Step 3. Find $r(T_{D_Q|D_Q}), r(T_{D_Q|D^*})$ and $r(T_{D_Q|D_C})$ (Eq. 1). Step 4. Find $KS(r(T_{D_Q|D_Q}), r(T_{D_Q|D^*})), KS(r(T_{D_Q|D_Q}), r(T_{D_Q|D_C}))$ (Eqs. 2, 3). Step 5. Find $\rho = \frac{KS(r(T_{D_Q|D_Q}), r(T_{D_Q|D^*}))}{KS(r(T_{D_Q|D_C}))}$ (Eq. 4). return If $\rho \ge 1$, the target model has forgotten D_Q . Otherwise ($\rho < 1$), has not.

3 Proposed method

We measure similarity of distributions to infer whether a model is trained with the query dataset D_Q , i.e. if D^* has D_Q or not. Note that this is not trivial since D^* is unknown. In addition, the possible overlap between D^* and D_Q introduces more challenges as we outlined above. To address both, we introduce another dataset D_C sampled from domain \mathbb{D} but without overlapping (in element-wise sense) D_Q . The method is summarized in Algorithm 1 with steps detailed below. Notation: We will consider x, a tensor, the input to the model e.g. an image. We will denote the *target* model as $f(\mathbf{x}; D^*, \theta^*)$ that is trained on dataset D^* parametrised by θ^* . Similarly, we define query $f(\mathbf{x}; D_Q, \theta_Q)$ and calibration models $f(\mathbf{x}; D_C, \theta_C)$, where both models share model design with the target model. If D_Q has N samples $(N \gg 1)$, we use $y^{\{i\}}, i = \{1, \dots, N\}$, a scalar, to denote the ground-truth label of data $\mathbf{x}^{\{i\}}$ of D_Q . We denote with $\mathbf{t}^{\{i\}}, i = \{1, \dots, N\}$, a *M*-dimensional vector, the *output* of a model with input $\mathbf{x}^{\{i\}}$. Hence, $\mathbf{t}^{\{i\}}[y^{\{i\}}]$ denotes the output of the model $\mathbf{t}^{\{i\}}$ for the ground-truth class $y^{\{i\}}$ (confidence score). We further define the $N \times 1$ vector $\mathbf{c}_{D_Q|D^*}$, which contains the values of $\mathbf{t}^{\{i\}}[y^{\{i\}}]$, $i = \{1, \dots, N\}$ in an increasing order. We define as $T_{D_Q|D^*}$ the $N \times M$ matrix that contains all outputs of the target model $f(\mathbf{x}; D^*, \theta^*)$ that is tested with D_Q . Similarly, we define $T_{D_Q|D_Q}$ and $T_{D_Q|D_C}$ for the outputs of $f(\mathbf{x}; D_Q, \theta_Q)$ and $f(\mathbf{x}; D_C, \theta_C)$, respectively, when both are tested with D_Q .

3.1 Assumptions on f()

We follow a gray-box scenario: access to the output of the model $f(\mathbf{x}; D^*, \theta^*)$ before any thresholds and knowledge of design (e.g. it is a neural network) but not of the parameters (e.g. weights) θ^* . (Model providers typically provide the best (un-thresholded) value as surrogate of (un)certainty.)

3.2 Kolmogorov-Smirnov distance between distributions

We denote as $r(T_{D_Q|D^*})$ (a $N \times 1$ vector) the empirical cumulative distribution (cdf) of the output of the target model when tested with D_Q . The n^{th} element $r_n(T_{D_Q|D^*})$ is defined as:

$$r_n(T_{D_Q|D^*}) = \frac{1}{N} \sum_{i=1}^N I_{[-\infty, \mathbf{c}_{D_Q|D^*}^{\{n\}}]}(\mathbf{c}_{D_Q|D^*}^{\{i\}}), \tag{1}$$

where $I_{[-\infty, \mathbf{c}_{D_Q|D^*}^{\{n\}}]}(\mathbf{c}_{D_Q|D^*}^{\{i\}})$ the indicator function, is equal to 1 if $\mathbf{c}_{D_Q|D^*}^{\{i\}} \leq \mathbf{c}_{D_Q|D^*}^{\{n\}}$ and equal to 0 otherwise.

Our next step is to create a proper empirical distribution as reference to compare against $r(T_{D_Q|D^*})$. Motivated by [20], we propose to train a query model $\tilde{f}(\mathbf{x}; D_Q, \theta_Q)$ with the same model design as the target model. We then obtain (as previously) the output cumulative distribution $r(T_{D_Q|D_Q})$. We propose to use $r(T_{D_Q|D_Q})$ as the reference distribution of dataset D_Q . Hence, measuring the similarity between $r(T_{D_Q|D^*})$ with $r(T_{D_Q|D_Q})$ can inform on the relationship between D_Q and D^* , which is extensively explored in the field of dataset bias [21]. Following [18], we use Kolmogorov–Smirnov (K-S) distance as a measure of the similarity between the two empirical distributions.

K-S distance was first used in [6] to compare a sample with a specific reference distribution or to compare two samples. For our purpose we will peruse the twosample K-S distance, which is given by:

$$KS(r(T_{D_Q|D_Q}), r(T_{D_Q|D^*})) = \sup |r(T_{D_Q|D_Q}) - r(T_{D_Q|D^*})|_1,$$
(2)

where sup denotes the largest value. K-S distance $\in [0, 1]$ with lower values pointing to greater similarity between $r(T_{D_Q|D^*})$ and $r(T_{D_Q|D_Q})$.

If D^* contains D_Q , the K-S distance between $r(T_{D_Q|D_Q})$ and $r(T_{D_Q|D^*})$ will be very small and ≈ 0 , i.e. $KS(r(T_{D_Q|D_Q}), r(T_{D_Q|D^*})) \approx 0$. On the contrary, if D^* has no samples from D_Q , then the value of $KS(r(T_{D_Q|D_Q}), r(T_{D_Q|D^*}))$ depends on the statistical overlap of D^* and D_Q . However, this overlap is challenging to measure because the training dataset D^* is assumed unknown (and we cannot approximate statistical overlap with element-wise comparisons).

3.3 Calibrating for data overlap

We assume the training data D^* of model $f(\mathbf{x}; D^*, \theta^*)$ are sampled from a domain \mathbb{D} . To calibrate against overlap between D^* and D_Q , we sample a *calibration* dataset D_C from \mathbb{D} but ensure that no samples in D_Q are included in D_C (by sample-wise comparisons) and train the calibration model $\tilde{f}(\mathbf{x}; D_C, \theta_C)$. Inference on model $\tilde{f}(\mathbf{x}; D_C, \theta_C)$ with D_Q , the output empirical cumulative distribution $r(T_{D_Q|D_C})$ can be calculated in a similar fashion as Eq. 1. To compare the similarity between D_Q and D_C , we calculate:

$$KS(r(T_{D_Q|D_Q}), r(T_{D_Q|D_C})) = sup |r(T_{D_Q|D_Q}) - r(T_{D_Q|D_C})|_1.$$
(3)

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We argue that Eq. 3 can be used to calibrate the overlap and inform if D^* contains D_Q .⁴ We discuss this for two scenarios:

1. D^* does not include any samples of D_Q . Since D_C and D_Q do not overlap in a element-wise sense by construction, then data of D^* should have higher probability of statistical overlap with D_C . In other words, the calibration model can be used to mimic the target model. By sampling more data from domain \mathbb{D} as D_C , there will be more statistical overlap between D_C and D_Q . Hence, the statistical overlap between a large-size dataset D_C and D_Q can be used to approximate the (maximum) possible overlap between D^* and D_Q i.e. $KS(r(T_{D_Q|D_Q}), r(T_{D_Q|D^*})) \geq KS(r(T_{D_Q|D_Q}), r(T_{D_Q|D_C}))$. 2. D^* contains samples of D_Q . In this case, samples in D_Q that are also in D^*

2. D^* contains samples of D_Q . In this case, samples in D_Q that are also in D^* will introduce statistical overlap between D_Q and D^* , which is well demonstrated in [18]. Based on our experiments, we find that the overlap between D_Q and D^* is consistently less than that of D_Q and D_C i.e. $KS(r(T_{D_Q|D_Q}), r(T_{D_Q|D^*})) < KS(r(T_{D_Q|D_Q}), r(T_{D_Q|D_C}))$.

Thus, we propose to use $KS(r(T_{D_Q|D_Q}), r(T_{D_Q|D_C}))$ as a data-driven indicator for detecting if D^* contains D_Q . To quantify how much the target model has forgotten about D_Q , we capture the two inequalities above in the ratio:

$$\rho = \frac{KS(r(T_{D_Q|D_Q}), r(T_{D_Q|D^*}))}{KS(r(T_{D_Q|D_Q}), r(T_{D_Q|D_C}))}.$$
(4)

In the first scenario i.e. D^* does not include any samples of D_Q , the inequality $KS(r(T_{D_Q|D_Q}), r(T_{D_Q|D^*})) \geq KS(r(T_{D_Q|D_Q}), r(T_{D_Q|D_C}))$ translates to $\rho \geq 1$. Instead, in the second scenario, $\rho < 1$. Thus, if $\rho \geq 1$, this implies the target model has forgotten D_Q . On the contrary, if $\rho < 1$, implies the target model has not forgotten D_Q .

4 Experiments

We first perform experiments on benchmark datasets such as MNIST [12], SVHN [15] and CIFAR-10 [11] to verify the effectiveness of the proposed method. Then, we test our method on the ACDC dataset using the pathology detection component of the challenge [3]. All classifiers in our experiments are well-trained (adequate training accuracy) such that the output statistics well approximate the input statistics. All model designs are included as Supplementary Material.

4.1 Benchmark datasets - MNIST, SVHN and CIFAR-10

Benchmark datasets: MNIST contains 60,000 images of 10 digits with image size 28×28 . Similar to MNIST, SVHN has over 600,000 digit images obtained from house numbers in Google Street view images. The image size of SVHN is

⁴ As mentioned previously we cannot measure statistical overlap between D^* and D_C (or D_Q) since D^* is unknown. Furthermore, statistical overlap in high-dimensional spaces is not trivial [2,7], and hence we want to avoid it.

Table 1. Benchmark datasets results. The query dataset D_Q is MNIST. The training dataset D^* has different variants as listed below. The K-S distances are calculated with Eq. 2. D_C for defining ρ is SVHN. Bold number in the table is the threshold.

Training dataset (D^*)	MNIST	%SVHN			SVH	1 + %	CIEAR 10	
		50%	75%	100%	10%	50%	100%	
$KS(r(t_{D_Q D_Q}), r(t_{D_Q D^*}))$	0.049	0.669	0.652	0.596	0.056	0.039	0.029	0.957
ρ	0.082	1.122	1.094	1.000	0.094	0.065	0.049	1.606

 32×32 . Since both datasets are for the task of digit recognition/classification we consider to belong to the same domain \mathbb{D} . We use CIFAR-10 as an *out of domain* \mathbb{D} dataset to validate the method. CIFAR-10 has 60,000 images (size 32×32) of 10-class objects (airplane, bird, etc). To train our models, we preprocess images of all three datasets to gray-scale and rescale to size 28×28 .

Results and discussion: We consider MNIST as D_Q . Experimental results are listed in Table 1. We first verify if our method can detect the simple case where the target model has been fully trained with D_Q . According to our findings (first column of Table 1) as expected $\rho = 0.082 \ll 1$. Other cases where we assume D^* is drawn only from D_C (SVHN) give $\rho > 1$. In the challenging setting when the target model was trained with a mix of data from SVHN and parts of MNIST, $\rho \sim 0$. We draw attention to the case of SVHN+%100MNIST. $\rho = 0.049$ indicates that the target definitely has not forgotten the query dataset.

For exploration purposes, we consider the extreme case of D^* not in the domain of D_Q . We train a target model with CIFAR-10 which has little statistical information overlap with MNIST. Hence, we would expect our method to return a $\rho > 1$, which agrees with the result $\rho = 1.606$ (highest in the table).

Next we assess the scenarios that D^* does not include any samples of D_Q . We consider SVHN as D_C and train target models with D^* , subsets of SVHN. The results ρ of 50% and 75% SVHN are 1.122 and 1.094 that both higher than $\rho = 1$. Higher ρ points to lower information overlap between D_Q and D^* . Hence, the large-size D_C i.e. %100SVHN has the largest statistical overlap with D_Q .

To check the sensitivity of our method when D^* contains part of D_Q , we train two models with SVHN and part of MNIST (10% and 50%). We observe that both have $\rho < 1$. Specifically, including only 10% of MNIST causes $(1-0.094) \div 1$ = 90.6% drop of ρ value. This suggests that if the training dataset contains part of D_Q , our method can still accurately detect these cases.

Some uncertainty may arise if ρ is hovering around 1. However, we did not observe such cases when D^* contains samples from D_Q in our extensive experiments. We advise to run multiple runs of experiments to statistically eliminate such uncertainty if needed.

4.2 Medical dataset - ACDC

ACDC dataset: Automated Cardiac Diagnosis Challenge (ACDC) dataset is composed of cardiac images of 150 different patients that are acquired with two



Fig. 2. Example images of ACDC 1.5T and ACDC 3T datasets. DCM: dilated cardiomyopathy. HCM: hypertrophic cardiomyopathy. MINF: myocardial infarction. NOR: normal subjects. RV: abnormal right ventricle.

MRI scanners of different magnetic strengths, i.e. 1.5T and 3.0T. For each patient, ACDC provides 4-D data (weight, height, diastolic phase instants, systolic phase instants). To simulate several data providers we split the original training data of ACDC into two datasets: we consider the images at 1.5T field strength as one source and those at 3T as the other. Overall, ACDC 1.5T has 17,408 images and ACDC 3T has 7,936 images (Examples are shown in Fig. 2.). For both datasets, four classes of pathology are provided and one for normal subjects (for a total of 5 classes). We randomly crop all images to 128×128 during training. **Results and discussion:** We consider ACDC 3T as D_Q and assume the domain D spans the space of cine images of the heart in 1.5T or 3T strength. ACDC 1.5T is the calibration dataset D_C . Results are shown in Table 2.

We first perform experiments to verify if our method can correctly detect whether D_Q is in D^* . Similar to benchmark datasets, all models trained with ACDC 3T or part of ACDC 3T have $\rho < 1$. Other models trained without any data from ACDC 3T have $\rho \geq 1$. Hence, for ACDC dataset, our method returns correct answers in all experiments.

According to %ACDC 1.5T results, a large-size dataset D_C (%100 ACDC 1.5T) achieves lower K-S distance compared to 50% and 75% ACDC 1.5T. This supports our discussion of the scenario when D^* does not include any samples of D_Q . Note that when mixing all data of ACDC 1.5T and only 10% of 3T, the K-S distance drops $(1 - 0.750) \div 1 = 25\%$. It suggests even if part of D_Q is in D^* , the proposed method is still sensitive to detect such case.

5 Conclusion

We introduced an approach that uses Kolmogorov-Smirnov (K-S) distance to detect if a model has used/forgotten a query dataset. Using the K-S distance we can obtain statistics about the output distribution of a target model without

Table 2. ACDC dataset results. The query dataset D_Q is ACDC 3T. The training dataset D^* has different variants as listed below. The K-S distances are calculated with Eq. 2. D_C for defining ρ is ACDC 1.5T. Bold number in the table is the threshold.

Training dataset (D^*)	ACDC 3T	%ACDC 1.5T			ACDC $1.5T + \% 3T$		
		50%	75%	100%	10%	50%	100%
$KS(r(T_{D_Q D_Q}), r(T_{D_Q D^*}))$	0.036	0.885	0.793	0.772	0.579	0.529	0.103
ρ	0.047	1.146	1.027	1.000	0.750	0.685	0.133

knowing the weights of the model. Since the model's training data are unknown we train new (shadow) models with the query dataset and another calibration dataset. By comparing the K-S values we can ascertain if the training data contain data from the query dataset even for the difficult case where data sources can overlap. We showed experiments in classical classification benchmarks but also classification problems in medical image analysis. We did not explore the effect of the query dataset's sample size but the cumulative distribution remains a robust measure even in small sample sizes. Finally, extensions to segmentation or regression tasks and further assessment whether differential-privacy techniques help protect data remain as future work.

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