Abstract—Deep Neural Networks (DNN) are typically tested for accuracy relying on a set of unlabelled real world data (operational dataset), from which a subset is selected, manually labelled and used as test suite. This subset is required to be small (due to manual labelling cost) yet to faithfully represent the operational context, with the resulting test suite containing roughly the same proportion of examples causing misprediction (i.e., failing test cases) as the operational dataset.

However, while testing to estimate accuracy, it is desirable to also learn as much as possible from the failing tests in the operational dataset, since they inform about possible bugs of the DNN. A smart sampling strategy may allow to intentionally include in the test suite many examples causing misprediction, thus providing this way more valuable inputs for DNN improvement while preserving the ability to get trustworthy unbiased estimates.

This paper presents a test selection technique (DeepEST) that actively looks for failing test cases in the operational dataset of a DNN, with the goal of assessing the DNN expected accuracy by a small and “informative” test suite (namely with a high number of mispredictions) for subsequent DNN improvement. Experiments with five subjects, combining four DNN models and three datasets, are described. The results show that DeepEST provides DNN accuracy estimates with precision close to (and often better than) those of existing sampling-based DNN testing techniques, while detecting from 5 to 30 times more mispredictions, with the same test suite size.

Index Terms—Software testing, Artificial neural networks

I. INTRODUCTION

Deep Neural Networks (DNN) are today integral part of many applications, including safety-critical software systems such as in the medical [1] and autonomous driving domains [2]. Testing is a crucial activity in the development of such systems, for both quality/safety reasons to avoid DNN-caused catastrophic failures [3], [4], and for cost as well – very large samples may be needed to reliably test a DNN [5], [6].

A significant research effort is currently being put on DNN testing. A primary goal is to find adversarial examples causing mispredictions, namely to expose as many failing behaviours as possible [7]–[11]. Several structural coverage criteria have been proposed to drive the automated generation of test inputs and assess the test adequacy – neuron coverage [7], k-multisection neuron coverage, neuron boundary coverage [8], combinatorial coverage [10]. It has been argued that these criteria may be misleading, because of the low correlation between the number of misclassified inputs in a test set and its coverage [12]. Wu et al. [13] and Kim et al. [14] recently considered discrepancy measures between the training/validation data and test data, to improve fault detection and to have coverage criteria better correlated to failure-inducing inputs.

The output of this type of failure-finding testing (and then debugging) process is an improved DNN, with higher accuracy. This resembles what is called debug testing in the traditional testing literature [15]. Beside differences in testing DNN-based and conventional software (e.g., the oracle definition), which make DNN testing problematic, a further issue is that the so-obtained testing results are not necessarily related to the accuracy actually experienced in operation. In fact, testing data may be not representative of the actual operational context. This may happen when test data are generated artificially (like in adversarial examples generation) or they differ significantly from input observed in the field. The number of mispredictions and/or the coverage achieved give only an “indirect” (and, for what said, inaccurate) measure of the expected accuracy in operation, and ultimately of the confidence that can be placed in the DNN.

A less investigated research path is testing a DNN with the explicit goal of providing a statistical estimate of its expected accuracy in the operational context. In software reliability engineering, this is a well established practice known as operational testing [16]. The objective is to assess how well a DNN will perform in the intended context using a small yet effective amount of test data. With a cost-effective accuracy estimate, testers can establish a threshold-based release criterion and correct or tune their artificial networks (e.g., adjust the DNN structure or hyper-parameters) until the criterion is met. The reference scenario is the following: a DNN model is trained with a set of data (training dataset) and is meant to operate in a given context (operational context). To test the model, an arbitrarily large set of operational data is available (operational dataset) containing examples whose correct label is unknown: the goal of the tester is to select a small subset of the operational data, to be manually labelled and used as test cases to estimate the accuracy of the model in operation. Due to manual labelling, testers typically look for a trade-off between a good estimate and the labelling cost. This problem has been recently faced by Li et al. [17]. Their selection criterion is to minimise cross-entropy between selected test data and the whole set of unlabelled operational data, so as to have a sample of tests representative of the operational context.

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As this approach does not explicitly look for failing examples, it does not give much room for improving the DNN accuracy. From this viewpoint, many tests selected are useless, as they are non-failing examples that serve only the purpose of having highly-accurate estimates.

Given the above, we either have a test suite exposing failures but not representative (adversarial-like testing), or representative tests but few failures exposed (operational testing). This paper targets the drawbacks of both strategies together. We present DeepEST (Deep neural networks Enhanced Sampler for operational Testing), an operational testing method for DNN that builds test suites from an operational dataset so as to provide a close and efficient estimate of the expected accuracy and to find, at the same time, a high number of failing examples. Paraphrasing a famous aphorism by O. Wilde (“Experience is the hardest kind of teacher. It gives you the test first and the lesson afterward.”), DeepEST aims to sample, from operational data, many failing tests to learn from while building a set of tests suited for the DNN accuracy estimate.

With respect to pure operational testing, DeepEST is expected to provide DNN accuracy estimates that are more precise and more efficient (for number of tests required), plus the practical advantage of enabling debug-testing-like scenarios (improving accuracy through debugging/re-training). With respect to adversarial-like testing, DeepEST is able to provide accuracy estimates, enabling evaluation of alternative designs or DNN fine tuning, and establishing a release criterion directly related to what will be observed in operation.

Experiments with various DNN and datasets are presented, which assess DeepEST effectiveness, sensitivity to the number of tests to select from the operational dataset, and dependency on the dataset. As DeepEST can exploit various types of auxiliary information to select tests, the experimental study considers four variants. The results show that all the variants produce accuracy estimates similar to those of the state-of-the-art technique, while detecting from 5 to 30 times more mispredictions, with the same sample size.

The paper is structured as follows. Section II introduces sampling-based operational testing concepts used by DeepEST. Section III describes the DeepEST algorithm. Sections IV and V report the experimentation. Section VI discusses related work. Section VII concludes the paper.

II. SAMPLING-BASED OPERATIONAL TESTING OF DNNs

A. Operational testing of DNNs

In the traditional testing literature, operational testing refers to the family of techniques that use an operational profile to test a system to estimate its expected reliability (i.e., probability of not failing) in operation. Likewise, the primary goal of DNN operational testing is to estimate the expected accuracy (i.e., probability of not having mispredictions) in a given operational context [17]. Two main challenges arise.

Data skew. The idea of operational testing is that testing should not just care about exposing possible failures, but should be able to spot those failure-causing inputs that are more likely to occur in operation. In case of a relevant mismatch between the pre-release test data and the post-release context, the system could be stimulated in operation with inputs never seen during testing, with unexpected failures.

Data skew is a concern for DNN, more than for traditional software systems. These are expected to work on a range of input data given to functionalities, and can be tested on a small carefully-selected sample of input data (e.g., obtained by input space partitioning). DNN are data-driven by nature; they are constructed around a training dataset, and generalising beyond what observed during training is hard. This data-driven approach is governed by a statistical process and, due also of the black-box nature of DNN, it is tricky to identify classes of inputs that homogeneously represent the expected behaviour in operation. For instance, white-box partitioning has been shown to be not clearly correlated to the failing behaviour [12]. Thus, a drift of the post-release operational context from the pre-release testing context is more likely to cause unexpected failures compared to traditional software.

The imitation bias of operational testing. The operational context drift is just a triggering condition for unexpected failures at runtime. The problem occurs because of the way in which operational testing is conducted. Operational testing selects a small data sample that can accurately represent the population; however, the mere imitation of the expected input can be inefficient, especially in highly reliable systems, because many failure-free tests are executed to get an acceptable estimate. A representative sample would roughly contain the same proportion of examples causing misprediction as the operational dataset. There are two problems with this: first, in highly accurate DNNs, the number of examples causing mispredictions is low, thus requiring other types of testing activities dedicated to detect mispredictions (e.g., through adversarial-like techniques) to possibly improve the accuracy. Second, just mimicking the expected usage is far from the estimation point of view as long as the imitation is faithful. If this is not the case, the risk of overestimation increases: if we only aim at having a representative sample of tests, the actual experienced accuracy may be significantly smaller than the estimated one if the operational context drifts, because of the occurrence of unexpected mispredictions.

Thus, a conventional approach for DNN operational testing allows to obtain the desired accuracy estimate but, since it reveals few mispredictions, can be ineffective (leading to wrong estimates) and inefficient (requiring further separate testing activities to detect mispredictions). Recent results in operational testing for traditional software show that exposing failures and estimating reliability are not contrasting objectives [18]. A strategy that actively looks for failures (rather than just mimicking the expected usage) can lead to accurate and stable estimates and, at the same time, expose many failures. Our aim is exactly to spot failing examples in a DNN operational dataset, while preserving the ability to yield effective (small error) and efficient (low variance) estimates.

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We use the terms misprediction and failure interchangeably for DNN.
B. Sampling-based testing

DeepEST uses statistical sampling, a natural way to cope with estimation problems: it serves to design sampling plans tailored for a population under study, providing effective and efficient estimators. In sampling-based testing [19], the sample is the set of \( n \) test cases \( T = \{ t_1, \ldots , t_n \} \), having a binary outcome (pass/fail). Test outcomes are a series of independent Bernoulli random variables \( z_{t_i} \), such that \( z_{t_i} = 1 \) if the DNN predicts the correct label for \( t_i \), \( z_{t_i} = 0 \) otherwise. The parameter of our interest is the DNN accuracy \( \theta \); we aim to an estimate \( \hat{\theta} \) with two desirable properties: unbiasedness – i.e., the expectation of the estimate \( \mathbb{E}[\hat{\theta}] \) should be equal to the true value \( \theta \) - and efficiency – for the given the sample size, the variance of \( \hat{\theta} \) should be as low as possible (for a highly confident, stable estimate). The probability that \( z_{t_i} = 1 \) corresponds to the true (unknown) proportion: \( \theta = \frac{\sum_{i=1}^{n} z_{t_i}}{N} \), with \( N \) being the population size (i.e., the size of the operational dataset).

Simple random sampling with replacement (SRSWR) is the baseline approach: an unbiased estimator of \( \theta \) is the observed proportion of correct predictions over the number of trials \( n \):

\[
\hat{\theta}_{\text{SRSWR}} = \frac{\sum_{i=1}^{n} z_{t_i}}{n}.
\]

Having assumed independent variables, the variance of \( \hat{\theta} \) is:

\[
V(\hat{\theta}_{\text{SRSWR}}) = \frac{\theta(1-\theta)}{n}.
\]

An improvement is represented by simple random sampling without replacement (SRSWOR), namely, the same test case is not selected twice: this reduces the variance to:

\[
V(\hat{\theta}_{\text{SRSWOR}}) = \frac{N-n}{N-1} \frac{\theta(1-\theta)}{n}.
\]

While SRS keeps the mathematical treatment simple, it is unable to exploit additional information a tester might have.

Exploiting auxiliary information to modify the sampling scheme is what is done in sampling theory to get more efficient estimators [20]. The sampling is made proportionally to an auxiliary observable variable assumed to be related to the (unknown) quantity to estimate; the estimator is then adjusted to account for the non-uniform sampling, so as to preserve unbiasedness. For instance, stratified sampling is a strategy which uses knowledge about which sample units are expected to have homogeneous values, and selects units contributing more to lower the estimate’s variance.

Li et al. [17] present two sampling strategies for DNN, which are, to our knowledge, the only attempt to DNN operational testing: Confidence-based Stratified Sampling (CSS) and Cross Entropy-based Sampling (CES) – the latter being the authors’ proposal. Both strategies exploit auxiliary information to drive the sampling task.

In CSS, sampling is proportional to the confidence value provided by classifiers when predicting a label: examples with higher confidence are more likely to be selected as part of the test suite. This works well when the classifier is reliable, namely if examples with higher confidence are actually those for which the prediction is more likely to be correct (in other words, the model is perfectly trained for the operation context). Whenever the operation context drifts from the training one, CSS exhibits poor performance.

CES attempts to overcome this limitation by using the output of the \( m \) neurons in the last hidden layer in the last hidden layer, assumed to be more robust to the operation context drift. It builds the test suite trying to minimize the average cross-entropy between the probability distribution of the \( m \)-dimensional representation of the output of neurons computed on the operational dataset and on the selected tests.

To pursue the double objective of sampling cases causing mispredictions and estimating accuracy efficiently, DeepEST adopts an adaptive and without-replacement sampling algorithm, described in Section III.

C. Auxiliary information

To look for mispredictions, the auxiliary information leveraged by DeepEST adaptive sampling should represent the belief about some factor(s) related to the model’s (in)accuracy. We consider two auxiliary variables, related to somehow opposite sources of information: the confidence value and the distance between the operational dataset example and the training dataset. The latter is based on the result by Kim et al. [14] that inputs more “distant” from training data are more likely to cause misprediction – they show that distance is correlated to mispredictions, which is what we look for. We borrow their distance metrics Distance-based Surprise Adequacy (DSA) and Likelihood-based Surprise Adequacy (LSA). These are computed by the Activation Trace (AT), namely a vector of activation values of each neuron of a certain layer corresponding to a certain input; we compute AT with reference to the last DNN activation layer. LSA is defined as the negative log of density (computed via Kernel Density Estimation). DSA is defined as the Euclidean distance between the AT of a new input and ATs observed during training. The variant using confidence is DeepEST\(_{CS}\); the variants using the distance metrics are DeepEST\(_{DSA}\) and DeepEST\(_{LSA}\).

Performance of an auxiliary variable can strongly depend on the DNN and on the training/operation dataset: for instance, for a distance metric the belief that examples far from the one in the training set are more likely to cause a misprediction is not an absolute truth. In particular, if an example is very similar to many others in the training set according to the distance metric, but has a different label, distance will be not a good metric to select it. Similarly, the confidence is a good proxy if the DNN is well trained for the operational context. In general, relying on a single auxiliary variable may work well in some settings and bad in others. Combining multiple beliefs is a choice that is expected to improve the stability of results across multiple settings. Based on this, we define a further variant of our algorithm, named DeepEST\(_{C}C\), considering as auxiliary variable the combination of confidence and distance:

\[
P = P_c \times (1 - P_d)
\]
where $P_c$ is the confidence value, $P_d$ is the DSA value normalized $[0, 1]$ and $P$ is probability of correct prediction. The intuition behind is that the probability of a correct prediction is related to both the confidence of the DNN and the drift of the example from what seen during training. In fact, $P_c$ is related to the probability that the DNN does a correct prediction according to what seen during training – hence due to the imperfection of training. If confidence $P_c$ is high and the example is close to the training dataset (i.e., $P_d$ is small), there is a high chance of correct prediction.

III. The DeepEST Algorithm

Both CES and CSS select a sample of tests representative of the operational context. DeepEST takes a different approach, favouring the sampling of failing (namely, misprediction-causing) tests, then used to estimate the DNN accuracy. While the estimation ability demands for probabilistic (hence “sampling-based”) selection of examples, the very idea of DeepEST is that, since mispredictions are usually rare compared to correct examples, looking for failing tests is well handled by adaptive sampling [21]: the examples progressively selected may give hints about the probability of finding other failing examples, so as to adapt the sampling process accordingly. Given a samples size, adaptive sampling implicitly assumes that inputs of interest are not uniformly spread across the input space, and adopts a disproportional selection to spot them, counterbalanced by the estimator to preserve unbiasedness. In software testing, this is expected to give smaller variance compared to conventional sampling (e.g., SRS), since the failing inputs are usually not uniformly distributed. The DeepEST sampling algorithm is inspired to the Adaptive Web Sampling [22], a flexible design for sampling rare populations.

The above auxiliary variables are used to define a weight $w_{i,j}$ between any pair of examples $i$ and $j$ of the operational dataset, used to explore the example space adaptively. For instance, let us assume to use the normalised DSA distance: if the example $i$ has distance $D_{di}$ (representing the belief that $i$ causes a misprediction), and a pre-defined threshold $\tau$ is exceeded (i.e., $i$ has a sufficiently high distance compared to the others), then all the $w_{i,j}$ values ($\forall j$ of the operational dataset) are set to their distance $D_{di}$; otherwise $w_{i,j} = 0$. This way, a strong-enough belief about example $i$ causing misprediction entails the activation of all the weights toward $i$. The latter are used, as explained hereafter, for sampling, and makes the algorithm follow the distance criterion to spot potential clusters of failing examples.

The DeepEST sampling strategy is sketched in Algorithm 1. Assuming $n$ examples to be selected from the operational dataset as test cases, the algorithm selects and executes one test case per step. The first input is selected by simple random sampling, namely, initially all examples have equal probability of being selected. Then, one of two sampling schemes is used to select next test: weight-based sampling (WBS), or simple random sampling (SRS). Example $i$ is selected from the operational dataset at step $k$ with probability $q_{k,i}$ given by:

$$
q_{k,i} = r \cdot \frac{\sum_{j \in s_k} w_{i,j}^{2}}{\sum_{h \notin s_k} w_{h,j}} + (1 - r) \cdot \frac{1}{N - n_{s_k}} \tag{5}
$$

where:

- $r$: probability of using WBS (hence, probability of using SRS = 1-$r$; $r$ is set to 0.8 in our implementation);
- $s_k$: current sample (all examples selected up to step $k$);
- $w_{i,j}^2$: weight relating example $j$ in $s_k$ to example $i$;
- $n_{s_k}$: size of the current sample $s_k$;
- $N$: size of the operational dataset.

For both WBS and SRS, the selection is without replacement.$^3$ WBS selects an example $i$ proportionally to the sum of weights $w_{i,j}$ of already selected examples toward $i$ – the chance of taking $i$ depends on the current sample, favouring the identification of clusters of failing examples if the auxiliary variable (hence, $w_{i,j}$) is well-correlated with mispredictions. As this is not always the case, the WBS “depth” exploration is balanced by SRS, chosen with probability (1-$r$), for a breadth exploration of the example space. This diversification in the search is useful to escape from unproductive cluster searches. The steps are repeated until the testing budget $n \leq N$ is over.

At step $1$, the probability that a randomly selected example will cause a misprediction is estimated as the outcome $y_1$ of the first test (1 in case of misprediction, 0 otherwise).

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$^2$In DeepEST, DSA is preferred to LSA since it has been shown to have better performance for the deeper layer [14].

$^3$Without replacement sampling schemes generally give smaller variance than their with replacement counterpart on the same sample size [20].
### TABLE I: Experimental DNN and datasets

| Subject | DNN | Dataset | Classes | Training set size | Test set size | True accuracy |
|---------|-----|---------|---------|------------------|--------------|--------------|
| S1      | CN5 | MNIST   | 10      | 60,000           | 10,000       | 0.9968       |
| S2      | LeNet5 | CIFAR10 | 10      | 50,000           | 10,000       | 0.9868       |
| S3      | CN12 | CIFAR10 | 10      | 50,000           | 10,000       | 0.8806       |
| S4      | VGG16 | CIFAR100 | 100    | 50,000          | 10,000       | 0.9359       |
| S5      | VGG16 | CIFAR100 | 100    | 50,000          | 10,000       | 0.7048       |

At step $k > 1$, example $i$, whose outcome is $y_i$, is selected with probability $q_{k,i}$ according to Eq. [5], and the estimator of the misprediction probability is that by Hansen-Hurwitz [23]:

$$z_k = \frac{1}{N} \left( \sum_{j \in S_k} y_j + \frac{y_i}{q_{k,i}} \right)$$

(6)

where the $y_j$ values are the outcome of the tests already selected. $z_k$ is an unbiased estimator of the expected misprediction probability at step $k$; the final estimator of the expected accuracy of the DNN is 1 minus the average of the $z_k$ values:

$$\hat{\theta} = 1 - \frac{1}{n} (y_1 + \sum_{k=2}^{n} z_k)$$

(7)

where $n$ is the number of tests run.

### IV. Evaluation

#### A. Experimental subjects

The four variants of DeepEST are evaluated against the SRSWR scheme as baseline and the mentioned state-of-the-art technique CES. Five experiments are conducted with four DNN models and three popular datasets.

The datasets are MNIST, a dataset of handwritten digits [24]; CIFAR10, for image processing systems; and CIFAR100, similar to the previous one but with 100 classes [25]. The chosen DNN models are ConvNet5 (here simply CN5) and LeNet5 for MNIST classification; ConvNet12 (simply, CN12) and VGG16 for CIFAR10 classification; VGG16 for CIFAR100. Table I lists the five subjects (DNN-dataset pairs); the true accuracy is in the last column.

#### B. Research questions and experiment design

The evaluation answers the following research questions.

**RQ1: Effectiveness.** How does DeepEST perform in finding inputs causing misprediction (i.e., failing examples) and simultaneously estimating a DNN operational accuracy?

To gauge DeepEST ability to provide effective DNN accuracy estimates with few examples, while spotting a high number of failing inputs, we set to 200 the number of tests to select (then varied to answer RQ2) and repeat 30 times the execution of the 6 compared techniques on the 5 subjects.

As for evaluation metrics, we compute:

- The accuracy $\hat{\theta}_i$ at the $i$-th repetition, and then compute the Mean Squared Error (MSE) as $MSE(\theta) = \frac{1}{n} \sum_{i=1}^{30} (\hat{\theta}_i - \theta)^2$, where $\theta$ is the true operational accuracy. Note that for unbiased estimators, MSE and variance can be considered indistinguishable. In fact, $MSE = \text{Variance} + \text{Bias}^2$ and $\text{Bias}(\hat{\theta}) = E[\hat{\theta}] - \theta = 0$.

  The precision of the estimator is: $\pi(\hat{\theta}) = \frac{1}{MSE(\theta)}$, and the relative precision (or relative efficiency) of estimator $A$ with respect to $B$ is: $\rho_{A,B} = \frac{\pi_A}{\pi_B}$. $1 > \rho_{A,B} > 0$ means that $A$ is better than $B$.

- The average number of failures ($\varphi = \text{Mean}(\varphi_i)$) with $\varphi_i$ being the number of failures in repetition $i$.

  For comparison purpose, we consider the relative number of failures of technique $A$ with respect to $B$: $\rho_{A,B} = \frac{\varphi_A}{\varphi_B}$. ($\rho_{A,B} > 1$ means that $A$ is better than $B$).

**RQ2: Sensitivity to sample size.** How does the performance of DeepEST vary with the sample size?

It is important to figure out how performance varies with the number of test cases to select from the operational dataset, namely with sample size. Indeed, DeepEST aims to perform well especially with a small sample size, so as to yield precise estimates with relatively few examples to be manually labelled.

**RQ3: Dataset influence.** How is DeepEST performance affected by the datasets?

In DNN testing, results are often heavily dependent on the (training and operational) datasets. This RQ aims to figure out how these may influence the ability of the auxiliary variables (confidence, DSA, LSA) to discriminate failing examples, affecting the performance of DeepEST. To answer RQ3, the test set is completely labeled, so as to identify all failures.

#### C. Implementation

DeepEST is implemented mostly in Java. The implementation of the distance metric in DeepEST$_{DSA}$ and DeepEST$_{LSA}$ is the same used by Kim et al. [14]: we used their Python scripts to compute DSA and LSA values. These are computed considering the last activation layer of each DNN. The threshold $\tau$ needed for the weights definition is set as follows:

- DeepEST$_{DSA}$: $\tau = \text{mean}(\text{DSA}) + 2 \times \text{Std}(\text{DSA})$;

- DeepEST$_{LSA}$: $\tau = \text{mean}(\text{LSA}) + \text{Var}(\text{LSA})$.

The threshold for confidence, used by DeepEST$_{CS}$ and DeepEST$_{C}$, is set to 0.7, assuming that lower confidence values are more related to misprediction (i.e., the weights are activated when the confidence is less than $\tau = 0.7$).

In CES, the selection exploits the output of the last hidden layer. We use the same configuration as the original article [17]. The size of the initial sample is $p=30$, enlarged by a group $Q^*$ of $q=5$ examples at each step. The number of random groups from which $Q^*$ is selected is $L = 300$. For CES and SRS, we used the Python scripts provided by Li et al. [17].

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4CN5 and CN12 are calibrated in the same way as Li et al. [17]; LeNet5 is calibrated as Kim et al. [14], for the VGG16 network we considered the weights at [https://github.com/geifmany/cifar-vgg](https://github.com/geifmany/cifar-vgg).
V. RESULTS

A. RQ1: effectiveness

Figure 1 plots the MSE of the estimated accuracy. The techniques exhibit comparable performances, with DeepEST\(_C\) being the best one for 3 of the 5 subjects. CES has good performance in terms of MSE, it is the second technique in 3 cases. Considering the single variables: confidence, DSA or LSA lead to results slightly more variable over the subjects – an aspect explored in RQ3. The SRS case is interesting, too: it is never the worst approach and is the best in one case.

Table II reports the average number and the standard deviation (\(\sigma\)) of the number of failing examples detected. All variants of DeepEST identify many more failing examples than SRS and CES, even up to a factor of 30x (DeepEST\(_{CS}\) vs CES for subject S2) and reaching in some cases (S1 and S2, DeepEST\(_{CS}\)) almost 50% of the total number of failing examples in the datasets (last column). The DeepEST algorithm leverages the adaptive sampling to spot clusters of failing examples with relatively few tests (set to 200 for RQ1). Its performance varies depending on the auxiliary information used, but it is always remarkably better than SRS and CES.

Among the DeepEST variants, confidence (DeepEST\(_{CS}\)) turns out to be the most effective auxiliary variable in detecting failures, showing the best performance for all datasets and models, followed by DSA. CES and SRS select the lowest number of mispredicted examples, and are close to each other. Considering both the failure detection ability and the estimate accuracy, DeepEST\(_C\) – that combines confidence and DSA, the two best auxiliary variables for failing examples detection - gives a good trade-off, since it provides stable (across subjects) and close-to-true estimates of the accuracy, with many more detected failing examples than CES and SRS.

Table II: RQ1 (effectiveness): Mean and standard deviation (\(\sigma\)) of the number of failing examples detected.

| Subject                   | SRS       | CES        | DeepEST\(_{CS}\) | DeepEST\(_{DSA}\) | DeepEST\(_{LSA}\) | DeepEST\(_C\) | Total failing examples |
|---------------------------|-----------|------------|-------------------|-------------------|-------------------|---------------|------------------------|
| S1 (CN5, MNIST)           | 1.90      | 2.23       | 44.57             | 0.68              | 30.17             | 4.03          | 3.29                   |
| S2 (LeNet5, MNIST)        | 2.77      | 2.03       | 65.13             | 0.86              | 57.57             | 5.12          | 4.74                   |
| S3 (CN12, CIFAR10)        | 37.93     | 33.77      | 106.10            | 7.26              | 98.33             | 4.62          | 6.61                   |
| S4 (VGG16, CIFAR10)       | 12.60     | 13.03      | 85.00             | 4.86              | 16.80             | 3.02          | 3.14                   |
| S5 (VGG16, CIFAR100)      | 57.33     | 55.77      | 131.17            | 7.35              | 78.20             | 6.62          | 3.68                   |

A replication package is at: https://github.com/dessertlab/DeepEST
TABLE III: Pairwise comparison of techniques. A value of $\rho_{R,C} > 1$ means the technique on the row has a greater precision than that on the column. Similarly for the relative number of detected failures $\pi_{R,C}$.

|        | CES | DSA | LSA | C  | SRS |
|--------|-----|-----|-----|----|-----|
| **DeepEST** |     |     |     |    |     |
| row vs col |     |     |     |    |     |
| CES | $\rho_{R,C}$ | $\pi_{R,C}$ | 0.0501 | 0.0340 | 0.0996 | 0.1872 | 1.1754 | 0.8929 |
| DSA | - | 1.4773 | 1.9866 | 3.7346 | 23.4561 | 15.8772 | 1.5150 |
| LSA | - | - | 1.8799 | 0.7187 | 11.8070 | 7.5621 |
| C  | - | - | - | 6.2807 | 7.8624 |
| **DeepEST** |     |     |     |    |     |
| row vs col |     |     |     |    |     |
| CES | $\rho_{R,C}$ | $\pi_{R,C}$ | 0.3183 | 0.3434 | 0.8870 | 0.4772 | 0.8902 |
| DSA | - | 1.0790 | 2.7872 | 1.4993 | 2.7970 |
| LSA | - | - | 2.5382 | 0.2265 | 0.5379 | 0.9954 |
| C  | - | - | - | 1.8858 | 2.8126 |
| **DeepEST** |     |     |     |    |     |
| row vs col |     |     |     |    |     |
| CES | $\rho_{R,C}$ | $\pi_{R,C}$ | 0.4252 | 0.7131 | 0.8266 | 0.5130 | 0.9727 |
| DSA | - | 1.6773 | 1.9442 | 1.2607 | 2.2878 |
| LSA | - | - | 1.1591 | 0.5056 | 1.1767 |
| C  | - | - | - | 1.8959 | 0.6892 |

Table III(a)–(e) show the results of the pairwise comparison of the techniques. For the four DeepEST variants, rows and columns headings list (in blue) the name of the auxiliary variables. The evaluation metrics are the ratio $\rho$ of the failing examples and the relative precision $\pi$ of the estimators. Values of $\rho$ or $\pi$ greater (lower) than 1 mean the technique on the row (column) has better performance. If a technique is better than the other in a pair for both metrics (values coloured in vs), it wins.

Table III(f) summarizes the number of wins (and losses). CES never wins over other approaches, while DeepESTC wins against CES 3 out of 5 times. SRS never wins over DeepEST, while it wins vs CES considering VGG16 on CIFAR10. DeepESTC never looses and collects the highest number of wins (10), showing the best trade-off among accuracy of the estimation and number of detected failing examples.

The choice of the DeepEST variant may be determined by which auxiliary information can be collected. We see that DeepESTC, exploiting a combination of two variables, has good and more stable results in terms of MSE than the other variants, at the expense of a slight decrease of detected failures. Single auxiliary variables are more sensitive to the specific dataset/model pair (e.g., confidence works well if the DNN is reliable), but expose more mispredictions. Confidence has the advantage of not requiring knowledge of the hidden layers and is easier to compute. When no information is available or easily computable, SRS could be a good low cost solution.

B. RQ2: sensitivity to sample size

To answer this RQ, experiments are run with the sample sizes 50, 100, 200, 400, 800, considering the subject with the highest accuracy (S1), and the one with the lowest accuracy.
(S5), so as to analyze how DeepEST performs when there are very few and many failing examples in the dataset, respectively. Figure 2 shows the mean values of the estimates’ accuracy over repetitions. Figures 3 and 4 plot the MSE and the mean number of detected failures, respectively. Expectedly, increasing the sample size all techniques exhibit a decreasing trend in MSE and an increasing trend in failing examples.

For the subject with highest accuracy (S1), we observe the following. DeepEST$_C$ shows very good performance for MSE (Fig. 3(a), Fig. 4), and it detects on average about six times the failures of SRS and CES (Table II). The advantage for MSE is more pronounced with the smallest sample sizes, which make DeepEST$_C$ particularly suited when the number of examples to select and label is very small. For larger sample sizes, the MSE is similar but the advantage of DeepEST$_C$ over SRS and CES is very pronounced for detected failures (Fig. 4(a)). DeepEST$_{CS}$ is the best among all techniques to detect failures for small sample sizes; for sizes 400 and 800, the best is DeepEST$_{DSA}$ (Fig. 4(a)). Although the estimates are unbiased (hence, they tend to the true value), if we look at the mean estimates over the repetitions (Figure 2(a)), CES shows bad performance with up to 200 test cases. SRS works well with low budget; its good results may be influenced by the very low number of failures: 18/30 repetitions show 0 failures and 100% accuracy. It is interesting to note that in most cases CES and SRS overestimate the true accuracy – an undesired
property, especially for critical systems. This is related to the low number of failures detected, as discussed in Section II.

For the subject with lowest accuracy (S5), CES and SRS outperform DeepEST$_{C}$ for small sample sizes (50 and 100); from 200 to 800, the estimation by CES starts diverging, while SRS and DeepEST keep good performance. The tendency to overestimate the accuracy by CES and SRS is confirmed. Performance in failing examples detection is always clearly in favour of all DeepEST variants. Performance in estimation accuracy is almost specular to what observed with the most accurate model. A reason is that DeepEST is a sampling techniques particularly suited for rare populations, which is not the case of S5. As for the ability to detect failing examples, confidence is the best auxiliary variable for DeepEST for subject S5: it presents the best values in all configurations.

C. RQ3: dataset influence

We have seen in the experiments for RQ1 and RQ2 that no single auxiliary variable performs best in all situations. For instance, we can consider the confidence a good auxiliary information for subject S1, and a bad choice for S5. This may depend on several factors: assuming a perfect training, the confidence could be affected by a bias in the training set; or, with a perfect training set, a wrong training phase (e.g., due to overfitting) could generate mispredictions with high confidence. In some cases the operational dataset could contain examples very similar (i.e., small distance) to those in the training set but with a different label, affecting the discriminative power of the DSA and LSA metrics. To analyze how the three datasets influence the ability of auxiliary variables to discriminate failing examples (impacting the performance observed in RQ1-RQ2), we consider the subjects S1 and S5, as for RQ2, plus the VGG16 DNN for CIFAR10.

Figures 5, 6, and 7 show the logistic regression for the three datasets. The curves fit the probabilities for the outcome fail/pass to the three predictors: confidence, DSA, and LSA. Consider MNIST, for which CN5 reaches the highest accuracy among all subjects; the probability for a test to fail is very low for values of confidence between 0.7 and 1 (Figure 5). This is clearly not the case for CIFAR10 (Figure 6 and, especially, CIFAR 100 (Figure 7). In the latter, there is a high chance of misprediction even with high values of confidence; this could be due to a high skew between training and test data.

The discriminating power of DSA and LSA is clearly greater for MNIST, as the high slope of the S-shaped curves in Figures 5(b)-(c) suggests, compared to corresponding ones in Figures 6 and 7. In this case, it actually happens that the farthest examples have highest probability to be related to a failure, with a sharp increase after DSA≈0.9 and LSA≈300. This means that if in operation there are (a lot of) examples far from what observed in training, re-training with a more representative dataset can be useful to improve the accuracy. This behaviour is not observed for CIFAR10 and CIFAR100, and DSA and LSA do not seem to be effective in dividing the two sets of examples. In CIFAR10, the DSA line is more horizontal, meaning that the DSA value does not reflect well the failure probability. The scarcely discriminative power of the auxiliary variables in CIFAR10 and CIFAR100 partially explains the smaller gain of DeepEST over CES and SRS (especially in terms of MSE); nevertheless, its adaptivity allows spotting many failing examples even in these conditions.

Finally, it is interesting to highlight the performance of DeepEST$_{CS}$ (based on confidence) on MNIST in RQ2. The saturation in its failure detection ability (Figure 4(a)) can be explained observing that, after a number of tests able to spot failures looking at low-confidence examples, the few remaining ones with high confidence are selected with low probability; the sharper discrimination made by DSA and LSA determines a high detection ability even when few failing examples remain. In summary, whenever a tester has good belief/evidence about the appropriateness of one of the above auxiliary variables, it is a good choice to select the specific DeepEST variant; if not, the combined variant DeepEST$_{C}$ has shown to give the best trade-offs in all five experimented cases.

D. Threats to validity

A threat to the internal validity comes from the selection of the experimental subjects. To favor the repeatability of the experiments under different possibly influencing factors, we have used publicly available networks and pre-trained models, to avoid incorrect implementation. The configuration of parameter $r$ in DeepEST and a different setting of thresholds may also affect the results (in terms of efficiency of the estimator), hence a fine-tuning is suggested before applying the method to other dataset-DNNs. Although the code developed was carefully inspected, a common threat is the correctness of the scripts to collect data and compute the results.

The choice of the sample size influences the effectiveness too. We ran a sensitivity analysis to show that DeepEST is still effective (compared to both CES and SRS) under five (from 50 to 800) values of the sample size, but different values could yield different results. Threats to external validity depend on both the number of models and datasets considered for experimentation. We strived to control this threat considering different widely used DNN and datasets. Although the results may change with different subjects, the diversity and significance of the chosen subjects give confidence to the general considerations. Replicability of the experiments on other subjects is to further mitigate this threat.

VI. RELATED WORK

Testing of DNN is a hot research topic. Zhang et al. [26] present a survey on Machine Learning testing, identifying three main families of techniques: mutation testing [27], metamorphic testing [9], [28], and cross referencing [7], [29]. The former two are for test generation: they generate adversarial examples causing mispredictions. Cross-referencing can highlight the most interesting test cases when different implementations of a system disagree. Most of these techniques are meant for fault detection, rather than for accuracy estimate.

Testing DNN for accuracy is the focus of Li et al. [17], who presented CES – which DeepEST has been compared against
DeepEST differs from CES in several aspects, the key ones being the sampling algorithm and the used auxiliary variables, which are conceived to improve failing examples detection while preserving the accuracy estimation power.

Operational testing, where tests are derived according to the operational profile, is an established practice to estimate software reliability. It was the core technique of Cleanroom software engineering and of the Software Reliability Engineering Test process. Cai et al. developed Adaptive Testing, still based on the operational profile, but foreseeing adaptation in assigning tests to partitions. Recently, Pietrantuno et al. stressed the use of unequal probability sampling to improve the estimation efficiency, to this aim formalizing several sampling schemes. Our proposal goes along this direction: we do not sample tests representative of the operational dataset, but we “alter” the selection and counterbalance the uneven selection in the estimator. Unequal probability, without-replacement and adaptive sampling are the key concepts we borrowed for operational testing of DNNs.

VII. CONCLUSION

Testing the accuracy of a DNN with operational data aims at precise estimates with small test suites, due to the cost for manually labelling the selected test cases. This effort motivates to pursue also the goal of exposing many mispredictions with the same test suite, so as to improve the DNN after testing. With these two concurrent goals, we presented DeepEST, a technique to select failing tests from an operational dataset, while ultimately yielding faithful estimates of the accuracy of a DNN under test.

We evaluated experimentally four variants of DeepEST, based on various types of auxiliary information that its adaptive sampling strategy can leverage. The results with four DNN models and three popular datasets show that all DeepEST variants provide accurate estimates, compared to existing sampling-based DNN testing techniques, while generally much outperforming them in exposing mispredictions. Practitioners may choose the appropriate variant, depending on the characteristics of their operational dataset and on which auxiliary information is available or they can collect.
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