

Detecting Errors in Databases with Bidirectional Recurrent Neural Networks

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ABSTRACT

In recent years we have seen an exponential growth of data. However, not only does the volume of data grow but often also the numbers of erroneous data values. Detecting and cleaning these errors is an important data management problem to enable high quality data science pipelines. To reduce the human effort and to achieve good results in detecting errors, it is important to label the tuples which give the best impact for the error detection system.

In this paper we introduce an architecture based on *bidirectional recurrent neural networks* to detect errors in databases. The experimental results with 6 different datasets demonstrate that our approach shows *similar performance to state-of-the-art error detection system* per dataset. When considering the average of the F1-scores over all datasets, our approach called Enriched Two-Stacked Bidirectional (ETSB-RNN) outperforms state-of-the-art systems. Moreover, our approach achieves a lower standard deviation than existing work, which shows that our system is more robust. Finally, our approach does *not require additional data augmentation techniques* to achieve high F1-scores. The system does not need any configuration or previous analysis of the data.

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1 INTRODUCTION

The amount of data is growing exponentially since it is collected by many different stakeholders such as people, companies and machines at many different locations across the globe – but also with varying degrees of quality. When data scientists want to analyze data, they often have to clean it first, which typically takes the most time of the whole data science pipeline [8]. Hence, finding ways of cleaning datasets (semi-)automatically with minimal user involvement is an important aspect for many companies and data science projects. However, due to the heterogeneity of different

datasets and the different ways of how data is collected and stored, data cleaning in practice is a hard problem.

The cleaning part can be split into two steps, first to *detect errors* and second to *repair errors*, i.e. to find the ground truth. Errors are values which deviate in the dirty dataset from the clean dataset (see highlighted values in Table 1) and can be categorized as *missing value* ('NaN'), *value/syntactic error* ('80,000', 'Romr', '12', 'BER', '850'), *integrity constraint violation* ('75000') and *data duplication*.

Table 1: Overview of a dirty and the corresponding clean dataset (A=Age, Sal=Salary).

A	Sal	ZIP	City	A	Sal	ZIP	City
21	80,000	8000	NaN	21	80000	8000	Zurich
45	98000	00100	Romr	45	98000	00100	Rome
30	92000	75000	Paris	30	92000	75000	Paris
12	99000	BER	Berlin	42	99000	10115	Berlin
26	850	75000	Vienna	26	85000	1010	Vienna

The focus of this work is on error detection. While related work have used different error detection strategies [5, 7, 16, 19, 21, 22], there is ample room for improvement. In our approach we use *bidirectional recurrent neural networks* [25] (RNN), which have received little attention for solving this kind of challenge. Our approach generates a model which tries to find the best parameter settings during learning the content of the data. For training the system, we need labelled data. To reduce the effort for the user to provide labels for the training data, we use only 20 tuples – similar to the state-of-the-art approaches. For selecting the data, we developed a novel algorithm to choose a diverse trainset, which gives our system the best impact and thus the most information content for learning error patterns in the data.

System in action: A typical data science pipeline that applies our error detection algorithm looks as follows: The user gives our system a dataset and chooses the number of tuples for training. The system is responsible for the data preparation step and uses our novel label sampling algorithm *DiverSet* to find the tuples with the most information content for the neural network architecture to learn. Afterwards, the user has to label the chosen tuples with either '0' (correct) or '1' (wrong) to produce the trainset. Finally, the system uses the trainset to learn which data points of the tuples are correct or wrong. The involvement of the user is only in the phase of labelling the data. Hence, the user does not need to understand the whole dataset but only the proposed tuples for labelling.

This paper has the following **contributions**:

- We introduce our end-to-end system architecture for error detection based on *bidirectional two-stacked RNNs*, which is divided in the two parts *data preparation* and *error detection*.

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- We developed a novel algorithm for *trainset selection* to choose the tuples which provide the most information content for our system to learn.
- The experimental results show that our approach achieves similar performance to *state-of-the-art error detection system* on 6 benchmark datasets. Over all datasets, our system shows a higher average F1-score and lower standard deviation, which demonstrates that our system is more robust than existing work. Finally, we do not require additional data augmentation techniques to achieve high F1-scores.

Outline: In Section 2 we discuss related work of error detection. For better understanding of our system, we give background information on character embedding and recurrent neural networks in Section 3. In the subsequent Section 4 we show the architecture of our novel system for error detection as well as the algorithms for automatically selecting training data. In Section 5 we introduce the results of our system in comparison with state-of-the-art error detection systems. At the end we give a summary about the insights in Section 6.

2 RELATED WORK

In this section we review the related work on error detection methods. A good overview can be found in [2].

Early approaches employ rules and heuristics for data cleaning [1, 4] or identify functional dependencies based on integrity constraints [3, 18, 24, 27]. Other approaches use outlier detection and statistical analysis [10, 12, 26].

More recent methods such as ActiveClean [19], HoloDetect [11], Raha [21] and Rotom [23] use machine learning approaches to tackle the problem. The main challenge of these approaches is how to design machine learning algorithms that work well with few training labels. We will discuss these machine learning-based methods in more detail.

ActiveClean [19] dynamically trains machine learning models by requesting new labels from users interactively. The approach supports convex loss models such as linear regressions and SVM and prioritizes cleaning those records that are most likely to effect the results.

Raha [21] uses a novel sampling and classification scheme to choose the most promising labels for training. In addition, information from knowledge bases such as DBPedia can be integrated to improve the detection of certain errors. Raha automatically configures error detection strategies such as outlier detection algorithms [22], pattern violation detection algorithms [16], rule violation detection algorithms [7] and knowledge base violation detection [5]. Moreover, Raha represents the results of various error detection algorithms as a feature vector for training. In order to select the most promising records to be labeled by humans, hierarchical agglomerative clustering is used. Raha outperforms other error detection systems like dBoost [22] (outlier detection tool), NADEEF [7] (rule-based data cleaning system), KATARA [5] (knowledge bases to detect errors) and ActiveClean [19].

HoloDetect [11] uses data augmentation – a weak supervision technique – to enrich the training data sets and to train high quality error detection models with minimal human involvement. The basic

idea is to learn data augmentation policies from potentially noisy input data.

Currently one of the most advanced methods is Rotom [23], which also uses data augmentation to improve the training of machine learning models. In particular, Rotom formulates data augmentation as a sequence to sequence task. The key idea is to use meta-learning to automatically learn policies for combining training examples from different data augmentation operators and to assemble them into high-quality training signals.

Our approach uses recurrent neural networks (RNN). The advantage of this approach is that no specific feature engineering is required and can be applied to a wide range of error detection problems. Unlike HoloDetect or Rotom, our approach does not require additional data augmentation techniques. Also, our approach does not need expensive GPU resources because it uses a relatively simple way to analyze sequential data. Compared to Long Short-Term Memory (LSTM) [14], gated recurrent unit (GRU) [6] or Bidirectional Encoder Representations from Transformers (BERT) [9], RNNs are less complex and therefore do not need as much time for training. To reduce the size of required training data, i.e. the number of labelled tuples, our system uses a novel algorithm to choose tuples which give the most information content for our algorithms to learn. This leads to better and more robust (lower standard deviation) results.

3 BACKGROUND

This section provides background information about character embedding [17] and recurrent neural networks that serves as the basis for our error detection.

3.1 Character Embedding

Typical values in databases are either of type numerical, date or text, i.e. character values. However, in order to train machine learning models for error detection, we need to convert all our input data into a vector space of numerical values between -1 and 1. The advantage in the comparison to the *one-hot representation* is that embedding considers the context of the occurring items, i.e. it groups similar items together in the vector space. This method typically yields better results than the *one-hot representation* because the low number of the different dimensions of the vector space typically prevent problems in the back propagation steps during training the model (for example, vanishing gradient [13]).

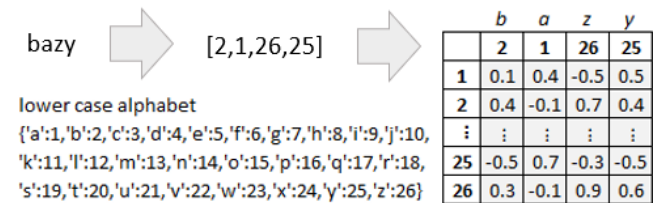


Figure 1: Character embedding for the example sequence 'bazy'.

The first step is to generate a dictionary where each unique character is encoded with an index ranging from 1 to the maximum

number of characters. As an example we use the lower case alphabet ('a':1,'b':2,...,'y':25,'z':26) consisting of 26 characters. This means that the character 'a' is encoded by the number 1. For example, if we want to encode a sequence of 4 characters, for instance 'bazy', we get the sequence 2,1,26,25. If we want to encode the above sequence ('bazy' -> [2,1,26,25]) using *character embedding*, we have to produce a matrix with dimension 4x26. The first dimension is of size 4 because our input sequence has a maximum length of 4. The second dimension is of size 26, which is due to the dimension space for the alphabet in our system – it could be higher or lower if desired (see example embedding in Figure 1). The values in the figure are fictitious and they are for visualization only.

In our error detection system we use this approach to transform a sequence of characters to a sequence of numbers in the data preparation step (see Section 4.1).

3.2 Recurrent Neural Networks

Recurrent Neural Networks (RNNs) are so-called sequence to sequence models that are typically used for natural language processing (NLP) tasks such as machine translation, language generation or question answering. The approach is helpful if you need the context from the previous input to make predictions about subsequent outputs, for instance, to predict the next word of a sentence. In these settings the inputs are character embeddings x_t to x_{t+n} – similar to the ones we discussed in the previous section.

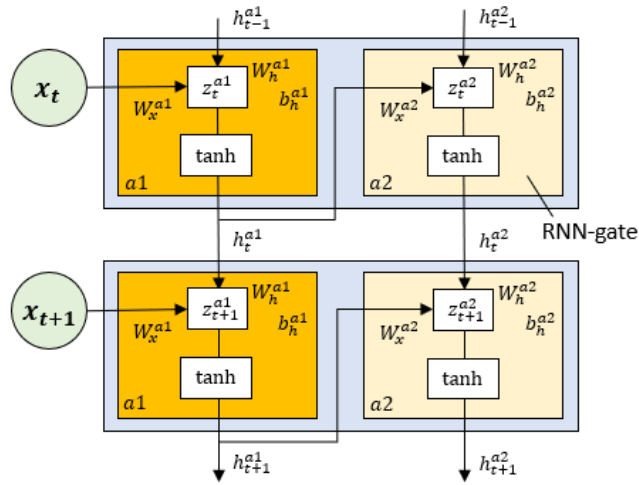


Figure 2: Two-stacked RNN.

Figure 2 shows a *forward two-stacked RNN* with the two levels $a1$ (left side, shown in orange) and $a2$ (right side, shown in yellow). The input to the RNN are the vectors of the characters, i.e. x_t and x_{t+1} . W_x and W_h are the weights and b_h are the biases of the RNN which we train. The weights and biases for all RNN-gates of the same level are equal. Moreover, the weights and biases of level $a2$ receive the computed hidden state h^{a1} as input. The hidden states h^{a1} and h^{a2} memorize the history of the previous step. Keeping in mind the

history of previous steps is important for learning dependencies between different characters, which is also the big difference to a fully connected neural network, which does not keep history. The parameter z^{a1} combines the information from the actual character x and the previous history h^{a1} . After this step the activation function \tanh is used to generate the history for the next RNN-gate.

The parameters z_t^{a1} , h_t^{a1} , z_t^{a2} and h_t^{a2} are calculated as follows:

$$z_t^{a1} = W_x^{a1} \cdot x_t + W_h^{a1} \cdot h_{t-1}^{a1} + b_h^{a1} \quad (1)$$

$$h_t^{a1} = \tanh(z_t^{a1}) \quad (2)$$

$$z_t^{a2} = W_x^{a2} \cdot h_t^{a1} + W_h^{a2} \cdot h_{t-1}^{a2} + b_h^{a2} \quad (3)$$

$$h_t^{a2} = \tanh(z_t^{a2}) \quad (4)$$

4 SYSTEM ARCHITECTURE

In this section we introduce two different neural network architectures for error detection given the clean and dirty version of a dataset. The main difference between the two architecture models is that the first model *Two-Stacked Bidirectional RNN* (TSB-RNN) receives input from one dataset (value_x, i.e. the actual data values), while the second model *Enriched Two-Stacked Bidirectional RNN* (ETSB-RNN) receives input from three datasets (value_x, i.e. the actual data values, length_norm, i.e. the standardized length of value_x, as well as metadata information about the attributes). Therefore, the second model is more complex and needs more time for training (see Section 5.6).

We will first discuss the data preparation step, then three different algorithms for selecting training data, and afterwards we describe the detailed system architecture for the models TSB-RNN and ETSB-RNN. The basic idea is that the data preparation step transforms the input data such that it can later be used by the neural network for error detection. Moreover, in order to enable effective learning with the minimal number of labeled training data samples, we discuss various methods of choosing those labeled training datasets.

4.1 Data Preparation

First we have to prepare the data so that we can produce a train- and testset, which in turn can be fed into a neural network to detect errors in the dataset and measure the results. Therefore, we perform the following steps (see Figure 3):

- (1) **Input:** We load both datasets (dirty and clean) as dirty_table and clean_table. Both are wide formats, i.e. each row is a tuple. In the example, cells with errors are marked in orange.
- (2) **Structure Transformation:** Next we remove preceding white spaces. Additionally, we add a column named 'id_' as a sequence number for every row to identify the tuple in later steps. Also we rename the different column names in the dirty_table to have identical names with the clean dataset. We need this renaming and the added column to combine the information of both datasets and create a new one (df).

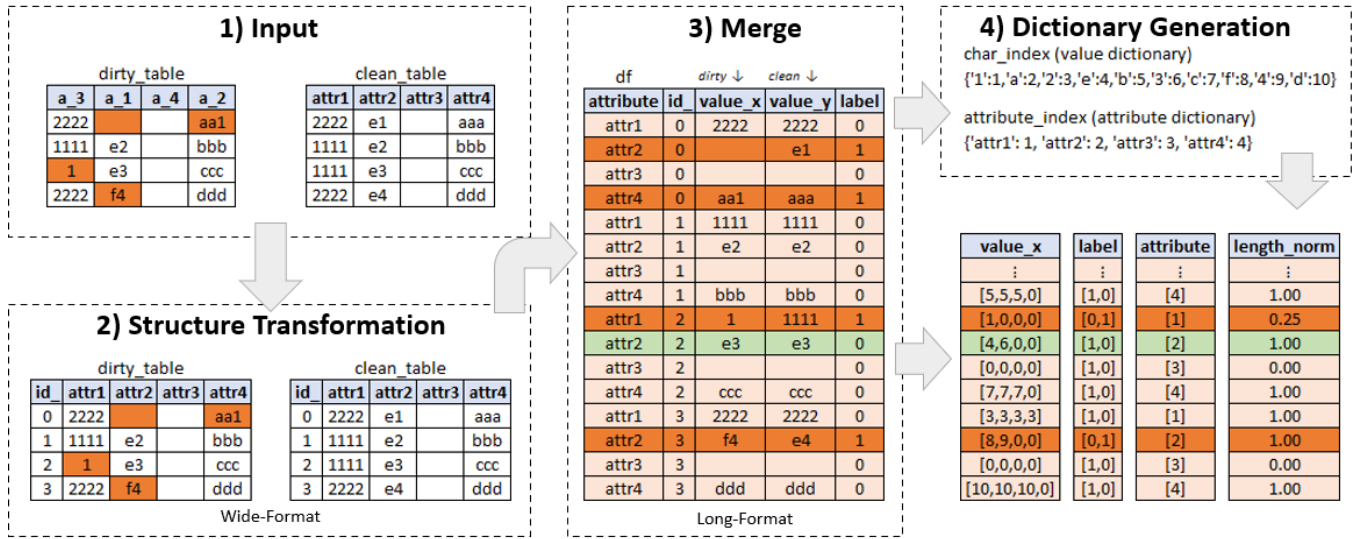


Figure 3: Data Preparation Process.

(3) **Merge:** Afterwards we combine the two tables to the dataset *df* where every cell of the *dirty_table* / *clean_table* is saved in the columns *value_x* / *value_y*, respectively. The new dataset has the long format, i.e. each tuple ('*id*') has the same number of rows (number of attributes). For the models we need a label, which includes '0' (correct value) or '1' (wrong value). We get this label when comparing *value_x* and *value_y*. If the value has more than 128 characters, e.g. in datasets hospital, movies and rayyan as described in Section 5, we cut them off. Our experiments showed that this approach achieves good F1-score results and reduced the training time.

For the Algorithm 3 (DiverSet) – see Section 4 – we need information if the *value_x* is empty (1) or not (0) and put it in a new column ('empty'). Also, the algorithm needs concatenated information about the attribute and the corresponding value ('*value_x*'). Therefore, we produce a new column ('concat').

Finally, we compute the length of *value_x* (i.e. the number of characters) in relation to the value with the longest length per attribute ('*length_norm*'). This information is used by our novel algorithm ETSB-RNN.

(4) **Dictionary Generation:** Before we can feed the data into a neural network, we need to transform the data types from character to numeric. We produce a *value dictionary* (*char_index*) which contains an index for each character in *value_x*. We use the dictionary to convert the sequence of characters to the appropriate indexes. To make sure all attribute values have the same length, we add the values 0 at the end. The character embedding is done inside the neural network architectures (see Section 4.3).

For the ETSB-RNN we also need an *attribute dictionary* (*attribute_index*) which includes an index for each attribute. The attribute dictionary contains metadata information about

attributes and shows from which attribute the content comes from.

Assume, for instance, that the model learns that all values of attribute 'age' only contain numeric values. Further assume that the model receives the value 'San Francisco' which comes from the attribute age. Since 'San Francisco' is not a numeric value, the model can recognize it as a wrong value for the attribute age.

Example (see bottom right side of Figure 3): The green highlighted row shows the transformed characters. For instance, the value 'e3' is encoded as the characters 'e' and '3' and is stored in the *value dictionary*. Moreover, since the value 'e3' refers to attribute 'attr2', this metadata information is stored in the *attribute dictionary*. Finally, since the value 'e3' is correct, the corresponding label is encoded as '0' (true). Therefore, the first value in column 'label' has the value 1.

In our simple running example we assume the length of our *value dictionary* is 10 characters and the longest sequence of characters in '*value_x*' is 4. Since the dimension of the sequence 'e3' is smaller than the longest value (e.g. '2222' or '1111'), the preparation step takes automatically the index 0 at the end. In other words, we pad short sequences of characters with the end-indicator.

4.2 Algorithms for Trainset Selection

After the preparation of the data, the system needs training and test data to train the system and then to test the trained machine learning model. However, real world datasets typically have very little or no labelled data. Hence, the user has to provide labels for the trainset. To reduce the human effort for generating training data, the system asks the user to provide labels for 20 tuples, that give our system the most impact during training. Hence, we analyze three different algorithms for choosing the training set.

For illustrating our algorithms, we use the dataset described in Section 4.1 in step (3) as input. Note: We do not use the *value_y* or

label for choosing tuples. We only consider the value_x from the dirty dataset.

Our baseline approach is Algorithm 1, which chooses 20 tuples at random. This means that every tuple ('id_') in the dataset has the same probability of being selected as labeled training data. Since we have a long-format dataset as input (see the dataset of the merge step in Figure 3), we have to delete duplicated id_s before we can choose tuples. The algorithm is very simple and does not consider the data content for choosing tuples.

Algorithm 1 RandomSet

Input: n_obs -> number of tuples (id_) for training
 df -> dataset with all data
Output: ID_train -> id_s of tuples for the trainset (size: n_obs)

```

1: function RANDOMSET( $n\_obs, df$ )
2:    $ID\_all \leftarrow \text{unique}(df['id\_'])$ 
3:    $ID\_train \leftarrow \text{random\_sample}(n\_obs, ID\_all)$ 
4:   return  $ID\_train$ 
5: end function
```

We also evaluate the label sampling approach of Raha [21] as shown in Algorithm 2. First, Raha generates and executes a set of error detection strategies. Then, it generates a vector which stores for each data cell and strategy a binary number referring to an error or no error. In the next step, Raha groups similar cells with the help of the previously created vectors. As the last step, Raha draws 20 samples from different clusters.

Algorithm 2 RahaSet

Input: n_obs -> number of tuples (id_) for training
 df -> dataset with all data
Output: ID_train -> id_s of tuples which are used for the trainset (size: n_obs)

```

1: function RAHASSET( $n\_obs, df$ )
2:    $ID\_train \leftarrow \text{RahaApproach}(n\_obs, df)$ 
3:   import raha
4:    $sampler\_list \leftarrow []$ 
5:   run\_strategies( $df$ )
6:   generate\_features( $df$ )
7:   build\_clusters( $df$ )
8:   while  $\text{len}(labeled\_tuples) < n\_obs$  do
9:     sample\_tuple( $df$ )
10:     $sampler\_list \leftarrow sampler\_list \cup sampler\_tuple$ 
11:  end while
12:  return( $sampler\_list$ )
13: }
14: return  $ID\_train$ 
15: end function
```

Finally we introduce a novel label sampling Algorithm 3, whose goal is to select a diverse trainset. The intuition is to select tuples with values that have not been seen previously. In other words, the newly selected tuples should increase the information content of our trainset the most. In case two candidate tuples contain the

same number of unseen attribute values, the tuple with the highest number of empty attribute values is chosen. Our hypothesis is that empty values give us more information for the system to learn – because if there are empty values in other attributes, we can learn if they should be empty or not. In case all candidate tuples have the same number of unseen attributes and empty attribute values, the tuples are chosen randomly.

Algorithm 3 DiverSet

Input: n_obs -> number of tuples (id_) for training
 df -> dataset with all data
Output: ID_train -> id_s of tuples which are used for the trainset (size: n_obs)

```

1: function DIVERSET( $n\_obs, df$ )
2:    $df\_rest \leftarrow df$ 
3:    $ID\_train \leftarrow []$ 
4:   for  $i=1$  do  $n\_obs$ 
5:      $\#unseenAttr \leftarrow \text{count}(df\_rest).groupby('id\_')$ 
6:      $\#empty \leftarrow \text{sum}(df\_rest['empty']).groupby('id\_')$ 
7:      $candidateID \leftarrow (id_, \#unseenAttr, \#empty)$ 
8:      $candidateID \leftarrow candidateID.max(\#unseenAttr)$ 
9:      $candidateID \leftarrow candidateID.max(\#empty)$ 
10:     $sampler\_ID \leftarrow \text{random\_sample}(1, candidateID)$ 
11:     $ID\_train \leftarrow ID\_train \cup sampler\_ID$ 
12:
13:     $seenAttr \leftarrow \text{unique}(df['concat'].where(id\_ = ID\_train))$ 
14:     $df\_rest \leftarrow df[df['concat'] \neq seenAttr]$ 
15:  end for
16:  return  $ID\_train$ 
17: end function
```

To show the algorithm in a more understandable way, consider Figure 4 as an example with two observations, i.e. $n_obs=2$. We used the same dataset (df) which we presented in Figure 3.

- **Start:** We initialize df_rest as df and ID_train as an empty list. Next we enter the loop, where the number of passes is given by n_obs .
- **i=1:** In the first step we produce the $candidateID$ with both attributes $\#unseenAttr$ (counts all occurring $id_$) and $\#empty$ (sums up all empty values per $id_$). Next, we keep only the tuples ($id_$) with the highest value in $\#unseenAttr$ and from this new result only the tuples with the highest value in $\#empty$. In this step we get $id_=0$ (bold framed) as a unique result and put it in ID_train . For $seenAttr$ we put all values from the column $concat$ where $id_=0$ in the dataset (df_rest). As the last step we delete all rows (marked in gray) which have occurring values in $seenAttr$ (yellow values) out of df_rest . The rows with $id_ \neq 0$ are deleted because the values in column $concat$ equals the values in the dataset $seenAttr$.
- **i=2:** The second step produces again the dataset $candidateIDs$ like the methods in $i=1$. Here, we do not receive a unique result because $id_=1$ and $id_=2$ have $\#unseenAttr=3$ and $\#empty=0$. From this $candidateID$ (bold framed) we randomly sample one of the two tuples ($id_$). We assume the random

df_rest				
attribute	id	value_x	concat	empty
attr1	0	2222	attr1_2222	0
attr2	0		attr2_	1
attr3	0		attr3_	1
attr4	0	aa1	attr4_aa1	0
attr1	1	1111	attr1_1111	0
attr2	1	e2	attr2_e2	0
attr3	1		attr3_	1
attr4	1	bbb	attr4_bbb	0
attr1	2	1	attr1_1	0
attr2	2	e3	attr2_e3	0
attr3	2		attr3_	1
attr4	2	ccc	attr4_ccc	0
attr1	3	2222	attr1_2222	0
attr2	3	f4	attr2_f4	0
attr3	3		attr3_	1
attr4	3	ddd	attr4_ddd	0

ID_train				
id_				
0				
1				

candidateID				
id	#unseenAttr	#empty		
0	4	2		
1	4	1		
2	4	1		
3	4	1		

ID_train				
id_				
0				

seenAttr				
concat				
attr1_2222				
attr2_				
attr3_				
attr4_aa1				

ID_train				
id_				
0				
2				

seenAttr				
concat				
...				
attr1_1				
attr2_e3				
attr4_ccc				

Figure 4: Example of Algorithm 3 for the first two loops to select a trainset.

sample chooses the tuple with $id_=2$ as result. We again append it to ID_train and append all values in column *concat* for this $id_$ to the dataset *seenAttr*. At the end we again delete all rows (marked in gray) which have occurring values in *seenAttr* (yellow values) out of *df_rest*.

- **Return:** As result the algorithm returns a list with all tuples ($id_$) which we will use as our trainset. This example gives us the tuples with $id_$ equal 0 and 1 as trainset.

Each algorithm returns a list of ' $id_$'s of all tuples that we take for training. With this we create the train- (X_train , $X_train_attribute$ and Y_train) and testsets (X_test , $X_test_attribute$ and Y_test) which are the input for our neural networks.

4.3 Neural Network Architectures

We now describe the two different neural network architectures that we use for error detection.

4.3.1 Two-Stacked Bidirectional RNN (TSB-RNN). As the first model we use a *two-stacked bidirectional RNN*, which has 64 units and as activation function the hyperbolic tangent (see top left part of Figure 5). Note that the orange parts refer again to the stacking level $a1$ and the yellow parts to the stacking level $a2$ as also shown previously in Figure 2. However, unlike the two-stacked RNN of Figure 2, our architecture shown in Figure 5 uses a *bidirectional RNN*, which means that in addition to the forward layer (highlighted in light blue), we also have a backward layer (highlighted in dark blue).

Therefore, we compute the settings forward and backward through the character sequences.

As input the model receives the datasets X_train and X_test , where every character is encoded as a sequence of numbers. For instance, in Figure 5, the input is the character string 'e3' represented as a sequence of numbers (7,9,0,0). First, in the embedding layer the system transforms the sequence of numbers into a vector with the dimension of the used value dictionary. The output of our model is a concatenation of the output from the forward path (dim 64) and the backward path (dim 64). In addition, we use a fully connected layer with dimension 32 and ReLu for activation. At the end there is a batch normalization [15] to standardize the input to the softmax. It reduce the generalization error and is a regularization technique.

Finally, the model has to decide if the sequence of characters has the value '0' (correct) or '1' (wrong). For this last step the model uses a fully connected layer with a softmax. During training the model needs the dataset Y_train to optimize the loss. The dataset Y_test is not used for training but for validating the trained model with it.

4.3.2 Enriched Two-Stacked Bidirectional RNN (ETSB-RNN). Our second architecture ETSB-RNN is an enriched version of TSB-RNN. In particular, we additionally provide the *attribute information* $X_train_attribute$ and $X_test_attribute$ as a second input. The attribute information tells the neural network that, for instance, the encoding of the character sequence 'e3' refers to attribute attr2 (see in the left lower part of Figure 5). Note that TSB-RNN, that we discussed previously, only has the information about the character embedding

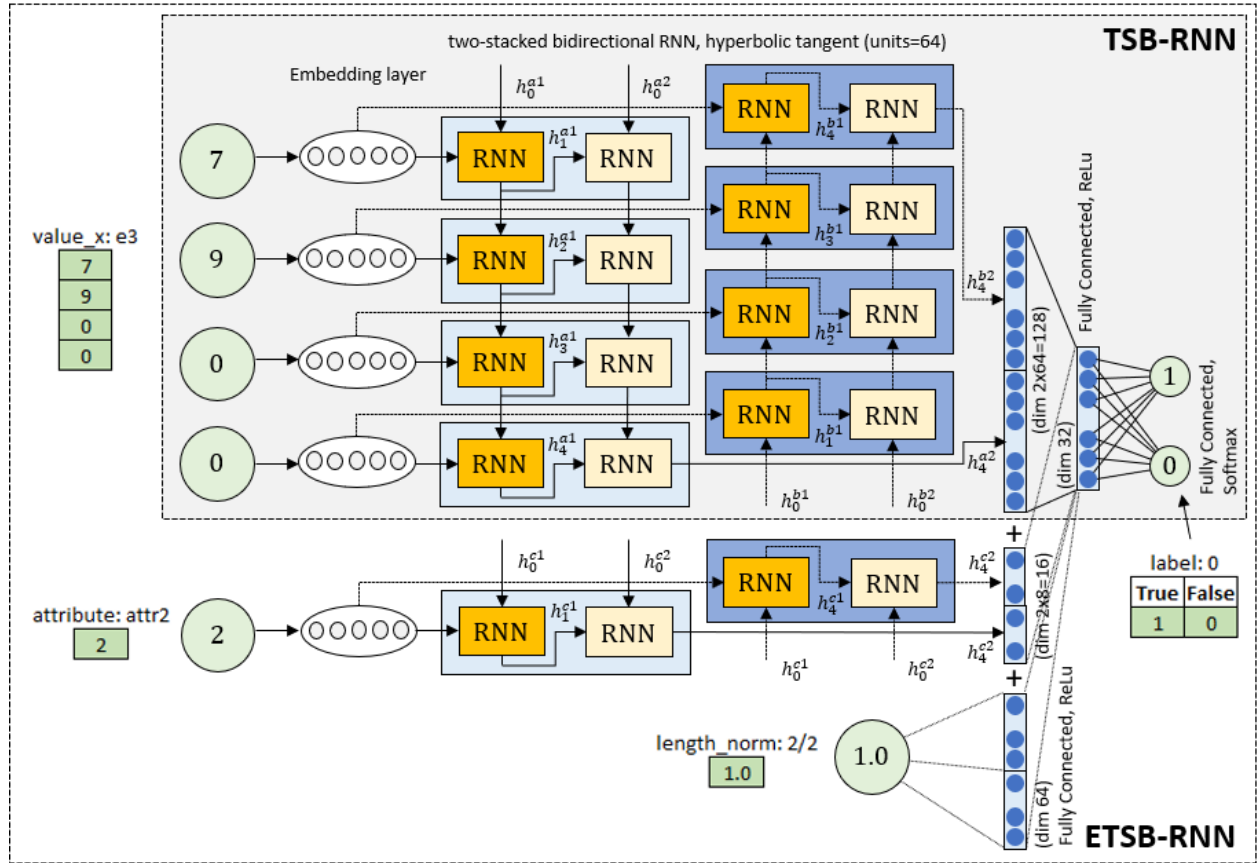


Figure 5: Overview of our bidirectional RNN architecture for error detection. Top part: TSB-RNN = Two-Stacked Bidirectional RNN. Bottom part: ETSB-RNN = Enriched Two-Stacked Bidirectional RNN.

but does not know to which attribute that value refers to. Also, we feed the system with the value about the ratio of the actual length of the input sequence to the maximum length of the attribute (length_norm) – see in the middle lower part of Figure 5.

As an additional layer the model uses an embedding layer and a two-stacked bidirectional RNN which has 8 units and as activation function the hyperbolic tangent. Also there is additionally a fully connected layer with dimension 64 and the activation function ReLu. We concatenate the output of these two new layers with the output of the previous two-stacked bidirectional RNNs. Afterwards we again use the two fully connected layers and the batch normalization described in TSB-RNN.

5 EXPERIMENTS

In this section we evaluate our two different architectures and measure how well they can detect errors in 6 commonly-used benchmark datasets. We also compare the results against the state-of-the-art-algorithms for error detection. Note that to enable the reproducibility of our experiments, we share our code on Github¹.

¹https://github.com/holzesev/E_TSB-RNN

5.1 Datasets

For our experiments we used 6 datasets that were also used by the state-of-the-art error detection systems, e.g. from Raha [21] and Rotom [23]. All datasets have a version of data with errors (dirty) and the corresponding ground truth (clean).

Table 2: Overview of datasets with error types. MV: Missing Value, T: Typo, FI: Formatting Issue, VAD: Violated Attribute Dependency.

Name	Size	Error Rate	Different Characters	Error Types
Beers	2,410x11	0.16	86	MV, FI, VAD
Flights	2,376x7	0.30	70	MV, FI, VAD
Hospital	1,000x20	0.03	46	T, VAD
Movies	7,390x17	0.06	135	MV, FI
Rayyan	1,000x10	0.09	101	MV, T, FI, VAD
Tax	200,000x15	0.04	69	T, FI, VAD

An overview of the datasets is given in Table 2. The table shows the size of each dataset including the number of records and attributes, the error rate, i.e. the fraction of data that is wrong, the number of different characters, i.e. the size of the value dictionary and finally, the error types² in the corresponding dataset.

Beers: This dataset is about different beers and contains information about the type of beer, the international bitterness unit (IBU), the alcohol by volume (ABV), the brewery and more. Typical data errors are, for instance, formatting issues (FI) in ounces ('12.0 oz' rather than '12.0'), ABV ('0.061%' rather than '0.061'), violated attribute dependencies (VAD) between city / state and also missing values (MV) in state ('NaN' rather than 'CA').

Flights: This dataset contains information about the same flights received from different data sources. Depending on the data sources, flights have violated attribute dependencies (VAD) in the departure and arrival times. In several cases the time information has missing values (MV) ('' rather than, e.g. '3:31 p.m.'), varies by few minutes ('9:00 a.m.' rather than '8:42 a.m.') or has a formatting issues (FI) ('12/02/2011 6:55 a.m.' rather than '6:55 a.m.').

Hospital: This dataset contains information about hospitals and the corresponding treatments (measures) for certain diseases, e.g. 'heart attack patients given aspirin at arrival'. In this dataset the hospital or measure information appears several times in the dataset but information varies, i.e. there are many violated attribute dependencies (VAD). Moreover, there are typos (T) in attributes, e.g. wrong city names such as 'Birmingxam' rather than 'Birmingham'. Note: For humans, the errors are not very hard to detect because they are marked with the character 'x'.

Movies: This dataset contains information about different movies and corresponding attributes like duration, genre etc. Data errors are formatting issues (FI) (e.g. RatingCount: '379,998' rather than '379998.0', e.g. name: 'Frankie & Johnny' rather than 'Frankie and Johnny' or RatingValue: '8.0' rather than '8'), missing values (MV) (e.g. duration: 'NaN' rather than '126 min'), several year indications instead of only one year or missing words.

Rayyan: This dataset includes scientific articles written by researchers, and the related attributes like title, author etc. Data errors are again formatting issues (FI) (e.g. journal_issn: 'Mar-22' rather than '22-Mar' or article_pagination: '70-6' rather than 'Jun-70'), missing values (MV) in article_jissue and typos (T) in journal_title or article_title.

Tax: This dataset contains information about people which is used for computing taxes, like name, gender, salary etc. Data errors are typos (T) in f_name (e.g. 'Jun"ichi' rather than 'Jun'ichi') or city (e.g. 'ARCHIE-*' rather than 'ARCHIE'), formatting issues (FI) in zip (e.g. '01907' rather than '1907') or rates (e.g. '7.0' rather than '7') and violated attribute dependencies (VAD) between state / city and marital_status / has_child.

5.2 Training and Testing Setup

We validated the models 10 times and used the Algorithm 3 (Diver-Set) for choosing the 20 tuples for training. For instance, for the dataset Beers we got a trainset of size 220, i.e. 20 tuples x 11 attributes, and a testset of size 26,290, i.e. 2,390 tuples x 11 attributes.

²We took the definition of the error types from Raha [21].

During training we used a model batch size of a quarter of the trainset, i.e. with Beers: $220 / 4 = 55$.

To find the algorithm which works best for our systems, we repeated the experiments several times with every algorithm described in Section 4.2. We reached the best results with our novel Algorithm 3. The basic idea of Algorithm 3 is to prevent, that the same values are selected for the trainset. We suppose that through the diverse trainset generated by Algorithm 3, the two neural networks algorithms *TSB-RNN* and *ETSB-RNN* received the most information content for the characters and hence could learn error patterns in the data.

The number of epochs was 120. After every epoch we saved the training weights (with a callback) if the computed loss of the trainset was less than in the previous epochs. For the loss-function we used the binary cross-entropy and as the optimizer RMSprop. We used the testset to measure precision, recall and F1-score for the best weights. Moreover, we computed the averages and standard deviations of these measures.

5.3 Comparison

In this section we compare our results with the Configuration Free Error Detection System (Raha, [21]), as well as with the Meta-Learned Data Augmentation Framework (Rotom, [23]), which is the state-of-the-art of error detection in databases. For both approaches we report the results of the original papers. The results are shown in Table 3.

Note that Rotom did not take the dataset Flights for validation. Therefore, in Table 4 we show the average (AVG) and standard deviation (S.D.) for the error detection systems for all datasets without Flights (1) and in addition for all datasets including Flights (2). Also both papers did not show the standard deviation from the 5 times experiments (we did 10 times experiments).

If we compare the experiments of both papers, we can see different results in F1-score for the datasets Beers (+/-0.01), Hospital (+/-0.02), Movies (+/-0.06) and Rayyan (+/-0.03) for Raha. However, neither the authors of Raha nor Rotom show the standard deviation. So we can not see how the state-of-the-art models do scatter. The dataset Tax was only used in the paper of Rotom. Moreover, they did not measure precision and recall, hence the value n/a in Table 3.

Additionally, Rotom and Rotom+SSL validated the results with 50, 100, 150 and 200 labelled cells and took the best results for comparison. However, in the real world we do not have a testset, therefore we do not know how many labelled cells would work best. So Rotom and Rotom+SSL would have to report the results from 200 labelled cells and not <200. Therefore, to have a direct comparison with both Raha and our approach, we reported the results for the models Rotom (Movies and Tax) and Rotom+SSL (Movies and Rayyan) using 200 labelled cells.

Let us now analyze the results. Our approach ETSB-RNN, which also uses the metadata information, outperforms the simpler model TSB-RNN on all datasets. Also for all datasets we got similar or less standard deviation, therefore we take the ETSB-RNN for comparing with Raha, Rotom and Rotom+SSL.

Our model ETSB-RNN outperform Raha on 3 out of 6 datasets for the F1-score (see Table 3). In particular, we got better results

Table 3: Comparison between the different models (20 labeled tuples). P = Precision, R = Recall, F1 = F1-Score, S.D. = Standard Deviation. Our approaches are TSB-RNN and ETSB-RNN.

Name	Beers			Flights			Hospital			Movies			Rayyan			Tax		
	P	R	F1	P	R	F1	P	R	F1	P	R	F1	P	R	F1	P	R	F1
Raha	0.99	0.99	0.99	0.82	0.81	0.81	0.94	0.59	0.72	0.85	0.88	0.86	0.81	0.78	0.79	n/a	n/a	0.91
Rotom	n/a	n/a	0.99	n/a	n/a	n/a	n/a	n/a	1.00	n/a	n/a	0.68	n/a	n/a	0.86	n/a	n/a	0.97
Rotom+SSL	n/a	n/a	0.99	n/a	n/a	n/a	n/a	n/a	1.00	n/a	n/a	0.54	n/a	n/a	0.76	n/a	n/a	1.00
TSB-RNN	0.99	0.94	0.96	0.77	0.63	0.69	0.98	0.95	0.97	0.96	0.79	0.87	0.83	0.73	0.78	0.83	0.90	0.85
S.D.	0.01	0.01	0.01	0.05	0.04	0.02	0.01	0.02	0.01	0.04	0.04	0.03	0.05	0.06	0.05	0.16	0.08	0.11
ETSB-RNN	1.00	0.96	0.98	0.81	0.68	0.74	0.98	0.95	0.97	0.96	0.81	0.88	0.87	0.83	0.85	0.82	0.92	0.86
S.D.	0.00	0.02	0.01	0.03	0.04	0.02	0.03	0.02	0.02	0.03	0.05	0.02	0.06	0.02	0.03	0.15	0.07	0.10

for Hospital (+0.25), for Movies (+0.02) and for Rayyan (+0.06). We achieved similar results for Beers (-0.01) and Tax (-0.04), only for Flights (-0.07) we got considerably worse results. When comparing the average F1-score (AVG) without Flights (1) (+0.06) and with Flights (1) (+0.01), our approach ETSB-RNN outperforms Raha (see Table 4). Also the standard deviation is smaller for our system.

In comparison to Rotom and Rotom+SSL (see Table 3), our model ETSB-RNN got a better result for Movies (+0.20 / +0.34). For Beers (-0.01 / -0.01), Hospital (-0.03 - 0.03) and Rayyan (-0.01 / +0.09) we achieved very similar results – only for Tax (-0.11 / -0.13) our approach performs considerably worse than Rotem+SSL. The average F1-score (AVG) is a little slightly higher (+0.01 / +0.05) than Rotom and Rotom+SSL (see Table 4). In addition, the standard deviation over all datasets is better.

Table 4: Average F1-score (AVG) and Standard Deviation (S.D.) for the different models of Table 3.

Name	Without Flights (1)		With Flights (2)	
	AVG	S.D.	AVG	S.D.
Raha	0.85	0.08	0.85	0.07
Rotom	0.90	0.10	n/a	n/a
Rotom+SSL	0.86	0.17	n/a	n/a
TSB-RNN	0.89	0.06	0.85	0.08
ETSB-RNN	0.91	0.05	0.88	0.06

5.4 Learning Analysis

We will now analyze the learning behavior of our approaches in more detail. First, we perform a direct comparison between our approaches TSB-RNN and ETSB-RNN. Afterwards, we compare the accuracy of our best approach ETSB-RNN during training and testing to see if our model does not show overfitting and generalizes well.

To show the different improvements of the accuracy over various epochs of our models, we plotted the average and the confidence intervals of the 10 experiments (see Figure 6) for the test dataset. We can see that TSB-RNN and ETSB-RNN show a gradual increase of the accuracy. There is no clear deterioration. As described we optimized the train loss and took the settings from the epoch which achieved the lowest loss. We can see the ETSB-RNN outperforms

TSB-RNN on all datasets, except for Tax, where both models show practically the same performance.

Finally, for ETSB-RNN, which performed best in our experiments, we also compared the average train- and test-loss (see Figure 7). The results show that the model performs well and does not suffer from overfitting. For all datasets our model achieves almost a perfect result for the train-accuracy. There are also gaps in the curves of the test-accuracy but these are not critical because the model chooses the epochs which do not have gaps and learns how to choose the optimal parameter setting based on the decreasing training loss.

The dataset flights, which reached the worst F1-score, has a big gap between the train- and test-accuracy and also the confidence interval is large. Therefore, our model does not work well for this dataset. The curve for test-accuracy for Hospital looks almost perfect, also because the confidence interval is very small. The datasets Beers, Movies and Rayyan also show a converging test-accuracy curve and the confidence interval is small. For Tax we got a curious result. After epoch 30 there are no wave movements in the test-accuracy. In addition, the model chooses the optimal parameter settings from training close to epoch 120, i.e. the green dots and blue triangles.

5.5 Error Analysis

We will now perform a more detailed error analysis of our algorithm ETSB-RNN for each of the datasets.

As described previously, the dataset **Flights** contains duplicate values due to different information sources for the same flight. ETSB-RNN reached an average F1-score of 0.81. However, it had some problems to detect errors due to varying departure and arrival times of different sources, e.g. flight 'UA-257-JFK-SFO' of source 'orbitz' has a departure time of '2:46 p.m.' while the same flight of source 'flightstats' has a departure time of '2:26 p.m.' The reason is that the model does not share the information about the name of the flight (i.e. the ID to recognize that flights are the same) and therefore it cannot detect duplicate records. Raha could find the dependencies in a better way, but there is also room for improving the result.

Detecting errors in the **Hospital** dataset is quite straightforward because the errors are marked with 'x' (e.g. 'hexrt fxilure'). The model TSB-RNN received an almost perfect result with an average F1-score of 0.97. These types of error are easy to detect and do not dependent on the attribute metadata information. Hence, the

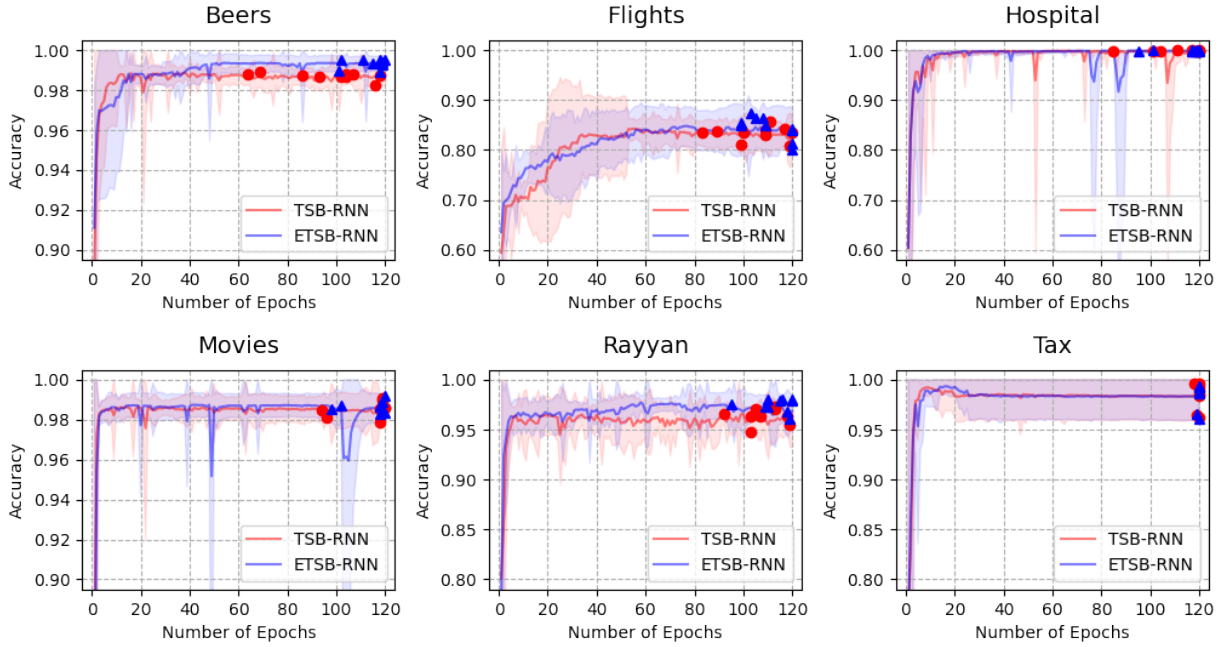


Figure 6: Analysis of the average test-accuracy during training (10-times experiment) and the confidence interval for the different epochs. The red dots (TSB-RNN) and blue triangles (ETSB-RNN) show the selected epoch for the best model and their corresponding test-accuracy.

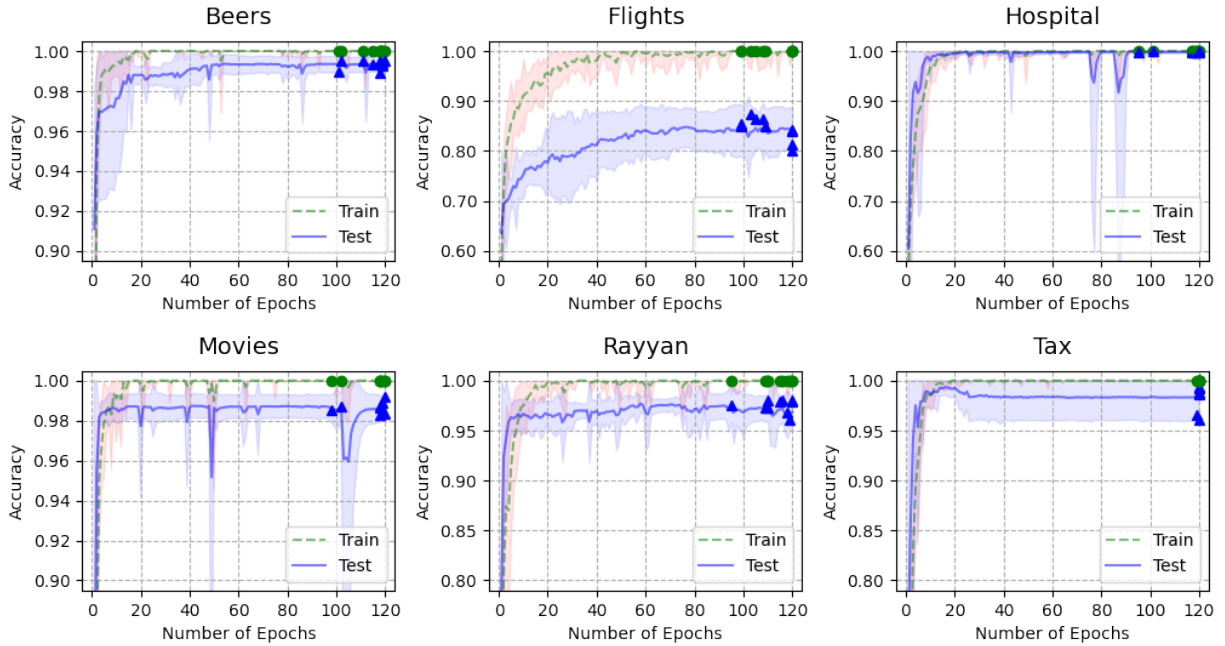


Figure 7: Comparison of average train- and test-accuracy (10-times experiment) and the confidence interval for the different epochs of ETSB-RNN. The green dots shows the epoch and corresponding train-accuracy with the lowest train-loss per experiment. The blue triangles show the corresponding test-accuracy.

models TSB-RNN and ETSB-RNN achieved very similar results. Rotom and Rotom+SSL outperform our results.

For the dataset **Movies** the model ETSB-RNN does not recognize errors in the attribute *Creator*. The reason is that some parts of the values are missing. (i.e. 'Roger Kumble' instead of 'Choderlos de Laclos, Roger Kumble'). For the attribute *Duration* the model ETSB-RNN cannot detect the value 'NaN' because in some rows the correct value is 'NaN' and in others the correct value is '96 min'. However, the model still achieved good results with an average F1-score of 0.88. The ETSB-RNN reached the best result for the dataset.

The errors of the dataset **Rayyan** are mostly due to non-recognized special characters. In this setting, the average F1-score for our model ETSB-RNN is 0.85. Rotom achieved a similar result to ETSB-RNN.

For the last dataset **Tax** we achieved a very high standard deviation. It means we got partly good and bad results on the 10-times experiment. This suggests that our algorithm *DiverSet* did not select the optimal data for the trainset. Rotom+SSL achieved almost a perfect result and therefore it clearly outperformed our result for ETSB-RNN.

5.6 Training Time

As we described previously, we ran the experiments 10 times. To see how fast our system is, we measured the training time per dataset and model. We used Colaboratory³ (Colab) which is a product from Google Research that allows using Python code in a Jupyter notebook in the browser. The access to the computing resource (incl. GPU and TPU) is for free.

"The types of GPUs that are available in Colab vary over time. This is necessary for Colab to be able to provide access to these resources for free. The GPUs available in Colab often include Nvidia K80s, T4s, P4s and P100s. There is no way to choose what type of GPU you can connect to in Colab at any given time."

Table 5: Training time [sec] for the different datasets using TSB-RNN and ETSB-RNN. AVG = Average, S.D. = Standard Deviation.

Name	TSB-RNN		ETSB-RNN	
	AVG	S.D.	AVG	S.D.
Beers	92	1	101	1
Flights	47	0	54	0
Hospital	283	3	287	2
Movies	302	3	312	3
Rayyan	199	2	209	2
Tax	176	1	183	1
AVG	183		191	

The training times of our systems TSB-RNN and ETSB-RNN are shown in Table 5. The time range is between less than one minute (Flights) until about 5 minutes (Movies). The different time between the datasets depends on the number of attributes, number

³<https://colab.research.google.com>

of different characters and the size of the longest value. The ETSB-RNN requires more time for training than TSB-RNN because it is the more complex neural network system. Note that the training times are similar to the results shown for Rotom. However, we use less powerful GPUs.

5.7 Improvements

We observed partially larger values in the confidence intervals for some datasets, therefore we achieved different F1-scores for different datasets. We suppose the training data was still too little heterogeneous. This suggests that there is room for improving our novel label-selection Algorithm *DiverSet* to reach even better results.

Moreover, our approach does not consider functional dependencies between different attributes, e.g. if there is an attribute age with value '12' and an attribute salary with value '99,000' (see Table 1). In other words, a 12 year old person with a such a high salary is most probably a wrong combination, therefore the value in age or salary should be detected as an error. Detecting such kind of data errors using *functional dependencies* is part of our future work.

Another point in which our model has weaknesses is the recognition of duplicates seen in the dataset **Flights**. To improve this, we should integrate a way to *identify primary keys* in the future work. This information allows us to identify identical records stemming from two different sources and our system would know that it has to fuse the values in one record.

6 CONCLUSION

We introduced a new way to detect errors in databases using two-stacked bidirectional recurrent neural networks. The prerequisite to use our approach is a clean and dirty dataset that is needed to train the models. During training our system reduces the train-loss and chooses the weights and biases with the lowest loss. We trained and tested our models with 6 publicly available datasets and compared the results with the state-of-the-art error detection systems Raha, Rotom and Rotom+SSL. Our approach shows similar F1-scores compared to state-of-the-art systems. When considering the average of the F1-scores over all datasets, our approach called Enriched Two-StackedBidirectional (ETSB-RNN) outperforms state-of-the-art systems.

Our experiments show promising results of using machine learning for identifying errors in databases, which is an important part of data management and thus a prerequisite for subsequent data science tasks. The ultimate goal, however, is not only to detect errors but also to correct them. An interesting avenue of future research is to integrate our approach with the data repair systems of HoloClean [24] and Baran [20].

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