Evaluation Platform for Artificial Intelligence Algorithms

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Abstract: Currently, artificial intelligence (AI) algorithms have been receiving a lot of attention from researchers as well as from commercial product developers. Hundreds of different AI algorithms aiming different kind of real-life problems have qualitatively different results, based on the nature of data, the nature of the problems and based on the context in which they are used. Choosing the most appropriate algorithm to solve a particular problem is not a trivial task. The goal of our research is to create a platform, which can be used in the early stage of problem solving. With this platform, the user could be able to quickly train, test and evaluate several artificial intelligence algorithms and also they will be able to find out which is the algorithm that performs best for a specific problem. Moreover, this platform will help developers to tune the parameters of the chosen algorithm in order to get better results on their problem. We will demonstrate our approach by running different types of algorithms initially in the case of breast cancer sample dataset and after that we will use the platform for solving an anomaly detection problem.

1 INTRODUCTION

In the last decade, because of some new data acquisition technologies (e.g. sensor networks, IoT, Industry 4.0) and the wide spread of Internet-based applications, huge amount of data is becoming available. Efficient and higher level processing of such "big data" requires intelligent algorithms, not just for academic purposes, but also for real software products. More and more companies are trying to use artificial intelligence to solve complex problems, which earlier required a human expert. The hardest thing in enterprise applications using artificial intelligence is to estimate the time needed to create a minimum viable product. The difficulty of the estimation arise from the fact that before starting the project it is very hard to know exactly which of the algorithms will be the best match for that particular problem and for that particular dataset.

Algorithms can have very different and surprising results based on the nature of the problem and the nature of the training data. A good approach to overcome this difficulty is to get a small part of the dataset, implement multiple artificial intelligence related algorithms and benchmark those algorithms using the small sample dataset. This can be a very time consuming process and usually for enterprise products you don't have such time. Our intention is to create a platform which can be used to test multiple algorithms and evaluate them using appropriate metrics, this way reducing the time to production.

Another problem which usually occur is the tidiness of data. Raw data can have multiple problems (e.g. noise, sample misses, etc.), which occur because of bugs in the data collecting process. Preprocessing of the data can be a challenging task, which may require multiple subsequent tasks; therefore our platform contains a set of preprocessing algorithms which can be applied by a domain specialist, without specific programming knowledge. The data can also be visualized in multiple way (2D-3D charts, histograms, FFT, etc.), which can help spotting problems within the raw data. Another important aspect of AI algorithms is that someone can combine multiple algorithms to improve the quality of the results. In order to combine different algorithms the platform should offer the possibility to pipeline the partial results in a reconfigurable manner.

As a practical result, in the following sections we will describe our platform, which can be used to evaluate different AI techniques with different types of datasets, configuring the most effective pipeline in different contexts. We will show results obtained with different algorithms with and without preprocessing, showing in this way the effectiveness of the preprocessing step. We will tune and evaluate multiple algorithms and the results will be compared. So the purpose of this article is to describe a platform that allows a user to select, tune and pipeline multiple AI algorithms as well as preprocessing and displaying tasks. All these operations may be done without writing any code and without programming skills. This tool is very useful in the early stage of a project when the feasibility of a given approach must be measured or decided. The tool will give an estimate of the quality level that may be obtain using different AI algorithms. This will reduce the time to production and it will increase the chance for a successful implementation. Furthermore, we will introduce a step-by-step process for choosing the right sequence of algorithms.

The rest of the paper is organized as follows. Section 2 shows a brief description of related works. Section 3 contains the theoretical classification of different algorithms and the decision tree which can be used to choose the proper algorithm for a given problem. The architecture of the platform is described in Section 4. Experiments can be found in Section 5 and Section 6 concludes the article.

2 RELATED WORK

Choosing the best algorithm for a specific problem domain, taking into consideration multiple factors (e.g. the type of the data involved, the nature of the problem etc.) is a general problem in developing intelligent products. There are multiple articles which are trying to classify AI algorithms, based on different points of view.

In (Dasgupta and Nath, 2016) we can find a typical classification of machine learn algorithms, based on the nature of the training data as follows:

- supervised, if the training data is labeled;

- unsupervised, if there are no labels;

- semi-supervised, if some of the class labels are missing.

The problem with this classification is the fact that it considers only the nature of the training data, without taking in consideration the context of the problem. This approach reduces the searching space, but each class has too many algorithms to be considered, tested and evaluated. Other articles describes specific problems, but only in the context of classification (Ilias et al., 2007), regression (Gulden and Nese, 2013) or only in case of one specific algorithm. There are lots of articles focusing on a specific problem and only one specific algorithm to resolve that problem. For example, in the case of anomaly detection, there are a plenty of articles discussing different scenarios and comparing results of different algorithms. In (Zareapoor et al., 2012) the authors make a comprehensive evaluation of the most popular algorithms used in the context of credit card fraud detection. This article contains a brief description of algorithms like Bayesian Networks, Neural Networks, SVM, etc. and at the end of the article there is a table comparing the results, using metrics like accuracy, speed and cost. Other types of articles has a different approach, creating a survey of all the algorithms which can be used for a specific problem.

In (Varun et al., 2009) there is a survey of all the algorithms which can be used to effectively detect anomalies. This article doesn't narrow down its view to only supervised or only unsupervised algorithms. On the contrary, the authors describe an extensive list of algorithms, which can be used in a larger context, in anomaly detection. There are plenty of research papers comparing two or more different algorithms, as in (Juan et al., 2004) (Murad et al., 2013) (Agrawal and Agrawal, 2016) (Gupta et al., 2014).

The papers mentioned above were trying to help engineers in finding the best algorithm for specific problems, but none of them managed to create a stepby-step process, a useful methodology on how to choose the right algorithm for any type of problem and any type of dataset. In this paper we will propose a step-by-step methodology that can be applied to choose an algorithm. Moreover we will present our platform, which can be used in order to evaluate the chosen algorithms.

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3 THE PROPOSED METHODOLOGY FOR CHOOSING AI ALGORITHMS

There are no universally good or bad algorithms, each algorithm is specific to a context, a problem or a type of dataset. This idea was demonstrated by David H. Wolpert and William G. Macready in (Wolpert and Macready, 1997) in the so called "No Free Lunch Theorems for Optimization". The NFLT are a set of mathematical proofs and general framework that explores the connection between general-purpose algorithms that are considered "black-box" and the problems they solve. This states that any algorithm that searches for an optimal cost or fitness solution is not universally superior to any other algorithm. Wolpert and Macready wrote in their paper that "If an algorithm performs better than random search on some class of problems then it must perform worse than random search on the remaining problems."

In the real world, we need to decide on engineer-



Figure 1: Algorithm Selection Decision Tree.

ing solutions to build practical models that solve real problems. For certain pattern recognition problems, we tend to find that certain algorithms perform better than others which is a consequence of the algorithm's fitness or cost function fit to the particular problem. For our particular given problem, in order to find the best algorithm, NFLT should remind us that we need to focus on the particular problem at hand, the assumptions, the priors (extra information), the data and the cost.

What the NFLT is trying to tell us is, we are generally not going to find off the shelf algorithms that fits perfectly to our data. We have to architect the algorithm to better fit the data. This means that there are no universally applicable algorithms, so in order to find optimized solutions, we must have a methodology to architect our model. Figure 1 shows a decision tree to help engineers in choosing the right algorithm based on the problem context and the nature of the data.

Based on Figure 1 the first step in choosing the appropriate AI algorithm is to determine the nature of the problem, establish the goal, what do we want to obtain. Based on this, the algorithms can be grouped in three categories: Classification, Clustering, Prediction or Visualization (or Dimensionality Reduction). After defining the nature of the problem, the next step is to focus on the nature of the training data. In this case we should ask questions about the dimension of the available training data, whether the dataset has labels or not or questions about the importance of the features, can we transform the features to get more relevant ones or should we keep the original features. Answering these questions can spectacularly reduce the search space, in some cases resulting in only one

algorithm, this way making the developer's job much easier. These questions are very important, because for example, if we don't have enough examples in the training set (Less than 100K), then Neural Networks can have poor results and linear SVC can do a better job. As another example, if the features are important, so we should preserve the original values, then PCA is not a good choice, because this algorithm modifies the features, using orthogonal transformation, so a better choice would be Kernel Approximation or GA.

4 ARCHITECTURE OF OUR PLATFORM

The platform developed by our team is meant to offer a pragmatic tool for specialists involved in some kind of artificial intelligence related projects. This tool is not meant as a final solution but as a starting point in the process of finding the best method that fits the quality and efficiency criteria of a given application domain. Therefore the platform contains those basic functionalities that allows a specialist to collect, process with AI algorithms and visualize data. Although the platform was meant for general use, most of our experiments were performed in the area of anomaly and outlier detection in big datasets. Here are the functionalities we considered necessary for such a platform:

 Data harvesting tools - that allows acquisition of data from a variety of datasets having different formats (Excel, CSV, ARFF) or from different physical sources (sensor networks, smart devices, IoT); a special treatment is given to real-time data



Figure 2: Application Diagram.

collections and datasets containing anomalies;

- Data preprocessing tools instruments meant to transform the raw data into a noise free and normalized data; typical methods included in this category are: parameterized filters, transforms (e.g. FFT, wavelet) or histograms; there are also methods for determining some statistical parameters of the input data such as: min-max, median value, standard deviation, etc.
- Artificial Intelligence algorithms a wide group of methods that try to cover the most representative nodes of the presented taxonomy; our goal is to offer a rich set of possibilities from were to choose and compare; the open nature of the platform allows new methods to be added to the existing library; methods can be tested without programming knowledge, because the platform offers an easy to use user interface.
- Visualization tools very important in the process of finding the best artificial intelligence methods because they offer a bi-dimensional representation of otherwise multidimensional data, much easier to understand for the human eye;
- Generators of datasets and artificial signals necessary in the process of validating or measuring the quality of some new artificial intelligence methods; special techniques are applied in order to combine "normal" data with artificially created anomalies.

The tools mentioned above are integrated into an open architecture platform allowing continuous extension with new methods. A unique internal predefined data format assures interoperability and interchangeability between the existing tools and functionalities.

The high level overview of the included functionalities is shown in Figure 2.

To create a more complex model, using multiple preprocessing techniques and/or multiple artificial intelligence algorithms, the platform supports creation of a pipeline. Each step of the pipeline is saved in .csv files, so the results of each step can be used separately. This way the pipeline is highly configurable, it can have multiple steps, some steps can be skipped, reordered or reused. Figure 3 shows a high level view of the pipeline, where steps can be reconfigured, replaced, skipped or reordered.

The most important library used to create this platform is (SciKit-Learn, 2017). Similar tool is Weka (Weka, 1997)

5 EXPERIMENTS

The purpose of this section is to show through some examples the usage possibilities of the artificial intelligence platform. Here we emphasize two aspects: the possibility to choose between different anomaly detection techniques and the ability to tune the parameters of a given method in order to obtain better results.

5.1 Experiments using the Breast Cancer Dataset

The next experiments were made on the same dataset that reflects information related to breast cancer



Figure 3: Overview of the pipeline.

(Lichman, 2013). This dataset contains 569 data points and each data point has 32 attributes. The dataset was labeled showing normal (e.g. benign) and abnormal (e.g. malign) data points. A part of the dataset was used in the training (learning) phase and the rest in the testing phase.

The next figures show how our platform can be used by someone that has no programming knowledge, only by choosing the options in the graphical user interface.

The first step is to load our dataset. At the beginning we select a given AI algorithm that can classify data in benign and malign samples. Our first choice is the SVM (support vector machine) algorithm. In the next step we will configure and train our SVM model using the loaded dataset, without preprocessing of the data. In the case of the SVM, we can configure the C parameter, which stands for penalty parameter of the error term, the value of gamma, which is the kernel coefficient for rbf, poly and sigmoid and the kernel type to be used in the algorithm. It must be one of linear, poly, rbf or sigmoid. Next step of the process is to evaluate the trained model. The result of the evaluation is shown in Figure 4. As we can see in this table, the scores of this model are not very promising, which is caused by the fact that no preprocessing was made.

In the last step we apply Min-Max scaler as a preprocessing algorithm and the configuration of the SVM model remains the same. From Figure 4 we can clearly see the effect of applying the correct preprocessing algorithm, the increase in the model's scores are spectacular. This demonstrates the importance of understanding our training dataset and the importance of using proper algorithms not just for classification, but also for preprocessing.

This process can be repeated using other types of algorithms, distance based, density based, hierarchical or even using neural networks. The following figures shows the usage of other algorithms.

In the case of KNN you can choose the number of neighbor nodes, the algorithm used to compute the nearest neighbors, the weight function used in prediction and the distance metric to use for the tree. The default metric is Minkowski, and with power=2 is equivalent to the standard Euclidean metric. Figure 4 shows the results in a table.

We can use the platform to configure more difficult models, like Neural Networks. Configuring the Neural Network model means that you can set the Alpha value or the L2 penalty (regularization term) parameter, set the solver for weight optimization (algorithm), choose the correct activation function, set the number of the hidden layer and the number of the nodes for the hidden layers. Figure 4 shows the results of the model as a confusion matrix.

In other cases, if you are not interested in the confusion matrix, but you want to run the training and the evaluation multiple times to see how does the algorithm behave, you can do it easily, using only the UI, with no programming knowledge.

To demonstrate this functionality, we ran the training and evaluation process 14 times. The accuracy diagram for this training and evaluation loop can be seen in Figure 5

To configure the Random Forest algorithm to be

Use Case	Accuracy	Precision	Recall	F1-Score
SVM without preprocessing	0.62	0.62	0.50	0.55
SVM using Min-Max Scaler as preprocessing algorithm	0.96	0.97	0.96	0.97
KNN	0.93	0.91	0.97	0.94
NN	0.93	0.91	0.98	0.95

Figure 4: Evaluation of the different algorithms.



Figure 5: Evaluation of the Random Forest model in a loop.



Figure 6: Scatter Plot.

evaluated it in a loop you can set the maximum number of evaluators (decision trees), choose the slicing method, configure a cutoff depth if needed and define the step size (the number of estimators will be increased with the step size in each step). Figure 5 contains a digram which shows the accuracy of the model in each step. This is very useful when you already know which algorithm do you want to use and you want to find the most optimal parameters to train the model with.

On the visualization part, we can see our data in a scatter plot (Figure 6), as a histogram (Figure 7) or as







time series (Figure 8), depending on the nature of the data.

As future development we aim to add some algorithms for visualization and dimensionality reduction (like PCA) and some preprocessing algorithms to cover most of the techniques and algorithms that a general purpose platform should offer.

5.2 Using the Proposed Methodology for Choosing Algorithms in the Context of Anomaly Detection

Anomaly detection finds data points that do not fit well with the rest of the data. It has a wide range of applications such as fraud detection, surveillance, diagnosis, data cleanup, and predictive maintenance.

Lets consider Thyroid Disease dataset, which has 3772 training instances and 3428 testing instances. It has 15 categorical and 6 real attributes. The problem is to determine whether a patient referred to the clinic is hypothyroid. Therefore three classes are built: normal (not hypothyroid), hyper function and subnormal functioning. For outlier detection, 3772 training instances are used, with only 6 real attributes. The hyper function class is treated as outlier class and other two classes are inliers, because hyper function is a clear minority class.

To find the appropriate algorithm for this particular dataset, the first question is about the nature of the problem. Because it is clear that we have a classification problem (classify examples as outliers or inliers), we can reduce the search space and go to the next step, to analyze the nature of the data. As we specified in the description of the thyroid disease dataset, we have labeled data, so we can step further and analyze the dimension of our dataset. The dataset contains 3772 instances, which is less than 100K, so we can try out the SVM algorithm.

After we found the appropriate algorithm, the next step is to use our platform to minimize the error and to maximize the precision, recall, accuracy and other evaluation results. We run the algorithm multiple times. Firstly we used C=1 and Gamma=0.1 with rbf kernel. Figure 9 shows the results of the experiment. The results are not too promising, because having a really small recall much of the outliers are omitted.

In the next step we increased both C and Gamma, which gave us a much better result, as we can see in Figure 10. From 0.08 we increased the recall to 0.6. Even if the precision has dropped, the overall result is much higher, the model behaves much better.

As we keep increasing the values of C and Gamma, we are getting better and better results until we reach a threshold, when the model starts to overfit. The best results we obtained using C=100 and Gamma=10 (see Figure 11)

In this case we still have a choice to make. Are the results good enough for our problem or should we test other algorithms too. If the answer is yes, then we can stop and use the SVM to solve our problem. If the answer is no, then we can still choose from plenty of algorithms. As we can see in the algorithm selection



Figure 9: SVM with C=1 and Gamma=0.1.



Figure 10: SVM with C=10 and Gamma=1.

tree (Figure 1) the next question is whether our data is text or not. In our case we don't have to deal with text, so we can choose between KNeighbors, Random Forest or we can use Ensemble Methods.

With this example for using our algorithm selection tree, we demonstrated the benefits of our algorithm selection methodology and the usefulness of our platform in finding the correct parameters for the chosen algorithm. This experiment clearly shows how can we reduce the search space when a developer have to choose between algorithms, this way increasing productivity and velocity of the developers and implicitly decreasing the time to production and also the costs of the project.



Figure 11: SVM with C=100 and Gamma=10.

6 CONCLUSIONS

The methodology and decision tree for choosing the proper algorithm based on the problem context and the nature of the data tries to show a coherent way allowing a developer to select the best methods that fits the requirements of a given application. This taxonomy was used as a theoretical background for the development of a platform that incorporate different artificial intelligence and data preprocessing techniques.

The platform implements the main functionalities needed by a developer in the process of finding the best strategy for a given domain that assure high quality problem solving in a reasonable execution time, increasing the productivity of the developers, increasing the probability of success and decreasing the costs and risk. It includes multiple data acquisition, preprocessing, anomaly detection and visualization functionalities that may be combined in a specific processing flow.

Every individual functionality is implemented as an autonomous service and multiple services may be orchestrated in a logical flow in order to obtain the final results. The richest group of functionalities is the one that contains the artificial intelligence methods. Here we tried to cover most of the development directions present today in the scientific literature, from statistical methods, towards, signal processing and neural networks. In this way the platform can be used as a training tool for those who must develop different solutions in different areas. For the same learning purposes we included a multitude of datasets that cover a diversity of application domains. The experimental part of the paper demonstrates some of the functionalities of the platform, including the possibility to compare the result obtained with different methods, to run the training and evaluation process multiple times and to visualize the data and the results in a meaningful way.

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