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Log Data Preparation for Predicting Critical Errors Occurrences

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Abstract. Failure anticipation is one of the key industrial research objectives with the advent of Industry 4.0. This paper presents an approach to predict high importance errors using log data emitted by machine tools. It uses the concept of bag to summarize events provided by remote machines, available within log files. The idea of bag is inspired by the Multiple Instance Learning paradigm. However, our proposal follows a different strategy to label bags, that we wanted as simple as possible. Three main setting parameters are defined to build the training set allowing the model to fine-tune the trade-off between early warning, historic informativeness and forecast accuracy. The effectiveness of the approach is demonstrated using a real industrial application where critical errors can be predicted up to seven days in advance thanks to a classification model.

Keywords: Predictive Maintenance (PdM), Machine Learning (ML), Classification, Log data, data preparation

1 Introduction

Predictive maintenance is of paramount importance for manufacturers ([6, 10]). Predicting failures allows anticipating machine maintenance and thus enabling financial saving thanks to shortening the off-period of a machine and anticipating the spare parts orders.

Nowadays, most modern machines in the manufacture domain are equipped with sensors which measure various physical properties such as oil pressure or coolant temperature. Any signal from these sensors is subsequently processed later in order to identify over time the indicators of abnormal operation ([12]). Once cleaned, historical data from sensors can be extremely valuable for prediction. Indeed, their richness makes it possible to visualize the state of the system over time in the form of a remaining useful life estimation ([5]).

In parallel, machines regularly deliver logs recording different events that enable tracking the conditions of use and abnormalities. Hence, event-based prediction is a key research topic in predictive maintenance ([13, 4]).

Our contribution in this paper is a data preparation approach whose output is used to learn a model for predicting errors/failures occurrences. It relies on the exploitation of historical log data associated to machines. The purpose is to predict sufficiently early the occurrence of a critical malfunction in the purpose of maintenance operations facilitation. The proposed framework is applied on a real industrial context and the empirical obtained performances show its effectiveness.

After introducing in the next section key definitions and parameters, we detail the data preparation process. Afterwards, we report on some of the experiments conducted in the context of our case study. Thereafter, we provide an overview of related work. Finally, we conclude and outline some indications for future work.

2 Dataset Collection and Preliminary Definitions

We suppose that we are given a log file \mathcal{F} emitted by a machine \mathcal{M} where the errors that have been observed by sensors associated to \mathcal{M} are reported. We assume that the errors are logged in regular time intervals. Moreover, we consider a set of errors $\mathcal{E} = \mathcal{L} \cup \mathcal{H}$ where \mathcal{L} is the set of low errors ℓ_j and \mathcal{H} is the set of critical errors h_j (high errors that will be called target in the following). Our aim is to predict high errors wrt other errors. Each record in \mathcal{F} is a pair $\langle i, \mathcal{E} \rangle$ where i is a timestamp expressed in days and $E \in \mathbb{N}^{|\mathcal{E}|}$ is a vector where E[j] represents the number of occurrences of error $e_j \in \mathcal{E}$ at timestamp i. Hence, \mathcal{F} is a temporal ordered sequence of such records. The following is the running example we use throughout the paper as an illustration of the approach.

Example 1. Let the errors set be $\mathcal{E} = \{\ell_1, \ell_2, \ell_3, \ell_4\} \cup \{h_1, h_2\}$ and consider the following sequence (Table 1):

Timestamp	ℓ_1	ℓ_2	ℓ_3	ℓ_4	h_1	h_2
1	0	12	6	1	0	0
2	0	0	3	2	0	0
3	0	1	4	1	1	1
4	1	0	1	2	0	0
5	0	1	1	2	0	0
6	0	1	1	1	0	0
7	0	1	1	0	0	1
8	1	0	1	8	0	1
9	0	0	6	1	1	0
10	1	0	7	1	1	0

Table 1: Example of a log event dataset with timestamp

For example, the first row/record whose timestamp is 1 says, among others, that at day 1, there have been 12 occurrences of low error ℓ_2 and no occurrence of high error h_1 .

To come up with a prediction model and collect the data, three parameters, depicted in Figure 1, are applied:

- Predictive Interval: it describes the history of data used to perform predictions and, its size in days is defined by the parameter PI. The information contained in that interval is gathered in a structure referred as a bag.
- Responsive Interval: since, from a practical point of view, making a prediction for the next day is of little interest, a common strategy is to constrain the model to forecast sufficiently early. The Responsive Interval is located immediately after the Predictive Interval and its size which controls how much early we want the model to forecast, is defined by the parameter RI.
- Error Interval: intuitively, this interval, whose size is defined by parameter
 EI, allows to make error occurrence prediction wrt a time interval instead of
 a time unit. It therefore introduces a degree of uncertainty of the prediction's
 temporality.



Fig. 1: The key parameters of model prediction

Example 2. Let PI = 3, RI = 4 and EI = 2. With this setting, we want to predict high errors wrt bags of size 3, and this prediction concerns a period whose duration is 2 days starting 4 days after the end of the observed bag. More precisely, the fact that EI = 2 means that a predicted error occurrence may happen either at the first, the second, or both days within the error interval. Consider the bag B_1 defined between timestamps i = 1 and i = 3 in the Table 1. Then the prediction period (Error Interval) concerned by B_1 is in timestamp [8, 9]. That is, at i = 3, we observe the bag B_1 and we should be able to say whether, or not, a high error will occur between i = 8 and i = 9.

3 Proposed Methodology

The current section describes how training and test sets are prepared. Indeed, in order to train a binary predictive model, a prerequisite is to build a training set where each sample is labeled YES/NO and corresponds to a situation where, respectively, a high error occurred or not.

Definition 1. Let B_i be a bag and given a target high error h_j . B_i is labeled YES iff time interval [i+PI+RI;i+PI+RI+EI-1] contains an occurrence of h_j . B_i is labeled NO otherwise.

Example 3. Let PI=3, RI=2 and EI=2. Consider again the running example and let us focus on error h_1 . To decide the label of bag B_1 , we need to check whether in [6; 7] error h_1 has been observed. Since target has not been observed, then B_1 takes label NO. Hence, referring to Table 1 and applying the definition above, the bags B_1 and B_2 are labeled as NO and the bags B_3 and B_4 are labeled as YES. Observe that since their corresponding EI period is outside the bounds of the available history, bags 5, 6, 7 and 8 cannot be labeled.

3.1 Summarization Strategy

The reader may notice that the above configuration (labeled bags) is close to the one encountered in the Multiple Instance Learning (MIL) [3]. However, our setting cannot be in conformance with MIL because the objective is a prediction based on the whole bag rather than individual samples. Thus, we choose an alternative approach, by summarizing every bag B_i by maximizing values and transform it into a single features vector³.

Example 4. Let PI = 3 and consider the bag B_1 starting at i = 1 and ending at i = 3. The bag is summarized by the vector (0, 12, 6, 2). That is, for each ℓ_j we take the maximum of its values in B_1 . So, with RI = 2 and EI = 2 and when considering the target error h_1 we obtain the data set presented in Figure 2.

Tim	estamp	ℓ_1	ℓ_2	ℓ_3	ℓ_4	h_1	h_2
	1	0	12	6	1	0	0
	2	0	0	3	2	0	0
	3	0	1	4	1	1	1
	4	1	0	1	2	0	0
	5	0	1	1	2	0	0
	6	0	1	1	1	0	0
	7	0	1	1	0	0	1
	8	1	0	1	8	0	1
	9	0	0	6	1	1	0
	10	1	0	7	1	1	0

Bag	ℓ_1	ℓ_2	ℓ_3	ℓ_4	h_2	Label h_1
B_1	0	12	6	2	1	NO
B_2	1	1	4	2	1	NO
B_3	1	1	4	2	1	YES
B_4	1	1	1	2	0	YES

Fig. 2: Bag summarization strategy by MAX() function

Each sample in the right pane of Figure 2 on one hand, is the total number of occurrences of every ℓ_j during a period of size PI=3. The last columns testify whether h_1 has occurred (YES) or not (NO) after this period at a precise time interval. The table on the left pane reminds the original data set.

³ Other methods of summarization can be used but in our use case, the MAX() function allowed us to obtain good prediction performances.

3.2 Overall Data Preparation Process

Figure 3 and Algorithm 1 depict our preprocessing stage. From a log file T, the consecutive records are gathered into bags using a sliding window of size PI. The label of each bag B_i depends on the content of interval EI_i . The starting point of EI_i is RI time units after the end point of B_i . Once the bags are identified and labeled, they are summarized by MAX() function to obtain the input data set T' for model learning. As it may be seen from Algorithm 1, the complexity of data preparation is linear with the size of the sequence: the outer most loop is executed O(|T|) times. At each iteration, PI data rows are summarized and the label is assigned after checking EI row records. Hence, the overall complexity is $O(|T| \times (PI + EI))$. From a computing point of view, the outer loop can be easily made parallel because the different iterations are independent of each others.

```
Algorithm 1: DATAPREP

Input: Sequence T, EI, PI, RI, Error h
Output: Table T' of labeled meta-instances.

begin

for i = PI + RI + 1 to |T| - EI + 1 do

s \leftarrow \text{Summarize}(T, i - RI - PI, i - RI - 1, \mathcal{S})

//A bag of size PI starting at index i - RI - PI is summarized using MAX() or another function \mathcal{S}

s.label \leftarrow False

for j = i to i + EI - 1 do

if T[j] contains an occurrence of h then

s.label \leftarrow True

Insert s into T'

return T'
```

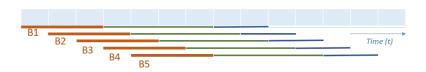


Fig. 3: Data Preparation Process. The bags of the training set are built by sliding windows with the following model parameters: PI=3; RI=4; EI=2

4 Experiments

4.1 Dataset Description

We collected raw data from a one-year history of several log files. Each file is associated to a specific machine. All machines belong to the same model⁴. However, they are not subject to the same conditions of use. Given a combination of parameters (PI;RI;EI), each log file is processed using Algorithm 1. Output files are then merged to constitute our input dataset. The latter is partitioned into a 80/20 ratio of training and test sets respectively. The partitioning relies on a stratified sampling to keep a similar proportion of positive samples both in the test and training sets.

The dataset contains 193 features (i.e., distinct low errors) and 26 high errors (our targets). For every high error, we designed several models, each of which is obtained by combination of three parameter values (PI;RI;EI). As part of our study, we were interested in the 11 most frequent target errors which occur in the dataset. Depending on the target, we ended up with an average of 1000 samples, among which, from 4% to 30% are positive. Although a balance correction technique is in general recommended in such a case, our experiments showed that subsampling did not provide any substantial accurracy gain⁵. Then, we decided not to make use of subsampling.

In the following experiments the default value for both parameters PI and RI is 7 days. The value of the RI parameter is chosen as constraint because the idea is that the model anticipates 7 days before the error target occurs.

4.2 Settings

All experiments were conducted with the RapidMiner software⁶. Several classifier algorithms were explored, e.g., SVM, Naive Bayes, Random Forest, Decision Tree and the multi-layer feed-forward artificial neural network (NN) algorithm of the H2o framework [1]. NN algorithm outperformed the others and was subsequentely chosen as the actual predictive model. Experiments were set with 4 layers of 100 neurons each, 1000 epochs and a Cross Entropy loss function for the learning phase.

To assess the performance of a prediction model, different metrics can be used. Among these, Accuracy and F1-score are suitable one. Accuracy is the ratio of correct predictions. F1-score wrt to the positive class is given by the harmonic mean of precision and recall⁷:

$$\frac{2(precision \times recall)}{(precision + recall)} = \frac{2 \times TP}{2 \times TP + FP + FN} \tag{1}$$

 $^{^4}$ Regretfully, for confidentiality reasons we cannot share the data used for this study

⁵ Due to lack of space the results will not be detailed in this paper.

 $^{^6}$ Rapidminer.com

⁷ precision= $\frac{TP}{TP+FP}$ and recall= $\frac{TP}{TP+FN}$.

TP, TN, FP, FN represent respectively, true positives, true negatives, false positives and false negatives. Because negative cases are much more frequent than positive one, accuracy alone is not sufficient to assess the performance of the model. F1-score allows to observe in more details the behavior of the model regarding the class of interest, in our case, the positive one.

4.3 Impact of the Error Interval Choice

Parameter EI introduces a flexibility wrt positive predictions as well as a rigidity for negative ones. Indeed, a positive prediction says that $at\ some\ point$ in a future time interval, a high error may occur while a negative prediction says, no occurrence $at\ any$ point in a future time interval. Intuitively and regarding positive predictions, one may expect that the higher EI, the more correct the model's predictions are. To analyse this hypothesis we varied EI from 1 to 4 days and for each target error, we computed F1-score wrt the positive class⁸ The obtained results are depicted in Figure 4.

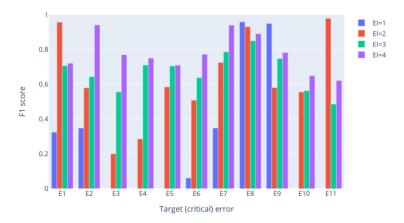


Fig. 4: F1-score evolution wrt EI

Overall, our initial assumption is confirmed. We observe that for about half of the errors, setting EI to 1 day leads to a significant degradation of the model's performance. In these cases the model yields to F1 score equal to zero (see errors E3; E4; E5; E10 and E11) explained by a null value of true positive prediction. For E8 and E9, the results show a different phenomenon from overall behaviour. Indeed, the largest F1 score for these errors is obtained when EI is set to 1.

F1-score gives a *global* picture of the predictive model behavior wrt positive class in that it combines precision and recall measures. From the application perspective, false alarms, i.e., false positives ratio (FPR) is a key indicator⁹.

 $^{^{8}}$ Due to space limitation, we do not present here the prediction for negative class.

⁹ $FPR = \frac{FP}{FP+TP} = 1 - precision$

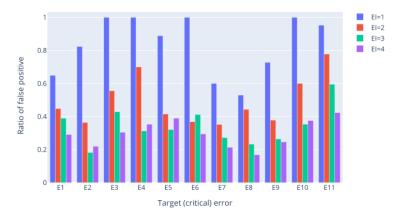


Fig. 5: False positive (false alarm) ratio wrt EI

Figure 5 shows the evolution of FPR with respect to EI. We observe that in general, increasing EI tends to decrease FPR, hence it increases the precision.

Increasing the size of the EI concomitantly increases the number of positives examples in the training dataset. That can explain that we obtained better performances with a greater EI. Although these performances may also in some cases mask an over-expression of false positives, the phenomenom is not present here, as proven in Figure 5.

Despite the general trend of F1-score and precision wrt EI, the model performance is not monotonic. For example, the learned model for error E1 has the highest F1-score when EI=2. This indicates that for every error, one needs to calibrate the most convenient EI regarding applications/F1-score needs.

5 Related Work

Fault anticipation for predictive maintenance has attracted a large body of research (please see e.g., recent surveys [9, 8, 14]) and several paradigms have been used for this purpose. Our solution belongs to the class of approaches based on supervised machine learning. As such, we focus our comparison to these techniques and more specifically on [11] and [7] that are, to our knowledge, the closest previous works.

Let us start with [11]. Despite some subtle differences, the task of tagging the bags as negative or positive instances is similar to that of [11]. However, in the latter, negative bags propagate their label to each individual instance belonging to them. By contrast, every positive bag is aggregated using the average measure and gives rise to a new positive meta-instance. Therefore, the training data set is a mix of meta-instances (positive in this case) and simple instances (negative ones). The rational behind the use of the average score is that the so obtained instance is considered as a representative of a daily behavior inside a bag. In fact, and unless having a small standard deviation inside positive bags, a daily

record can be very different from the average of the bag it belongs to. Hence learning from bag averages and predicting from single instances may present some discrepancy. The data for our case study present high deviations inside the bags, which makes it hopeless to identify positive records using averaged records. In our case, the data preparation process considers positive and negative bags equally. In that way, consider the occurrence or not of a high error/failure is explained by the whole bag associated to it, and not just some instances belonging to that bag. Moreover, in contrast to what has been suggested in the work of [11], this way of proceeding reduces the issue of unbalanced classes that we are confronted with.

Now let us consider the prediction phase. In [11], prediction is instance based (i.e., daily records) and these predictions are then propagated to bags. More precisely, during the prediction phase, successive instances belonging to the unclassified bag, are provided to the learned model. If one of them is classified as positive then the whole bag is positive. It is negative otherwise¹⁰. So fundamentally, predictions are daily based in the sense that the real explanation of a possible error/failure occurrence depends actually on what happens during a day while ours needs are to combine events of larger intervals. This is also the reason why, unlike [11], our work is not in conformance with Multi-Instance Learning paradigm (MIL). Indeed, the MIL main hypothesis states that a bag is positive iff one of its instances is positive [3, 2], while our approach does not impose any condition on individual instances. We have applied the methodology of [11] over our use case, but could not obtain F1-score better than 0.2.

A more recent work with a similar data preparation process as [11] is [7]. The main difference is that its learned model is a regression that estimates the probability of a future error occurrence. If this probability is higher than some fixed threshold, then an alert is fired. Hence, during data preparation, the labels associated to bags are probabilities instead of YES/NO labels. Intuitively, the closer a bag is to an error, the higher is its probability. To handle this property, the authors used a sigmoid function. The reported results show that the proposed techniques outperform those of [11] by a large margin. We believe that this is mainly due to the data nature defining the use case and in general there is no absolute winner. As it has been shown in the experiments, our targeted use case does not need such complex techniques to achieve high prediction accuracy.

6 Conclusion and Future Works

We described an approach to exploit log data for predicting critical errors that may cause costly failures in the industry manufacture domain. The approach is based on aggregation of temporal intervals that precede these errors. Combined with a neural network, our solution turns to be highly accurate. Its main advantage, compared to other techniques, is its simplicity. Even if we do not claim that

Notice that in [11], during data preparation, bags are first labeled and then, their labels are propagated to instances while for bag label prediction, the labels of its instances are first predicted and then the bag label is deduced.

it should work for every similar setting (log based prediction), we argue that it could be considered as a baseline before trying more sophisticated options. We applied our solution to a real industrial use case where a dozen of highly critical errors were to be predicted. The obtained accuracy and F1-score belong to respectively [0.8; 0.95], and [0.6; 0.9] which can be considered as highly effective compared to the reported results on log based failures prediction litterature.

So far, we designed one model per target error. In the future, we plan to analyze in more depth the errors to see whether it would be possible to combine different predictions in order to reduce the number of models. In addition, we are interested in automating the parameters setting: given a (set of) target(s), to find the optimal values of PI, EI and RI such that the learned model performances is maximized.

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