Handling Distribution Shift in Tire Design

Antoine de Mathelin^{1,2} Francois Deheeger¹ Mathilde Mougeot^{2,3} Nicolas Vayatis² ¹Manufacture Française des Pneumatiques Michelin, Clermont-Ferrand, France ²Université Paris-Saclay, CNRS, ENS Paris-Saclay, Centre Borelli, Gif-sur-Yvette, France ³ ENSIIE, Évry-Courcouronnes, France antoine.de-mathelin-de-papigny@michelin.com

Abstract

The recent success of machine learning methods in the industrial sector open new perspectives for the design of innovative products. However, these promising results are often challenged when it comes to industrial model deployment. Indeed, it frequently appears that the performance of the model is degraded when used on application data due to the distribution shift between the training and the targeted data. This issue is even more critical for model dedicated to the research of innovative designs as the model is mainly used on unseen regions of the design space. In this work, we present, on a real application of tire design, how distribution shifts impact the model performance and what can be expected from several domain adaptation methods. In an objective of industrial model deployment, we conduct this benchmark with the use of unsupervised evaluation metrics that considerably help the model selection.

1 Introduction

The design of complex industrial products such as tires requires a time consuming and expensive process. Designers need to explore for multiple design choices to find novel products with desired properties. In this context, the use of machine learning models able to learn the relationship between the design of a product and its properties can substantively improve the process [25].

Recent works in design space exploration show that machine learning methods can efficiently learn the mapping between a product components and its properties [3, 11, 12, 16]. However, it has been observed that the learned models are somehow limited to the domain defined by the training data and provide degraded performances on unseen regions of the design space [13]. These observations have been confirmed in our experiments (cf Section 2.1). This is an important drawback as designers aim to explore for innovative designs to find new products with competitive properties.

To tackle this issue, one solution is to correct the shift between the training distribution and the distribution on which the model is applied. To this end, the learner can use domain adaptation methods which either transform the feature space to find more robust representations [8, 19, 17, 26] or reweight the training instances to correct the sample bias [5, 7, 10, 14, 21]. Both strategies have been applied for design space exploration with promising results [1, 13, 15]. However, many questions still need to be addressed before deploying such methods. Because of the huge number of existing methods, practitioners are in particular wondering how to make a relevant choice between them and how to select the right hyper-parameters? They are also wondering what benefit a few labels can bring in the target domain?

In this paper we tackle these questions by studying on a real scenario of technological drift in tire design, the impact of several domain adaptation methods on the target score. We also explore the influence of several unsupervised metrics on the choice of method and hyper-parameters. The organization of the paper is as follows: first we present the issue of distribution shift in tire design and

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the impact of this shift on the model performances. We then present the metrics that are considered to make the model selection. We finally present the results obtained for real tire design data with the main domain adaptation methods provided in the ADAPT¹ library [6].



Figure 1: Domain adaptation for tire design: to correct the shift that exists between already deployed tires and potential innovative designs, instance-based or feature-based strategies can be considered.

2 Motivation and Evaluation metrics

2.1 Motivation

We consider the problem of tire performance prediction where the learner has access to a set of *source* labeled tires data $S = \{(x_1, y_1), ..., (x_m, y_m)\} \in \mathbb{R}^p \times \mathbb{R}$ where each x_i is a vector of size p corresponding to the features of one particular tire's design (i.e. the height and width of the different rubber layers, the materials properties etc...) and y_i is one recorded performance of interest. S is referred as the source data set which contains the design of already deployed tires. The learner has also access to an unlabeled set of tires $\mathcal{T} = \{x'_1, ..., x'_n\} \in \mathbb{R}^p$ which corresponds to possible new tire's designs. The goal is to find the design x' in \mathcal{T} with the best corresponding performance y'.

The performance y' corresponding to each $x' \in \mathcal{T}$ can be approximated by using a machine learning model $f : \mathbb{R}^p \to \mathbb{R}$. The standard approach consists to train f using only the data set S. We perform this experiment on a set of real tire designs. We first split S in a train and test sets and fit f on the first set. We then compute the prediction errors on the source train and test sets as well as on the target set \mathcal{T} . The results are reported on Figure 2.B and 2.C, we can observe that the errors provided by f on the target set are far above the validation error computed on the source test set. We observe, besides, on Figure 2.A, a clear distribution shift in the input space between S and \mathcal{T} . This shift certainly causes the degraded performances of the model on the target set.

Facing this issue, one solution would be to use a domain adaptation method to correct the shift in the input space. However, as in practice no or only a few labels are available for the data from \mathcal{T} , finding the right method to use is not straightforward. Furthermore, many domain adaptation methods are very sensitive to negative transfer [20] and require a careful model selection. For these reasons, we introduce, in the next section, evaluations metrics that will be considered to help the model selection.

2.2 Evaluation Metrics

When dealing with a domain adaptation problem, two main scenarios can be considered: the supervised and unsupervised domain adaptation. In the first one, the learner has access to a small

¹https://adapt-python.github.io/adapt/



Figure 2: From the left to the right: t-SNE visualization [24] of the input space, validation plot and distribution of the errors for a model trained on source data only.

sample of labeled target data and can then estimate a target score. In the second scenario, no label are available on the target domain. Thus, to help selecting an adequate model, one can only consider unsupervised metrics. In this work, we consider the metrics below, some of them have already been used for hyper-parameter selection as the J-score and the Reverse Validation, others are measures of divergence between distributions:

- **J-score** (js): is proposed in the method KLIEP [21] for selecting the bandwidth of the kernel, it is correlated to the KL divergence between source and target distributions.
- **Domain Classifier (dc)**: characterizes the shift between domains by the ability of a classifier to discriminate between data from different domains [2, 8].
- Frechet Distance (fd): is inspired from the FID score [9] used for GANs to compare the real distribution to the generaed distribution.
- Linear Discrepancy (ls): is one of the first unsupervised metric used for domain adaptation, it can be computed for linear classifiers using linear algebra [14].
- **Reverse Validation** (rev): is computed as a source error by inversing the role of the source and target domains, this requires pseudo target labels which are computed using the domain adaptation model [27, 8]

According to [8] a good feature representation is where the source and target distributions are matching and the source task can be learned. Considering only the source error for model selection is not enough as we have shown in Section 2.1 but considering only unsupervised metric may also be inefficient. Indeed, we could produce a representation where all data are regrouped in one localization, thus, the two distributions will match but it will not be possible to learn the task anymore. Thus, to take into account this two objectives, we would like to consider, for model evaluation, a combination of the source error relatively to the metric is difficult. We then propose the following evaluation process: first training several models using the domain adaptation method and then evaluate the source validation error and the unsupervised metric for each run. The source errors and the metrics are then rescaled with standard scaling. An evaluation score is then given by summing them together. We thus give the same importance to the learning of the source task and the matching of distributions.

3 Experiments

3.1 Setup

We perform the experiments on a tire design data set composed of around 7500 tire design vectors of size 470 describing the features of the tire designs. In this data set, we select all data belonging to a same tire line (around 1500 data) which act as the set of potential new tire's designs \mathcal{T} . The remaining 6000 data define the set of previous tire's designs \mathcal{S} .

We consider the following domain adaptation methods available in the ADAPT library [6]: the two unsupervised instance-based methods KMM [10] and KLIEP [21] and the four unsupervised feature-based methods CORAL [22], DeepCORAL [23], DANN [8] and mSDA [4]. We also consider the supervised method TrAdaBoostR2 [18].

We perform model selection using the metrics describe in Section 2.2. For each domain adaptation method, several models are trained with different set of hyper-parameters and random seed. For each trained model the evaluation metrics are recorded along with a source validation error. The evaluation scores described in Section 2.2 are then computed. We finally select the model which returns the smallest score. The list of hyper-parameters considered for each method as well as the networks architecture and optimization parameters are presented in Appendix A.

3.2 Results and Discussion

The results for the unsupervised experiments are reported in Table 1. We compare the target MAE obtained using the selected model according to each of the unsupervised metrics. The average target MAE (avg tgt) over all trained models is reported along with the MAE of the best model (min tgt).

Table 1: Target MAE across different methods and evaluation metrics. The values below "src+rev" correspond to the target MAEs of the models selected according to the scores computed with the combination of the "rev" metric and the source validation error of each method (cf Section 2.2).

method	src+rev	src+fd	src+ld	src+js	src+dc	avg tgt	min tgt
Src-Only	0.243	0.243	0.243	0.243	0.243	0.243	0.243
KMM	0.236	0.239	0.224	0.236	0.233	0.235	0.217
KLIEP	0.223	0.244	0.228	0.633	0.228	0.350	0.223
mSDA	0.349	0.451	0.410	0.559	0.684	0.486	0.296
CORAL	0.254	0.314	0.232	0.266	0.229	0.344	0.221
DeepCORAL	0.577	0.629	0.606	0.539	0.608	0.570	0.406
DANN	0.228	0.331	0.331	0.276	0.642	0.391	0.222

Table 1 first shows that, for any metric, the MAE obtained by selecting the model with the smallest score is most of the time below the average target error (avg tgt). This means that using an unsupervised metric to select the model is in general improving the performances one can expect from domain adaptation against using a model with arbitrary parameters. We observe in particular that the two metrics Reverse Validation (rev) and Linear Discrepancy (ld) help to select better model than other metrics on our tire data set. We finally observe that the two instance-based methods KMM and KLIEP provide the best scores.

We then visualize the impact of adding labels in the target domain. We now select the score according to a validation error obtained on 20% of the labeled target, the 80% others are used during training along with sources. In this category, TrAdaBoostR2 specifically exploits the labeled target data. In this experiment, this last method improves substantively the model performances (cf Table 2).

Table 2:	Target N	MAEs of	the	models	selected	with	target	validation	scores

method	5 tgt	10 tgt	15 tgt	20 tgt	30 tgt	50 tgt
No Adapt	0.226	0.212	0.223	0.215	0.219	0.212
KMM	0.226	0.255	0.230	0.197	0.204	0.191
KLIEP	0.229	0.241	0.246	0.219	0.208	0.183
TrAdaBoostR2	0.216	0.188	0.198	0.176	0.191	0.176

4 Conclusion

In this work, we address the challenge of distribution shift in tire design. We show, in particular, that instance-based domain adaptation methods are useful to improve the model performances in both supervised and unsupervised scenarios. We conduct our work in an industrial model deployment perspective and propose, for this purpose, an unsupervised evaluation process to make accurate

model selection. We are convinced that our study will give the practitioners more confidence in using domain adaptation for industrial design.

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Appendices

A Architectures and hyper-parameters

Layer type	Layer size	Activation
Fully-connected	100	ReLU
Fully-connected	100	ReLU
Fully-connected	100	ReLU
Fully-connected	10	ReLU
Fully-connected	1	Linear

Table 4: Architecture for the networks used in DANN and DeepCORAL

Network	Layer type	Layer size	Activation
Encoder	Fully-connected	100	ReLU
	Fully-connected	10	ReLU
Task	Fully-connected	100	ReLU
	Fully-connected	10	ReLU
	Fully-connected	1	Linear
Discriminator	Fully-connected	100	ReLU
	Fully-connected	10	ReLU
	Fully-connected	1	Sigmoid

Table 5: Architecture for the networks used in mSDA

Network	Layer type	Layer size	Activation
Encoder	Fully-connected	100	ReLU
	Fully-connected	100	ReLU
	Fully-connected	latent size	ReLU
Decoder	Fully-connected	100	ReLU
	Fully-connected	100	ReLU
	Fully-connected	470	Linear

Table 6: Classifier used for Domain Classifer metric

Layer type	Layer size	Activation
Fully-connected	10	ReLU
Fully-connected	10	ReLU
Fully-connected	1	Sigmoid

Table 7: Hyper-parameters

Method	Hyper-parameters
КММ	$ sigma = \{2^i\}_{i \in [1-5,5]}$
KLIEP	max centers = $\{100, 200, 300, 500\}$ (sigma is selected via the LCV procedure)
mSDA	latent size = $\{10, 20, 50, 100\}$, noise level = $\{2^i\}_{i \in [[-5, 5]]}$
CORAL	lambda = $\{2^i\}_{i \in [-10,10]}$
DeepCORAL	$ \text{lambda} = \{2^i\}_{i \in [-10,10]}$
DANN	$ \text{ lambda} = \{2^i\}_{i \in [-5,5]}$



B Visualization of the true target error in function of the unsupervised scores

Figure 3: Blue dots show the score evaluated for a model with a particular set of hyper-parameters and random seed. The red dot in each figure is the model selected, i.e. the model with the smallest score. We can see that this model has, in general, a target error below the average target error.