

# Journal of Statistical Software

November 2024, Volume 111, Issue 10.

doi: 10.18637/jss.v111.i10

## Estimating Conditional Distributions with Neural Networks Using R Package deeptrafo

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#### Abstract

Contemporary empirical applications frequently require flexible regression models for complex response types and large tabular or non-tabular, including image or text, data. Classical regression models either break down under the computational load of processing such data or require additional manual feature extraction to make these problems tractable. Here, we present **deeptrafo**, a package for fitting flexible regression models for conditional distributions using a **tensorflow** back end with numerous additional processors, such as neural networks, penalties, and smoothing splines. Package deeptrafo implements deep conditional transformation models (DCTMs) for binary, ordinal, count, survival, continuous, and time series responses, potentially with uninformative censoring. Unlike other available methods, DCTMs do not assume a parametric family of distributions for the response. Further, the data analyst may trade off interpretability and flexibility by supplying custom neural network architectures and smoothers for each term in an intuitive formula interface. We demonstrate how to set up, fit, and work with DCTMs for several response types. We further showcase how to construct ensembles of these models, evaluate models using inbuilt cross-validation, and use other convenience functions for DCTMs in several applications. Lastly, we discuss DCTMs in light of other approaches to regression with non-tabular data.

Keywords: deep learning, distributional regression, neural networks, transformation models.

### 1. Introduction

Regression analysis aims to characterize the conditional distribution of a response Y given a set of covariates X, thereby describing how changes in the covariates propagate to the conditional distribution of Y given X (Fahrmeir, Kneib, Lang, and Marx 2013). In this paper, we present deeptrafo (Kook, Baumann, and Rügamer 2024), an R package for estimating a broad class of distributional regression models for various types of responses (continuous, survival, count, ordinal, binary) using tabular or non-tabular (e.g., image or text) data or both. Package deeptrafo is available from the Comprehensive R Archive Network (CRAN) at https://CRAN.R-project.org/package=deeptrafo. Due to a flexible tensorflow (Allaire and Tang 2024) back end and mini-batch optimization, deeptrafo does not only scale well with non-tabular (imaging, text) data but also large tabular data sets. Many well-known models fall into the class of transformation models (TMs), such as normal linear regression (Lm), Cox proportional hazards models (CoxPH), and proportional odds logistic regression (Polr, Hothorn, Möst, and Bühlmann 2018). In the following, we review existing software for fitting these models.

Existing software packages. TMs for tabular data are implemented in tram (Hothorn, Barbanti, and Siegfried 2024) using mlt (Hothorn 2020a) and fitted via maximum likelihood, relying on alabama (Varadhan 2023) and BB (Varadhan and Gilbert 2009) for optimization. Package tram provides an intuitive interface for fitting a multitude of distributional regression models, ranging from shift and shift-scale (Siegfried, Kook, and Hothorn 2024) to tensor-product (or "conditional") transformation models (Hothorn, Kneib, and Bühlmann 2014). Several extensions of transformation models exist. For instance, **cotram** for count TMs (Siegfried and Hothorn 2020), tramME for mixed effects TMs and TMs including smoothing splines (Tamási and Hothorn 2021), and tramnet as well as tramvs for regularized TMs (Kook and Hothorn 2021; Kook 2024). Transformation boosting machines (Hothorn 2020b) and transformation trees and random forests (Hothorn 2023) offer extensions to classical machine learning models. Table 1 summarizes the commonalities and differences between the packages implementing different (extensions of) transformation models in terms of model classes, support for mgcv-based splines and tensorflow-based neural networks and scalable optimization (via mini-batch training, see Appendix H). The **deeptrafo** package is currently the only package implementing transformation models which supports neural network architectures enabling direct handling of text, image, and other deep learning-related data without requiring feature engineering.

Neural network-based transformation models. With the advent of (deep) neural networks and the routine collection of non-tabular data, the idea to combine deep learning and distributional regression approaches was adopted in several ways. For instance, Rügamer, Kolb, and Klein (2024) parameterize distributional regression models via neural networks, Sick, Hothorn, and Dürr (2021) describe flexible deep transformation models for continuous responses. Kook, Herzog, Hothorn, Dürr, and Sick (2022) focus on semi-structured regression for ordinal responses, and Rügamer, Baumann, Kneib, and Hothorn (2023a) extend the DCTM approach to distributional autoregressive models for time series responses. Alternative approaches to combining regression with neural networks including generalized additive models for location, scale, and shape have been implemented in deepregression (Rügamer

Package	Model class	Nonlinear	Splines	Neural networks	Scalable optimization
tram	Linear TMs	Х	Х	Х	Х
$\operatorname{cotram}$	Count TMs	X	Х	Х	Х
tramnet	$L_1/L_2$ -penalized TMs	X	Х	Х	Х
$\mathbf{tramvs}$	$L_0$ -penalized TMs	X	Х	X	Х
$\mathbf{tbm}$	Additive TMs	✓	Х	X	Х
$\mathbf{trtf}$	Transformation forests	✓	Х	Х	Х
$\mathbf{tramME}$	Additive mixed TMs	✓	✓	X	Х
${f deeptrafo}$	Additive TMs	✓	✓	✓	✓

Table 1: Overview of packages for estimating different classes of transformation models. Packages **tramME**, **tbm**, **trtf**, and **deeptrafo** support estimation of nonlinear TMs. Specifically, **tramME** supports splines from **mgcv**, **tbm** fits nonlinear model components via score-based boosting, **trtf** fits nonlinear effects by aggregating trees with TMs in the leaves, and **deeptrafo** supports both splines from **mgcv** and neural networks from **tensorflow**. Package **deeptrafo** allows for scalable optimization via mini-batch training.

et al. 2023c). In this paper, we present **deeptrafo**, which unifies the above DCTM approaches in a single R package.

Comparison to existing packages. Combining distributional regression with neural network-based estimation has many advantages, such as modularity (data analysts can easily use well-established problem-specific neural network architectures), and easy handling of big datasets (e.g., through mini-batch gradient descent with adaptive learning rates). Like tram, deeptrafo relies on maximizing a likelihood function. However, stochastic first-order optimization, such as stochastic gradient descent and the ability to deal with non-tabular data distinguishes the two packages (Table 1). Further, deeptrafo covers and extends models implemented in cotram. Like tramME, deeptrafo also allows the specification of smoothing splines via mgcv (Wood 2023). However, the focus of our package does not lie on random effects. Penalization as in tramnet is also available for deeptrafo. Lastly, unlike models in deepregression, DCTMs do not require specification of a parametric family of distributions for the response given covariates.

The rest of this paper is organized as follows. Section 1.1 introduces the statistical theory behind TMs and DCTMs. The inner workings of **deeptrafo** are described in Section 2, where several case studies on how to setup up, fit, validate, and interpret DCTMs are presented. We present an application to binary classification with tabular and text modalities, and an application to time series modeling via autoregressive TMs (Rügamer *et al.* 2023a). The appendix contains information on advanced usage of the package, e.g., how censored responses are handled (Appendix B) or how to warm-start or fix parameters of certain predictors (Appendix C). In Appendix H, we demonstrate the package for large tabular datasets and factors with many levels, which cannot be handled by standard implementations of classical regression models.

#### 1.1. Deep conditional transformation models

Transformation models (Hothorn et al. 2014, 2018) estimate the conditional cumulative dis-

tribution function (CDF) of a response  $Y \in \mathcal{Y} \subseteq \mathbb{R}$  given a realization  $\boldsymbol{x}$  of covariates  $\boldsymbol{X} \in \mathcal{X}$ ,

$$F_{Y|X=x}(y) := P(Y \le y \mid X = x)$$

without committing to a particular parametric family of distributions for  $F_{Y|X=x}$ . Instead of estimating the CDF directly, transformation models estimate how to transform the response (conditional on covariates) to a latent variable  $Z := h(Y \mid x)$  with fixed and user-defined CDF  $F_Z : \mathbb{R} \to [0,1]$ , using the transformation function  $h : \mathcal{Y} \times \mathcal{X} \to \mathbb{R}$ , which is constrained to be monotonically non-decreasing in  $y \in \mathcal{Y}$  for all  $x \in \mathcal{X}$ . Then, the conditional CDF of the outcome given covariates can be evaluated using the latent CDF  $F_Z$  and the transformation function h:

$$P(Y \le y \mid \boldsymbol{X} = \boldsymbol{x}) = P(h(Y \mid \boldsymbol{x}) \le h(y \mid \boldsymbol{x}) \mid \boldsymbol{X} = \boldsymbol{x}) = P(Z \le h(y \mid \boldsymbol{x})) = F_Z(h(y \mid \boldsymbol{x})).$$

For continuous responses, h is continuous and for discrete responses, h is discrete (see Figure 1). Expressing the conditional CDF in terms of  $F_Z$  and h yields simple expressions for probability density and mass functions and thus also the log-likelihood.

Depending on the choice of  $F_Z$  and restrictions on the functional form and parameterization of h, TMs cover a wide range of well-known models with varying complexity.

**Example 1 (Beyond normal linear regression)** Choosing  $F_Z = \Phi$  and  $h(y \mid \boldsymbol{x}) = \sigma^{-1}(y - \alpha - \boldsymbol{x}^{\top}\boldsymbol{\beta})$ , with standard deviation  $\sigma > 0$ , and intercept  $\alpha \in \mathbb{R}$ , is equivalent to a normal linear regression model, since  $P(Y \leq y \mid \boldsymbol{X} = \boldsymbol{x}) = \Phi(\sigma^{-1}(y - \alpha - \boldsymbol{x}^{\top}\boldsymbol{\beta}))$ . Fixing the transformation function to be linear will always result in conditionally normal outcome distributions. However, this restriction can be lifted by using a nonlinear increasing transformation,  $h_Y : \mathcal{Y} \to \mathbb{R}$ , i.e.,  $h(y \mid \boldsymbol{x}) = h_Y(y) - \boldsymbol{x}^{\top}\boldsymbol{\beta}$ , which now assumes that the transformed response  $h_Y(Y)$  (instead of the original response) is normal with mean  $\boldsymbol{x}^{\top}\boldsymbol{\beta}$ .

Example 2 (Beyond Weibull regression) Choosing  $F_Z(z) = 1 - \exp(-\exp(z))$  with  $h(y \mid \boldsymbol{x}) = a + b \log y + \boldsymbol{x}^\top \boldsymbol{\beta}$ , with intercept a and slope b > 0, is equivalent to a Weibull regression model, since  $P(Y \leq y \mid \boldsymbol{X} = \boldsymbol{x}) = 1 - \exp(-\exp(a + b \log y + \boldsymbol{x}^\top \boldsymbol{\beta})) = 1 - \exp(-\tilde{a}y^b \exp(\boldsymbol{x}^\top \boldsymbol{\beta}))$ , where  $\tilde{a} \coloneqq \exp(a)$ . Also in this example, log-linearity of the transformation function fixes the conditional outcome to be Weibull distributed. Allowing an arbitrary increasing function,  $h_Y : \mathcal{Y} \to \mathbb{R}$ , instead, i.e.,  $h(y \mid \boldsymbol{x}) = h_Y(y) + \boldsymbol{x}^\top \boldsymbol{\beta}$ , results in the Cox proportional hazards model, since the survivor function equals  $P(Y \geq y \mid \boldsymbol{X} = \boldsymbol{x}) = \exp(-\exp(h_Y(y)) \exp(\boldsymbol{x}^\top \boldsymbol{\beta}))$  and  $\exp(h_Y(y))$  is the cumulative baseline hazards.

Thus, TMs contain both normal linear and Weibull regression but also extend both to a more flexible counterpart that does not assume a parametric family of conditional outcome distributions.

Parameterizing the transformation function. In semi-structured regression, we have access to J input modalities  $X_1, \ldots, X_J$ , such as tabular data, images, or text, from which we construct structured (e.g., linear, sparse, or smooth) or unstructured (e.g., neural network) predictors. These inputs may be non-tabular, i.e., there may be a j for which  $X_j \in \mathcal{X}_j \not\subseteq \mathbb{R}^d$ . By  $\mathcal{X} := \mathcal{X}_1 \times \cdots \times \mathcal{X}_J$  we denote the entire input space. In DCTMs, restrictions on the functional form of h, i.e., the way predictors are constructed based on the input data, lead

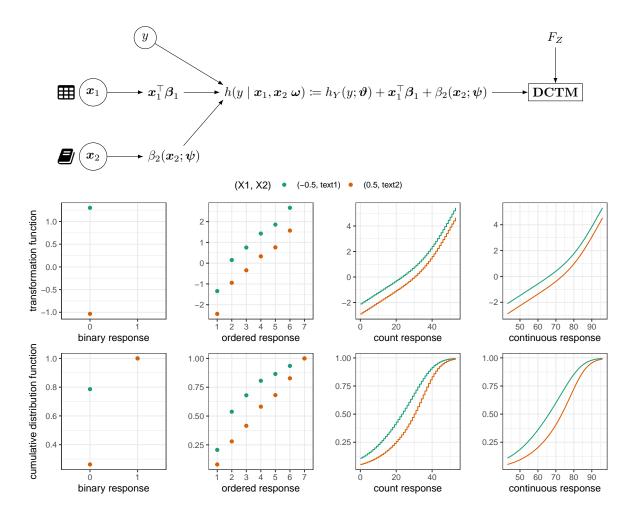


Figure 1: Example of a DCTM with transformation function  $h(y \mid x_1, x_2)$  depending on a tabular modality  $X_1$  and a text modality  $X_2$ , which both enter as an additive shift term. The tabular modality enters as a simple linear predictor  $x_1^{\top}\beta_1$  and the text data via the output of a neural network  $\beta_2$ , which is specified by the user. Together with a baseline transformation  $h_Y$ , whose parameterization is discussed later, and the latent distribution  $F_Z$ , the DCTM is fully specified. On the bottom, the transformation function h and cumulative distribution function  $F_{Y|X_1=x_1,X_2=x_2}=F_Z\circ h$  are depicted for a binary, ordered, count, and continuous response for two realizations of the tabular (X1, X2) and text modalities (text1, text2). For binary and ordered responses with K levels, the transformation function contains one and K-1 parameters, respectively, because the CDF is constrained to one for the largest class.

to varying degrees of interpretability and flexibility of the model. We begin with an example before introducing h in its most flexible form. Consider a problem with a single tabular  $(X_1 \in \mathcal{X}_1 \subseteq \mathbb{R}^p)$  and a single text modality  $(X_2 \in \mathcal{X}_2)$ . Data analysts commonly assume additivity in the effects the separate modalities, which can be realized by modeling the effect of both modalities as shift terms,

$$h(y \mid \boldsymbol{x}_1, \boldsymbol{x}_2; \boldsymbol{\omega}) = h_Y(y; \boldsymbol{\vartheta}) + \boldsymbol{x}_1^{\top} \boldsymbol{\beta}_1 + \beta_2(\boldsymbol{x}_2; \boldsymbol{\psi}), \quad y \in \mathcal{Y},$$
(1)

where  $h_Y: \mathcal{Y} \to \mathbb{R}$  denotes the baseline transformation (i.e., the transformation function obtained when  $\boldsymbol{x}_1^{\top}\boldsymbol{\beta}_1 + \beta_2(\boldsymbol{x}_2) = 0$ , which is parameterized in terms of  $\boldsymbol{\vartheta} \in \mathbb{R}^M$ ). Further,  $\boldsymbol{\beta}_1$  denotes the coefficients of the linear predictor and  $\beta_2: \mathcal{X}_2 \to \mathbb{R}$  denotes the unstructured predictor for the text data, which are typically controlled by a neural network with weights  $\boldsymbol{\psi}$ . By  $\boldsymbol{\omega} := (\boldsymbol{\vartheta}, \boldsymbol{\beta}, \boldsymbol{\psi})$ , we denote the collection of all parameters, including the neural network weights. A DCTM with (1) is distribution-free because for any constellation of covariates for which the shifting predictor is zero, i.e., for all  $(\boldsymbol{x}_1^0, \boldsymbol{x}_2^0) \in S^0 := \{(\boldsymbol{x}_1, \boldsymbol{x}_2) \in \mathcal{X}_1 \times \mathcal{X}_2 \mid \boldsymbol{x}_1^{\top}\boldsymbol{\beta}_1 + \beta_2(\boldsymbol{x}_2) = 0\}$ , and all conditional distributions  $Y \mid \boldsymbol{X}_1 = \boldsymbol{x}_1^0, \boldsymbol{X}_2 = \boldsymbol{x}_2^0$ , there exists a unique baseline transformation given by  $h_Y = F_Z^{-1} \circ F_{Y|X_1 = \boldsymbol{x}_1^0, \boldsymbol{X}_2 = \boldsymbol{x}_2^0}$ . In (1), covariate effects are assumed to enter additively on the scale of the transformation function, thus restricting distributions that can be modeled for  $(\boldsymbol{x}_1, \boldsymbol{x}_2) \in \mathcal{X} \setminus S^0$ . This argument can be extended to more complex DCTMs (for shift-scale see, e.g., Siegfried *et al.* 2024). The example in (1) is depicted in Figure 1 for typical types of responses and standard logistic latent distribution.

In **deeptrafo**, the most general transformation function h is parameterized in terms of  $\omega := (\vartheta, \beta, \phi, \psi) \in \mathbb{R}^{Md} \times \mathbb{R}^p \times \mathbb{R}^q \times \mathbb{R}^s$  which serves as the collection of parameters for basis expansions (potentially including neural networks) of the response and input modalities,

$$h(y \mid \boldsymbol{x}; \boldsymbol{\omega}) = (\boldsymbol{a}(y) \otimes \boldsymbol{b}(\boldsymbol{x}; \boldsymbol{\phi}))^{\top} \boldsymbol{\vartheta} + \boldsymbol{s}(\boldsymbol{x}; \boldsymbol{\psi})^{\top} \boldsymbol{\beta}, \quad y \in \mathcal{Y}, \ \boldsymbol{x} \in \mathcal{X},$$
(2)

where  $\otimes$  denotes the Kronecker product and  $\boldsymbol{a}:\mathcal{Y}\to\mathbb{R}^M,\boldsymbol{b}:\mathcal{X}\to\mathbb{R}^d,\boldsymbol{s}:\mathcal{X}\to\mathbb{R}^p$  denote the bases for the response, and the J predictors, which either interact  $(\boldsymbol{b}(\cdot;\boldsymbol{\phi}))$  with the response or simply shift  $(\boldsymbol{s}(\cdot;\boldsymbol{\psi}))$  the transformation function. The dimensions of the neural network weights  $\boldsymbol{\phi}$  and  $\boldsymbol{\psi}$  depend on the complexity of the neural network architectures which the user has full control over. In **deeptrafo**, the basis for the response is not data-dependent and thus contains no parameters. The interacting and shifting basis, however, depend on the covariates and may include splines or neural networks, whose parameters are collected in  $\boldsymbol{\phi}$  and  $\boldsymbol{\psi}$ , respectively.

The transformation function is required to be monotonically non-decreasing for all  $x \in \mathcal{X}$ . Hence, depending on the choice of basis, the parameters  $\vartheta$  in (2) need to fulfill positivity or monotonicity constraints (Hothorn *et al.* 2014), which can be enforced by appropriate reparameterizations. Without interacting predictors, Bernstein polynomials and discrete bases require  $\vartheta_1 \leq \vartheta_2 \leq \cdots \leq \vartheta_M$  and linear and log-linear bases require positive slopes. For more complex interacting predictors, the positivity of  $b(\cdot; \phi)$  has to be enforced together with more complex constraints on  $\vartheta$  to ensure a monotonically non-decreasing transformation function (for details, see Baumann, Hothorn, and Rügamer 2021).

Shift effects are constant across all values of the response, i.e., the transformation h can only shift up- or downwards (see Figure 1). The effect of interacting predictors may vary with the response and thus the shape of h may change for different predictor values. For instance, an interacting binary predictor leads to two separate transformations for each level, much like stratum variables in survival analysis allow for separate hazard functions (Collett 2015). However, in its general form, interacting predictors may also include neural networks and thus unstructured predictors, making them extremely versatile. Scale effects as introduced in Siegfried et al. (2024) are a special case of interacting predictors, which are included in deeptrafo by using  $b: \mathcal{X} \to \mathbb{R}_+$  with  $x \mapsto \sqrt{\exp(\gamma(x))}$  and, e.g., a neural network  $\gamma: \mathcal{X} \to \mathbb{R}$ . With a linear basis a in y,  $a(y) = (1, y)^{\top}$ , this is equivalent to location-scale regression with error distribution  $F_Z$ .

Supported response types. Several types of univariate, potentially censored, responses can be handled. This includes continuous  $(\mathcal{Y} \subseteq \mathbb{R})$ , survival  $(\mathcal{Y} \subseteq \mathbb{R}_+)$ , count  $(\mathcal{Y} = \mathbb{N})$ , and ordered  $(\mathcal{Y} = \{y_1, \dots, y_K\})$  responses. For continuous responses, the basis for Y is a smooth function parameterized via polynomials in Bernstein form of order M-1, denoted by  $\mathbf{a}_{\mathrm{Bs},M-1}(y)$ . For count responses (M=K), the polynomials in Bernstein form are evaluated only at the integers, i.e.,  $\mathbf{a}_{\mathrm{Bs},M-1}(\lfloor y \rfloor)$  (Siegfried and Hothorn 2020). For ordered responses, a dummy-encoding is used, i.e., for  $k=1,\dots,K$ ,  $\mathbf{b}(y_k)=\mathbf{e}_k$ , where  $\mathbf{e}_k$  denotes the k-th unit vector. Linear and log-linear bases are supported as well. In Appendix D, we describe how the user can supply custom basis functions.

Fitting transformation models. Finally, transformation models can be fitted by minimizing the negative average log-likelihood over the class of transformation functions  $h(y \mid x; \omega)$  with parameters  $\omega$ ,

$$\mathrm{NLL}(\boldsymbol{\omega}; y_i, \boldsymbol{x}_i) \coloneqq -\frac{1}{n} \sum_{i=1}^n \ell(\boldsymbol{\omega}; y_i, \boldsymbol{x}_i),$$

where the observations  $\{(y_i, x_i)\}_{i=1}^n$  are assumed to be independent. In **deeptrafo**, the default optimizer is (stochastic) gradient descent using Adam (Kingma and Ba 2015). However, any **keras** (Allaire and Chollet 2024) or **tensorflow** optimizer or a custom optimization routine can be used instead. For a single observation (y, x), the log-likelihood contribution  $\ell(h; y, x)$  depends on the type of censoring of the observed response. Exact responses y contribute  $\log f_Z(h(y \mid x))h'(y \mid x)$  to the log-likelihood. Interval-censored responses  $(y, \bar{y}]$  contribute  $\log(F_Z(h(\bar{y} \mid x)) - F_Z(h(\bar{y} \mid x)))$ . Left- and right-censored observations follow from the interval-censored contribution as a special case, by letting  $y \to -\infty$  and  $y \to +\infty$ , respectively (Hothorn *et al.* 2014). In **deeptrafo**, the log-likelihood contributions are implemented in terms of mathematical operations implemented in **tensorflow**, which call their Python (Van Rossum *et al.* 2011) counterpart via **reticulate** and allow efficient computation of the log-likelihood, its gradients and weight updates during optimization.

#### 1.2. Autoregressive transformation models

Time series data pose one particular case where the independence assumption between observations is not tenable and needs to be taken into account. Formally, the joint distribution of a time series  $(Y_t)_{t \in \mathcal{T}}$  with  $\mathcal{T} \subseteq \mathbb{N}_0$  can always be factorized in its conditional distributions, i.e., by conditioning  $Y_t$  on its full history  $\mathcal{F}_{t,1} := (Y_{t-1}, \dots, Y_1)$ . A simplification is to impose a Markov property of order  $p \geq 1$  which implies that the conditional distribution of  $Y_t$  only depends on the history up to and including t - p, that is  $\mathcal{F}_{t,t-p} := (Y_{t-1}, \dots, Y_{t-p})$  rather than the entire history  $\mathcal{F}_{t,1}$ .

Package **deeptrafo** offers three ways on how to model time series data assuming the Markov property. The naive way is given by classical transformation models where  $\mathcal{F}_{t,t-p}$  is regarded in the basis expansion of the transformation function shown in (2) where elements of  $\mathcal{F}_{t,t-p}$  may interact with the response  $Y_t$  and simultaneously shift the transformation function. Furthermore, Rügamer *et al.* (2023a) proposed the class of autoregressive transformation models (ATMs) which differ from the naive approach (i.e., classical transformation models) in two perspectives. First, the transformation function  $h_t$  in ATMs can be time-varying which may result in different transformations for different sub-periods. Second, the same  $h_t$  is applied to

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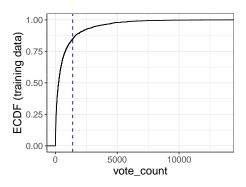


Figure 2: Left: The pre-processed movie review of a picked instance of the movies dataset, in which stop words and punctuation have already been removed. Right: The empirical CDF of vote\_count over all movies in the training, where the picked instance has a vote\_count of 1400 as indicated by the dashed line (placing it above the top quartile). The used tabular input data comprise popularity (4.08 for the picked instance) and revenue (400 million US dollars for the picked instance).

 $Y_t$  and each element of  $\mathcal{F}_{t,t-p}$  simultaneously, resulting in a shared transformation between  $Y_t$  and its lags.

A special subclass of ATMs are AT(p) models which do not allow for interacting elements of  $\mathcal{F}_{t,t-p}$  with  $Y_t$  through  $\boldsymbol{b}$  but restrict to a linear shift impact of the transformed values of  $\mathcal{F}_{t,t-p}$  on the scale of h. The class of AT(p) models is closely related to a well-known class of time series models, i.e., autoregressive models of order p (AR(p), Hamilton 2020). In fact, AT(p) models are equivalent to AR(p) models for  $\boldsymbol{a}(y) = (1,y)^{\top}$ ,  $\boldsymbol{s}(\boldsymbol{x}) \equiv \boldsymbol{x}$  and the independent white noise follows the distribution  $F_Z$  (for details, see Rügamer et al. 2023a). Learning the transformation simultaneously for the response and its lags as it is done in ATMs is particularly important for ordinal time series, for which the dimensionality of the model can thereby be reduced. Instead of modeling each level of the lagged response, the one-dimensional transformed lagged response is included. It also allows for a more consistent interpretation in the sense of autoregression because we model  $h(Y_t)$  at the current step (auto)regress the next time point  $h(Y_{t+1})$  on the likewise transformed response  $h(Y_t)$ , not on the untransformed  $Y_t$ . We showcase the practical differences between linear transformation models, AT(p) and ATM models in Section 4.

#### 1.3. Application datasets

Movies data. In Section 2, we will illustrate the features of deeptrafo using the movies dataset (Kaggle 2017). The dataset contains information on 45,000 movies released prior to July 2017, including number of ratings, budget, revenue, popularity, run time, and genre. In addition, non-tabular reviews of the movies are available as text data. In Section 2, we will focus on estimating the conditional distribution of vote\_count given whether a movie is an action movie, its budget, its popularity score, and the text review. In Section 3, we will switch to the binary classification task of deciding whether a movie falls into the action genre or not. This way, we can showcase how to apply DCTMs for a wider range of outcome types. We pre-process budget, revenue, and popularity using  $\log(1+x)$ , due to their skewed nature. In Figure 2, we show the empirical CDF of the variable vote\_count of the movies dataset

and provide more information on the used variables for one specific movie. For the text data, we use a text\_tokenizer with a 1,000 word vocabulary, convert text to sequence and pad sequences to a maximum length of 100 and truncate the end of a review. We use such a simple embedding to illustrate the key steps of the analysis and make the computations feasible on a standard laptop with 8 gigabytes of RAM. We additionally present results with a pre-trained embedding that performs comparably in terms of test NLL in Appendix E.

Temperature data. An application of autoregressive transformation models to a time series of monthly mean maximum temperature in Melbourne (Australia) in degrees Celsius between January 1971 and December 1990 (240 records) is presented in Section 1.2. The temperature time series was recorded by the Australian Bureau of Meteorology and later provided in Hyndman and Yang (2022).

## 2. The package

Package deeptrafo builds upon tensorflow as a fitting engine and deepregression for setting up structured model terms such as linear effects or splines within a neural network. In contrast to deepregression, which implements models with parametric families and individual additive predictors, deeptrafo supports more complicated computations such as in (2). This is exposed to the user via **deeptrafo**'s formula interface. In **deeptrafo**, response, interacting, and shifting terms are represented as 'formula' objects and correspond to the bases in (2). Internally, a processor is defined for each model term, which evaluates its basis functions and optional penalties via deeptrafo internal, mgcv, or keras/tensorflow functions. For instance, for a continuous response, a polynomial basis in Bernstein form and its derivatives are set up by default (cf. Table 2). The corresponding basis functions are implemented in **deeptrafo**. Package deeptrafo can include terms modeled by user-specified neural network architectures for the interacting and shifting terms (see Figure 1 and Figure 3). When initializing the model using such a formula-based call, the model is internally translated to tensorflow computations using a computational graph. In the end, a single end-to-end trainable neural network is set up, which may contain different neural network components for different terms in the interacting or shifting predictor. Together with the supplied latent\_distr  $F_Z$ , the DCTM is fully specified and its parameters can be estimated by minimizing the NLL via stochastic gradient descent (SGD). Since the DCTM has internally been translated to a model from tensorflow, the optimization can be done using the keras API, which implements the SGD routine with many choices for adaptive learning rates while providing training metrics without requiring users to define training loops for parameter updates. An appropriate last-layer transformation ensures monotonicity constraints of the interacting model term in the response.

Workflow. Typical workflows around deeptrafo, including the illustration in Section 2 and both applications on binary classification (Section 3) and distributional time series (Section 4), are structured as follows: First, a model formula is set up. The 'formula' object encodes in which way each feature enters the model. If neural network components are used, the corresponding architectures have to be defined beforehand. Next, the latent distribution  $F_Z$  is chosen and decides which scale the partial effects of components in the formula are interpreted. Although the formula together with the latent distribution formally specify the

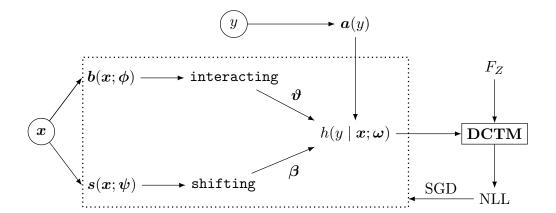


Figure 3: Schematic depiction of setting up and fitting DCTMs. Bases for input predictors x and response y (circles) are evaluated and enter the two neural network components interacting and shifting according to (3). The components' outputs make up the transformation function h which is parameterized in terms of  $\omega$ . Together with the latent distribution  $F_Z$ , the loss, e.g., NLL, and its gradients can be evaluated and used to update parameters  $\omega$ . Since  $F_Z$  is parameter-free, all trainable parameters are in the transformation function, as indicated by the dotted box.

Model function	Model name	Default basis	Default latent distribution
BoxCoxNN	Transformed normal	Bernstein	Standard normal
ColrNN	Continuous outcome logistic	Bernstein	Standard logistic
cotramNN	Count transformation	Bernstein	Standard logistic
CoxphNN	Cox proportional hazards	Bernstein	Standard minimum extreme value
LehmannNN	Lehmann-type	Bernstein	Standard maximum extreme value
LmNN	Normal linear	Linear	Standard normal
PolrNN	Proportional odds logistic	Discrete	Standard logistic
SurvregNN	Weibull	Log-linear	Standard minimum extreme value

Table 2: Supported models together with the default choice of basis function and latent distribution. Model functions summarized here are implemented with a specific choice of basis and latent distribution that define commonly applied regression models.

TM completely (Figure 3), the data and optimizer have to be supplied at this stage. For deep learning models (as opposed to statistical models), it is common to separate model building from model fitting, in order to supply more arguments (such as callbacks) to the optimization routine. Now, hyperparameters can be tuned based on cross-validation. Finally, with the chosen hyperparameters, either a single instance of the DCTM or an ensemble is fitted and can be used for downstream prediction tasks. In Section 2.1, we describe each step of the workflow in more detail using the movies data.

Each step in the **deeptrafo** workflow is highly customizable, e.g., custom functions for basis evaluation (Appendix D), custom last-layer transformations, and general-purpose optimization routines (Section 3), such as SGD with adaptive learning rates (Appendix F), can be supplied.

Effect / Processor	Example formula
Linear Smooth Tensor product splines Lasso Group lasso Row-wise tensor product	$y \sim x$ $y \sim s(x,)$ $y \sim [te ti t2](x,)$ $y \sim lasso(x)$ $y \sim grlasso(x)$ $y \sim rwt(x)$
Neural network	$y \sim nn(x)$
Processor	*_processor e.g., fac_processor

Table 3: Implemented choices of interacting and shift processors. All splines from mgcv are supported. Custom neural networks can be supplied as functions via  $list_of_deep_models$ . Additional processors, for example, for faster processing of large factors or interactions from safareg, can be included via additional\_processors (Rügamer  $et\ al.\ 2023b$ ; Rügamer 2024). All terms can also be included as interacting effects on the left-hand side of the formula, e.g., y | term(x, ...)  $\sim 1$ .

#### 2.1. Main components

We describe the main components of **deeptrafo** below by showing how to use the formula interface, set up a DCTM, and fit the model. In this section, all steps are illustrated with the movies dataset. In the following examples, we assign non-default values to some of the arguments that can be supplied to functions and methods for building and fitting **keras**-based neural networks. This is not because the models have been tuned extensively, but rather to illustrate the most important hyperparameters that are involved in building and fitting DCTMs.

#### Formula interface

Models can be specified via a formula interface akin to the one used in **tram** (Hothorn *et al.* 2024), where covariates interacting with the response are supplied on the left-hand side, and shift effects are supplied on the right-hand side of the formula, as illustrated below.

response | interacting ~ shifting

Thus, the formula interface mimics the transformation function as introduced in (2):

$$\left(\underbrace{\boldsymbol{a}(y)}_{\text{response}} \otimes \underbrace{\boldsymbol{b}(\boldsymbol{x}; \boldsymbol{\phi})}^{\text{interacting}}\right)^{\top} \boldsymbol{\vartheta} + \underbrace{\boldsymbol{s}(\boldsymbol{x}; \boldsymbol{\psi})}_{\text{shifting}}^{\top} \boldsymbol{\beta}. \tag{3}$$

Case study: Formula interface. We begin by modeling the conditional distribution of vote\_count given a binary indicator of whether the movie is categorized as an action movie or not (genreAction), the movie's budget and its popularity. The below formula allows

for separate baseline transformations of the response for action movies vs. all other genres, a smooth effect for budget and a linear effect for popularity. Here, we use the standard spline basis representation implemented in mgcv. A non-exhaustive list of smoothers and other processors is given in Table 3. Processors are specialized functions for handling predictors which can speed up computation. For instance, fac\_processor() from safareg evaluates factors online and thus facilitates computation for large factor models (Rügamer et~al.~2023b, also see the illustration in Appendix H).

```
R> fm <- vote_count | genreAction ~ 0 + s(budget, df = 3) + popularity
```

In the above formula we exclude an additional intercept in the shift term by specifying 0 + ..., because the interacting basis already contains an intercept.

Setting up DCTMs

DCTMs can be generically set up using the deeptrafo() function.

```
deeptrafo(formula = response | interacting ~ shifting, data = ...)
```

The data can be supplied as a data.frame or list. The function returns a 'deeptrafo' object, whose methods are described in Section 2.2.

Special cases of DCTMs coincide with well-known models and are given their own function in **deeptrafo**. The naming conventions in **deeptrafo** follow the **tram** package (Hothorn *et al.* 2024) and add the suffix NN. For instance, the proportional odds logistic regression model (ordinal response and  $F_Z = \exp{it}$ ) is implemented as Polr() in **tram** and PolrNN() in **deeptrafo** (see Table 2 for an overview).

Case study: Setting up DCTMs. For the movies data, we set up a count transformation model with standard logistic latent distribution. The logistic distribution is chosen, so that the partial effects of the features are interpretable as log-odds ratios. Example interpretations are given in Section 3. We supply the Adam optimizer (the default, see Appendix F) for SGD with learning rate of 0.1 decaying with a rate of  $4 \cdot 10^{-4}$  (Kingma and Ba 2015). The training data train is the result of the pre-processing steps described in Section 1.3. The code for reproducing all output and figures can be found in the online supplement.

```
R> opt <- optimizer_adam(learning_rate = 0.1, decay = 4e-4)
R> (m_fm <- cotramNN(formula = fm, data = train, optimizer = opt))</pre>
```

Untrained count outcome deep conditional transformation model

```
Call:
cotramNN(formula = fm, data = train, optimizer = opt)
Interacting: vote_count | genreAction
Shifting: ~0 + s(budget, df = 6) + popularity
```

```
Shift coefficients:

s(budget, df = 6)1 s(budget, df = 6)2 s(budget, df = 6)3 s(budget, df = 6)4

0.557 -0.702 0.760 -0.181

s(budget, df = 6)5 s(budget, df = 6)6 s(budget, df = 6)7 s(budget, df = 6)8

-0.201 -0.687 0.670 0.671

s(budget, df = 6)9 popularity

-0.377 -0.888
```

Printing the model to the console shows the model specification and shift coefficients. Note that the model has only been randomly initialized and not yet fitted, as indicated by "Untrained count outcome deep conditional transformation model" in the print() call. Upon calling fit(), ensemble(), or cv(), the model's history will be non-empty and it will be considered "trained" when printed again.

### Fitting DCTMs

For fitting DCTMs the user calls fit(), which calls the model internal mod\$fit\_fun(), per default a wrapper around fit.keras.engine.training.Model(), with the supplied arguments (for instance epochs, batch\_size). All functionalities of fitting keras models carry over to fitting DCTMs, including callbacks (i.e., custom operations applied after every iteration or mini-batch update).

Case study: Fitting DCTMs. The 'deeptrafo' object returned by cotramNN is fitted for 1,000 epochs, with a batch size of 64, and a 10% validation split. The validation split is used during training to judge whether overfitting occurs (Goodfellow, Bengio, and Courville 2016). Below, we print the (now trained) model.

```
R> m_fm_hist <- fit(m_fm, epochs = 1e3, validation_split = 0.1,
     batch_size = 64, verbose = FALSE)
R> unlist(coef(m_fm, which = "shifting"))
s(budget, df = 6)1 s(budget, df = 6)2 s(budget, df = 6)3 s(budget, df = 6)4
                             -0.28824
                                                 -0.04608
           0.38339
s(budget, df = 6)5 s(budget, df = 6)6 s(budget, df = 6)7 s(budget, df = 6)8
                             -0.02692
           0.00616
                                                 -0.00511
                                                                     0.01355
s(budget, df = 6)9
                           popularity
          -0.36587
                             -0.82771
```

Figure 4A depicts the training and validation loss trajectory for inspecting convergence and overfitting, which can be generated with plot(m\_fm\_hist). The learning curves indicate that the model is not fully trained after 1000 epochs and there is no evidence for overfitting. Figure 4B shows the estimated transformation function. In Section 2.2, we describe how to produce plots of the transformation function and density. Since genreAction is included as a response-varying effect, the two transformation functions are allowed to cross.

### Working with neural networks

The **deeptrafo** package allows to directly model effects of, for instance, text or image data via neural networks. In DCTMs, neural networks map from a complex input space, such as text

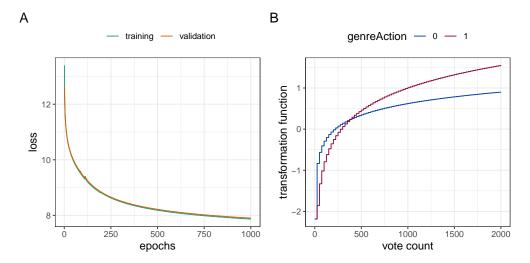


Figure 4: A: Training and validation loss trajectory for m\_fm. B: Estimated transformation functions for both levels of genreAction with popularity and budget fixed at their mean in the training data.

or images, to Euclidean space. When the neural network enters as a shift term, the output of the network is a real number which is interpretable on the latent scale  $F_Z^{-1}$ , i.e., the scale of the transformation function. Custom neural networks can be supplied to deeptrafo as functions or 'keras\_model's via the list\_of\_deep\_models argument.

Case study: Working with neural networks. In our running example, we use the following architecture to model the contribution of the movie reviews provided as textual descriptions. In Section 3, we present an application with further downstream analysis of the text embedding and how this simple embedding compares against using larger pre-trained embeddings.

```
R> embd_mod <- function(x) x |>
+ layer_embedding(input_dim = nr_words, output_dim = embedding_size) |>
+ layer_lstm(units = 50, return_sequences = TRUE) |>
+ layer_lstm(units = 50, return_sequences = FALSE) |>
+ layer_dropout(rate = 0.1) |> layer_dense(25) |>
+ layer_dropout(rate = 0.2) |> layer_dense(5) |>
+ layer_dropout(rate = 0.3) |> layer_dense(1)
```

The neural network  $embd_mod$  maps movie ratings to a real value (for more details see the case study in Section 3). The interpretational scale of output depends on the choice of latent distribution. Here, the logistic distribution ( $F_Z = expit$ ) renders the output of  $embd_mod$  interpretable on the log-odds scale. In turn, differences in the output of  $embd_mod$  can be interpreted as log odds-ratios when changing, for instance, a single word in a sentence and leaving everything else constant. In our deeptrafo model, we can now supply a named list list(deep =  $embd_mod$ ) and use  $embd_mod$  in the formula.

```
R> fm_deep <- update(fm, . ~ . + deep(texts))
R> m_deep <- cotramNN(fm_deep, data = train,</pre>
```

```
+ list_of_deep_models = list(deep = embd_mod))
R> fit(m_deep, epochs = 50, validation_split = 0.1, batch_size = 32,
+ callbacks = list(callback_early_stopping(patience = 5)),
+ verbose = FALSE)
```

The default optimization routine may not produce optimization paths as smooth as when omitting the neural network component. However, adaptively scheduled learning rates for SGD often work well out-of-the-box, e.g., using optimizer = optimizer\_adam() as an argument when initializing the 'deeptrafo' model. Sometimes also different learning schedules are needed for the different modalities (see Section 3).

#### Ensembling DCTMs

A simple and popular method to improve prediction performance and to quantify training stability (i.e., uncertainty from random initialization and stochastic optimization) are deep ensembles (Lakshminarayanan, Pritzel, and Blundell 2017). In a deep ensemble, a neural network model is trained B times using the same training and validation data, but different initial weights. Training via SGD may then converge to different (local) minima and the members may yield different predictions. However, averaging the predicted densities of the B ensemble members is guaranteed to improve upon the average individual performance (e.g., in terms of NLL). In **deeptrafo**, an ensemble of a model can be fitted via **ensemble()**. Besides classical deep ensembling, **deeptrafo** implements transformation ensembles (Kook, Götschi, Baumann, Hothorn, and Sick 2022). Transformation ensembles are specifically tailored towards DCTMs and preserve their additive structure and thus (partial) interpretability by averaging the predicted transformation functions instead of the predicted densities.

Case study: Ensembling DCTMs. Below, we fit five instances of m\_deep. Then, we combine their predictions on the scale of the transformation function and can investigate uncertainty in the effects of the shifting predictors and prediction performance on the test set.

Figure 5 shows the estimated smooth effect of budget with training stability indicated by the shaded area. Investigating the out-of-sample prediction performance, we see that the transformation ensemble performs better than the members do on average (see Proposition 3 in Kook *et al.* 2022).

```
R> unlist(logLik(ens_deep, convert_fun = \(x) - mean(x)))
members1 members2 members3 mean ensemble
-8.28 -8.50 -8.32 -8.37 -8.35
```

Cross-validating DCTMs for hyperparameter tuning

With cv(), deeptrafo provides a cross-validation function for 'deeptrafo' objects. When supplying an integer to cv\_folds, the data is split into cv\_folds number of folds. Alternatively, the user can specify a list with two elements indicating data indices for training

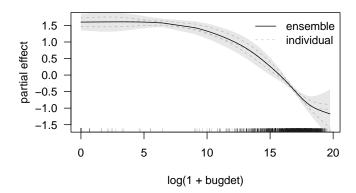


Figure 5: Training stability in the estimated smooth partial effect of budget on vote\_count obtained via transformation ensembling.

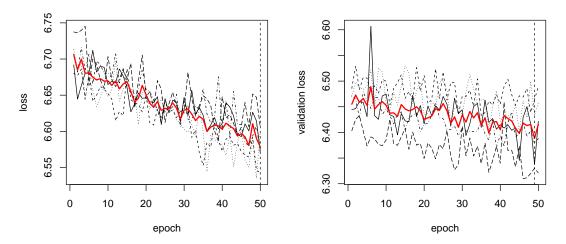


Figure 6: Default plot generated by cv.deeptrafo(). The vertical lines indicate the epoch with minimal average train/validation loss.

and validation. The output of cv() can be used for tuning smoothing hyperparameters, choosing between including a predictor as interacting or shifting, or different neural network architectures.

Case study: Cross-validating DCTMs. The following call to cv() performs 5-fold cross validation while fitting each instance of m\_deep for 50 epochs. Train and validation loss trajectories are shown in Figure 6. The vertical bars indicate the epoch with the best average train/validation loss.

#### 2.2. Methods overview

In the following, we briefly describe S3 methods for 'deeptrafo' and 'dtEnsemble' objects.

### Methods for 'deeptrafo' objects

- coef() returns coefficients for the interacting or shifting terms (controllable via which\_param = c("shifting", "interacting", "autoregressive")).
- predict() returns in-sample predictions when newdata is not supplied. The supported types are "trafo", "pdf", "cdf", "interaction", "shift", "terms". When newdata is supplied, predictions are evaluated at the response, if it is contained in newdata. The response can be omitted from newdata to predict the whole conditional distribution. Then, predictions are evaluated on a grid of length K, which is automatically generated based on the response's support in the training data set. A custom grid of response values can be supplied via q, which overwrites K.
- logLik() evaluates in- or out-of-sample log-likelihoods. This can be useful for model criticism and evaluating predictive performance, respectively. The argument convert\_fun controls how the individual NLL contributions are summarized. The default is function(x) = -sum(x) to compute the log-likelihood. Other common choices include identity to obtain the individual NLL contributions, or mean to get the average NLL.
- plot() by default plots smooth components in the shifting formula part. Data for plotting can be obtained by setting only\_data = TRUE. Smooth terms in interacting can be plotted by setting which\_param = "interacting". In the same manner as in predict, densities evaluated in-sample (type = "pdf"), CDFs (or probability integral transforms, with type = "cdf"), and transformation functions (type = "trafo") can be obtained. When omitting the response from newdata, the whole density, cumulative distribution, or transformation function can be plotted.
- print() prints a brief summary of the DCTM including coefficients of additive linear and smooth terms in shifting. Setting with\_baseline = TRUE also prints coefficients of linear and smooth terms in interacting. The print\_model argument toggles whether the keras summary of the DCTM should be printed in addition.

#### Methods for 'dtEnsemble' objects

Methods coef() and predict() of 'deeptrafo' objects take the same arguments as their 'deeptrafo' counterparts. The output is returned for all ensemble members. Likewise, logLik() returns the processed NLL contributions for individual ensemble members, their average, and the transformation ensemble.

## 3. Application: Binary classification

In this application, we use the movies dataset and fit four different models with the goal to predict the binary response action (0: non-action movie, 1: action movie, defined in the next code chunk), which encodes whether a movie is an action movie or not. The model m\_0 is unconditional; m\_tab uses only one tabular predictor, popularity, as linear shift predictor; m\_text uses only texts as an unstructured shift predictor; m\_semi is a semi-structured model which uses both modalities as shift predictors. The purpose of the analysis is to show the

potential gains in prediction performance that can be achieved when including the text data and learning an embedding, for which conventional statistical models would require extensive feature engineering. The models that do not include the text data could, in principle, also be fitted using glm() from the MASS (Ripley 2024) package and yield virtually the same results as deeptrafo.

First, we encode the binary response as an ordered factor allowing us to use the framework of ordinal neural network transformation models (Kook *et al.* 2022). This step is necessary because unordered factors are not supported by **deeptrafo**.

```
R> train$action <- ordered(train$genreAction)
R> test$action <- ordered(test$genreAction, levels = levels(train$action))</pre>
```

We then set up the formulas for the four models. The unconditional model is specified without any predictors and 1 on the right-hand side. Later, we will restrict this additional intercept to zero for identification (see warmstart\_weights in the definition of  $m_0$ ). For all other models, we remove the intercept directly by specifying  $0 + \ldots$  on the right-hand side.

```
R> fm_0 <- action ~ 1
R> fm_tab <- action ~ 0 + popularity
R> fm_text <- action ~ 0 + deep(texts)
R> fm_semi <- action ~ 0 + popularity + deep(texts)</pre>
```

Here, deep is the same neural network architecture with text embedding as in Section 2.1. We use a custom 'keras\_model' to which we can refer to by the name "embd" and wrap it in a function, which allows us to create multiple instances of the same model. This may be necessary in applications to make the resulting Python objects point to different copies in memory and avoid unintended re-use of already trained or modified models.

```
R> make_keras_model <- function() {
    return(keras_model_sequential(name = "embd") |>
    layer_embedding(input_dim = nr_words, output_dim = embedding_size) |>
    layer_lstm(units = 50, return_sequences = TRUE) |>
    layer_lstm(units = 50, return_sequences = FALSE) |>
    layer_dropout(rate = 0.1) |> layer_dense(25) |>
    layer_dropout(rate = 0.2) |> layer_dense(5, name = "penultimate") |>
    layer_dropout(rate = 0.3) |> layer_dense(1))
    layer_dropout(rate = 0.3) |> layer_dense(1))
```

Next, we use Polrnn() to set up the different models with a standard logistic latent distribution. Models including text data are trained for ten epochs with early stopping and a patience of two, and the weights from the epoch with the best validation loss are restored. The unconditional and tabular-only models are trained full-batch and without validation split until converging to the minimum since convexity of the problem implies a unique solution.

Besides the simple text embedding that is trained from scratch, we also present how to use a pre-trained word2vec embedding in Appendix E.

#### 3.1. Unconditional model

For the unconditional model, the intercept is fixed to zero via warmstart\_weights to ensure identification. The details explaining the next code chunk can be found in Appendix C.

The unconditional model  $m_0$  has one parameter which estimates the log-odds of a movie belonging to a non-action genre without any predictors. The estimated intercept parameter, given by  $coef(m_0)$ , which = "interacting"), corresponds to the single (fix) value of the transformation function h for a binary response (see Figure 1). The code chunk below shows that the estimated intercept agrees with the observed log-odds of a movie belonging to a non-action genre up to numerical inaccuracies.

We can obtain the unconditional log-odds also using predict() with type = "trafo". From the estimated log-odds we can determine the probability for a movie to belong to a non-action genre which matches the prevalence of non-action movies in the train set. The prevalence of non-action movies can also be computed directly by using the predict function and setting the argument type = "pdf" and supplying action = 0 in newdata.

#### 3.2. Tabular-only model

[1] TRUE

Next, we set up and fit m\_tab including popularity as a linear shift predictor.

```
R> m_tab <- PolrNN(fm_tab, data = train, optimizer = optimizer_adam(
+ learning_rate = 0.1, decay = 1e-4))
R> fit(m_tab, epochs = 1e3, batch_size = length(train$action),
+ validation_split = 0, verbose = FALSE)
```

We obtain the estimated linear shift parameter  $\hat{\beta}$  of m\_tab by coef(m\_tab, which\_param = "shifting"). Here, the odds for a movie to belong to genre action change by the factor  $\exp(-\hat{\beta})$ 

```
R> exp(-unlist(coef(m_tab, which = "shifting")))
popularity
    1.54
```

when the predictor popularity increases by one unit. Without flipping the sign, the coefficient  $\hat{\beta}$  represents a log-odds ratio for a movie belonging to a non-action genre compared to genre action upon a one-unit change in popularity. Thus, the interpretation of  $\hat{\beta}$  depends on the parameterization of the model, in particular, the sign of the shifting predictor. In deeptrafo, the shifting predictor is consistently parameterized with a plus sign for all models, which may differ from other implementations of the same model type (e.g., generalized linear models in MASS, or TMs in tram).

### 3.3. Text-only model

We now define and fit m\_text including only the tokenized movie reviews.

Analogously to smooth partial effects, the differences between two shift estimates resulting from two different text inputs can still be interpreted as log odds-ratios.

We now have a closer look at what the embd\_mod has learned. The network takes as input the words (encoded as indices). Here, we use a vocabulary (all words in the data set) of 10000 words and limit each review text to a size of 100 words. Review texts which are shorter are prepended with zeros, longer movie descriptions are cut after 100 words. All punctuation is removed. The layer\_embedding learns to embed the word indices into an embedding\_size-dimensional representation. The resulting word embeddings of the text are the input sequence to an LSTM layer with a 50-dimensional memory state. The second LSTM layer outputs the 50-dimensional state after the last word in the text, which is then further processed by a fully connected neural network including dropout to prevent overfitting.

We can now use the trained embd to extract and analyze the derived latent features of the embedding of single words or whole texts. We can obtain the embedding of a single word as the output of layer\_embedding(). If we use a whole review as input, the latent features in the layer "penultimate" correspond to a five-dimensional representation of the text embedding processed by subsequent layers.

Figure 7 shows the first two components of a principle component analysis (PCA) applied to the word embedding (left) and to the features learned in the penultimate layer for whole reviews (right). The left plot reveals, that words hinting at an action movie, have a similar embedding, and are separated from words that are rather representative of a romance movie. The plot on the right of Figure 7 confirms that the features derived from the embedding are tailored to discriminate action movies from other genres since latent features of action movies cluster together and are fairly well separated from romantic movies.

#### 3.4. Semi-structured model

Finally, we set up the most complex model m\_semi which takes both data modalities as input. To achieve efficient training of the tabular part and avoid overfitting of the embedding network emdb\_semi we use two different learning rates for the structured and unstructured

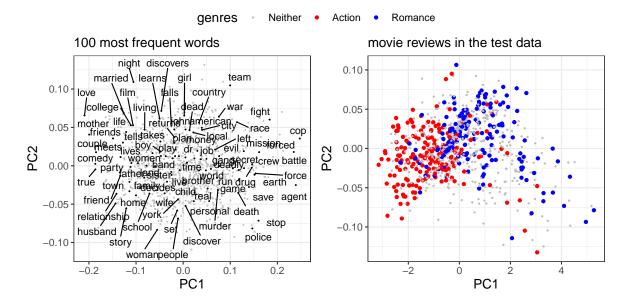


Figure 7: The first two principal components of the "embedding" layer for single words (left) and the lower-dimensional representation learned in "penultimate" (right) for whole movie reviews in the test data. On the left, the PCA is computed on the word embedding of the 1,000 most frequent words (we display only the 100 most frequent, black dots), and on the right based on the low-dimensional representation of the embedding of the 888 full movie reviews contained in the test data.

part of the model. Specifically, we optimize the intercept (with name "ia\_1\_\_2") and tabular shift predictor (with name "popularity\_3") with a higher learning rate, than the embedding model ("embd"). In the embedding model, some layers are named explicitly, the names for the other components can be obtained from the 'keras\_model' summary by initializing and calling print(m\_semi, print\_model = TRUE).

#### 3.5. Model comparison

Comparing the prediction performance of the models (measured in terms of NLL) indicates that mainly the text modality contains information for separating action movies from other genres. However, for a more reliable assessment of this statement, the training schedule should be optimized further. We compute 95% bootstrap confidence intervals as a simple uncertainty measure for the test NLL. In Appendix E, we illustrate how to use pre-trained embeddings with a shallow and deeper neural network architecture and obtain comparable results in terms of out-of-sample NLL. Using pre-trained embeddings may reduce computation times and yield comparable predictions, especially when the training sample size is small (Goodfellow et al. 2016).

```
R> bci <- function(mod) {</pre>
     11i <- logLik(mod, newdata = test, convert_fun = identity)</pre>
     bt <- boot(lli, statistic = (x, d) mean(x[d]), R = 1e4)
     btci <- boot.ci(bt, conf = 0.95, type = "perc")$percent[1, 4:5]
     c("nll" = mean(lli), "lwr" = btci[1], "upr" = btci[2])
+ }
R> mods <- list("unconditional" = m_0, "tabular only" = m_tab,</pre>
    "text only" = m_text, "semi-structured" = m_semi)
R> do.call("cbind", lapply(mods, bci))
    unconditional tabular only text only semi-structured
nll
            0.531
                          0.516
                                     0.437
                                                      0.423
            0.501
                          0.486
                                     0.390
                                                      0.372
lwr
upr
            0.562
                          0.549
                                     0.486
                                                      0.478
```

Like m\_tab the model m\_semi estimates a linear shift parameter for popularity which can also be interpreted as a (conditional) log odds-ratio. The parameter goes in the same direction as in m\_tab but has a reduced absolute value and is now interpretable as a conditional log-odds ratio because the text information that is now additionally accounted for.

The presented case study is meant to showcase some functionality of the package **deeptrafo** for binary responses. A PolrNN model for an ordinal response that has K levels and yields K-1 values for a discrete transformation function (see Figure 1) can be interpreted analogously, e.g., linear shift terms are still interpreted as log odds-ratios (for details and more examples, see Kook *et al.* 2022).

## 4. Application: Autoregressive transformation models

We now return to ATMs, first discussed in Section 1.2. One special form of ATMs are AT(p) models. AT(p) models assume a linear impact of the transformed values of  $\mathcal{F}_{t,t-p}$  on the scale of h. Because the transformation is the same as for the response, AT(p) models thus learn a joint transformation of the response and its lags. For an illustration of transformation models applied to time series data, the temperature dataset is used. We aim to estimate the conditional distribution of the monthly mean maximum temperature in degrees Celsius (°C) in Melbourne (Australia) between January 1971 and December 1990. A descriptive analysis of the time series shows a strong seasonal pattern. This motivates the application of a flexible approach that allows modeling the quickly changing moments of the conditional distribution over time.

In the following, we compare three different forms of autoregressive transformation models. The most flexible model (ATM) includes the lags as interacting predictors and transformed lags in the shift term. The AT(3) model only includes the transformed lags in the shift term. Lastly, the naive ColrNN model (Colr) includes the lags as an additive linear term. In addition, every model contains a shift effect for month. The ATM and AT(3) model can currently only be fitted using deeptrafo, whereas the other two models could be fitted using conventional TMs implemented in tram. We compare the three models based on their estimated transformation functions and conditional densities. We start by creating a factor variable month for the calendar month as well as the lags  $Y_{t-p}$ , p = 1, 2, 3 denoted by y\_lag\_ for including raw additive lags. AT(p) lags are included using the internal atplag() processor.

```
R > lags <- c(paste0("y_lag_", 1:p, collapse = "+"))
```

The formula for the ATM model is given as follows. We include all three lags as interacting predictors on the left-hand side of the formula and specify the atplags on the right-hand side.

```
R> (fm_atm <- as.formula(paste0("y |", lags, "~ 0 + month + atplag(1:p)")))
y | y_lag_1 + y_lag_2 + y_lag_3 ~ 0 + month + atplag(1:p)</pre>
```

ATP lags can be conveniently included in the formula by specifying the lags inside atplag(). For the AT(3) model, we include the transformed lags in the shift but not in the interacting term.

```
R> (fm_atp <- y ~ 0 + month + atplag(1:p))
y ~ 0 + month + atplag(1:p)</pre>
```

The third model (Colr) we compare is a ColrNN model which includes the raw lags in an additive shift term.

```
R> (fm_colr <- as.formula(paste0("y ~ 0 + month + ", lags)))
y ~ 0 + month + y_lag_1 + y_lag_2 + y_lag_3</pre>
```

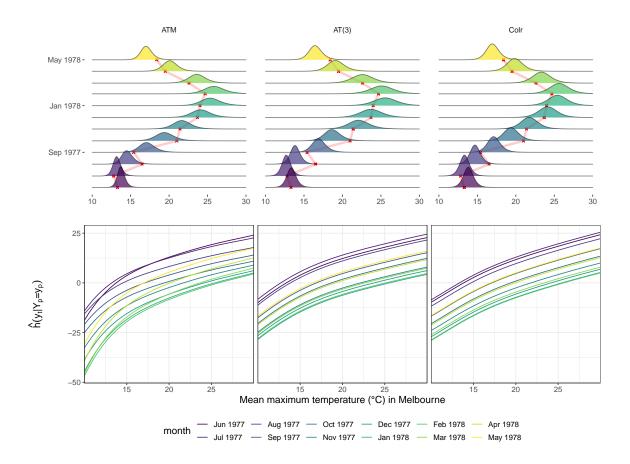


Figure 8: Estimated conditional densities (top row) of monthly temperature records between June 1983 and May 1984, based on the ATM model (left), the AT(3) model (center) and the Colr model (right). The observed values across this time span are depicted in red. The plots in the bottom row show the corresponding estimated conditional transformation functions.

After pre-processing, the temperature dataset is saved in d\_ts. We fix the support of the response to  $min\_supp = 10$  and  $max\_supp = 30$  and specify Bernstein polynomials of order P = 6. We use Colrnn() to specify all models. ATM and AT(3) include atplags and the third model, Colr, does not.

```
R> mod_fun <- function(fm, d) ColrNN(fm, data = d,
+ trafo_options = trafo_control(order_bsp = P,
+ support = c(min_supp, max_supp)), tf_seed = 1,
+ optimizer = optimizer_adam(learning_rate = 0.01))
R> mods <- lapply(list(fm_atm, fm_atp, fm_colr), mod_fun)</pre>
```

After defining the models, we proceed with training all three models. In addition, we include callbacks to reduce the learning rate upon encountering a plateau in the training loss, to ensure convergence of the optimization procedure.

```
+ callback_reduce_lr_on_plateau(patience = 5)), batch_size = nrow(d_ts_lag),
+ verbose = FALSE)
R> lapply(mods, fit_fun)
```

We compare the in-sample log-likelihood for the three models for the subset of data between June 1977 and May 1978 in t\_idx.

The comparison shows that the Colr and the ATM model fit similarly well compared to the slightly less favorable fit of the AT(3) model. A visual inspection of the estimated conditional densities depicted in Figure 8 shows similar results for all three estimation methods. In summary, the ATM class may be favored over naive TMs (Colr) in the time series domain for its autoregressive structural assumption, i.e., lags entering in a transformed way, identical to the transformation of  $y_t$  (see Rügamer  $et\ al.\ 2023a$ ).

#### 5. Conclusion

With deeptrafo, we introduce the first R package for fitting a broad class of distributional regression models with a neural network back-end. Package deeptrafo combines the advantages of transformation models, i.e., flexible distribution-free, yet interpretable models for conditional distributions, with the advantages of neural network-based machine learning, which scales well for large or non-tabular datasets. The intuitive formula interface allows users familiar with packages such as stats (R Core Team 2024), MASS, tram, survival (Therneau 2024), mgcv, and others to easily adapt their workflow to neural networks and more complex datasets out-of-the-box.

Users can supply custom basis functions, loss functions, optimization routines and neural network architectures to adapt and extend functionalities from **deeptrafo** to problems in which the goal is learning a conditional cumulative distribution function. We illustrate **deeptrafo** with tabular and text, as well as time series data with count, discrete, and continuous outcomes, which are all handled in a unified way. We demonstrate how custom neural network architectures and optimizers can be used, and how to tune, evaluate, and interpret DCTMs.

Applying neural network-based models to analyze text or image data typically comes with higher flexibility but also larger computational costs compared to more conventional statistical models. We demonstrate how pre-trained text embeddings can be used to obtain competitive results to training an embedding from scratch and reduce the computational and the environmental burden.

## Acknowledgments

We thank Sandra Siegfried for her valuable comments on the manuscript. The research of LK was supported by the Swiss National Science Foundation (Grant No. 214457). LK conducted part of this work at the University of Copenhagen, University of Zurich and Zurich University of Applied Sciences. The research of DR has been partially supported by the German Federal Ministry of Education and Research (BMBF) under Grant No. 01IS18036A. The research of LK and BS was supported by Novartis Research Foundation (FreeNovation 2019) and by the Swiss National Science Foundation (Grant No. S-86013-01-01 and S-42344-04-01). The research of OD has been partially supported by BMBF under Grant No. 01IS19083A. The authors of this work take full responsibilities for its content.

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#### A. Additional details

In the appendix, we describe how **deeptrafo** handles censored responses (Appendix B), how the user can warmstart and fix weights of interacting and shifting terms (Appendix C), and how to include custom basis functions (Appendix D). We demonstrate how to use pre-trained embeddings (Appendix E) and give details on the most commonly used options for optimization (Appendix F). In addition, we describe an alternative formula interface (Appendix G) and show how to use **deeptrafo** for large tabular datasets (Appendix H).

## B. Handling censored responses

Package **deeptrafo** detects the type of response automatically. However, the user may specify the type explicitly via response\_type in deeptrafo() and all alias/wrapper functions. Allowed types of responses are continuous, count, survival, ordered (including binary). Censored responses can be supplied as 'Surv' objects. Internally, ordered and count responses are treated as censored. For instance, the two observations c(OL, 1L) with response\_type = "count" are internally represented as left- and interval-censored, respectively.

## C. Warmstarting and fixing weights

Warmstarting and fixing weights may be important in numerical experiments, for fine-tuning parts of the models, or transfer learning (Goodfellow et al. 2016). In **deeptrafo**, the user can supply a 'keras\_model', as returned, for instance, by keras\_model\_sequential(). When defining the model, keras specific arguments for controlling weight initialization can be used, as shown below.

```
R> nn <- keras_model_sequential() |>
+ layer_dense(input_shape = 1L, units = 3L, activation = "relu",
+ use_bias = FALSE, kernel_initializer = initializer_constant(
+ value = 1))
R> unlist(get_weights(nn))
[1] 1 1 1
```

To warmstart or fix coefficients of the interacting or shifting part of a DCTM, the weight\_options argument in deeptrafo() can supplied with the output of weight\_control(), which, in addition to others, takes the same arguments as the keras layers above.

```
R> args(weight_control)

function (specific_weight_options = NULL, general_weight_options = list(
    activation = NULL, use_bias = FALSE, trainable = TRUE,
    kernel_initializer = "glorot_uniform", bias_initializer = "zeros",
    kernel_regularizer = NULL, bias_regularizer = NULL,
    activity_regularizer = NULL, kernel_constraint = NULL,
    bias_constraint = NULL), warmstart_weights = NULL,
    shared_layers = NULL)
```

Below, we warmstart the shift coefficient for a PolrNN model. Here, warmstart\_weights takes a list with three components, of which the first two control the weights of the interacting predictor and the last the weights of the shift predictor. The weights can be referred to by the name of the covariate, i.e., "temp" = 0.

The three lists correspond to the three formula components response, interacting, and shifting. The list corresponding to the response is always empty, since it does not contain any parameters. In case there is no interacting predictor, the second list corresponds to the parameters of the basis function of the response, i.e., the intercept function. In case there is no shift term, an intercept is set up which can be referred to as "1" and frozen as illustrated in the main text (Section 3). In the example above, we warmstart weights of a component in the shift term and supply two empty lists for the other components.

## D. Including custom basis functions

Linear, log-linear, and Bernstein bases, as used by **deeptrafo**, require (linear) inequality constraints on their parameters. Internally, these constraints are handled in trafo\_control(), by supplying an **keras** layer, which transforms the weights for the interacting predictor appropriately. In **deeptrafo**, the implemented bases are "bernstein", "ordered", and "shiftscale". The former two require  $\vartheta_{jP+1} \leq \vartheta_{jP+2} \leq \cdots \leq \vartheta_{jP+P}, \ l=0,\ldots,L-1$  for  $\boldsymbol{b}(\boldsymbol{x}) \in \mathbb{R}^L$  and degree P-1 Bernstein basis or ordered response with P+1 levels. The shift-scale basis requires only  $\vartheta_1 > 0$  in  $y \mapsto \vartheta_0 + \vartheta_1 y$ .

The user can now supply custom basis functions as shown below. First, the basis (linear\_basis) and its derivative (linear\_basis\_prime) are defined. Afterwards, the constraints on the parameters are defined using Python- and tensorflow-specified constructs (tf\$...).

```
R> linear_basis <- function(y) {</pre>
     ret <- cbind(1, y)</pre>
     if (NROW(ret) == 1)
       return(as.vector(ret))
+ }
R> linear_basis_prime <- function(y) {</pre>
     ret <- cbind(0, rep(1, length(y)))</pre>
     if (NROW(ret) == 1)
       return(as.vector(ret))
     ret
+
R> constraint <- function(w, bsp_dim) {</pre>
     w_res <- tf$reshape(w, shape = list(bsp_dim, as.integer(nrow(w) /</pre>
     bsp_dim()))
     w1 \leftarrow tfslice(w_res, c(OL, OL), size = c(1L, ncol(w_res)))
     wrest <- tfmathsoftplus(tfslice(w_res, c(1L, 0L), size = c(
     as.integer(nrow(w_res) - 1), ncol(w_res))))
     w_w_cons <- k_concatenate(list(w1, wrest), axis = 1L)</pre>
     tf$reshape(w_w_cons, shape = list(nrow(w), 1L))
+ }
R> tfc <- trafo_control(</pre>
     order_bsp = 1L,
     y_basis_fun = linear_basis,
     y_basis_fun_prime = linear_basis_prime,
     basis = constraint
+ )
```

We can now compare our re-implementation of a transformation model with linear basis against Lm() from tram. To efficiently fit DCTMs for small tabular datasets, we recommend full-batch (i.e., batch size n) training with a large learning rate (0.01) and either decay or callbacks for reducing the learning on validation loss plateaus.

## E. Binary classification with pre-trained embeddings

As large pre-trained language models become more and more practice in natural language processing, we show an alternative way to fit a DCTM using a pre-trained embedding called word2vec (Mikolov, Chen, Corrado, and Dean 2013). The embedding is provided by Google and can be downloaded from their servers. Due to its corpus size, the embedding file is multiple Gigabytes large. After storing the embedding in the ./Data/ folder, we can load the embedding using the gensim Python library (Rehurek and Sojka 2011) and transform every word in the training dataset into a vector in the embedding space.

```
R> embedding_dim <- 300
R> if (file.exists("word2vec_embd_matrix.RDS")) {
    embedding_matrix <- readRDS("word2vec_embd_matrix.RDS")</pre>
    vocab_size <- nrow(embedding_matrix)</pre>
R>
R> } else {
   gensim <- import("gensim")</pre>
R>
    model <- gensim$models$KeyedVectors$load_word2vec_format(</pre>
+
     ".../Data/GoogleNews-vectors-negative300.bin", binary = TRUE)
   vocab_size <- length(words$word)</pre>
R>
    embedding_matrix <- matrix(0, nrow = vocab_size, ncol = embedding_dim)</pre>
R>
R.>
    names_model <- names(model$key_to_index)</pre>
    for (i in 1:vocab_size) {
R.>
R>
      word <- words$word[i]</pre>
      if (word %in% names_model) {
R>
        embedding_matrix[i, ] <- model[[word]]</pre>
R>
R>
      }
R>
R> saveRDS(embedding matrix, file = "word2vec embd matrix.RDS")
R> }
```

Having transformed the text data into vectors in the embedding space, we can use these in an embedding layer to define our model. We start with a shallow neural network that flattens the vectors for each word and learns a linear model for the resulting data matrix.

```
R> w2v_mod <- function(x) x |>
+ layer_embedding(input_dim = vocab_size, output_dim = embedding_dim,
+ weights = list(embedding_matrix), trainable = FALSE) |>
+ layer_flatten() |>
+ layer_dense(units = 1)
R> fm_w2v <- action ~ 0 + shallow(texts)
R> m_w2v <- deeptrafo(fm_w2v, data = train,
+ list_of_deep_models = list(shallow = w2v_mod),
+ optimizer = optimizer_adam(learning_rate = 1e-5))
R> dhist <- fit(m_w2v, epochs = 200, validation_split = 0.1,
+ batch_size = 32, callbacks = list(
+ callback_early_stopping(patience = 5)), verbose = FALSE)
R> bci(m_w2v)
```

```
nll lwr upr
0.523 0.494 0.553
```

While stopping the training later may result in further model improvement, the negative log-likelihood values already indicate similar performance to the tabular-only model in Section 3.5, which yielded a test NLL of 0.52. We can improve this model by learning a deep neural network on top of the pre-trained embedding using 1D convolutions as follows.

```
R> w2v2\_mod <- function(x) x >
     layer_embedding(input_dim = vocab_size, output_dim = embedding_dim,
     weights = list(embedding_matrix), trainable = FALSE) |>
    layer_conv_1d(filters = 128, kernel_size = 5, activation = 'relu') |>
     layer_max_pooling_1d(pool_size = 5) |>
     layer_conv_1d(filters = 128, kernel_size = 5, activation = 'relu') |>
     layer_global_max_pooling_1d() |>
     layer_dense(units = 128, activation = 'relu') |>
     layer_dropout(rate = 0.5) |>
    layer_dense(units = 1)
R> fm_w2v2 <- action ~ 0 + deep(texts)
R > m_w2v2 < - deeptrafo(fm_w2v2, data = train,
     list_of_deep_models = list(deep = w2v2_mod),
     optimizer = optimizer_adam(learning_rate = 1e-5))
R> dhist <- fit(m_w2v2, epochs = 200, validation_split = 0.1,
     batch_size = 32, callbacks = list(
     callback_early_stopping(patience = 5)), verbose = FALSE)
R> bci(m_w2v2)
 n11
        lwr
              upr
0.510 0.486 0.534
```

The test NLL resulting from the deeper architecture is lower than the one obtained from the shallow architecture above, but not as low as the one obtained using the text-only model in Section 3.5 of 0.44 (95% bootstrap confidence interval from 0.390 to 0.486).

## F. Options for optimization

In deep learning, selecting an appropriate optimizer is crucial for model performance. If **deeptrafo** specifies a model with a deep predictor, exact optimization is not possible anymore and routines that require second- or higher-order derivatives of the objective are too expensive. Optimization is therefore done using first-order methods, in particular variations of stochastic gradient descent (SGD).

Adam (Kingma and Ba 2015) is widely used due to its effectiveness across various
applications, offering adaptive learning rates that handle sparse gradients efficiently. It
is by far the most common choice and hence our default option. While Adam's default
learning rate and momentum parameters can be changed, this must be done with care.

- Another option is SGD with momentum, preferred for optimizing large CNNs, with its momentum term accelerating gradients for faster convergence. In contrast to Adam, SGD with momentum does not come with a well-working default and hence often requires hyperparameter tuning for the momentum.
- Other notable options include RMSprop (Tieleman and Hinton 2012), designed for nonstationary objectives and noisy gradients, and Nadam (Dozat 2016), which combines elements of Adam and Nesterov accelerated gradient. These optimizers are, however, typically chosen for specific applications and should be used only after careful hyperparameter tuning.

Irrespective of the choice of the optimizer, semi-structured models such as DCTM typically have an imbalance in their optimization dynamic when including deep neural networks due to the large difference in the number of parameters for structured and unstructured model components. This can, in particular, lead to slow convergence of the structured model part. To mitigate this problem, users can use warm-starts as described in Appendix C, or use optimizers with different learning rates for the different model components as described in Section 3.4.

#### G. Alternative formula interface

Following ONTRAMS (ordinal neural network transformation models), introduced in Kook & Herzog et al. (2022), deeptrafo offers an alternative formula interface. Here, the user supplies a separate formula for the intercepts (before: interacting) and for the shift (before: shifting) and avoids using the pipe | on the left-hand-side of the formula. Internally, the formula is translated back into the form in (2). All other functionalities in the article carry over to ONTRAMS as well. The same interface for other than ordinal responses is implemented in dctm().

```
R> dord <- data.frame(Y = ordered(sample.int(6, 100, TRUE)),
+ X = rnorm(100), Z = rnorm(100))
R> ontram(response = ~ Y, intercept = ~ X, shift = ~ 0 + s(Z, df = 3),
+ data = dord)
```

Untrained ordinal outcome deep conditional transformation model

## H. Large factor models

We consider a large factor model with  $10^6$  observations and a factor variable with  $10^3$  levels. The standard implementation of lm() and lmNN() fail to process the data, due to evaluating the large model matrix. However, we can use  $fac_processor()$  from safareg to circumvent this issue and use mini-batch stochastic gradient descent to fit the model on a standard machine. Now, deeptrafo can fit large factor models for arbitrary types of responses and censoring.

```
R> set.seed(0)
R> library("safareg")
R> n <- 1e6
R> nlevs <- 1e3
R> X <- factor(sample.int(nlevs, n, TRUE))</pre>
R > Y < -(X == 2) - (X == 3) + rnorm(n)
R > d < -data.frame(Y = Y, X = X)
R>m \leftarrow LmNN(Y \sim 0 + fac(X), data = d, additional_processor = list(
     fac = fac_processor))
R> fit(m, batch_size = 1e4, epochs = 20, validation_split = 0,
     callbacks = list(callback_early_stopping("loss", patience = 3),
     callback_reduce_lr_on_plateau("loss", 0.9, 2)))
R> bl <- unlist(coef(m, which = "interacting"))</pre>
R > - (unlist(coef(m))[1:5] + bl[1]) / bl[2]
fac(X)1 fac(X)2 fac(X)3 fac(X)4 fac(X)5
-0.0204 0.9986 -1.0156 -0.0249 0.0477
```

To compute the log-likelihood in models with vast amounts of data, specifying batch-wise computation avoids memory issues.

```
R> logLik(m, batch_size = 1e4)
[1] -1.42
```

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Submitted: 2024-01-17

 $Accepted \hbox{: } 2024\hbox{-}05\hbox{-}28$