

# Package ‘RemixAutoML’

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**Title** Remix Automated Machine Learning

**Version** 0.5.0

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**Maintainer** Adrian Antico <adrianantico@gmail.com>

**Description** Automates and ensures high quality output for most of your machine learning and data science tasks. The package contains high quality functions that run at efficient speed with minimal memory constraints for supervised learning, unsupervised learning, feature engineering, model evaluation and interpretation, along with some helper functions for graphing. `AutoCatBoostClassifier()`, `AutoCatBoostRegression()`, and `AutoCatBoostMultiClass()` have a dependency to the `catboost` package which isn't part of the CRAN repository at the time of this writing. The link to the `catboost` URL to download the package for use is in the `Additional_repositories` field below, which has the installation instructions. You need to install that package to make use of the `AutoCatBoost_` functions.

**License** MPL-2.0

**URL** <https://github.com/AdrianAntico/RemixAutoML>

**BugReports** <https://github.com/AdrianAntico/RemixAutoML/issues>

**Depends** R (≥ 3.5.0)

**Imports** `catboost`, `caTools`, `data.table`, `doParallel`, `foreach`, `forecast`, `ggplot2`, `grid`, `h2o`, `itertools`, `lubridate`, `Matrix`, `magick`, `methods`, `monreg`, `nortest`, `parallel`, `pROC`, `RColorBrewer`, `recommenderlab`, `ROCR`, `scatterplot3d`, `stats`, `stringr`, `tm`, `tsoutliers`, `utils`, `wordcloud`, `xgboost`, `zoo`

**Suggests** `knitr`, `rmarkdown`, `sde`, `testthat`

**VignetteBuilder** `knitr`

**Additional\_repositories**

<https://github.com/catboost/catboost/tree/master/catboost/R-package>

**Contact** Adrian Antico

**Encoding** UTF-8

**Language** en-US

**LazyData** true

**NeedsCompilation** no

**RoxygenNote** 6.1.1

**SystemRequirements** Java ( $i=7.0$ )

**Author** Adrian Antico [aut, cre],  
Douglas Pestana [aut]

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AutoCatBoostCARMA	<i>AutoCatBoostCARMA Automated CatBoost Calendar, ARMA, and Trend Variables Forecasting</i>
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## Description

AutoCatBoostCARMA Automated CatBoost Calendar, ARMA, and Trend Variables Forecasting. Create hundreds of thousands of time series forecasts using this function.

## Usage

```
AutoCatBoostCARMA(data, TargetColumnName = "Target",
  DateColumnName = "DateTime", GroupVariables = NULL,
  FC_Periods = 30, TimeUnit = "week", TargetTransformation = FALSE,
  Lags = c(1:5), MA_Periods = c(1:5), CalendarVariables = FALSE,
  TimeTrendVariable = FALSE, DataTruncate = FALSE,
  SplitRatios = c(0.7, 0.2, 0.1), TaskType = "GPU",
  EvalMetric = "MAPE", GridTune = FALSE, GridEvalMetric = "mape",
  ModelCount = 1, NTrees = 1000, PartitionType = "timeseries",
  Timer = TRUE)
```

## Arguments

<code>data</code>	Supply your full series data set here
<code>TargetColumnName</code>	List the column name of your target variables column. E.g. "Target"
<code>DateColumnName</code>	List the column name of your date column. E.g. "DateTime"
<code>GroupVariables</code>	Defaults to NULL. Use NULL when you have a single series. Add in GroupVariables when you have a series for every level of a group or multiple groups.
<code>FC_Periods</code>	Set the number of periods you want to have forecasts for. E.g. 52 for weekly data to forecast a year ahead

TimeUnit	List the time unit your data is aggregated by. E.g. "hour", "day", "week", "year"
TargetTransformation	Run AutoTransformationCreate() to find best transformation for the target variable. Tests YeoJohnson, BoxCox, and Asigh (also Asin and Logit for proportion target variables).
Lags	Select the periods for all lag variables you want to create. E.g. c(1:5,52)
MA_Periods	Select the periods for all moving average variables you want to create. E.g. c(1:5,52)
CalendarVariables	Set to TRUE to have calendar variables created. The calendar variables are numeric representations of second, minute, hour, week day, month day, year day, week, isoweek, quarter, and year
TimeTrendVariable	Set to TRUE to have a time trend variable added to the model. Time trend is numeric variable indicating the numeric value of each record in the time series (by group). Time trend starts at 1 for the earliest point in time and increments by one for each success time point.
DataTruncate	Set to TRUE to remove records with missing values from the lags and moving average features created
SplitRatios	E.g c(0.7,0.2,0.1) for train, validation, and test sets
TaskType	Default is "GPU" but you can also set it to "CPU"
EvalMetric	Select from "RMSE", "MAE", "MAPE", "Poisson", "Quantile", "LogLin-Quantile", "Lq", "NumErrors", "SMAPE", "R2", "MSLE", "MedianAbsoluteError"
GridTune	Set to TRUE to run a grid tune
GridEvalMetric	This is the metric used to find the threshold 'poisson', 'mae', 'mape', 'mse', 'msle', 'kl', 'cs', 'r2'
ModelCount	Set the number of models to try in the grid tune
NTrees	Select the number of trees you want to have built to train the model
PartitionType	Select "random" for random data partitioning "time" for partitioning by time frames
Timer	Set to FALSE to turn off the updating print statements for progress

## Value

Returns a data.table of original series and forecasts, the catboost model objects (everything returned from AutoCatBoostRegression()), a time series forecast plot, and transformation info if you set TargetTransformation to TRUE. The time series forecast plot will plot your single series or aggregate your data to a single series and create a plot from that.

## See Also

Other Time Series: [AutoH2oDRFCARMA](#), [AutoH2oGBMCARMA](#), [AutoTS](#), [AutoXGBoostCARMA](#)

## Examples

```
Results <- AutoCatBoostCARMA(data,
                             TargetColumnName = "Weekly_Sales",
                             DateColumnName = "Date",
                             GroupVariables = c("Store", "Dept"),
                             FC_Periods = 52,
                             TimeUnit = "week",
                             TargetTransformation = FALSE,
                             Lags = c(1:5, 52),
                             MA_Periods = c(1:5, 52),
                             CalendarVariables = TRUE,
                             TimeTrendVariable = TRUE,
                             DataTruncate = FALSE,
                             SplitRatios = c(1-2*30/143, 30/143, 30/143),
                             TaskType = "GPU",
                             EvalMetric = "MAE",
                             GridTune = FALSE,
                             GridEvalMetric = "mae",
                             ModelCount = 1,
                             NTrees = 1000,
                             PartitionType = "timeseries",
                             Timer = TRUE)

Results$TimeSeriesPlot
Results$Forecast
Results$ModelInformation$...
```

---

### AutoCatBoostClassifier

*AutoCatBoostClassifier is an automated catboost model grid-tuning classifier and evaluation system*

---

## Description

AutoCatBoostClassifier is an automated modeling function that runs a variety of steps. First, a stratified sampling (by the target variable) is done to create train, validation, and test sets (if not supplied). Then, the function will run a random grid tune over N number of models and find which model is the best (a default model is always included in that set). Once the model is identified and built, several other outputs are generated: validation data with predictions (on test data), an ROC plot, evaluation plot, evaluation metrics, variable importance, partial dependence calibration plots, partial dependence calibration box plots, and column names used in model fitting. You can download the catboost package using devtools, via: `devtools::install_github('catboost/catboost', subdir = 'catboost/R-package')`

## Usage

```
AutoCatBoostClassifier(data, ValidationData = NULL, TestData = NULL,
                       TargetColumnName = NULL, FeatureColNames = NULL,
                       PrimaryDateColumn = NULL, ClassWeights = NULL, IDcols = NULL,
                       task_type = "GPU", eval_metric = "AUC", Trees = 50,
                       GridTune = FALSE, grid_eval_metric = "f", MaxModelsInGrid = 10,
                       model_path = NULL, ModelID = "FirstModel", NumOfParDepPlots = 3,
```

```
ReturnModelObjects = TRUE, SaveModelObjects = FALSE,
PassInGrid = NULL)
```

### Arguments

<b>data</b>	This is your data set for training and testing your model
<b>ValidationData</b>	This is your holdout data set used in modeling either refine your hyperparameters. Catboost using both training and validation data in the training process so you should evaluate out of sample performance with this data set.
<b>TestData</b>	This is your holdout data set. Catboost using both training and validation data in the training process so you should evaluate out of sample performance with this data set.
<b>TargetColumnName</b>	Either supply the target column name OR the column number where the target is located, but not mixed types. Note that the target column needs to be a 0 — 1 numeric variable.
<b>FeatureColNames</b>	Either supply the feature column names OR the column number where the target is located, but not mixed types. Also, not zero-indexed.
<b>PrimaryDateColumn</b>	Supply a date or datetime column for catboost to utilize time as its basis for handling categorical features, instead of random shuffling
<b>ClassWeights</b>	Supply a vector of weights for your target classes. E.g. <code>c(0.25, 1)</code> to weight your 0 class by 0.25 and your 1 class by 1.
<b>IDcols</b>	A vector of column names or column numbers to keep in your data but not include in the modeling.
<b>task_type</b>	Set to "GPU" to utilize your GPU for training. Default is "CPU".
<b>eval_metric</b>	This is the metric used inside catboost to measure performance on validation data during a grid-tune. "AUC" is the default, but other options include "Logloss", "CrossEntropy", "Precision", "Recall", "F1", "BalancedAccuracy", "BalancedErrorRate", "MCC", "Accuracy", "CtrFactor", "AUC", "BrierScore", "HingeLoss", "HammingLoss", "ZeroOneLoss", "Kappa", "WKappa", "LogLikelihoodOfPrediction"
<b>Trees</b>	The maximum number of trees you want in your models
<b>GridTune</b>	Set to TRUE to run a grid tuning procedure. Set a number in <code>MaxModelsInGrid</code> to tell the procedure how many models you want to test.
<b>grid_eval_metric</b>	This is the metric used to find the threshold "f", "auc", "tpr", "fnr", "fpr", "tnr", "prbe", "f", "odds"
<b>MaxModelsInGrid</b>	Number of models to test from grid options.
<b>model_path</b>	A character string of your path file to where you want your output saved
<b>ModelID</b>	A character string to name your model and output
<b>NumOfParDepPlots</b>	Tell the function the number of partial dependence calibration plots you want to create. Calibration boxplots will only be created for numerical features (not dummy variables)

**ReturnModelObjects**

Set to TRUE to output all modeling objects. E.g. plots and evaluation metrics

**SaveModelObjects**

Set to TRUE to return all modeling objects to your environment

**PassInGrid**

Defaults to NULL. Pass in a single row of grid from a previous output as a data.table (they are collected as data.tables)

**Value**

Saves to file and returned in list: VariableImportance.csv, Model (the model), ValidationData.csv, ROC\_Plot.png, EvaluationPlot.png, EvaluationMetrics.csv, ParDepPlots.R a named list of features with partial dependence calibration plots, GridCollect, and GridList

**Author(s)**

Adrian Antico

**See Also**

Other Supervised Learning: [AutoCatBoostMultiClass](#), [AutoCatBoostRegression](#), [AutoCatBoostScoring](#), [AutoCatBoostHurdleModel](#), [AutoH2OMLScoring](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFClassifier](#), [AutoH2oDRFMultiClass](#), [AutoH2oDRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#), [AutoXGBoostScoring](#)

**Examples**

```
Correl <- 0.85
N <- 1000
data <- data.table::data.table(Target = runif(N))
data[, x1 := qnorm(Target)]
data[, x2 := runif(N)]
data[, Independent_Variable1 := log(pnorm(Correl * x1 +
                                         sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable2 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable3 := exp(pnorm(Correl * x1 +
                                         sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable4 := exp(exp(pnorm(Correl * x1 +
                                              sqrt(1-Correl^2) * qnorm(x2))))]
data[, Independent_Variable5 := sqrt(pnorm(Correl * x1 +
                                           sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable6 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.10]
data[, Independent_Variable7 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.25]
data[, Independent_Variable8 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.75]
data[, Independent_Variable9 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^2]
data[, Independent_Variable10 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^4]
data[, Independent_Variable11 := as.factor(
  ifelse(Independent_Variable2 < 0.20, "A",
```

```

        ifelse(Independent_Variable2 < 0.40, "B",
              ifelse(Independent_Variable2 < 0.6, "C",
                    ifelse(Independent_Variable2 < 0.8, "D", "E")))))]
data[, 'x1' := (x1 = NULL, x2 = NULL)]
data[, Target := ifelse(Target < 0.5, 1, 0)]
TestModel <- AutoCatBoostClassifier(data,
                                   ValidationData = NULL,
                                   TestData = NULL,
                                   TargetColumnName = "Target",
                                   FeatureColNames = c(2:12),
                                   PrimaryDateColumn = NULL,
                                   ClassWeights = NULL,
                                   IDcols = NULL,
                                   MaxModelsInGrid = 3,
                                   task_type = "GPU",
                                   eval_metric = "AUC",
                                   grid_eval_metric = "auc",
                                   Trees = 50,
                                   GridTune = FALSE,
                                   model_path = NULL,
                                   ModelID = "ModelTest",
                                   NumOfParDepPlots = 15,
                                   ReturnModelObjects = TRUE,
                                   SaveModelObjects = FALSE,
                                   PassInGrid = NULL)

```

---

#### AutoCatBoostdHurdleModel

*AutoCatBoostdHurdleModel is a Retrain Function for the Regression Models for the Subsetted Data in P6*

---

### Description

AutoCatBoostdHurdleModel is a Retrain Function for the Regression Models for the Sub-setted Data in P6

### Usage

```

AutoCatBoostdHurdleModel(data, ValidationData = NULL, TestData = NULL,
  Buckets = c(1, 5, 10, 20), TargetColumnName = "Target",
  FeatureColNames = 4:ncol(data), PrimaryDateColumn = NULL,
  IDcols = NULL, TransformNumericColumns = NULL, ClassWeights = NULL,
  SplitRatios = c(0.7, 0.2, 0.1), task_type = "GPU",
  ModelID = "ModelTest", Paths = NULL, SaveModelObjects = TRUE,
  Trees = 15000, GridTune = TRUE, MaxModelsInGrid = 1,
  NumOfParDepPlots = 10, PassInGrid = NULL)

```

### Arguments

<b>data</b>	Source training data. Do not include a column that has the class labels for the buckets as they are created internally.
<b>ValidationData</b>	Source validation data. Do not include a column that has the class labels for the buckets as they are created internally.



TestData	Source test data. Do not include a column that has the class labels for the buckets as they are created internally.
Buckets	A numeric vector of the buckets used for subsetting the data. NOTE: the final Bucket value will first create a subset of data that is less than the value and a second one thereafter for data greater than the bucket value.
TargetColumnName	Supply the column name or number for the target variable
FeatureColNames	Supply the column names or number of the features (not included the PrimaryDateColumn)
PrimaryDateColumn	Supply a date column if the data is functionally related to it
IDcols	Includes PrimaryDateColumn and any other columns you want returned in the validation data with predictions
TransformNumericColumns	Transform numeric column inside the AutoCatBoostRegression() function
ClassWeights	Utilize these for the classifier model
SplitRatios	Supply vector of partition ratios. For example, c(0.70,0.20,0.10).
task_type	Set to "GPU" or "CPU"
ModelID	Define a character name for your models
Paths	A character vector of the path file strings. EITHER SUPPLY 1 file path or N file paths for N models
SaveModelObjects	Set to TRUE to save the model objects to file in the folders listed in Paths
Trees	Default 15000
GridTune	Set to TRUE if you want to grid tune the models
NumOfParDepPlots	Set to pull back N number of partial dependence calibration plots.
PassInGrid	Pass in a grid for changing up the parameter settings for catboost
NumberModelsInGrid	Set to a numeric value for the number of models to try in grid tune

## Value

Returns AutoCatBoostRegression() model objects: VariableImportance.csv, Model, ValidationData.csv, EvaluationPlot.png, EvaluationBoxPlot.png, EvaluationMetrics.csv, ParDepPlots.R a named list of features with partial dependence calibration plots, ParDepBoxPlots.R, GridCollect, and catboostgrid

## See Also

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoCatBoostRegression](#), [AutoCatBoostScoring](#), [AutoH2OMLScoring](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFClassifier](#), [AutoH2oDRFMultiClass](#), [AutoH2oDRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#), [AutoXGBoostScoring](#)

## Examples

```
Output <- RemixAutoML::AutoCatBoostHurdleModel(
  data,
  ValidationData = NULL,
  TestData = NULL,
  Buckets = c(1, 5, 10, 20),
  TargetColumnName = "PLND_LABOR_UNITS",
  FeatureColNames = 4:ncol(data),
  PrimaryDateColumn = "PLND_STRT_DT",
  IDcols = c(1,3),
  TransformNumericColumns = NULL,
  ClassWeights = NULL,
  SplitRatios = c(0.7, 0.2, 0.1),
  task_type = "GPU",
  ModelID = "P6",
  Paths = c(paste0(getwd()),"/P6-Buckets")),
  SaveModelObjects = TRUE,
  Trees = 5000,
  GridTune = FALSE,
  MaxModelsInGrid = 1,
  NumOfParDepPlots = 10,
  PassInGrid = grid)
```

---

### AutoCatBoostMultiClass

*AutoCatBoostMultiClass is an automated catboost model grid-tuning multinomial classifier and evaluation system*

---

## Description

AutoCatBoostMultiClass is an automated modeling function that runs a variety of steps. First, a stratified sampling (by the target variable) is done to create train and validation sets. Then, the function will run a random grid tune over N number of models and find which model is the best (a default model is always included in that set). Once the model is identified and built, several other outputs are generated: validation data with predictions, evaluation metrics, variable importance, and column names used in model fitting. You can download the catboost package using devtools, via: `devtools::install_github('catboost/catboost', subdir = 'catboost/R-package')`.

## Usage

```
AutoCatBoostMultiClass(data, ValidationData = NULL, TestData = NULL,
  TargetColumnName = NULL, FeatureColNames = NULL,
  PrimaryDateColumn = NULL, ClassWeights = NULL, IDcols = NULL,
  task_type = "GPU", eval_metric = "MultiClassOneVsAll", Trees = 50,
  GridTune = FALSE, grid_eval_metric = "Accuracy",
  MaxModelsInGrid = 10, model_path = NULL, ModelID = "FirstModel",
  ReturnModelObjects = TRUE, SaveModelObjects = FALSE,
  PassInGrid = NULL)
```

**Arguments**

<code>data</code>	This is your data set for training and testing your model
<code>ValidationData</code>	This is your holdout data set used in modeling either refine your hyperparameters. Catboost using both training and validation data in the training process so you should evaluate out of sample performance with this data set.
<code>TestData</code>	This is your holdout data set. Catboost using both training and validation data in the training process so you should evaluate out of sample performance with this data set.
<code>TargetColumnName</code>	Either supply the target column name OR the column number where the target is located, but not mixed types.
<code>FeatureColNames</code>	Either supply the feature column names OR the column number where the target is located, but not mixed types. Also, not zero-indexed.
<code>PrimaryDateColumn</code>	Supply a date or datetime column for catboost to utilize time as its basis for handling categorical features, instead of random shuffling
<code>ClassWeights</code>	Supply a vector of weights for your target classes. E.g. <code>c(0.25, 1)</code> to weight your 0 class by 0.25 and your 1 class by 1.
<code>IDcols</code>	A vector of column names or column numbers to keep in your data but not include in the modeling.
<code>task_type</code>	Set to "GPU" to utilize your GPU for training. Default is "CPU".
<code>eval_metric</code>	This is the metric used inside catboost to measure performance on validation data during a grid-tune. "MultiClass" or "MultiClassOneVsAll"
<code>Trees</code>	The maximum number of trees you want in your models
<code>GridTune</code>	Set to TRUE to run a grid tuning procedure. Set a number in <code>MaxModelsInGrid</code> to tell the procedure how many models you want to test.
<code>grid_eval_metric</code>	This is the metric used to find the threshold "auc", "accuracy"
<code>MaxModelsInGrid</code>	Number of models to test from grid options. 1080 total possible options
<code>model_path</code>	A character string of your path file to where you want your output saved
<code>ModelID</code>	A character string to name your model and output
<code>ReturnModelObjects</code>	Set to TRUE to output all modeling objects. E.g. plots and evaluation metrics
<code>SaveModelObjects</code>	Set to TRUE to return all modeling objects to your environment
<code>PassInGrid</code>	Defaults to NULL. Pass in a single row of grid from a previous output as a data.table (they are collected as data.tables)

**Value**

Saves to file and returned in list: `VariableImportance.csv`, `Model` (the model), `ValidationData.csv`, `EvaluationMetrics.csv`, `GridCollect`, and `GridList`

**Author(s)**

Adrian Antico

**See Also**

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostRegression](#), [AutoCatBoostScoring](#), [AutoCatBoostHurdleModel](#), [AutoH2OMLScoring](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFClassifier](#), [AutoH2oDRFMultiClass](#), [AutoH2oDRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#), [AutoXGBoostScoring](#)

**Examples**

```
Correl <- 0.85
N <- 1000
data <- data.table::data.table(Target = runif(N))
data[, x1 := qnorm(Target)]
data[, x2 := runif(N)]
data[, Independent_Variable1 := log(pnorm(Correl * x1 +
                                         sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable2 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable3 := exp(pnorm(Correl * x1 +
                                         sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable4 := exp(exp(pnorm(Correl * x1 +
                                             sqrt(1-Correl^2) * qnorm(x2))))]
data[, Independent_Variable5 := sqrt(pnorm(Correl * x1 +
                                           sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable6 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.10]
data[, Independent_Variable7 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.25]
data[, Independent_Variable8 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.75]
data[, Independent_Variable9 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^2]
data[, Independent_Variable10 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^4]
data[, Target := as.factor(
  ifelse(Independent_Variable2 < 0.20, "A",
    ifelse(Independent_Variable2 < 0.40, "B",
      ifelse(Independent_Variable2 < 0.6, "C",
        ifelse(Independent_Variable2 < 0.8, "D", "E"))))]
data[, ':= ' (x1 = NULL, x2 = NULL)]
TestModel <- AutoCatBoostMultiClass(data,
  ValidationData = NULL,
  TestData = NULL,
  TargetColumnName = "Target",
  FeatureColNames = c(2:11),
  PrimaryDateColumn = NULL,
  ClassWeights = NULL,
  IDcols = NULL,
  MaxModelsInGrid = 1,
  task_type = "GPU",
  eval_metric = "MultiClass",
```

```

grid_eval_metric = "Accuracy",
Trees = 50,
GridTune = FALSE,
model_path = NULL,
ModelID = "ModelTest",
ReturnModelObjects = TRUE,
SaveModelObjects = FALSE,
PassInGrid = NULL)

```

---

## AutoCatBoostRegression

*AutoCatBoostRegression is an automated catboost model grid-tuning classifier and evaluation system*

---

## Description

AutoCatBoostRegression is an automated modeling function that runs a variety of steps. First, the function will run a random grid tune over N number of models and find which model is the best (a default model is always included in that set). Once the model is identified and built, several other outputs are generated: validation data with predictions, evaluation plot, evaluation boxplot, evaluation metrics, variable importance, partial dependence calibration plots, partial dependence calibration box plots, and column names used in model fitting. You can download the catboost package using devtools, via: `devtools::install_github('catboost/catboost', subdir = 'catboost/R-package')`

## Usage

```

AutoCatBoostRegression(data, ValidationData = NULL, TestData = NULL,
  TargetColumnName = NULL, FeatureColNames = NULL,
  PrimaryDateColumn = NULL, IDcols = NULL,
  TransformNumericColumns = NULL, task_type = "GPU",
  eval_metric = "RMSE", Alpha = NULL, Trees = 50, GridTune = FALSE,
  grid_eval_metric = "mae", MaxModelsInGrid = 10, model_path = NULL,
  ModelID = "FirstModel", NumOfParDepPlots = 3,
  ReturnModelObjects = TRUE, SaveModelObjects = FALSE,
  PassInGrid = NULL)

```

## Arguments

<b>data</b>	This is your data set for training and testing your model
<b>ValidationData</b>	This is your holdout data set used in modeling either refine your hyperparameters. Catboost using both training and validation data in the training process so you should evaluate out of sample performance with this data set.
<b>TestData</b>	This is your holdout data set. Catboost using both training and validation data in the training process so you should evaluate out of sample performance with this data set.
<b>TargetColumnName</b>	Either supply the target column name OR the column number where the target is located (but not mixed types).

<b>FeatureColNames</b>	Either supply the feature column names OR the column number where the target is located (but not mixed types)
<b>PrimaryDateColumn</b>	Supply a date or datetime column for catboost to utilize time as its basis for handling categorical features, instead of random shuffling
<b>IDcols</b>	A vector of column names or column numbers to keep in your data but not include in the modeling.
<b>TransformNumericColumns</b>	Set to NULL to do nothing; otherwise supply the column names of numeric variables you want transformed
<b>task_type</b>	Set to "GPU" to utilize your GPU for training. Default is "CPU".
<b>eval_metric</b>	This is the metric used inside catboost to measure performance on validation data during a grid-tune. "RMSE" is the default, but other options include: "MAE", "MAPE", "Poisson", "Quantile", "LogLinQuantile", "Lq", "NumErrors", "SMAPE", "R2", "MSLE", "MedianAbsoluteError".
<b>Alpha</b>	This is the quantile value you want to use for quantile regression. Must be a decimal between 0 and 1.
<b>Trees</b>	The maximum number of trees you want in your models
<b>GridTune</b>	Set to TRUE to run a grid tuning procedure. Set a number in MaxModelsInGrid to tell the procedure how many models you want to test.
<b>grid_eval_metric</b>	This is the metric used to find the threshold 'poisson', 'mae', 'mape', 'mse', 'msle', 'kl', 'cs', 'r2'
<b>MaxModelsInGrid</b>	Number of models to test from grid options (1080 total possible options)
<b>model_path</b>	A character string of your path file to where you want your output saved
<b>ModelID</b>	A character string to name your model and output
<b>NumOfParDepPlots</b>	Tell the function the number of partial dependence calibration plots you want to create. Calibration boxplots will only be created for numerical features (not dummy variables)
<b>ReturnModelObjects</b>	Set to TRUE to output all modeling objects (E.g. plots and evaluation metrics)
<b>SaveModelObjects</b>	Set to TRUE to return all modeling objects to your environment
<b>PassInGrid</b>	Defaults to NULL. Pass in a single row of grid from a previous output as a data.table (they are collected as data.tables)

## Value

Saves to file and returned in list: VariableImportance.csv, Model, ValidationData.csv, EvaluationPlot.png, EvaluationBoxPlot.png, EvaluationMetrics.csv, ParDepPlots.R a named list of features with partial dependence calibration plots, ParDepBoxPlots.R, GridCollect, catboostgrid, and a transformation details file.

**Author(s)**

Adrian Antico

**See Also**

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoCatBoostScoring](#), [AutoCatBoostHurdleModel](#), [AutoH2OMLScoring](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFClassifier](#), [AutoH2oDRFMultiClass](#), [AutoH2oDRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#), [AutoXGBoostScoring](#)

**Examples**

```
Correl <- 0.85
N <- 1000
data <- data.table::data.table(Target = runif(N))
data[, x1 := qnorm(Target)]
data[, x2 := runif(N)]
data[, Independent_Variable1 := log(pnorm(Correl * x1 +
                                         sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable2 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable3 := exp(pnorm(Correl * x1 +
                                         sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable4 := exp(exp(pnorm(Correl * x1 +
                                              sqrt(1-Correl^2) * qnorm(x2))))]
data[, Independent_Variable5 := sqrt(pnorm(Correl * x1 +
                                          sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable6 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.10]
data[, Independent_Variable7 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.25]
data[, Independent_Variable8 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.75]
data[, Independent_Variable9 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^2]
data[, Independent_Variable10 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^4]
data[, Independent_Variable11 := as.factor(
  ifelse(Independent_Variable2 < 0.20, "A",
    ifelse(Independent_Variable2 < 0.40, "B",
      ifelse(Independent_Variable2 < 0.6, "C",
        ifelse(Independent_Variable2 < 0.8, "D", "E"))))]
data[, ':= ' (x1 = NULL, x2 = NULL)]
TestModel <- AutoCatBoostRegression(data,
  ValidationData = NULL,
  TestData = NULL,
  TargetColumnName = "Target",
  FeatureColNames = c(2:12),
  PrimaryDateColumn = NULL,
  IDcols = NULL,
  TransformNumericColumns = NULL,
  MaxModelsInGrid = 1,
  task_type = "GPU",
  eval_metric = "RMSE",
```

```
Alpha = NULL,
grid_eval_metric = "r2",
Trees = 50,
GridTune = FALSE,
model_path = NULL,
ModelID = "ModelTest",
NumOfParDepPlots = 3,
ReturnModelObjects = TRUE,
SaveModelObjects = FALSE,
PassInGrid = NULL)
```

---

AutoCatBoostScoring	<i>AutoCatBoostScoring is an automated scoring function that compliments the AutoCatBoost model training functions.</i>
---------------------	---

---

## Description

AutoCatBoostScoring is an automated scoring function that compliments the AutoCatBoost model training functions. This function requires you to supply features for scoring. It will run ModelDataPrep() to prepare your features for catboost data conversion and scoring.

## Usage

```
AutoCatBoostScoring(TargetType = NULL, ScoringData = NULL,
  FeatureColumnNames = NULL, IDcols = NULL, ModelObject = NULL,
  ModelPath = NULL, ModelID = NULL, ReturnFeatures = TRUE,
  TransformNumeric = FALSE, BackTransNumeric = FALSE,
  TargetColumnName = NULL, TransformationObject = NULL,
  TransID = NULL, TransPath = NULL, MDP_Impute = TRUE,
  MDP_CharToFactor = TRUE, MDP_RemoveDates = TRUE,
  MDP_MissFactor = "0", MDP_MissNum = -1)
```

## Arguments

TargetType	Set this value to "regression", "classification", or "multiclass" to score models built using AutoCatBoostRegression(), AutoCatBoostClassify() or AutoCatBoostMultiClass().
ScoringData	This is your data.table of features for scoring. Can be a single row or batch.
FeatureColumnNames	Supply either column names or column numbers used in the AutoCatBoostRegression() function
IDcols	Supply ID column numbers for any metadata you want returned with your predicted values
ModelObject	Supply the model object directly for scoring instead of loading it from file. If you supply this, ModelID and ModelPath will be ignored.
ModelPath	Supply your path file used in the AutoCatBoost__() function
ModelID	Supply the model ID used in the AutoCatBoost__() function



<b>ReturnFeatures</b>	Set to TRUE to return your features with the predicted values.
<b>TransformNumeric</b>	Set to TRUE if you have features that were transformed automatically from an <code>Auto__Regression()</code> model AND you haven't already transformed them.
<b>BackTransNumeric</b>	Set to TRUE to generate back-transformed predicted values. Also, if you return features, those will also be back-transformed.
<b>TargetColumnName</b>	Input your target column name used in training if you are utilizing the transformation service
<b>TransformationObject</b>	Set to NULL if you didn't use transformations or if you want the function to pull from the file output from the <code>Auto__Regression()</code> function. You can also supply the transformation <code>data.table</code> object with the transformation details versus having it pulled from file.
<b>TransID</b>	Set to the ID used for saving the transformation <code>data.table</code> object or set it to the <code>ModelID</code> if you are pulling from file from a build with <code>Auto__Regression()</code> .
<b>TransPath</b>	Set the path file to the folder where your transformation <code>data.table</code> detail object is stored. If you used the <code>Auto__Regression()</code> to build, set it to the same path as <code>ModelPath</code> .
<b>MDP_Impute</b>	Set to TRUE if you did so for modeling and didn't do so before supplying <code>ScoringData</code> in this function
<b>MDP_CharToFactor</b>	Set to TRUE to turn your character columns to factors if you didn't do so to your <code>ScoringData</code> that you are supplying to this function
<b>MDP_RemoveDates</b>	Set to TRUE if you have date of timestamp columns in your <code>ScoringData</code>
<b>MDP_MissFactor</b>	If you set <code>MDP_Impute</code> to TRUE, supply the character values to replace missing values with
<b>MDP_MissNum</b>	If you set <code>MDP_Impute</code> to TRUE, supply a numeric value to replace missing values with

## Value

A `data.table` of predicted values with the option to return model features as well.

## See Also

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoCatBoostRegression](#), [AutoCatBoostHurdleModel](#), [AutoH2OMLScoring](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFClassifier](#), [AutoH2oDRFMultiClass](#), [AutoH2oDRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#), [AutoXGBoostScoring](#)

## Examples

```
Preds <- AutoCatBoostScoring(TargetType = "regression",
                             ScoringData = data,
```

```

FeatureColumnNames = 2:12,
IDcols = NULL,
ModelObject = NULL,
ModelPath = "home",
ModelID = "ModelTest",
ReturnFeatures = TRUE,
TransformNumeric = FALSE,
BackTransNumeric = FALSE,
TargetColumnName = NULL,
TransformationObject = NULL,
TransID = NULL,
TransPath = NULL,
MDP_Impute = TRUE,
MDP_CharToFactor = TRUE,
MDP_RemoveDates = TRUE,
MDP_MissFactor = "0",
MDP_MissNum = -1)

```

---

AutoDataPartition

*The AutoDataPartition function*


---

## Description

This function will take your ratings matrix and model and score your data in parallel.

## Usage

```

AutoDataPartition(data, NumDataSets = 3, Ratios = c(0.7, 0.2, 0.1),
  PartitionType = "random", StratifyColumnNames = NULL,
  StratifyNumericTarget = NULL, StratTargetPrecision = 3,
  TimeColumnName = NULL)

```

## Arguments

- |                            |  |
|----------------------------|--|
| <b>data</b>                | Source data to do your partitioning on   |
| <b>NumDataSets</b>         | The number of total data sets you want built   |
| <b>Ratios</b>              | A vector of values for how much data each data set should get in each split. E.g. c(0.70, 0.20, 0.10)  |
| <b>PartitionType</b>       | Set to either "random", "timeseries", or "time". With "random", your data will be partitioned randomly (with stratified sampling if column names are supplied). With "timeseries", you can partition by time with a stratify option (so long as you have an equal number of records for each strata). With "time" you will have data sets generated so that the training data contains the earliest records in time, validation data the second earliest, test data the third earliest, etc. |
| <b>StratifyColumnNames</b> | Supply column names of categorical features to use in a stratified sampling procedure for partitioning the data. Partition type must be "random" to use this option  |

**StratifyNumericTarget**

Supply a column name that is numeric. Use for "random" PartitionType, you can stratify your numeric variable by splitting up based on percRank to ensure a proper allocation of extreme values in your created data sets.

**StratTargetPrecision**

For "random" PartitionType and when StratifyNumericTarget is not null, precision will be the number of decimals used in the percentile calculation. If you supply a value of 1, deciles will be used. For a value of 2, percentiles will be used. Larger values are supported.

**TimeColumnName** Supply a date column name or a name of a column with an ID for sorting by time such that the smallest number is the earliest in time.

**Value**

Returns a list of data.tables

**Author(s)**

Adrian Antico and Douglas Pestana

**See Also**

Other Feature Engineering: [AutoTransformationCreate](#), [AutoTransformationScore](#), [AutoWord2VecModeler](#), [CreateCalendarVariables](#), [DT.GDL.Feature\\_Engineering](#), [DummifyDT](#), [GDL.Feature\\_Engineering](#), [ModelDataPrep](#), [Partial\\_DT.GDL.Feature\\_Engineering2](#), [Partial\\_DT.GDL.Feature\\_Engineering](#), [Scoring.GDL.Feature\\_Engineering](#)

**Examples**

```
dataSets <- AutoDataPartition(data,
                              NumDataSets = 3,
                              Ratios = c(0.70,0.20,0.10),
                              PartitionType = "random",
                              StratifyColumnNames = NULL,
                              StratifyNumericTarget = NULL,
                              StratTargetPrecision = 1,
                              TimeColumnName = NULL)
```

---

AutoH2oDRFCARMA

*AutoH2oDRFCARMA Automated CatBoost Calendar, ARMA, and Trend Variables Forecasting*

---

**Description**

AutoH2oDRFCARMA Automated CatBoost Calendar, ARMA, and Trend Variables Forecasting. Create hundreds of thousands of time series forecasts using this function.

**Usage**

```
AutoH2oDRFCARMA(data, TargetColumnName = "Target",
  DateColumnName = "DateTime", GroupVariables = NULL,
  FC_Periods = 30, TimeUnit = "week", TargetTransformation = FALSE,
  Lags = c(1:5), MA_Periods = c(1:5), CalendarVariables = FALSE,
  TimeTrendVariable = FALSE, DataTruncate = FALSE,
  SplitRatios = c(0.7, 0.2, 0.1), EvalMetric = "MAE",
  GridTune = FALSE, ModelCount = 1, NTrees = 1000,
  PartitionType = "timeseries", MaxMem = "32G", NThreads = max(1,
  parallel::detectCores() - 2), Timer = TRUE)
```

**Arguments**

<b>data</b>	Supply your full series data set here
<b>TargetColumnName</b>	List the column name of your target variables column. E.g. "Target"
<b>DateColumnName</b>	List the column name of your date column. E.g. "DateTime"
<b>GroupVariables</b>	Defaults to NULL. Use NULL when you have a single series. Add in GroupVariables when you have a series for every level of a group or multiple groups.
<b>FC_Periods</b>	Set the number of periods you want to have forecasts for. E.g. 52 for weekly data to forecast a year ahead
<b>TimeUnit</b>	List the time unit your data is aggregated by. E.g. "hour", "day", "week", "year"
<b>TargetTransformation</b>	Run AutoTransformationCreate() to find best transformation for the target variable. Tests YeoJohnson, BoxCox, and Asigh (also Asin and Logit for proportion target variables).
<b>Lags</b>	Select the periods for all lag variables you want to create. E.g. c(1:5,52)
<b>MA_Periods</b>	Select the periods for all moving average variables you want to create. E.g. c(1:5,52)
<b>CalendarVariables</b>	Set to TRUE to have calendar variables created. The calendar variables are numeric representations of second, minute, hour, week day, month day, year day, week, isoweek, quarter, and year
<b>TimeTrendVariable</b>	Set to TRUE to have a time trend variable added to the model. Time trend is numeric variable indicating the numeric value of each record in the time series (by group). Time trend starts at 1 for the earliest point in time and increments by one for each success time point.
<b>DataTruncate</b>	Set to TRUE to remove records with missing values from the lags and moving average features created
<b>SplitRatios</b>	E.g c(0.7,0.2,0.1) for train, validation, and test sets
<b>EvalMetric</b>	Select from "RMSE", "MAE", "MAPE", "Poisson", "Quantile", "LogLin-Quantile", "Lq", "NumErrors", "SMAPE", "R2", "MSLE", "MedianAbsoluteError"
<b>GridTune</b>	Set to TRUE to run a grid tune
<b>ModelCount</b>	Set the number of models to try in the grid tune

NTrees	Select the number of trees you want to have built to train the model
PartitionType	Select "random" for random data partitioning "time" for partitioning by time frames
MaxMem	Set to the maximum amount of memory you want to allow for running this function. Default is "32G".
NThreads	Set to the number of threads you want to dedicate to this function.
Timer	Set to FALSE to turn off the updating print statements for progress

### Value

Returns a data.table of original series and forecasts, the catboost model objects (everything returned from AutoCatBoostRegression()), a time series forecast plot, and transformation info if you set TargetTransformation to TRUE. The time series forecast plot will plot your single series or aggregate your data to a single series and create a plot from that.

### See Also

Other Time Series: [AutoCatBoostCARMA](#), [AutoH2oGBMCARMA](#), [AutoTS](#), [AutoXGBoostCARMA](#)

### Examples

```
Results <- AutoH2oDRFCARMA(data,
  TargetColumnName = "Weekly_Sales",
  DateColumnName = "Date",
  GroupVariables = c("Store", "Dept"),
  FC_Periods = 52,
  TimeUnit = "week",
  TargetTransformation = FALSE,
  Lags = c(1:5, 52),
  MA_Periods = c(1:5, 52),
  CalendarVariables = TRUE,
  TimeTrendVariable = TRUE,
  DataTruncate = FALSE,
  SplitRatios = c(1-2*30/143, 30/143, 30/143),
  EvalMetric = "MAE",
  GridTune = FALSE,
  ModelCount = 1,
  NTrees = 1000,
  PartitionType = "timeseries",
  MaxMem = "32G",
  NThreads = max(1, parallel::detectCores() - 2),
  Timer = TRUE)

Results$TimeSeriesPlot
Results$Forecast
Results$ModelInformation$...
```

---

AutoH2oDRFClassifier	<i>AutoH2oDRFClassifier is an automated H2O modeling framework with grid-tuning and model evaluation</i>
----------------------	--

---

## Description

AutoH2oDRFClassifier is an automated H2O modeling framework with grid-tuning and model evaluation that runs a variety of steps. First, a stratified sampling (by the target variable) is done to create train and validation sets. Then, the function will run a random grid tune over N number of models and find which model is the best (a default model is always included in that set). Once the model is identified and built, several other outputs are generated: validation data with predictions, evaluation plot, evaluation metrics, variable importance, partial dependence calibration plots, and column names used in model fitting.

## Usage

```
AutoH2oDRFClassifier(data, ValidationData = NULL, TestData = NULL,
  TargetColumnName = NULL, FeatureColNames = NULL,
  eval_metric = "auc", Trees = 50, GridTune = FALSE,
  MaxMem = "32G", MaxModelsInGrid = 2, model_path = NULL,
  ModelID = "FirstModel", NumOfParDepPlots = 3,
  ReturnModelObjects = TRUE, SaveModelObjects = FALSE,
  IfSaveModel = "mojo")
```

## Arguments

<code>data</code>	This is your data set for training and testing your model
<code>ValidationData</code>	This is your holdout data set used in modeling either refine your hyper-parameters.
<code>TestData</code>	This is your holdout data set. Catboost using both training and validation data in the training process so you should evaluate out of sample performance with this data set.
<code>TargetColumnName</code>	Either supply the target column name OR the column number where the target is located (but not mixed types). Note that the target column needs to be a 0 — 1 numeric variable.
<code>FeatureColNames</code>	Either supply the feature column names OR the column number where the target is located (but not mixed types)
<code>eval_metric</code>	This is the metric used to identify best grid tuned model. Choose from "AUC" or "logloss"
<code>Trees</code>	The maximum number of trees you want in your models
<code>GridTune</code>	Set to TRUE to run a grid tuning procedure. Set a number in MaxModelsInGrid to tell the procedure how many models you want to test.
<code>MaxMem</code>	Set the maximum amount of memory you'd like to dedicate to the model run. E.g. "32G"
<code>MaxModelsInGrid</code>	Number of models to test from grid options (1080 total possible options)
<code>model_path</code>	A character string of your path file to where you want your output saved

ModelID	A character string to name your model and output
NumOfParDepPlots	Tell the function the number of partial dependence calibration plots you want to create.
ReturnModelObjects	Set to TRUE to output all modeling objects (E.g. plots and evaluation metrics)
SaveModelObjects	Set to TRUE to return all modeling objects to your environment
IfSaveModel	Set to "mojo" to save a mojo file, otherwise "standard" to save a regular H2O model object

### Value

Saves to file and returned in list: VariableImportance.csv, Model, ValidationData.csv, EvaluationPlot.png, EvaluationMetrics.csv, ParDepPlots.R a named list of features with partial dependence calibration plots, GridCollect, and GridList

### Author(s)

Adrian Antico

### See Also

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoCatBoostRegression](#), [AutoCatBoostScoring](#), [AutoCatBoostHurdleModel](#), [AutoH2OMLScoring](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFMultiClass](#), [AutoH2oDRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#), [AutoXGBoostScoring](#)

### Examples

```
Correl <- 0.85
N <- 1000
data <- data.table::data.table(Target = runif(N))
data[, x1 := qnorm(Target)]
data[, x2 := runif(N)]
data[, Independent_Variable1 := log(pnorm(Correl * x1 +
                                          sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable2 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable3 := exp(pnorm(Correl * x1 +
                                         sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable4 := exp(exp(pnorm(Correl * x1 +
                                              sqrt(1-Correl^2) * qnorm(x2))))]
data[, Independent_Variable5 := sqrt(pnorm(Correl * x1 +
                                           sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable6 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.10]
data[, Independent_Variable7 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.25]
data[, Independent_Variable8 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.75]
data[, Independent_Variable9 := (pnorm(Correl * x1 +
```

```

                                sqrt(1-Correl^2) * qnorm(x2)))^2]
data[, Independent_Variable10 := (pnorm(Correl * x1 +
                                sqrt(1-Correl^2) * qnorm(x2)))^4]
data[, Independent_Variable11 := as.factor(
  ifelse(Independent_Variable2 < 0.20, "A",
    ifelse(Independent_Variable2 < 0.40, "B",
      ifelse(Independent_Variable2 < 0.6, "C",
        ifelse(Independent_Variable2 < 0.8, "D", "E")))))]
data[, ':= ' (x1 = NULL, x2 = NULL)]
data[, Target := as.factor(ifelse(Independent_Variable2 < 0.5, 1, 0))]
TestModel <- AutoH2oDRFClassifier(data,
  ValidationData = NULL,
  TestData = NULL,
  TargetColumnName = "Target",
  FeatureColNames = 2:ncol(data),
  eval_metric = "auc",
  Trees = 50,
  GridTune = FALSE,
  MaxMem = "32G",
  MaxModelsInGrid = 10,
  model_path = NULL,
  ModelID = "FirstModel",
  NumOfParDepPlots = 3,
  ReturnModelObjects = TRUE,
  SaveModelObjects = FALSE,
  IfSaveModel = "mojo")

```

---

AutoH2oDRFMultiClass    *AutoH2oDRFMultiClass is an automated H2O modeling framework with grid-tuning and model evaluation*

---

## Description

AutoH2oDRFMultiClass is an automated H2O modeling framework with grid-tuning and model evaluation that runs a variety of steps. First, a stratified sampling (by the target variable) is done to create train and validation sets. Then, the function will run a random grid tune over N number of models and find which model is the best (a default model is always included in that set). Once the model is identified and built, several other outputs are generated: validation data with predictions, evaluation metrics, confusion matrix, and variable importance.

## Usage

```

AutoH2oDRFMultiClass(data, ValidationData = NULL, TestData = NULL,
  TargetColumnName = NULL, FeatureColNames = NULL,
  eval_metric = "logloss", Trees = 50, GridTune = FALSE,
  MaxMem = "32G", MaxModelsInGrid = 2, model_path = NULL,
  ModelID = "FirstModel", ReturnModelObjects = TRUE,
  SaveModelObjects = FALSE, IfSaveModel = "mojo")

```



**Arguments**

<code>data</code>	This is your data set for training and testing your model
<code>ValidationData</code>	This is your holdout data set used in modeling either refine your hyper-parameters.
<code>TestData</code>	This is your holdout data set. Catboost using both training and validation data in the training process so you should evaluate out of sample performance with this data set.
<code>TargetColumnName</code>	Either supply the target column name OR the column number where the target is located (but not mixed types).
<code>FeatureColNames</code>	Either supply the feature column names OR the column number where the target is located (but not mixed types)
<code>eval_metric</code>	This is the metric used to identify best grid tuned model. Choose from "logloss", "r2", "RMSE", "MSE"
<code>Trees</code>	The maximum number of trees you want in your models
<code>GridTune</code>	Set to TRUE to run a grid tuning procedure. Set a number in <code>MaxModelsInGrid</code> to tell the procedure how many models you want to test.
<code>MaxMem</code>	Set the maximum amount of memory you'd like to dedicate to the model run. E.g. "32G"
<code>MaxModelsInGrid</code>	Number of models to test from grid options (1080 total possible options)
<code>model_path</code>	A character string of your path file to where you want your output saved
<code>ModelID</code>	A character string to name your model and output
<code>ReturnModelObjects</code>	Set to TRUE to output all modeling objects (E.g. plots and evaluation metrics)
<code>SaveModelObjects</code>	Set to TRUE to return all modeling objects to your environment
<code>IfSaveModel</code>	Set to "mojo" to save a mojo file, otherwise "standard" to save a regular H2O model object

**Value**

Saves to file and returned in list: `VariableImportance.csv`, `Model`, `ValidationData.csv`, `EvaluationMetrics.csv`, `GridCollect`, and `GridList`

**Author(s)**

Adrian Antico

**See Also**

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoCatBoostRegression](#), [AutoCatBoostScoring](#), [AutoCatBoostHurdleModel](#), [AutoH2OMLScoring](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFClassifier](#), [AutoH2oDRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#), [AutoXGBoostScoring](#)

## Examples

```

Correl <- 0.85
N <- 1000
data <- data.table::data.table(Target = runif(N))
data[, x1 := qnorm(Target)]
data[, x2 := runif(N)]
data[, Independent_Variable1 := log(pnorm(Correl * x1 +
                                          sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable2 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable3 := exp(pnorm(Correl * x1 +
                                          sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable4 := exp(exp(pnorm(Correl * x1 +
                                              sqrt(1-Correl^2) * qnorm(x2))))]
data[, Independent_Variable5 := sqrt(pnorm(Correl * x1 +
                                           sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable6 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.10]
data[, Independent_Variable7 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.25]
data[, Independent_Variable8 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.75]
data[, Independent_Variable9 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^2]
data[, Independent_Variable10 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^4]
data[, Independent_Variable11 := as.factor(
  ifelse(Independent_Variable2 < 0.20, "A",
    ifelse(Independent_Variable2 < 0.40, "B",
      ifelse(Independent_Variable2 < 0.6, "C",
        ifelse(Independent_Variable2 < 0.8, "D", "E"))))]
data[, Target :=
  ifelse(Independent_Variable2 < 0.25, "A",
    ifelse(Independent_Variable2 < 0.45, "B",
      ifelse(Independent_Variable2 < 0.65, "C",
        ifelse(Independent_Variable2 < 0.85, "D", "E"))))]
data[, ':= ' (x1 = NULL, x2 = NULL)]
TestModel <- AutoH2oDRFMultiClass(data,
  ValidationData = NULL,
  TestData = NULL,
  TargetColumnName = "Target",
  FeatureColNames = 2:ncol(data),
  eval_metric = "logloss",
  Trees = 50,
  GridTune = FALSE,
  MaxMem = "32G",
  MaxModelsInGrid = 10,
  model_path = NULL,
  ModelID = "FirstModel",
  ReturnModelObjects = TRUE,
  SaveModelObjects = FALSE,
  IfSaveModel = "mojo")

```

---

AutoH2oDRFRegression	<i>AutoH2oDRFRegression is an automated H2O modeling framework with grid-tuning and model evaluation</i>
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## Description

AutoH2oDRFRegression is an automated H2O modeling framework with grid-tuning and model evaluation that runs a variety of steps. First, the function will run a random grid tune over N number of models and find which model is the best (a default model is always included in that set). Once the model is identified and built, several other outputs are generated: validation data with predictions, evaluation plot, evaluation boxplot, evaluation metrics, variable importance, partial dependence calibration plots, partial dependence calibration box plots, and column names used in model fitting.

## Usage

```
AutoH2oDRFRegression(data, ValidationData = NULL, TestData = NULL,
  TargetColumnName = NULL, FeatureColNames = NULL,
  TransformNumericColumns = NULL, eval_metric = "RMSE", Trees = 50,
  GridTune = FALSE, MaxMem = "32G", MaxModelsInGrid = 2,
  model_path = NULL, ModelID = "FirstModel", NumOfParDepPlots = 3,
  ReturnModelObjects = TRUE, SaveModelObjects = FALSE,
  IfSaveModel = "mojo", StopH2O = TRUE)
```

## Arguments

<b>data</b>	This is your data set for training and testing your model
<b>ValidationData</b>	This is your holdout data set used in modeling either refine your hyper-parameters.
<b>TestData</b>	This is your holdout data set. Catboost using both training and validation data in the training process so you should evaluate out of sample performance with this data set.
<b>TargetColumnName</b>	Either supply the target column name OR the column number where the target is located (but not mixed types).
<b>FeatureColNames</b>	Either supply the feature column names OR the column number where the target is located (but not mixed types)
<b>TransformNumericColumns</b>	Set to NULL to do nothing; otherwise supply the column names of numeric variables you want transformed
<b>eval_metric</b>	This is the metric used to identify best grid tuned model. Choose from "MSE", "RMSE", "MAE", "RMSLE"
<b>Trees</b>	The maximum number of trees you want in your models
<b>GridTune</b>	Set to TRUE to run a grid tuning procedure. Set a number in MaxModelsInGrid to tell the procedure how many models you want to test.
<b>MaxMem</b>	Set the maximum amount of memory you'd like to dedicate to the model run. E.g. "32G"

MaxModelsInGrid	Number of models to test from grid options (1080 total possible options)
model_path	A character string of your path file to where you want your output saved
ModelID	A character string to name your model and output
NumOfParDepPlots	Tell the function the number of partial dependence calibration plots you want to create. Calibration boxplots will only be created for numerical features (not dummy variables)
ReturnModelObjects	Set to TRUE to output all modeling objects (E.g. plots and evaluation metrics)
SaveModelObjects	Set to TRUE to return all modeling objects to your environment
IfSaveModel	Set to "mojo" to save a mojo file, otherwise "standard" to save a regular H2O model object
StopH2O	For use in other functions.

### Value

Saves to file and returned in list: VariableImportance.csv, Model, ValidationData.csv, EvaluationPlot.png, EvaluationBoxPlot.png, EvaluationMetrics.csv, ParDepPlots.R a named list of features with partial dependence calibration plots, ParDepBoxPlots.R, GridCollect, GridList, and Transformation metadata

### Author(s)

Adrian Antico

### See Also

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoCatBoostRegression](#), [AutoCatBoostScoring](#), [AutoCatBoostHurdleModel](#), [AutoH2OMLScoring](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFClassifier](#), [AutoH2oDRFMultiClass](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#), [AutoXGBoostScoring](#)

### Examples

```
Correl <- 0.85
N <- 1000
data <- data.table::data.table(Target = runif(N))
data[, x1 := qnorm(Target)]
data[, x2 := runif(N)]
data[, Independent_Variable1 := log(pnorm(Correl * x1 +
                                         sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable2 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable3 := exp(pnorm(Correl * x1 +
                                         sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable4 := exp(exp(pnorm(Correl * x1 +
                                             sqrt(1-Correl^2) * qnorm(x2))))]
data[, Independent_Variable5 := sqrt(pnorm(Correl * x1 +
```

```

                                sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable6 := (pnorm(Correl * x1 +
                                sqrt(1-Correl^2) * qnorm(x2)))^0.10]
data[, Independent_Variable7 := (pnorm(Correl * x1 +
                                sqrt(1-Correl^2) * qnorm(x2)))^0.25]
data[, Independent_Variable8 := (pnorm(Correl * x1 +
                                sqrt(1-Correl^2) * qnorm(x2)))^0.75]
data[, Independent_Variable9 := (pnorm(Correl * x1 +
                                sqrt(1-Correl^2) * qnorm(x2)))^2]
data[, Independent_Variable10 := (pnorm(Correl * x1 +
                                sqrt(1-Correl^2) * qnorm(x2)))^4]
data[, Independent_Variable11 := as.factor(
  ifelse(Independent_Variable2 < 0.20, "A",
    ifelse(Independent_Variable2 < 0.40, "B",
      ifelse(Independent_Variable2 < 0.6, "C",
        ifelse(Independent_Variable2 < 0.8, "D", "E")))))]
data[, ':= ' (x1 = NULL, x2 = NULL)]
TestModel <- AutoH2oDRFRegression(data,
  ValidationData = NULL,
  TestData = NULL,
  TargetColumnName = "Target",
  FeatureColNames = 2:ncol(data),
  TransformNumericColumns = NULL,
  eval_metric = "RMSE",
  Trees = 50,
  GridTune = FALSE,
  MaxMem = "32G",
  MaxModelsInGrid = 10,
  model_path = NULL,
  ModelID = "FirstModel",
  NumOfParDepPlots = 3,
  ReturnModelObjects = TRUE,
  SaveModelObjects = FALSE,
  IfSaveModel = "mojo",
  StopH2O = TRUE)

```

AutoH2oGBMCARMA

*AutoH2oGBMCARMA Automated CatBoost Calendar, ARMA,  
and Trend Variables Forecasting*

## Description

AutoH2oDRFCARMA Automated CatBoost Calendar, ARMA, and Trend Variables Forecasting. Create hundreds of thousands of time series forecasts using this function.

## Usage

```

AutoH2oGBMCARMA(data, TargetColumnName = "Target",
  DateColumnName = "DateTime", GroupVariables = NULL,
  FC_Periods = 30, TimeUnit = "week", TargetTransformation = FALSE,
  Lags = c(1:5), MA_Periods = c(1:5), CalendarVariables = FALSE,
  TimeTrendVariable = FALSE, DataTruncate = FALSE,
  SplitRatios = c(0.7, 0.2, 0.1), EvalMetric = "MAE",

```

```
GridTune = FALSE, ModelCount = 1, NTrees = 1000,
PartitionType = "timeseries", MaxMem = "32G", NThreads = max(1,
parallel::detectCores() - 2), Timer = TRUE)
```

### Arguments

<b>data</b>	Supply your full series data set here
<b>TargetColumnName</b>	List the column name of your target variables column. E.g. "Target"
<b>DateColumnName</b>	List the column name of your date column. E.g. "DateTime"
<b>GroupVariables</b>	Defaults to NULL. Use NULL when you have a single series. Add in GroupVariables when you have a series for every level of a group or multiple groups.
<b>FC_Periods</b>	Set the number of periods you want to have forecasts for. E.g. 52 for weekly data to forecast a year ahead
<b>TimeUnit</b>	List the time unit your data is aggregated by. E.g. "hour", "day", "week", "year"
<b>TargetTransformation</b>	Run AutoTransformationCreate() to find best transformation for the target variable. Tests YeoJohnson, BoxCox, and Asigh (also Asin and Logit for proportion target variables).
<b>Lags</b>	Select the periods for all lag variables you want to create. E.g. c(1:5,52)
<b>MA_Periods</b>	Select the periods for all moving average variables you want to create. E.g. c(1:5,52)
<b>CalendarVariables</b>	Set to TRUE to have calendar variables created. The calendar variables are numeric representations of second, minute, hour, week day, month day, year day, week, isoweek, quarter, and year
<b>TimeTrendVariable</b>	Set to TRUE to have a time trend variable added to the model. Time trend is numeric variable indicating the numeric value of each record in the time series (by group). Time trend starts at 1 for the earliest point in time and increments by one for each success time point.
<b>DataTruncate</b>	Set to TRUE to remove records with missing values from the lags and moving average features created
<b>SplitRatios</b>	E.g c(0.7,0.2,0.1) for train, validation, and test sets
<b>EvalMetric</b>	Select from "RMSE", "MAE", "MAPE", "Poisson", "Quantile", "LogLin-Quantile", "Lq", "NumErrors", "SMAPE", "R2", "MSLE", "MedianAbsoluteError"
<b>GridTune</b>	Set to TRUE to run a grid tune
<b>ModelCount</b>	Set the number of models to try in the grid tune
<b>NTrees</b>	Select the number of trees you want to have built to train the model
<b>PartitionType</b>	Select "random" for random data partitioning "time" for partitioning by time frames
<b>MaxMem</b>	Set to the maximum amount of memory you want to allow for running this function. Default is "32G".
<b>NThreads</b>	Set to the number of threads you want to dedicate to this function.
<b>Timer</b>	Set to FALSE to turn off the updating print statements for progress

**Value**

Returns a data.table of original series and forecasts, the catboost model objects (everything returned from AutoCatBoostRegression()), a time series forecast plot, and transformation info if you set TargetTransformation to TRUE. The time series forecast plot will plot your single series or aggregate your data to a single series and create a plot from that.

**See Also**

Other Time Series: [AutoCatBoostCARMA](#), [AutoH2oDRFCARMA](#), [AutoTS](#), [AutoXGBoostCARMA](#)

**Examples**

```
Results <- AutoH2oGBMCARMA(data,
  TargetColumnName = "Weekly_Sales",
  DateColumnName = "Date",
  GroupVariables = c("Store", "Dept"),
  FC_Periods = 52,
  TimeUnit = "week",
  TargetTransformation = FALSE,
  Lags = c(1:5, 52),
  MA_Periods = c(1:5, 52),
  CalendarVariables = TRUE,
  TimeTrendVariable = TRUE,
  DataTruncate = FALSE,
  SplitRatios = c(1-2*30/143, 30/143, 30/143),
  EvalMetric = "MAE",
  GridTune = FALSE,
  ModelCount = 1,
  NTrees = 1000,
  PartitionType = "timeseries",
  MaxMem = "32G",
  NThreads = max(1, parallel::detectCores() - 2),
  Timer = TRUE)

Results$TimeSeriesPlot
Results$Forecast
Results$ModelInformation$...
```

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AutoH2oGBMClassifier	<i>AutoH2oGBMClassifier is an automated H2O modeling framework with grid-tuning and model evaluation</i>
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**Description**

AutoH2oGBMClassifier is an automated H2O modeling framework with grid-tuning and model evaluation that runs a variety of steps. First, a stratified sampling (by the target variable) is done to create train and validation sets. Then, the function will run a random grid tune over N number of models and find which model is the best (a default model is always included in that set). Once the model is identified and built, several other outputs are generated: validation data with predictions, evaluation plot, evaluation metrics, variable importance, partial dependence calibration plots, and column names used in model fitting.

**Usage**

```
AutoH2oGBMClassifier(data, ValidationData = NULL, TestData = NULL,
  TargetColumnName = NULL, FeatureColNames = NULL,
  eval_metric = "auc", Trees = 50, GridTune = FALSE,
  MaxMem = "32G", MaxModelsInGrid = 2, model_path = NULL,
  ModelID = "FirstModel", NumOfParDepPlots = 3,
  ReturnModelObjects = TRUE, SaveModelObjects = FALSE,
  IfSaveModel = "mojo")
```

**Arguments**

<b>data</b>	This is your data set for training and testing your model
<b>ValidationData</b>	This is your holdout data set used in modeling either refine your hyper-parameters.
<b>TestData</b>	This is your holdout data set. Catboost using both training and validation data in the training process so you should evaluate out of sample performance with this data set.
<b>TargetColumnName</b>	Either supply the target column name OR the column number where the target is located (but not mixed types). Note that the target column needs to be a 0 — 1 numeric variable.
<b>FeatureColNames</b>	Either supply the feature column names OR the column number where the target is located (but not mixed types)
<b>eval_metric</b>	This is the metric used to identify best grid tuned model. Choose from "AUC" or "logloss"
<b>Trees</b>	The maximum number of trees you want in your models
<b>GridTune</b>	Set to TRUE to run a grid tuning procedure. Set a number in MaxModelsInGrid to tell the procedure how many models you want to test.
<b>MaxMem</b>	Set the maximum amount of memory you'd like to dedicate to the model run. E.g. "32G"
<b>MaxModelsInGrid</b>	Number of models to test from grid options (1080 total possible options)
<b>model_path</b>	A character string of your path file to where you want your output saved
<b>ModelID</b>	A character string to name your model and output
<b>NumOfParDepPlots</b>	Tell the function the number of partial dependence calibration plots you want to create.
<b>ReturnModelObjects</b>	Set to TRUE to output all modeling objects (E.g. plots and evaluation metrics)
<b>SaveModelObjects</b>	Set to TRUE to return all modeling objects to your environment
<b>IfSaveModel</b>	Set to "mojo" to save a mojo file, otherwise "standard" to save a regular H2O model object

**Value**

Saves to file and returned in list: VariableImportance.csv, Model, ValidationData.csv, EvaluationPlot.png, EvaluationMetrics.csv, ParDepPlots.R a named list of features with partial dependence calibration plots, GridCollect, and GridList



**Author(s)**

Adrian Antico

**See Also**

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoCatBoostRegression](#), [AutoCatBoostScoring](#), [AutoCatBoostHurdleModel](#), [AutoH2OMLScoring](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFClassifier](#), [AutoH2oDRFMultiClass](#), [AutoH2oDRFRegression](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#), [AutoXGBoostScoring](#)

**Examples**

```
Correl <- 0.85
N <- 1000
data <- data.table::data.table(Target = runif(N))
data[, x1 := qnorm(Target)]
data[, x2 := runif(N)]
data[, Independent_Variable1 := log(pnorm(Correl * x1 +
                                         sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable2 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable3 := exp(pnorm(Correl * x1 +
                                         sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable4 := exp(exp(pnorm(Correl * x1 +
                                              sqrt(1-Correl^2) * qnorm(x2))))]
data[, Independent_Variable5 := sqrt(pnorm(Correl * x1 +
                                           sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable6 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.10]
data[, Independent_Variable7 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.25]
data[, Independent_Variable8 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.75]
data[, Independent_Variable9 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^2]
data[, Independent_Variable10 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^4]
data[, Independent_Variable11 := as.factor(
  ifelse(Independent_Variable2 < 0.20, "A",
    ifelse(Independent_Variable2 < 0.40, "B",
      ifelse(Independent_Variable2 < 0.6, "C",
        ifelse(Independent_Variable2 < 0.8, "D", "E"))))]
data[, ':= ' (x1 = NULL, x2 = NULL)]
data[, Target := as.factor(ifelse(Independent_Variable2 < 0.5, 1, 0))]
TestModel <- AutoH2oGBMClassifier(data,
  ValidationData = NULL,
  TestData = NULL,
  TargetColumnName = "Target",
  FeatureColNames = 2:ncol(data),
  eval_metric = "auc",
  Trees = 50,
  GridTune = FALSE,
  MaxMem = "32G",
  MaxModelsInGrid = 10,
```

```

model_path = NULL,
ModelID = "FirstModel",
NumOfParDepPlots = 3,
ReturnModelObjects = TRUE,
SaveModelObjects = FALSE,
IfSaveModel = "mojo")

```

---

**AutoH2oGBMMultiClass**    *AutoH2oGBMMultiClass is an automated H2O modeling framework with grid-tuning and model evaluation*

---

## Description

AutoH2oGBMMultiClass is an automated H2O modeling framework with grid-tuning and model evaluation that runs a variety of steps. First, a stratified sampling (by the target variable) is done to create train and validation sets. Then, the function will run a random grid tune over N number of models and find which model is the best (a default model is always included in that set). Once the model is identified and built, several other outputs are generated: validation data with predictions, evaluation metrics, confusion matrix, and variable importance.

## Usage

```

AutoH2oGBMMultiClass(data, ValidationData = NULL, TestData = NULL,
  TargetColumnName = NULL, FeatureColNames = NULL,
  eval_metric = "logloss", Trees = 50, GridTune = FALSE,
  MaxMem = "32G", MaxModelsInGrid = 2, model_path = NULL,
  ModelID = "FirstModel", ReturnModelObjects = TRUE,
  SaveModelObjects = FALSE, IfSaveModel = "mojo")

```

## Arguments

<b>data</b>	This is your data set for training and testing your model
<b>ValidationData</b>	This is your holdout data set used in modeling either refine your hyper-parameters.
<b>TestData</b>	This is your holdout data set. Catboost using both training and validation data in the training process so you should evaluate out of sample performance with this data set.
<b>TargetColumnName</b>	Either supply the target column name OR the column number where the target is located (but not mixed types).
<b>FeatureColNames</b>	Either supply the feature column names OR the column number where the target is located (but not mixed types)
<b>eval_metric</b>	This is the metric used to identify best grid tuned model. Choose from "logloss", "r2", "RMSE", "MSE"
<b>Trees</b>	The maximum number of trees you want in your models
<b>GridTune</b>	Set to TRUE to run a grid tuning procedure. Set a number in MaxModelsInGrid to tell the procedure how many models you want to test.

MaxMem	Set the maximum amount of memory you'd like to dedicate to the model run. E.g. "32G"
MaxModelsInGrid	Number of models to test from grid options (1080 total possible options)
model_path	A character string of your path file to where you want your output saved
ModelID	A character string to name your model and output
ReturnModelObjects	Set to TRUE to output all modeling objects (E.g. plots and evaluation metrics)
SaveModelObjects	Set to TRUE to return all modeling objects to your environment
IfSaveModel	Set to "mojo" to save a mojo file, otherwise "standard" to save a regular H2O model object

### Value

Saves to file and returned in list: VariableImportance.csv, Model, ValidationData.csv, EvaluationMetrics.csv, GridCollect, and GridList

### Author(s)

Adrian Antico

### See Also

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoCatBoostRegression](#), [AutoCatBoostScoring](#), [AutoCatBoostHurdleModel](#), [AutoH2OMLScoring](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFClassifier](#), [AutoH2oDRFMultiClass](#), [AutoH2oDRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#), [AutoXGBoostScoring](#)

### Examples

```
Correl <- 0.85
N <- 1000
data <- data.table::data.table(Target = runif(N))
data[, x1 := qnorm(Target)]
data[, x2 := runif(N)]
data[, Independent_Variable1 := log(pnorm(Correl * x1 +
                                         sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable2 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable3 := exp(pnorm(Correl * x1 +
                                         sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable4 := exp(exp(pnorm(Correl * x1 +
                                             sqrt(1-Correl^2) * qnorm(x2))))]
data[, Independent_Variable5 := sqrt(pnorm(Correl * x1 +
                                           sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable6 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.10]
data[, Independent_Variable7 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.25]
data[, Independent_Variable8 := (pnorm(Correl * x1 +
```

```

                                sqrt(1-Correl^2) * qnorm(x2)))^0.75]
data[, Independent_Variable9 := (pnorm(Correl * x1 +
                                sqrt(1-Correl^2) * qnorm(x2)))^2]
data[, Independent_Variable10 := (pnorm(Correl * x1 +
                                sqrt(1-Correl^2) * qnorm(x2)))^4]
data[, Independent_Variable11 := as.factor(
  ifelse(Independent_Variable2 < 0.20, "A",
    ifelse(Independent_Variable2 < 0.40, "B",
      ifelse(Independent_Variable2 < 0.6, "C",
        ifelse(Independent_Variable2 < 0.8, "D", "E")))))]
data[, Target :=
  ifelse(Independent_Variable2 < 0.25, "A",
    ifelse(Independent_Variable2 < 0.45, "B",
      ifelse(Independent_Variable2 < 0.65, "C",
        ifelse(Independent_Variable2 < 0.85, "D", "E"))))]
data[, ':= ' (x1 = NULL, x2 = NULL)]
TestModel <- AutoH2oGBMMultiClass(data,
                                ValidationData = NULL,
                                TestData = NULL,
                                TargetColumnName = "Target",
                                FeatureColNames = 2:ncol(data),
                                eval_metric = "logloss",
                                Trees = 50,
                                GridTune = FALSE,
                                MaxMem = "32G",
                                MaxModelsInGrid = 10,
                                model_path = NULL,
                                ModelID = "FirstModel",
                                ReturnModelObjects = TRUE,
                                SaveModelObjects = FALSE,
                                IfSaveModel = "mojo")

```

---

AutoH2oGBMRegression     *AutoH2oGBMRegression is an automated H2O modeling framework with grid-tuning and model evaluation*

---

## Description

AutoH2oGBMRegression is an automated H2O modeling framework with grid-tuning and model evaluation that runs a variety of steps. First, the function will run a random grid tune over N number of models and find which model is the best (a default model is always included in that set). Once the model is identified and built, several other outputs are generated: validation data with predictions, evaluation plot, evaluation boxplot, evaluation metrics, variable importance, partial dependence calibration plots, partial dependence calibration box plots, and column names used in model fitting.

## Usage

```

AutoH2oGBMRegression(data, ValidationData, TestData = NULL,
  TargetColumnName = NULL, FeatureColNames = NULL,
  TransformNumericColumns = NULL, Alpha = NULL,
  Distribution = "poisson", eval_metric = "RMSE", Trees = 50,
  GridTune = FALSE, MaxMem = "32G", MaxModelsInGrid = 2,

```

```

model_path = NULL, ModelID = "FirstModel", NumOfParDepPlots = 3,
ReturnModelObjects = TRUE, SaveModelObjects = FALSE,
IfSaveModel = "mojo", StopH2O = TRUE)

```

### Arguments

<code>data</code>	This is your data set for training and testing your model
<code>ValidationData</code>	This is your holdout data set used in modeling either refine your hyper-parameters.
<code>TestData</code>	This is your holdout data set. Catboost using both training and validation data in the training process so you should evaluate out of sample performance with this data set.
<code>TargetColumnName</code>	Either supply the target column name OR the column number where the target is located (but not mixed types).
<code>FeatureColNames</code>	Either supply the feature column names OR the column number where the target is located (but not mixed types)
<code>TransformNumericColumns</code>	Set to NULL to do nothing; otherwise supply the column names of numeric variables you want transformed
<code>Alpha</code>	This is the quantile value you want to use for quantile regression. Must be a decimal between 0 and 1.
<code>Distribution</code>	Choose from "gaussian", "poisson", "gamma", "tweedie", "laplace", "quantile", "huber"
<code>eval_metric</code>	This is the metric used to identify best grid tuned model. Choose from "MSE", "RMSE", "MAE", "RMSLE"
<code>Trees</code>	The maximum number of trees you want in your models
<code>GridTune</code>	Set to TRUE to run a grid tuning procedure. Set a number in <code>MaxModelsInGrid</code> to tell the procedure how many models you want to test.
<code>MaxMem</code>	Set the maximum amount of memory you'd like to dedicate to the model run. E.g. "32G"
<code>MaxModelsInGrid</code>	Number of models to test from grid options (1080 total possible options)
<code>model_path</code>	A character string of your path file to where you want your output saved
<code>ModelID</code>	A character string to name your model and output
<code>NumOfParDepPlots</code>	Tell the function the number of partial dependence calibration plots you want to create. Calibration boxplots will only be created for numerical features (not dummy variables)
<code>ReturnModelObjects</code>	Set to TRUE to output all modeling objects (E.g. plots and evaluation metrics)
<code>SaveModelObjects</code>	Set to TRUE to return all modeling objects to your environment
<code>IfSaveModel</code>	Set to "mojo" to save a mojo file, otherwise "standard" to save a regular H2O model object
<code>StopH2O</code>	Set to FALSE to keep H2O running after you build your model



```

FeatureColNames = 2:ncol(data),
TransformNumericColumns = NULL,
Alpha = NULL,
Distribution = "poisson",
eval_metric = "RMSE",
Trees = 50,
GridTune = FALSE,
MaxMem = "32G",
MaxModelsInGrid = 10,
model_path = NULL,
ModelID = "FirstModel",
NumOfParDepPlots = 3,
ReturnModelObjects = TRUE,
SaveModelObjects = FALSE,
IfSaveModel = "mojo",
StopH2O = TRUE)

```

---

AutoH2OMLScoring	<i>AutoH2OMLScoring is an automated scoring function that compliments the AutoH2o model training functions.</i>
------------------	---

---

## Description

AutoH2OMLScoring is an automated scoring function that compliments the AutoH2oGBM\_\_() and AutoH2oDRF\_\_() models training functions. This function requires you to supply features for scoring. It will run ModelDataPrep() to prepare your features for H2O data conversion and scoring.

## Usage

```

AutoH2OMLScoring(ScoringData = NULL, ModelObject = NULL,
  ModelType = "mojo", H2OShutdown = TRUE, MaxMem = "28G",
  JavaOptions = "-Xmx1g -XX:ReservedCodeCacheSize=256m",
  ModelPath = NULL, ModelID = NULL, ReturnFeatures = TRUE,
  TransformNumeric = FALSE, BackTransNumeric = FALSE,
  TargetColumnName = NULL, TransformationObject = NULL,
  TransID = NULL, TransPath = NULL, MDP_Impute = TRUE,
  MDP_CharToFactor = TRUE, MDP_RemoveDates = TRUE,
  MDP_MissFactor = "0", MDP_MissNum = -1)

```

## Arguments

ScoringData	This is your data.table of features for scoring. Can be a single row or batch.
ModelObject	Supply a model object from AutoH2oDRF__()
ModelType	Set to either "mojo" or "standard" depending on which version you saved
H2OShutdown	Set to TRUE is you are scoring a "standard" model file and you aren't planning on continuing to score.
MaxMem	Set to you dedicated amount of memory. E.g. "28G"
JavaOptions	Change the default to your machines specification if needed. Default is '-Xmx1g -XX:ReservedCodeCacheSize=256m',

ModelPath	Supply your path file used in the AutoH2o__() function
ModelID	Supply the model ID used in the AutoH2o__() function
ReturnFeatures	Set to TRUE to return your features with the predicted values.
TransformNumeric	Set to TRUE if you have features that were transformed automatically from an Auto__Regression() model AND you haven't already transformed them.
BackTransNumeric	Set to TRUE to generate back-transformed predicted values. Also, if you return features, those will also be back-transformed.
TargetColumnName	Input your target column name used in training if you are utilizing the transformation service
TransformationObject	Set to NULL if you didn't use transformations or if you want the function to pull from the file output from the Auto__Regression() function. You can also supply the transformation data.table object with the transformation details versus having it pulled from file.
TransID	Set to the ID used for saving the transformation data.table object or set it to the ModelID if you are pulling from file from a build with Auto__Regression().
TransPath	Set the path file to the folder where your transformation data.table detail object is stored. If you used the Auto__Regression() to build, set it to the same path as ModelPath.
MDP_Impute	Set to TRUE if you did so for modeling and didn't do so before supplying ScoringData in this function
MDP_CharToFactor	Set to TRUE to turn your character columns to factors if you didn't do so to your ScoringData that you are supplying to this function
MDP_RemoveDates	Set to TRUE if you have date of timestamp columns in your ScoringData
MDP_MissFactor	If you set MDP_Impute to TRUE, supply the character values to replace missing values with
MDP_MissNum	If you set MDP_Impute to TRUE, supply a numeric value to replace missing values with

### Value

A data.table of predicted values with the option to return model features as well.

### See Also

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoCatBoostRegression](#), [AutoCatBoostScoring](#), [AutoCatBoostHurdleModel](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFClassifier](#), [AutoH2oDRFMultiClass](#), [AutoH2oDRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#), [AutoXGBoostScoring](#)



## Examples

```
Preds <- AutoH2OMLScoring(ScoringData = data,
                           ModelObject = NULL,
                           ModelType = "mojo",
                           H2OShutdown = TRUE,
                           MaxMem = "28G",
                           JavaOptions = '-Xmx1g -XX:ReservedCodeCacheSize=256m',
                           ModelPath = NULL,
                           ModelID = "ModelTest",
                           ReturnFeatures = TRUE,
                           TransformNumeric = FALSE,
                           BackTransNumeric = FALSE,
                           TargetColumnName = NULL,
                           TransformationObject = NULL,
                           TransID = NULL,
                           TransPath = NULL,
                           MDP_Impute = TRUE,
                           MDP_CharToFactor = TRUE,
                           MDP_RemoveDates = TRUE,
                           MDP_MissFactor = "0",
                           MDP_MissNum = -1)
```

---

AutoH2OModeler

*An Automated Machine Learning Framework using H2O*


---

## Description

Steps in the function include: See details below for information on using this function.

## Usage

```
AutoH2OModeler(Construct, max_memory = "28G", ratios = 0.8,
               BL_Trees = 500, nthreads = 1, model_path = NULL,
               MaxRuntimeSeconds = 3600, MaxModels = 30, TrainData = NULL,
               TestData = NULL, SaveToFile = FALSE, ReturnObjects = TRUE)
```

## Arguments

Construct	Core instruction file for automation (see Details below for more information on this)
max_memory	The ceiling amount of memory H2O will utilize
ratios	The percentage of train samples from source data (remainder goes to validation set)
BL_Trees	The number of trees to build in baseline GBM or RandomForest
nthreads	Set the number of threads to run function
model_path	Directory path for where you want your models saved
MaxRuntimeSeconds	Number of seconds of run time for grid tuning
MaxModels	Number of models you'd like to have returned

TrainData	Set to NULL or supply a data.table for training data
TestData	Set to NULL or supply a data.table for validation data
SaveToFile	Set to TRUE to save models and output to model_path
ReturnObjects	Set to TRUE to return objects from function

## Details

1. Logic: Error checking in the modeling arguments from your Construction file
2. ML: Build grid-tuned models and baseline models for comparison and checks which one performs better on validation data
3. Evaluation: Collects the performance metrics for both
4. Evaluation: Generates calibration plots (and boxplots for regression) for the winning model
5. Evaluation: Generates partial dependence calibration plots (and boxplots for regression) for the winning model
6. Evaluation: Generates variable importance tables and a table of non-important features
7. Production: Creates a storage file containing: model name, model path, grid tune performance, baseline performance, and threshold (if classification) and stores that file in your model\_path location

The Construct file must be a data.table and the columns need to be in the correct order (see examples). Character columns must be converted to type "Factor". You must remove date columns or convert them to "Factor". For classification models, your target variable needs to be a (0,1) of type "Factor." See the examples below for help with setting up the Construct file for various modeling target variable types. There are examples for regression, classification, multinomial, and quantile regression. For help on which parameters to use, look up the r/h2o documentation. If you misspecify the construct file, it will produce an error and outputfile of what was wrong and suggestions for fixing the error.

Let's go over the construct file, column by column. The Targets column is where you specify the column number of your target variable (in quotes, e.g. "c(1)").

The Distribution column is where you specify the distribution type for the modeling task. For classification use bernoulli, for multilabel use multinomial, for quantile use quantile, and for regression, you can choose from the list available in the H2O docs, such as gaussian, poisson, gamma, etc. It's not set up to handle tweedie distributions currently but I can add support if there is demand.

The Loss column tells H2O which metric to use for the loss metrics. For regression, I typically use "mse", quantile regression, "mae", classification "auc", and multinomial "logloss". For deeplearning models, you need to use "quadratic", "absolute", and "crossentropy".

The Quantile column tells H2O which quantile to use for quantile regression (in decimal form).

The ModelName column is the name you wish to give your model as a prefix.

The Algorithm column is the model you wish to use: gbm, randomForest, deeplearning, AutoML, XGBoost, LightGBM.

The dataName column is the name of your data.

The TargetCol column is the column number of your target variable.

The FeatureCols column is the column numbers of your features.

The CreateDate column is for tracking your model build dates.

The GridTune column is a TRUE / FALSE column for whether you want to run a grid tune model for comparison.

The ExportValidData column is a TRUE / FALSE column indicating if you want to export the validation data.

The ParDep column is where you put the number of partial dependence calibration plots you wish to generate.

The PD\_Data column is where you specify if you want to generate the partial dependence plots on "All" data, "Validate" data, or "Train" data.

The ThreshType column is for classification models. You can specify "f1", "f2", "f0point5", or "CS" for cost sensitive.

The FSC column is the feature selection column. Specify the percentage importance cutoff to create a table of "unimportant" features.

The tpProfit column is for when you specify "CS" in the ThreshType column. This is your true positive profit.

The tnProfit column is for when you specify "CS" in the ThreshType column. This is your true negative profit.

The fpProfit column is for when you specify "CS" in the ThreshType column. This is your false positive profit.

The fnProfit column is for when you specify "CS" in the ThreshType column. This is your false negative profit.

The SaveModel column is a TRUE / FALSE indicator. If you are just testing out models, set this to FALSE.

The SaveModelType column is where you specify if you want a "standard" model object saved or a "mojo" model object saved.

The PredsAllData column is a TRUE / FALSE column. Set to TRUE if you want all the predicted values returns (for all data).

The TargetEncoding column lets you specify the column number of features you wish to run target encoding on. Set to NA to not run this feature.

The SupplyData column lets you supply the data names for training and validation data. Set to NULL if you want the data partitioning to be done internally.

## Value

Returns saved models, corrected Construct file, variable importance tables, evaluation and partial dependence calibration plots, model performance measure, and a file called grid\_tuned\_paths.Rdata which contains paths to your saved models for operationalization.

## Author(s)

Adrian Antico

## See Also

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoCatBoostRegression](#), [AutoCatBoostScoring](#), [AutoCatBoostHurdleModel](#), [AutoH2OMLScoring](#), [AutoH2OScoring](#), [AutoH2oDRFClassifier](#), [AutoH2oDRFMultiClass](#), [AutoH2oDRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#), [AutoXGBoostScoring](#)



```

TargetEncoding = rep(NA,3),
SupplyData      = rep(FALSE,3))

AutoH2OModeler(Construct,
  max_memory = "28G",
  ratios = 0.75,
  BL_Trees = 500,
  nthreads = 5,
  model_path = NULL,
  MaxRuntimeSeconds = 3600,
  MaxModels = 30,
  TrainData = NULL,
  TestData = NULL,
  SaveToFile = FALSE,
  ReturnObjects = TRUE)

# Multinomial Example
Correl <- 0.85
aa <- data.table::data.table(target = runif(1000))
aa[, x1 := qnorm(target)]
aa[, x2 := runif(1000)]
aa[, Independent_Variable1 := log(pnorm(Correl * x1 +
  sqrt(1-Correl^2) * qnorm(x2)))]
aa[, Independent_Variable2 := (pnorm(Correl * x1 +
  sqrt(1-Correl^2) * qnorm(x2)))]
aa[, Independent_Variable3 := exp(pnorm(Correl * x1 +
  sqrt(1-Correl^2) * qnorm(x2)))]
aa[, Independent_Variable4 := exp(exp(pnorm(Correl * x1 +
  sqrt(1-Correl^2) * qnorm(x2)))]
aa[, Independent_Variable5 := sqrt(pnorm(Correl * x1 +
  sqrt(1-Correl^2) * qnorm(x2)))]
aa[, Independent_Variable6 := (pnorm(Correl * x1 +
  sqrt(1-Correl^2) * qnorm(x2)))^0.10]
aa[, Independent_Variable7 := (pnorm(Correl * x1 +
  sqrt(1-Correl^2) * qnorm(x2)))^0.25]
aa[, Independent_Variable8 := (pnorm(Correl * x1 +
  sqrt(1-Correl^2) * qnorm(x2)))^0.75]
aa[, Independent_Variable9 := (pnorm(Correl * x1 +
  sqrt(1-Correl^2) * qnorm(x2)))^2]
aa[, Independent_Variable10 := (pnorm(Correl * x1 +
  sqrt(1-Correl^2) * qnorm(x2)))^4]

aa[, ':= ' (x1 = NULL, x2 = NULL)]
aa[, target := as.factor(ifelse(target < 0.33, "A", ifelse(target < 0.66, "B", "C")))]
Construct <- data.table::data.table(Targets = rep("target",3),
  Distribution = c("multinomial",
    "multinomial",
    "multinomial"),
  Loss = c("auc", "logloss", "accuracy"),
  Quantile = rep(NA,3),
  ModelName = c("GBM", "DRF", "DL"),
  Algorithm = c("gbm",
    "randomForest",
    "deeplearning"),
  dataName = rep("aa",3),
  TargetCol = rep(c("1"),3),
  FeatureCols = rep(c("2:11"),3),
  CreateDate = rep(Sys.time(),3),
  GridTune = rep(FALSE,3),

```



```

                                "gaussian"),
Loss                           = c("MSE", "MSE", "Quadratic"),
Quantile                       = rep(NA, 3),
ModelName                     = c("GBM", "DRF", "DL"),
Algorithm                      = c("gbm",
                                "randomForest",
                                "deeplearning"),
dataName                      = rep("aa", 3),
TargetCol                     = rep(c("1"), 3),
FeatureCols                   = rep(c("2:11"), 3),
CreateDate                   = rep(Sys.time(), 3),
GridTune                      = rep(FALSE, 3),
ExportValidData              = rep(TRUE, 3),
ParDep                        = rep(2, 3),
PD_Data                      = rep("All", 3),
ThreshType                   = rep("f1", 3),
FSC                          = rep(0.001, 3),
tpProfit                     = rep(NA, 3),
tnProfit                     = rep(NA, 3),
fpProfit                     = rep(NA, 3),
fnProfit                     = rep(NA, 3),
SaveModel                    = rep(FALSE, 3),
SaveModelType                = c("Mojo", "standard", "mojo"),
PredsAllData                 = rep(TRUE, 3),
TargetEncoding               = rep(NA, 3),
SupplyData                   = rep(FALSE, 3))

AutoH2OModeler(Construct,
  max_memory = "28G",
  ratios = 0.75,
  BL_Trees = 500,
  nthreads = 5,
  model_path = NULL,
  MaxRuntimeSeconds = 3600,
  MaxModels = 30,
  TrainData = NULL,
  TestData = NULL,
  SaveToFile = FALSE,
  ReturnObjects = TRUE)

# Quantile Regression Example
Correl <- 0.85
aa <- data.table::data.table(target = runif(1000))
aa[, x1 := qnorm(target)]
aa[, x2 := runif(1000)]
aa[, Independent_Variable1 := log(pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))]
aa[, Independent_Variable2 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))]
aa[, Independent_Variable3 := exp(pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))]
aa[, Independent_Variable4 := exp(exp(pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))]
aa[, Independent_Variable5 := sqrt(pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))]
aa[, Independent_Variable6 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.10]
aa[, Independent_Variable7 := (pnorm(Correl * x1 +

```

```

                                sqrt(1-Correl^2) * qnorm(x2)))^0.25]
aa[, Independent_Variable8 := (pnorm(Correl * x1 +
                                sqrt(1-Correl^2) * qnorm(x2)))^0.75]
aa[, Independent_Variable9 := (pnorm(Correl * x1 +
                                sqrt(1-Correl^2) * qnorm(x2)))^2]
aa[, Independent_Variable10 := (pnorm(Correl * x1 +
                                sqrt(1-Correl^2) * qnorm(x2)))^4]
aa[, ':= ' (x1 = NULL, x2 = NULL)]
Construct <- data.table::data.table(Targets = rep("target",3),
                                Distribution = c("quantile",
                                                  "quantile"),
                                Loss = c("MAE", "Absolute"),
                                Quantile = rep(0.75,2),
                                ModelName = c("GBM", "DL"),
                                Algorithm = c("gbm",
                                              "deeplearning"),
                                dataName = rep("aa",2),
                                TargetCol = rep(c("1"),2),
                                FeatureCols = rep(c("2:11"),2),
                                CreateDate = rep(Sys.time(),2),
                                GridTune = rep(FALSE,2),
                                ExportValidData = rep(TRUE,2),
                                ParDep = rep(4,2),
                                PD_Data = rep("All",2),
                                ThreshType = rep("f1",2),
                                FSC = rep(0.001,2),
                                tpProfit = rep(NA,2),
                                tnProfit = rep(NA,2),
                                fpProfit = rep(NA,2),
                                fnProfit = rep(NA,2),
                                SaveModel = rep(FALSE,2),
                                SaveModelType = c("Mojo", "mojo"),
                                PredsAllData = rep(TRUE,2),
                                TargetEncoding = rep(NA,2),
                                SupplyData = rep(FALSE,2))

AutoH20Modeler(Construct,
               max_memory = "28G",
               ratios = 0.75,
               BL_Trees = 500,
               nthreads = 5,
               model_path = NULL,
               MaxRuntimeSeconds = 3600,
               MaxModels = 30,
               TrainData = NULL,
               TestData = NULL,
               SaveToFile = FALSE,
               ReturnObjects = TRUE)

```

---

AutoH2OScoring

*AutoH2OScoring is the complement of AutoH20Modeler.*


---

## Description

AutoH2OScoring is the complement of AutoH20Modeler. Use this for scoring models. You can score regression, quantile regression, classification, multinomial, clustering, and



text models (built with the `Word2VecModel` function). You can also use this to score multioutcome models so long as there are two models: one for predicting the count of outcomes (a count outcome in character form) and a multinomial model on the label data. You will want to ensure you have a record for each label in your training data in (0,1) as factor form.

## Usage

```
AutoH2OScoring(Features = data, GridTuneRow = c(1:3),
  ScoreMethod = "Standard", TargetType = rep("multinomial", 3),
  ClassVals = rep("probs", 3), TextType = "individual",
  TextNames = NULL, NThreads = 6, MaxMem = "28G",
  JavaOptions = "-Xmx1g -XX:ReservedCodeCacheSize=256m",
  SaveToFile = FALSE, FilePath = NULL, H2OShutDown = rep(FALSE, 3))
```

## Arguments

Features	This is a <code>data.table</code> of features for scoring.
GridTuneRow	Numeric. The row numbers of <code>grid_tuned_paths</code> , <code>KMeansModelFile</code> , or <code>StoreFile</code> containing the model you wish to score
ScoreMethod	"Standard" or "Mojo": Mojo is available for supervised models; use standard for all others
TargetType	"Regression", "Classification", "Multinomial", "MultiOutcome", "Text", "Clustering". MultiOutcome must be two multinomial models, a count model (the count of outcomes, as a character value), and the multinomial model predicting the labels.
ClassVals	Choose from "p1", "Probs", "Label", or "All" for classification and multinomial models.
NThreads	Number of available threads for H2O
MaxMem	Amount of memory to dedicate to H2O
JavaOptions	Modify to your machine if the default doesn't work
SaveToFile	Set to TRUE if you want your model scores saved to file.
FilePath	Set this to the folder where your models and model files are saved
H2OShutDown	TRUE to shutdown H2O after the run. Use FALSE if you will be repeatedly scoring and shutdown somewhere else in your environment.

## Value

Returns a list of predicted values. Each list element contains the predicted values from a single model predict call.

## Author(s)

Adrian Antico

## See Also

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoCatBoostRegression](#), [AutoCatBoostScoring](#), [AutoCatBoostHurdleModel](#), [AutoH2OMLScoring](#), [AutoH2OModeler](#), [AutoH2oDRFClassifier](#), [AutoH2oDRFMultiClass](#), [AutoH2oDRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#), [AutoXGBoostScoring](#)



```

                                TargetEncoding = rep(NA,3),
                                SupplyData      = rep(FALSE,3))

AutoH2OModeler(Construct,
  max_memory = "28G",
  ratios = 0.75,
  BL_Trees = 500,
  nthreads = 5,
  model_path = NULL,
  MaxRuntimeSeconds = 3600,
  MaxModels = 30,
  TrainData = NULL,
  TestData  = NULL,
  SaveToFile = FALSE,
  ReturnObjects = TRUE)

N <- 3
data <- AutoH2OScoring(Features      = aa,
  GridTuneRow = c(1:N),
  ScoreMethod = "standard",
  TargetType  = rep("multinomial",N),
  ClassVals   = rep("Probs",N),
  NThreads    = 6,
  MaxMem       = "28G",
  JavaOptions  = '-Xmx1g -XX:ReservedCodeCacheSize=256m',
  SaveToFile   = FALSE,
  FilesPath    = NULL,
  H2OShutDown  = rep(FALSE,N))

```

---

AutoH2OTextPrepScoring

*AutoH2OTextPrepScoring is for NLP scoring*


---

## Description

This function returns prepared tokenized data for H2O Word2VecModeler scoring

## Usage

```
AutoH2OTextPrepScoring(data, string = NULL, MaxMem = NULL,
  NThreads = NULL, StartH2O = TRUE)
```

## Arguments

<code>data</code>	The text data
<code>string</code>	The name of the string column to prepare
<code>MaxMem</code>	Amount of memory you want to let H2O utilize
<code>NThreads</code>	The number of threads you want to let H2O utilize
<code>StartH2O</code>	Set to TRUE to have H2O start inside this function

## Author(s)

Adrian Antico

**See Also**

Other Misc: [ChartTheme](#), [PrintObjectsSize](#), [RemixTheme](#), [SimpleCap](#), [multiplot](#), [percRank](#), [tempDatesFun](#), [tokenizeH2O](#)

**Examples**

```
data <- AutoH2OTextPrepScoring(data = x,
                               string = "text_column",
                               MaxMem = "28G",
                               NThreads = 8,
                               StartH2O = TRUE)
```

AutoKMeans

*AutoKMeans Automated row clustering for mixed column types***Description**

AutoKMeans adds a column to your original data with a cluster number identifier. Uses glrm (grid tune-able) and then k-means to find optimal k.

**Usage**

```
AutoKMeans(data, nthreads = 8, MaxMem = "28G", SaveModels = NULL,
            PathFile = NULL, GridTuneGLRM = TRUE, GridTuneKMeans = TRUE,
            glrmCols = c(1:5), IgnoreConstCols = TRUE, glrmFactors = 5,
            Loss = "Absolute", glrmMaxIters = 1000, SVDMethod = "Randomized",
            MaxRunTimeSecs = 3600, KMeansK = 50, KMeansMetric = "totss")
```

**Arguments**

<code>data</code>	is the source time series data.table
<code>nthreads</code>	set based on number of threads your machine has available
<code>MaxMem</code>	set based on the amount of memory your machine has available
<code>SaveModels</code>	Set to "standard", "mojo", or NULL (default)
<code>PathFile</code>	Set to folder where you will keep the models
<code>GridTuneGLRM</code>	If you want to grid tune the glrm model, set to TRUE, FALSE otherwise
<code>GridTuneKMeans</code>	If you want to grid tune the KMeans model, set to TRUE, FALSE otherwise
<code>glrmCols</code>	the column numbers for the glrm
<code>IgnoreConstCols</code>	tell H2O to ignore any columns that have zero variance
<code>glrmFactors</code>	similar to the number of factors to return from PCA
<code>Loss</code>	set to one of "Quadratic", "Absolute", "Huber", "Poisson", "Hinge", "Logistic", "Periodic"
<code>glrmMaxIters</code>	max number of iterations
<code>SVDMethod</code>	choose from "Randomized", "GramSVD", "Power"
<code>MaxRunTimeSecs</code>	set the timeout for max run time
<code>KMeansK</code>	number of factors to test out in k-means to find the optimal number
<code>KMeansMetric</code>	pick the metric to identify top model in grid tune c("totss", "betweeness", "withinss")

**Value**

Original data.table with added column with cluster number identifier

**Author(s)**

Adrian Antico

**See Also**

Other Unsupervised Learning: [GenTSAnomVars](#), [ResidualOutliers](#)

**Examples**

```
data <- data.table::as.data.table(iris)
data <- AutoKMeans(data,
  nthreads = 8,
  MaxMem = "28G",
  SaveModels = NULL,
  PathFile = NULL,
  GridTuneGLRM = TRUE,
  GridTuneKMeans = TRUE,
  glrmCols = 1:(ncol(data)-1),
  IgnoreConstCols = TRUE,
  glrmFactors = 2,
  Loss = "Absolute",
  glrmMaxIters = 1000,
  SVDMethod = "Randomized",
  MaxRunTimeSecs = 3600,
  KMeansK = 5,
  KMeansMetric = "totss")
unique(data[["Species"]])
unique(data[["ClusterID"]])
temp <- data[, mean(ClusterID), by = "Species"]
Setosa <- round(temp[Species == "setosa", V1][[1]],0)
Versicolor <- round(temp[Species == "versicolor", V1][[1]],0)
Virginica <- round(temp[Species == "virginica", V1][[1]],0)
data[, Check := "a"]
data[ClusterID == eval(Setosa), Check := "setosa"]
data[ClusterID == eval(Virginica), Check := "virginica"]
data[ClusterID == eval(Versicolor), Check := "versicolor"]
data[, Acc := as.numeric(ifelse(Check == Species, 1, 0))]
data[, mean(Acc)][[1]]
```

---

AutoMarketBasketModel	<i>AutoMarketBasketModel function runs a market basket analysis automatically</i>
-----------------------	---

---

**Description**

AutoMarketBasketModel function runs a market basket analysis automatically. It will convert your data, run the algorithm, and add on additional significance values not originally contained within.

## Usage

```
AutoMarketBasketModel(data, OrderIDColumnName, ItemIDColumnName,  
  LHS_Delimiter = ",", Support = 0.001, Confidence = 0.1,  
  MaxLength = 2, MinLength = 2, MaxTime = 5)
```

## Arguments

<code>data</code>	This is your transactions data set
<code>OrderIDColumnName</code>	Supply your column name for the Order ID Values
<code>ItemIDColumnName</code>	Supply your column name for the Item ID Values
<code>LHS_Delimiter</code>	Default delimiter for separating multiple ItemID's is a comma.
<code>Support</code>	Threshold for inclusion using support
<code>Confidence</code>	Threshold for inclusion using confidence
<code>MaxLength</code>	Maximum combinations of Item ID (number of items in basket to consider)
<code>MinLength</code>	Minimum length of combinations of ItemID (number of items in basket to consider)
<code>MaxTime</code>	Max run time per iteration (default is 5 seconds)

## Author(s)

Adrian Antico and Douglas Pestana

## See Also

Chi-sq statistics and p-values based on this paper: <http://www.cs.bc.edu/~alvarez/ChiSquare/chi2tr.pdf>

Other Recommender Systems: [AutoRecomDataCreate](#), [AutoRecommenderScoring](#), [AutoRecommender](#)

## Examples

```
rules_data <- AutoMarketBasketModel(  
  data,  
  OrderIDColumnName = "OrderNumber",  
  ItemIDColumnName = "ItemNumber",  
  LHS_Delimiter = ",",  
  Support = 0.001,  
  Confidence = 0.1,  
  MaxLength = 2,  
  MinLength = 2,  
  MaxTime = 5)
```

AutoNLS

*AutoNLS is a function for automatically building nls models*

## Description

This function will build models for 9 different nls models, along with a non-parametric monotonic regression and a polynomial regression. The models are evaluated, a winner is picked, and the predicted values are stored in your data table.

## Usage

```
AutoNLS(data, y, x, monotonic = TRUE)
```

## Arguments

<code>data</code>	Data is the data table you are building the modeling on
<code>y</code>	Y is the target variable name in quotes
<code>x</code>	X is the independent variable name in quotes
<code>monotonic</code>	This is a TRUE/FALSE indicator - choose TRUE if you want monotonic regression over polynomial regression

## Value

A list containing "PredictionData" which is a data table with your original column replaced by the nls model predictions; "ModelName" the model name; "ModelObject" The winning model to later use; "EvaluationMetrics" Model metrics for models with ability to build.

## Author(s)

Adrian Antico

## See Also

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoCatBoostRegression](#), [AutoCatBoostScoring](#), [AutoCatBoostHurdleModel](#), [AutoH2OMLScoring](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFClassifier](#), [AutoH2oDRFMultiClass](#), [AutoH2oDRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#), [AutoXGBoostScoring](#)

## Examples

```
# Create Growth Data
data <-
  data.table::data.table(Target = seq(1, 500, 1),
                        Variable = rep(1, 500))
for (i in as.integer(1:500)) {
  if (i == 1) {
    var <- data[i, "Target"][[1]]
    data.table::set(data,
                    i = i,
                    j = 2L,
                    value = var * (1 + runif(1) / 100))
  }
}
```

```

    } else {
      var <- data[i - 1, "Variable"][[1]]
      data.table::set(data,
        i = i,
        j = 2L,
        value = var * (1 + runif(1) / 100))
    }
  }

# Add jitter to Target
data[, Target := jitter(Target,
  factor = 0.25)]

# To keep original values
data1 <- data.table::copy(data)

# Merge and Model data
data11 <- AutoNLS(
  data = data,
  y = "Target",
  x = "Variable",
  monotonic = TRUE
)

# Join predictions to source data
data2 <- merge(
  data1,
  data11$PredictionData,
  by = "Variable",
  all = FALSE
)

# Plot output
ggplot2::ggplot(data2, ggplot2::aes(x = Variable)) +
  ggplot2::geom_line(ggplot2::aes(y = data2[["Target.x"]],
    color = "Target")) +
  ggplot2::geom_line(ggplot2::aes(y = data2[["Target.y"]],
    color = "Predicted")) +
  RemixAutoML::ChartTheme(Size = 12) +
  ggplot2::ggtitle(paste0("Growth Models AutoNLS: ",
    data11$ModelName)) +
  ggplot2::ylab("Target Variable") +
  ggplot2::xlab("Independent Variable") +
  ggplot2::scale_colour_manual("Values",
    breaks = c("Target",
      "Predicted"),
    values = c("red",
      "blue"))

summary(data11$ModelObject)
data11$EvaluationMetrics

```



## Description

Convert transactional data.table to a binary ratings matrix

## Usage

```
AutoRecomDataCreate(data, EntityColName = "CustomerID",  
  ProductColName = "StockCode", MetricColName = "TotalSales",  
  ReturnMatrix = FALSE)
```

## Arguments

<code>data</code>	This is your transactional data.table. Must include an Entity (typically customer), ProductCode (such as SKU), and a sales metric (such as total sales).
<code>EntityColName</code>	This is the column name in quotes that represents the column name for the Entity, such as customer
<code>ProductColName</code>	This is the column name in quotes that represents the column name for the product, such as SKU
<code>MetricColName</code>	This is the column name in quotes that represents the column name for the metric, such as total sales
<code>ReturnMatrix</code>	Set to FALSE to coerce the object (desired route) or TRUE to return a matrix

## Value

A BinaryRatingsMatrix

## Author(s)

Adrian Antico and Douglas Pestana

## See Also

Other Recommender Systems: [AutoMarketBasketModel](#), [AutoRecommenderScoring](#), [AutoRecommender](#)

## Examples

```
RatingsMatrix <- AutoRecomDataCreate(data,  
  EntityColName = "CustomerID",  
  ProductColName = "StockCode",  
  MetricColName = "TotalSales",  
  ReturnMatrix = TRUE)
```

---

AutoRecommender	<i>Automatically build the best recommender model among models available.</i>
-----------------	---

---

## Description

This function returns the winning model that you pass onto AutoRecommenderScoring

## Usage

```
AutoRecommender(data, Partition = "Split", KFold = 1, Ratio = 0.75,
  Given = 1, RatingType = "TopN", RatingsKeep = 20,
  SkipModels = "AssociationRules", ModelMetric = "TPR")
```

## Arguments

data	This is your BinaryRatingsMatrix. See function RecomDataCreate
Partition	Choose from "split", "cross-validation", "bootstrap". See evaluation-Scheme in recommenderlab for details.
KFold	Choose 1 for traditional train and test. Choose greater than 1 for the number of cross validations
Ratio	The ratio for train and test. E.g. 0.75 for 75 percent data allocated to training
Given	The number of products you would like to evaluate. Negative values implement all-but schemes.
RatingType	Choose from "TopN", "ratings", "ratingMatrix"
RatingsKeep	The total ratings you wish to return. Default is 20.
SkipModels	AssociationRules runs the slowest and may crash your system. Choose from: "AssociationRules", "ItemBasedCF", "UserBasedCF", "PopularItems", "RandomItems"
ModelMetric	Choose from "Precision", "Recall", "TPR", or "FPR"

## Value

The winning model used for scoring in the AutoRecommenderScoring function

## Author(s)

Adrian Antico and Douglas Pestana

## See Also

Other Recommender Systems: [AutoMarketBasketModel](#), [AutoRecomDataCreate](#), [AutoRecommenderScoring](#)

## Examples

```
WinningModel <- AutoRecommender(RatingsMatrix,
                                Partition = "Split",
                                KFold = 1,
                                Ratio = 0.75,
                                Given = 1,
                                RatingType = "TopN",
                                RatingsKeep = 20,
                                SkipModels = "AssociationRules",
                                ModelMetric = "TPR")
```

---

### AutoRecommenderScoring

*The AutoRecomScoring function scores recommender models from AutoRecommender()*

---

## Description

This function will take your ratings matrix and model and score your data in parallel.

## Usage

```
AutoRecommenderScoring(data, WinningModel, EntityColName = "CustomerID",
                        ProductColName = "StockCode", NumItemsReturn = 1)
```

## Arguments

data	The binary ratings matrix from RecomDataCreate()
WinningModel	The winning model returned from AutoRecommender()
EntityColName	Typically your customer ID
ProductColName	Something like "StockCode"
NumItemsReturn	Number of items to return on scoring

## Value

Returns the prediction data

## Author(s)

Adrian Antico and Douglas Pestana

## See Also

Other Recommender Systems: [AutoMarketBasketModel](#), [AutoRecomDataCreate](#), [AutoRecommender](#)

## Examples

```
# F(G(Z(x))): AutoRecommenderScoring(AutoRecommender(AutoRecomDataCreate(TransactionData)))
Results <- AutoRecommenderScoring(
  data = AutoRecomDataCreate(
    data,
    EntityColName = "CustomerID",
    ProductColName = "StockCode",
    MetricColName = "TotalSales"),
  WinningModel = AutoRecommender(
    AutoRecomDataCreate(
      data,
      EntityColName = "CustomerID",
      ProductColName = "StockCode",
      MetricColName = "TotalSales"),
    Partition = "Split",
    KFold = 2,
    Ratio = 0.75,
    RatingType = "TopN",
    RatingsKeep = 20,
    SkipModels = "AssociationRules",
    ModelMetric = "TPR"),
  EntityColName = "CustomerID",
  ProductColName = "StockCode")
```

---

### AutoTransformationCreate

*AutoTransformationCreate is a function for automatically identifying the optimal transformations for numeric features and transforming them once identified.*

---

## Description

AutoTransformationCreate is a function for automatically identifying the optimal transformations for numeric features and transforming them once identified. This function will loop through your selected transformation options (YeoJohnson, BoxCox, Asinh, Asin, and Logit) and find the one that produces data that is the closest to normally distributed data. It then makes the transformation and collects the metadata information for use in the AutoTransformationScore() function, either by returning the objects (always) or saving them to file (optional).

## Usage

```
AutoTransformationCreate(data, ColumnNames = NULL,
  Methods = c("BoxCox", "YeoJohnson", "Asinh", "Asin", "Logit"),
  Path = NULL, TransID = "ModelID", SaveOutput = FALSE)
```

## Arguments

data	This is your source data
ColumnNames	List your columns names in a vector, for example, c("Target", "IV1")

Methods	Choose from "YeoJohnson", "BoxCox", "Asinh", "Asin", and "Logit". "Identity" is always run.
Path	Set to the directly where you want to save all of your modeling files
TransID	Set to a character value that corresponds with your modeling project
SaveOutput	Set to TRUE to save necessary file to run AutoTransformationScore()

**Value**

data with transformed columns and the transformation object for back-transforming later

**Author(s)**

Adrian Antico

**See Also**

Other Feature Engineering: [AutoDataPartition](#), [AutoTransformationScore](#), [AutoWord2VecModeler](#), [CreateCalendarVariables](#), [DT\\_GDL\\_Feature\\_Engineering](#), [DummifyDT](#), [GDL\\_Feature\\_Engineering](#), [ModelDataPrep](#), [Partial\\_DT\\_GDL\\_Feature\\_Engineering2](#), [Partial\\_DT\\_GDL\\_Feature\\_Engineering](#), [Scoring\\_GDL\\_Feature\\_Engineering](#)

**Examples**

```
Correl <- 0.85
N <- 1000
data <- data.table::data.table(Adrian = runif(N))
data[, x1 := qnorm(Adrian)]
data[, x2 := runif(N)]
data[, Adrian1 := log(pnorm(Correl * x1 +
                           sqrt(1-Correl^2) * qnorm(x2)))]
data <- AutoTransformationCreate(data,
                                ColumnNames = "Sample",
                                Methods = c("BoxCox", "Asin"),
                                Path = NULL,
                                TransID = "Trans",
                                SaveOutput = FALSE)
```

---

**AutoTransformationScore**

*AutoTransformationScore()* is a the complimentary function to *AutoTransformationCreate()*

---

**Description**

`AutoTransformationScore()` is a the compliment function to `AutoTransformationCreate()`. Automatically apply or inverse the transformations you identified in `AutoTransformationCreate()` to other data sets. This is useful for applying transformations to your validation and test data sets for modeling. It's also useful for back-transforming your target and prediction columns after you have build and score your models so you can obtain statistics on the original features.

**Usage**

```
AutoTransformationScore(ScoringData, FinalResults, Type = "Inverse",
  TransID = "TestModel", Path = NULL)
```

**Arguments**

ScoringData	This is your source data
FinalResults	This is the FinalResults output object from AutoTransformationCreate().
Type	Set to "Inverse" to back-transfrom or "Apply" for applying the transformation.
TransID	Set to a character value that corresponds with your modeling project
Path	Set to the directly where you want to save all of your modeling files

**Value**

data with transformed columns

**Author(s)**

Adrian Antico

**See Also**

Other Feature Engineering: [AutoDataPartition](#), [AutoTransformationCreate](#), [AutoWord2VecModeler](#), [CreateCalendarVariables](#), [DT\\_GDL\\_Feature\\_Engineering](#), [DummifyDT](#), [GDL\\_Feature\\_Engineering](#), [ModelDataPrep](#), [Partial\\_DT\\_GDL\\_Feature\\_Engineering2](#), [Partial\\_DT\\_GDL\\_Feature\\_Engineering](#), [Scoring\\_GDL\\_Feature\\_Engineering](#)

**Examples**

```
Correl <- 0.85
N <- 1000
data <- data.table::data.table(Adrian = runif(N))
data[, x1 := qnorm(Adrian)]
data[, x2 := runif(N)]
data[, Adrian1 := log(pnorm(Correl * x1 +
  sqrt(1-Correl^2) * qnorm(x2)))]
data <- AutoTransformationScore(ScoringData,
  FinalResults,
  Path = NULL,
  TransID = "Trans")
```

## Description

Step 1 is to build all the models and evaluate them on the number of HoldOutPeriods periods you specify. Step 2 is to pick the winner and rebuild the winning model on the full data set. Step 3 is to generate forecasts with the final model for FCPeriods that you specify. AutoTS builds the best time series models for each type, using optimized box-cox transformations and using a user-supplied frequency for the ts data conversion along with a model-based frequency for the ts data conversion, compares all types, selects the winner, and generates a forecast. Models include:

## Usage

```
AutoTS(data, TargetName = "Target", DateName = "DateTime",
       FCPeriods = 30, HoldOutPeriods = 30, EvaluationMetric = "MAPE",
       TimeUnit = "day", Lags = 25, SLags = 2, NumCores = 4,
       SkipModels = NULL, StepWise = TRUE, TSClean = TRUE,
       ModelFreq = TRUE, PrintUpdates = FALSE)
```

## Arguments

<code>data</code>	is the source time series data as a <code>data.table</code> - or a data structure that can be converted to a <code>data.table</code>
<code>TargetName</code>	is the name of the target variable in your <code>data.table</code>
<code>DateName</code>	is the name of the date column in your <code>data.table</code>
<code>FCPeriods</code>	is the number of periods into the future you wish to forecast
<code>HoldOutPeriods</code>	is the number of periods to use for validation testing
<code>EvaluationMetric</code>	Set this to either "MAPE", "MSE", or "MAE". Default is "MAPE"
<code>TimeUnit</code>	is the level of aggregation your dataset comes in
<code>Lags</code>	is the number of lags you wish to test in various models (same as moving averages)
<code>SLags</code>	is the number of seasonal lags you wish to test in various models (same as moving averages)
<code>NumCores</code>	is the number of cores available on your computer
<code>SkipModels</code>	Don't run specified models - e.g. exclude all models "DSHW" "ARFIMA" "ARIMA" "ETS" "NNET" "TBATS" "TSLM"
<code>StepWise</code>	Set to TRUE to have ARIMA and ARFIMA run a stepwise selection process. Otherwise, all models will be generated in parallel execution, but still run much slower.
<code>TSClean</code>	Set to TRUE to have missing values interpolated and outliers replaced with interpolated values: creates separate models for a larger comparison set
<code>ModelFreq</code>	Set to TRUE to run a separate version of all models where the time series frequency is chosen algorithmically
<code>PrintUpdates</code>	Set to TRUE for a print to console of function progress

## Details

DSHW: Double Seasonal Holt Winters

ARFIMA: Auto Regressive Fractional Integrated Moving Average

ARIMIA: Stepwise Auto Regressive Integrated Moving Average with specified max lags, seasonal lags, moving averages, and seasonal moving averages

ETS: Additive and Multiplicative Exponential Smoothing and Holt Winters

NNetar: Auto Regressive Neural Network models automatically compares models with 1 lag or 1 seasonal lag compared to models with up to N lags and N seasonal lags

TBATS: Exponential smoothing state space model with Box-Cox transformation, ARMA errors, Trend and Seasonal components

TSLM: Time Series Linear Model - builds a linear model with trend and season components extracted from the data

## Value

Returns a list containing 1: A data.table object with a date column and the forecasted values; 2: The model evaluation results; 3: The champion model for later use if desired; 4: The name of the champion model; 5: A time series ggplot with historical values and forecasted values.

## Author(s)

Adrian Antico and Douglas Pestana

## See Also

Other Time Series: [AutoCatBoostCARMA](#), [AutoH2oDRFCARMA](#), [AutoH2oGBMCARMA](#), [AutoXGBoostCARMA](#)

## Examples

```
data <- data.table::data.table(DateTime = as.Date(Sys.time()),
  Target = stats::filter(rnorm(100,
    mean = 50,
    sd = 20),
    filter=rep(1,10),
    circular=TRUE))
data[, temp := seq(1:100)][, DateTime := DateTime - temp][, temp := NULL]
data <- data[order(DateTime)]
output <- AutoTS(data,
  TargetName      = "Target",
  DateName        = "DateTime",
  FCPeriods       = 1,
  HoldOutPeriods  = 1,
  EvaluationMetric = "MAPE",
  TimeUnit        = "day",
  Lags            = 1,
  SLags           = 1,
  NumCores        = 4,
  SkipModels      = c("NNET", "TBATS", "ETS", "TSLM", "ARFIMA", "DSHW"),
  StepWise        = TRUE,
  TSClean         = FALSE,
  ModelFreq       = TRUE,
  PrintUpdates    = FALSE)
```



```
ForecastData <- output$Forecast
ModelEval    <- output$EvaluationMetrics
WinningModel <- output$TimeSeriesModel
```

---

AutoWord2VecModeler      *Automated word2vec data generation via H2O*

---

## Description

This function allows you to automatically build a word2vec model and merge the data onto your supplied dataset

## Usage

```
AutoWord2VecModeler(data, BuildType = "Combined",
  stringCol = c("Text_Col1", "Text_Col2"), KeepStringCol = FALSE,
  model_path = NULL, vects = 100, SaveStopWords = FALSE,
  MinWords = 1, WindowSize = 12, Epochs = 25, StopWords = NULL,
  SaveModel = "standard", Threads = max(1, parallel::detectCores() -
  2), MaxMemory = "28G", SaveOutput = FALSE)
```

## Arguments

<code>data</code>	Source data table to merge vects onto
<code>stringCol</code>	A string name for the column to convert via word2vec
<code>KeepStringCol</code>	Set to TRUE if you want to keep the original string column that you convert via word2vec
<code>model_path</code>	A string path to the location where you want the model and metadata stored
<code>vects</code>	The number of vectors to retain from the word2vec model
<code>SaveStopWords</code>	Set to TRUE to save the stop words used
<code>MinWords</code>	For H2O word2vec model
<code>WindowSize</code>	For H2O word2vec model
<code>Epochs</code>	For H2O word2vec model
<code>StopWords</code>	For H2O word2vec model
<code>SaveModel</code>	Set to "standard" to save normally; set to "mojo" to save as mojo. NOTE: while you can save a mojo, I haven't figured out how to score it in the AutoH2OScoring function.
<code>Threads</code>	Number of available threads you want to dedicate to model building
<code>MaxMemory</code>	Amount of memory you want to dedicate to model building
<code>SaveOutput</code>	Set to TRUE to save your models to file

## Author(s)

Adrian Antico

## See Also

Other Feature Engineering: [AutoDataPartition](#), [AutoTransformationCreate](#), [AutoTransformationScore](#), [CreateCalendarVariables](#), [DT\\_GDL\\_Feature\\_Engineering](#), [DummifyDT](#), [GDL\\_Feature\\_Engineering](#), [ModelDataPrep](#), [Partial\\_DT\\_GDL\\_Feature\\_Engineering2](#), [Partial\\_DT\\_GDL\\_Feature\\_Engineering](#), [Scoring\\_GDL\\_Feature\\_Engineering](#)

## Examples

```
data <- AutoWord2VecModeler(data,
  BuildType = "individual",
  stringCol = c("Text_Col1",
    "Text_Col2"),
  KeepStringCol = FALSE,
  model_path = NULL,
  vects = 100,
  SaveStopWords = FALSE,
  MinWords = 1,
  WindowSize = 1,
  Epochs = 25,
  StopWords = NULL,
  SaveModel = "standard",
  Threads = max(1,parallel::detectCores()-2),
  MaxMemory = "28G",
  SaveOutput = TRUE)
```

---

AutoWordFreq

*Automated Word Frequency and Word Cloud Creation*


---

## Description

This function builds a word frequency table and a word cloud. It prepares data, cleans text, and generates output.

## Usage

```
AutoWordFreq(data, TextColName = "DESCR",
  GroupColName = "ClusterAllNoTarget", GroupLevel = 0,
  RemoveEnglishStopwords = TRUE, Stemming = TRUE,
  StopWords = c("bla", "bla2"))
```

## Arguments

<code>data</code>	Source data table
<code>TextColName</code>	A string name for the column
<code>GroupColName</code>	Set to NULL to ignore, otherwise set to Cluster column name (or factor column name)
<code>GroupLevel</code>	Must be set if GroupColName is defined. Set to cluster ID (or factor level)
<code>RemoveEnglishStopwords</code>	Set to TRUE to remove English stop words, FALSE to ignore

Stemming            Set to TRUE to run stemming on your text data  
 StopWords          Add your own stopwords, in vector format

### Author(s)

Adrian Antico

### See Also

Other EDA: [ProblematicFeatures](#), [ProblematicRecords](#)

### Examples

```
data <- data.table::data.table(
  DESCR = c("Gru, Gru, Gru, Gru, Gru, Gru, Gru, Gru, Gru, Gru, Gru, Gru, Gru, Gru,
    Urkle, Urkle, Urkle, Urkle, Urkle, Urkle, Urkle, Urkle, Gru, Gru, Gru,
    bears, bears, bears, bears, bears, bears, bears, smug, smug, smug, smug,
    smug, smug, smug, smug, smug, smug, smug, smug, smug, smug, smug,
    eats, eats, eats, eats, eats, eats, beats, beats, beats, beats,
    beats, beats, beats, beats, beats, beats, beats, beats, science, science,
    Dwig, Dwig, Dwig, Dwig, Dwig, Dwig, Dwig, Dwig, Dwig, Dwig, Dwig, Dwig,
    Schrute, Schrute, Schrute, Schrute, Schrute, Schrute, Schrute,
    James, James, James, James, James, James, James, James, James, James,
    Halpert, Halpert, Halpert, Halpert, Halpert, Halpert, Halpert, Halpert"))
data <- AutoWordFreq(data,
  TextColName = "DESCR",
  GroupColName = NULL,
  GroupLevel = NULL,
  RemoveEnglishStopwords = FALSE,
  Stemming = FALSE,
  StopWords = c("Bla"))
```

---

AutoXGBoostCARMA	<i>AutoXGBoostCARMA Automated XGBoost Calendar, ARMA, and Trend Variables Forecasting</i>
------------------	---

---

### Description

AutoXGBoostCARMA Automated XGBoost Calendar, ARMA, and Trend Variables Forecasting. Create hundreds of thousands of time series forecasts using this function.

### Usage

```
AutoXGBoostCARMA(data, TargetColumnName = "Target",
  DateColumnName = "DateTime", GroupVariables = NULL,
  FC_Periods = 30, TimeUnit = "week", TargetTransformation = FALSE,
  Lags = c(1:5), MA_Periods = c(1:5), CalendarVariables = FALSE,
  TimeTrendVariable = FALSE, DataTruncate = FALSE,
  SplitRatios = c(0.7, 0.2, 0.1), TreeMethod = "hist",
  NThreads = max(1, parallel::detectCores() - 2), EvalMetric = "MAPE",
  GridTune = FALSE, GridEvalMetric = "mape", ModelCount = 1,
  NTrees = 1000, PartitionType = "timeseries", Timer = TRUE)
```

**Arguments**

<code>data</code>	Supply your full series data set here
<code>TargetColumnName</code>	List the column name of your target variables column. E.g. "Target"
<code>DateColumnName</code>	List the column name of your date column. E.g. "DateTime"
<code>GroupVariables</code>	Defaults to NULL. Use NULL when you have a single series. Add in GroupVariables when you have a series for every level of a group or multiple groups.
<code>FC_Periods</code>	Set the number of periods you want to have forecasts for. E.g. 52 for weekly data to forecast a year ahead
<code>TimeUnit</code>	List the time unit your data is aggregated by. E.g. "hour", "day", "week", "year"
<code>TargetTransformation</code>	Run <code>AutoTransformationCreate()</code> to find best transformation for the target variable. Tests YeoJohnson, BoxCox, and Asigh (also Asin and Logit for proportion target variables).
<code>Lags</code>	Select the periods for all lag variables you want to create. E.g. <code>c(1:5,52)</code>
<code>MA_Periods</code>	Select the periods for all moving average variables you want to create. E.g. <code>c(1:5,52)</code>
<code>CalendarVariables</code>	Set to TRUE to have calendar variables created. The calendar variables are numeric representations of second, minute, hour, week day, month day, year day, week, isoweek, quarter, and year
<code>TimeTrendVariable</code>	Set to TRUE to have a time trend variable added to the model. Time trend is numeric variable indicating the numeric value of each record in the time series (by group). Time trend starts at 1 for the earliest point in time and increments by one for each success time point.
<code>DataTruncate</code>	Set to TRUE to remove records with missing values from the lags and moving average features created
<code>SplitRatios</code>	E.g <code>c(0.7,0.2,0.1)</code> for train, validation, and test sets
<code>TreeMethod</code>	Choose from "hist", "gpu_hist"
<code>NThreads</code>	Set the maximum number of threads you'd like to dedicate to the model run. E.g. 8
<code>EvalMetric</code>	Select from "r2", "RMSE", "MSE", "MAE"
<code>GridTune</code>	Set to TRUE to run a grid tune
<code>GridEvalMetric</code>	This is the metric used to find the threshold 'poisson', 'mae', 'mape', 'mse', 'msle', 'kl', 'cs', 'r2'
<code>ModelCount</code>	Set the number of models to try in the grid tune
<code>NTrees</code>	Select the number of trees you want to have built to train the model
<code>PartitionType</code>	Select "random" for random data partitioning "time" for partitioning by time frames
<code>Timer</code>	= TRUE

**Value**

Returns a data.table of original series and forecasts, the catboost model objects (everything returned from AutoCatBoostRegression()), a time series forecast plot, and transformation info if you set TargetTransformation to TRUE. The time series forecast plot will plot your single series or aggregate your data to a single series and create a plot from that.

**See Also**

Other Time Series: [AutoCatBoostCARMA](#), [AutoH2oDRFCARMA](#), [AutoH2oGBMCARMA](#), [AutoTS](#)

**Examples**

```
Results <- AutoXGBoostCARMA(data,
                             TargetColumnName = "Target",
                             DateColumnName = "DateTime",
                             GroupVariables = NULL,
                             FC_Periods = 30,
                             TimeUnit = "week",
                             TargetTransformation = FALSE,
                             Lags = c(1:5),
                             MA_Periods = c(1:5),
                             CalendarVariables = FALSE,
                             TimeTrendVariable = FALSE,
                             DataTruncate = FALSE,
                             SplitRatios = c(0.7, 0.2, 0.1),
                             TreeMethod = "hist",
                             NThreads = max(1, parallel::detectCores()-2),
                             EvalMetric = "MAE",
                             GridTune = FALSE,
                             GridEvalMetric = "mape",
                             ModelCount = 1,
                             NTrees = 1000,
                             PartitionType = "timeseries",
                             Timer = TRUE)

Results$TimeSeriesPlot
Results$Forecast
Results$ModelInformation$...
```

---

AutoXGBoostClassifier *AutoXGBoostClassifier is an automated XGBoost modeling framework with grid-tuning and model evaluation*

---

**Description**

AutoXGBoostClassifier is an automated XGBoost modeling framework with grid-tuning and model evaluation that runs a variety of steps. First, a stratified sampling (by the target variable) is done to create train and validation sets. Then, the function will run a random grid tune over N number of models and find which model is the best (a default model is always included in that set). Once the model is identified and built, several other outputs are generated: validation data with predictions, evaluation plot, evaluation box-plot, evaluation metrics, variable importance, partial dependence calibration plots, partial dependence calibration box plots, and column names used in model fitting.

**Usage**

```
AutoXGBoostClassifier(data, ValidationData = NULL, TestData = NULL,
  TargetColumnName = NULL, FeatureColNames = NULL, IDcols = NULL,
  eval_metric = "auc", Trees = 50, GridTune = FALSE,
  grid_eval_metric = "auc", TreeMethod = "hist",
  MaxModelsInGrid = 10, NThreads = 8, model_path = NULL,
  ModelID = "FirstModel", NumOfParDepPlots = 3, Verbose = 0,
  ReturnModelObjects = TRUE, SaveModelObjects = FALSE,
  PassInGrid = NULL)
```

**Arguments**

<b>data</b>	This is your data set for training and testing your model
<b>ValidationData</b>	This is your holdout data set used in modeling either refine your hyper-parameters.
<b>TestData</b>	This is your holdout data set. Catboost using both training and validation data in the training process so you should evaluate out of sample performance with this data set.
<b>TargetColumnName</b>	Either supply the target column name OR the column number where the target is located (but not mixed types). Note that the target column needs to be a 0 — 1 numeric variable.
<b>FeatureColNames</b>	Either supply the feature column names OR the column number where the target is located (but not mixed types)
<b>IDcols</b>	A vector of column names or column numbers to keep in your data but not include in the modeling.
<b>eval_metric</b>	This is the metric used to identify best grid tuned model. Choose from "logloss", "error", "aucpr", "auc"
<b>Trees</b>	The maximum number of trees you want in your models
<b>GridTune</b>	Set to TRUE to run a grid tuning procedure. Set a number in MaxModelsInGrid to tell the procedure how many models you want to test.
<b>grid_eval_metric</b>	Set to "f", "auc", "tpr", "fnr", "fpr", "tnr", "prbe", "f", "odds"
<b>TreeMethod</b>	Choose from "hist", "gpu_hist"
<b>MaxModelsInGrid</b>	Number of models to test from grid options (243 total possible options)
<b>NThreads</b>	Set the maximum number of threads you'd like to dedicate to the model run. E.g. 8
<b>model_path</b>	A character string of your path file to where you want your output saved
<b>ModelID</b>	A character string to name your model and output
<b>NumOfParDepPlots</b>	Tell the function the number of partial dependence calibration plots you want to create.
<b>Verbose</b>	Set to 0 if you want to suppress model evaluation updates in training
<b>ReturnModelObjects</b>	Set to TRUE to output all modeling objects (E.g. plots and evaluation metrics)

SaveModelObjects	Set to TRUE to return all modeling objects to your environment
PassInGrid	Default is NULL. Provide a data.table of grid options from a previous run.

### Value

Saves to file and returned in list: VariableImportance.csv, Model, ValidationData.csv, EvaluationPlot.png, EvaluationMetrics.csv, ParDepPlots.R a named list of features with partial dependence calibration plots, GridCollect, and GridList

### Author(s)

Adrian Antico

### See Also

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoCatBoostRegression](#), [AutoCatBoostScoring](#), [AutoCatBoostdHurdleModel](#), [AutoH2OMLScoring](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFClassifier](#), [AutoH2oDRFMultiClass](#), [AutoH2oDRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#), [AutoXGBoostScoring](#)

### Examples

```
Correl <- 0.85
N <- 10000
data <- data.table::data.table(Target = runif(N))
data[, x1 := qnorm(Target)]
data[, x2 := runif(N)]
data[, Independent_Variable1 := log(pnorm(Correl * x1 +
                                         sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable2 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable3 := exp(pnorm(Correl * x1 +
                                         sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable4 := exp(exp(pnorm(Correl * x1 +
                                             sqrt(1-Correl^2) * qnorm(x2))))]
data[, Independent_Variable5 := sqrt(pnorm(Correl * x1 +
                                          sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable6 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.10]
data[, Independent_Variable7 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.25]
data[, Independent_Variable8 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.75]
data[, Independent_Variable9 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^2]
data[, Independent_Variable10 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^4]
data[, Independent_Variable11 := as.factor(
  ifelse(Independent_Variable2 < 0.20, "A",
    ifelse(Independent_Variable2 < 0.40, "B",
      ifelse(Independent_Variable2 < 0.6, "C",
        ifelse(Independent_Variable2 < 0.8, "D", "E")))))]
```

```

data[, ']:= (x1 = NULL, x2 = NULL)]
data[, Target := ifelse(Target > 0.5, 1, 0)]
TestModel <- AutoXGBoostClassifier(data,
                                   ValidationData = NULL,
                                   TestData = NULL,
                                   TargetColumnName = 1,
                                   FeatureColNames = 2:12,
                                   IDcols = NULL,
                                   eval_metric = "auc",
                                   Trees = 50,
                                   GridTune = TRUE,
                                   grid_eval_metric = "auc",
                                   MaxModelsInGrid = 10,
                                   NThreads = 8,
                                   TreeMethod = "hist",
                                   model_path = getwd(),
                                   ModelID = "FirstModel",
                                   NumOfParDepPlots = 3,
                                   ReturnModelObjects = TRUE,
                                   SaveModelObjects = FALSE,
                                   PassInGrid = NULL)

```

---

**AutoXGBoostMultiClass** *AutoXGBoostMultiClass is an automated XGBoost modeling framework with grid-tuning and model evaluation*

---

## Description

AutoXGBoostMultiClass is an automated XGBoost modeling framework with grid-tuning and model evaluation that runs a variety of steps. First, a stratified sampling (by the target variable) is done to create train and validation sets. Then, the function will run a random grid tune over N number of models and find which model is the best (a default model is always included in that set). Once the model is identified and built, several other outputs are generated: validation data with predictions, evaluation metrics, variable importance, and column names used in model fitting.

## Usage

```

AutoXGBoostMultiClass(data, ValidationData = NULL, TestData = NULL,
  TargetColumnName = NULL, FeatureColNames = NULL, IDcols = NULL,
  eval_metric = "merror", Trees = 50, GridTune = FALSE,
  grid_eval_metric = "merror", TreeMethod = "hist",
  MaxModelsInGrid = 10, NThreads = 8, model_path = NULL,
  ModelID = "FirstModel", Verbose = 0, ReturnModelObjects = TRUE,
  SaveModelObjects = FALSE, PassInGrid = NULL)

```

## Arguments

<b>data</b>	This is your data set for training and testing your model
<b>ValidationData</b>	This is your holdout data set used in modeling either refine your hyper-parameters.



TestData	This is your holdout data set. Catboost using both training and validation data in the training process so you should evaluate out of sample performance with this data set.
TargetColumnName	Either supply the target column name OR the column number where the target is located (but not mixed types). Target should be in factor or character form.
FeatureColNames	Either supply the feature column names OR the column number where the target is located (but not mixed types)
IDcols	A vector of column names or column numbers to keep in your data but not include in the modeling.
eval_metric	This is the metric used to identify best grid tuned model. Choose from "merror", "mlogloss"
Trees	The maximum number of trees you want in your models
GridTune	Set to TRUE to run a grid tuning procedure. Set a number in MaxModelsInGrid to tell the procedure how many models you want to test.
grid_eval_metric	Set to "accuracy" (only option currently)
TreeMethod	Choose from "hist", "gpu_hist"
MaxModelsInGrid	Number of models to test from grid options (243 total possible options)
NThreads	Set the maximum number of threads you'd like to dedicate to the model run. E.g. 8
model_path	A character string of your path file to where you want your output saved
ModelID	A character string to name your model and output
Verbose	Set to 0 if you want to suppress model evaluation updates in training
ReturnModelObjects	Set to TRUE to output all modeling objects (E.g. plots and evaluation metrics)
SaveModelObjects	Set to TRUE to return all modeling objects to your environment
PassInGrid	Default is NULL. Provide a data.table of grid options from a previous run.

**Value**

Saves to file and returned in list: VariableImportance.csv, Model, ValidationData.csv, EvaluationMetrics.csv, GridCollect, and GridList

**Author(s)**

Adrian Antico

**See Also**

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoCatBoostRegression](#), [AutoCatBoostScoring](#), [AutoCatBoostHurdleModel](#), [AutoH2OMLScoring](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFClassifier](#), [AutoH2oDRFMultiClass](#), [AutoH2oDRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoXGBoostClassifier](#), [AutoXGBoostRegression](#), [AutoXGBoostScoring](#)

## Examples

```

Correl <- 0.85
N <- 10000
data <- data.table::data.table(Target = runif(N))
data[, x1 := qnorm(Target)]
data[, x2 := runif(N)]
data[, Independent_Variable1 := log(pnorm(Correl * x1 +
                                         sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable2 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable3 := exp(pnorm(Correl * x1 +
                                         sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable4 := exp(exp(pnorm(Correl * x1 +
                                             sqrt(1-Correl^2) * qnorm(x2))))]
data[, Independent_Variable5 := sqrt(pnorm(Correl * x1 +
                                           sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable6 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.10]
data[, Independent_Variable7 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.25]
data[, Independent_Variable8 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.75]
data[, Independent_Variable9 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^2]
data[, Independent_Variable10 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^4]
data[, Target := as.factor(
  ifelse(Independent_Variable2 < 0.20, "A",
    ifelse(Independent_Variable2 < 0.40, "B",
      ifelse(Independent_Variable2 < 0.6, "C",
        ifelse(Independent_Variable2 < 0.8, "D", "E"))))]
data[, Independent_Variable11 := as.factor(
  ifelse(Independent_Variable2 < 0.25, "A",
    ifelse(Independent_Variable2 < 0.35, "B",
      ifelse(Independent_Variable2 < 0.65, "C",
        ifelse(Independent_Variable2 < 0.75, "D", "E"))))]
data[, ':= ' (x1 = NULL, x2 = NULL)]
TestModel <- AutoXGBoostMultiClass(data,
  ValidationData = NULL,
  TestData = NULL,
  TargetColumnName = 1,
  FeatureColNames = 2:12,
  IDcols = NULL,
  eval_metric = "merror",
  Trees = 50,
  GridTune = TRUE,
  grid_eval_metric = "accuracy",
  MaxModelsInGrid = 10,
  NThreads = 8,
  TreeMethod = "hist",
  model_path = getwd(),
  ModelID = "FirstModel",
  ReturnModelObjects = TRUE,
  SaveModelObjects = FALSE,
  PassInGrid = NULL)

```

---

**AutoXGBoostRegression** *AutoXGBoostRegression is an automated XGBoost modeling framework with grid-tuning and model evaluation*

---

## Description

AutoXGBoostRegression is an automated XGBoost modeling framework with grid-tuning and model evaluation that runs a variety of steps. First, the function will run a random grid tune over N number of models and find which model is the best (a default model is always included in that set). Once the model is identified and built, several other outputs are generated: validation data with predictions, evaluation plot, evaluation boxplot, evaluation metrics, variable importance, partial dependence calibration plots, partial dependence calibration box plots, and column names used in model fitting.

## Usage

```
AutoXGBoostRegression(data, ValidationData = NULL, TestData = NULL,
  TargetColumnName = NULL, FeatureColNames = NULL, IDcols = NULL,
  ReturnFactorLevels = FALSE, TransformNumericColumns = NULL,
  eval_metric = "RMSE", Trees = 50, GridTune = FALSE,
  grid_eval_metric = "mae", TreeMethod = "hist",
  MaxModelsInGrid = 10, NThreads = max(1, parallel::detectCores() - 2),
  model_path = NULL, ModelID = "FirstModel", NumOfParDepPlots = 3,
  Verbose = 0, ReturnModelObjects = TRUE, SaveModelObjects = FALSE,
  PassInGrid = NULL)
```

## Arguments

<b>data</b>	This is your data set for training and testing your model
<b>ValidationData</b>	This is your holdout data set used in modeling either refine your hyper-parameters.
<b>TestData</b>	This is your holdout data set. Catboost using both training and validation data in the training process so you should evaluate out of sample performance with this data set.
<b>TargetColumnName</b>	Either supply the target column name OR the column number where the target is located (but not mixed types).
<b>FeatureColNames</b>	Either supply the feature column names OR the column number where the target is located (but not mixed types)
<b>IDcols</b>	A vector of column names or column numbers to keep in your data but not include in the modeling.
<b>ReturnFactorLevels</b>	Set to TRUE to have the factor levels returned with the other model objects
<b>TransformNumericColumns</b>	Set to NULL to do nothing; otherwise supply the column names of numeric variables you want transformed

<code>eval_metric</code>	This is the metric used to identify best grid tuned model. Choose from "r2", "RMSE", "MSE", "MAE"
<code>Trees</code>	The maximum number of trees you want in your models
<code>GridTune</code>	Set to TRUE to run a grid tuning procedure. Set a number in <code>MaxModelsInGrid</code> to tell the procedure how many models you want to test.
<code>grid_eval_metric</code>	Choose from "poisson", "mae", "mape", "mse", "msle", "kl", "cs", "r2"
<code>TreeMethod</code>	Choose from "hist", "gpu_hist"
<code>MaxModelsInGrid</code>	Number of models to test from grid options (243 total possible options)
<code>NThreads</code>	Set the maximum number of threads you'd like to dedicate to the model run. E.g. 8
<code>model_path</code>	A character string of your path file to where you want your output saved
<code>ModelID</code>	A character string to name your model and output
<code>NumOfParDepPlots</code>	Tell the function the number of partial dependence calibration plots you want to create.
<code>Verbose</code>	Set to 0 if you want to suppress model evaluation updates in training
<code>ReturnModelObjects</code>	Set to TRUE to output all modeling objects (E.g. plots and evaluation metrics)
<code>SaveModelObjects</code>	Set to TRUE to return all modeling objects to your environment
<code>PassInGrid</code>	Default is NULL. Provide a data.table of grid options from a previous run.

### Value

Saves to file and returned in list: `VariableImportance.csv`, `Model`, `ValidationData.csv`, `EvaluationPlot.png`, `EvaluationBoxPlot.png`, `EvaluationMetrics.csv`, `ParDepPlots.R` a named list of features with partial dependence calibration plots, `ParDepBoxPlots.R`, `GridCollect`, and `GridList`

### Author(s)

Adrian Antico

### See Also

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoCatBoostRegression](#), [AutoCatBoostScoring](#), [AutoCatBoostHurdleModel](#), [AutoH2OMLScoring](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFClassifier](#), [AutoH2oDRFMultiClass](#), [AutoH2oDRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#), [AutoXGBoostScoring](#)

### Examples

```
Correl <- 0.85
N <- 10000
data <- data.table::data.table(Target = runif(N))
```

```

data[, x1 := qnorm(Target)]
data[, x2 := runif(N)]
data[, Independent_Variable1 := log(pnorm(Correl * x1 +
                                         sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable2 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable3 := exp(pnorm(Correl * x1 +
                                         sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable4 := exp(exp(pnorm(Correl * x1 +
                                              sqrt(1-Correl^2) * qnorm(x2))))]
data[, Independent_Variable5 := sqrt(pnorm(Correl * x1 +
                                           sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable6 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.10]
data[, Independent_Variable7 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.25]
data[, Independent_Variable8 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.75]
data[, Independent_Variable9 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^2]
data[, Independent_Variable10 := (pnorm(Correl * x1 +
                                        sqrt(1-Correl^2) * qnorm(x2)))^4]
data[, Independent_Variable11 := as.factor(
  ifelse(Independent_Variable2 < 0.20, "A",
    ifelse(Independent_Variable2 < 0.40, "B",
      ifelse(Independent_Variable2 < 0.6, "C",
        ifelse(Independent_Variable2 < 0.8, "D", "E")))))]
data[, ':= ' (x1 = NULL, x2 = NULL)]
TestModel <- AutoXGBoostRegression(data,
  ValidationData = NULL,
  TestData = NULL,
  TargetColumnName = 1,
  FeatureColNames = 2:12,
  IDcols = NULL,
  ReturnFactorLevels = FALSE,
  TransformNumericColumns = NULL,
  eval_metric = "RMSE",
  Trees = 50,
  GridTune = TRUE,
  grid_eval_metric = "mae",
  MaxModelsInGrid = 10,
  NThreads = max(1, parallel::detectCores()-2),
  TreeMethod = "hist",
  model_path = getwd(),
  ModelID = "FirstModel",
  NumOfParDepPlots = 3,
  ReturnModelObjects = TRUE,
  SaveModelObjects = FALSE,
  PassInGrid = NULL)

```

## Description

AutoXGBoostScoring is an automated scoring function that compliments the AutoCatBoost model training functions. This function requires you to supply features for scoring. It will run ModelDataPrep() and the DummifyDT() function to prepare your features for xgboost data conversion and scoring.

## Usage

```
AutoXGBoostScoring(TargetType = NULL, ScoringData = NULL,
  FeatureColumnNames = NULL, IDcols = NULL, FactorLevelsList = NULL,
  OneHot = FALSE, ModelObject = NULL, ModelPath = NULL,
  ModelID = NULL, ReturnFeatures = TRUE, TransformNumeric = FALSE,
  BackTransNumeric = FALSE, TargetColumnName = NULL,
  TransformationObject = NULL, TransID = NULL, TransPath = NULL,
  MDP_Impute = TRUE, MDP_CharToFactor = TRUE, MDP_RemoveDates = TRUE,
  MDP_MissFactor = "0", MDP_MissNum = -1)
```

## Arguments

TargetType	Set this value to "regression", "classification", or "multiclass" to score models built using AutoCatBoostRegression(), AutoCatBoostClassify() or AutoCatBoostMultiClass().
ScoringData	This is your data.table of features for scoring. Can be a single row or batch.
FeatureColumnNames	Supply either column names or column numbers used in the AutoXGBoost__() function
IDcols	Supply ID column numbers for any metadata you want returned with your predicted values
FactorLevelsList	Supply the factor variables' list from DummifyDT()
OneHot	Set to TRUE to have one-hot-encoding run. Otherwise, N columns will be made for N levels of a factor variable
ModelObject	Supply a model for scoring, otherwise it will have to search for it in the file path you specify
ModelPath	Supply your path file used in the AutoXGBoost__() function
ModelID	Supply the model ID used in the AutoXGBoost__() function
ReturnFeatures	Set to TRUE to return your features with the predicted values.
TransformNumeric	Set to TRUE if you have features that were transformed automatically from an Auto__Regression() model AND you haven't already transformed them.
BackTransNumeric	Set to TRUE to generate back-transformed predicted values. Also, if you return features, those will also be back-transformed.
TargetColumnName	Input your target column name used in training if you are utilizing the transformation service

**TransformationObject**

Set to NULL if you didn't use transformations or if you want the function to pull from the file output from the `Auto__Regression()` function. You can also supply the transformation data.table object with the transformation details versus having it pulled from file.

**TransID** Set to the ID used for saving the transformation data.table object or set it to the ModelID if you are pulling from file from a build with `Auto__Regression()`.

**TransPath** Set the path file to the folder where your transformation data.table detail object is stored. If you used the `Auto__Regression()` to build, set it to the same path as ModelPath.

**MDP\_Impute** Set to TRUE if you did so for modeling and didn't do so before supplying ScoringData in this function

**MDP\_CharToFactor** Set to TRUE to turn your character columns to factors if you didn't do so to your ScoringData that you are supplying to this function

**MDP\_RemoveDates** Set to TRUE if you have date of timestamp columns in your ScoringData

**MDP\_MissFactor** If you set MDP\_Impute to TRUE, supply the character values to replace missing values with

**MDP\_MissNum** If you set MDP\_Impute to TRUE, supply a numeric value to replace missing values with

**Value**

A data.table of predicted values with the option to return model features as well.

**See Also**

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoCatBoostRegression](#), [AutoCatBoostScoring](#), [AutoCatBoostHurdleModel](#), [AutoH2OMLScoring](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2ODRFClassifier](#), [AutoH2ODRFMultiClass](#), [AutoH2ODRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#)

**Examples**

```
Preds <- AutoXGBoostScoring(TargetType = "regression",
                             ScoringData = data,
                             FeatureColumnNames = 2:12,
                             IDcols = NULL,
                             FactorLevelsList = NULL,
                             OneHot = FALSE,
                             ModelObject = NULL,
                             ModelPath = "home",
                             ModelID = "ModelTest",
                             ReturnFeatures = TRUE,
                             MDP_Impute = TRUE,
                             MDP_CharToFactor = TRUE,
                             MDP_RemoveDates = TRUE,
                             MDP_MissFactor = "0",
                             MDP_MissNum = -1)
```

---

ChartTheme	<i>ChartTheme function is a ggplot theme generator for ggplots</i>
------------	--

---

## Description

This function helps your ggplots look professional with the choice of the two main colors that will dominate the theme

## Usage

```
ChartTheme(Size = 12)
```

## Arguments

Size                      The size of the axis labels and title

## Value

An object to pass along to ggplot objects following the "+" sign

## Author(s)

Adrian Antico

## See Also

Other Misc: [AutoH20TextPrepScoring](#), [PrintObjectsSize](#), [RemixTheme](#), [SimpleCap](#), [multiplot](#), [percRank](#), [tempDatesFun](#), [tokenizeH20](#)

## Examples

```
data <- data.table::data.table(DateTime = as.Date(Sys.time()),
  Target = stats::filter(rnorm(1000,
    mean = 50,
    sd = 20),
    filter=rep(1,10),
    circular=TRUE))
data[, temp := seq(1:1000)][, DateTime := DateTime - temp][, temp := NULL]
data <- data[order(DateTime)]
p <- ggplot2::ggplot(data, ggplot2::aes(x = DateTime, y = Target)) + ggplot2::geom_line()
p <- p + ChartTheme(Size = 12)
```



---

**CreateCalendarVariables***CreateCalendarVariables Create Calendar Variables*

---

**Description**

CreateCalendarVariables Rapidly creates calendar variables based on the date column you provide

**Usage**

```
CreateCalendarVariables(data, DateCols = c("Date", "Date2"),
  AsFactor = FALSE, TimeUnits = "wday")
```

**Arguments**

<b>data</b>	This is your data
<b>DateCols</b>	Supply either column names or column numbers of your date columns you want to use for creating calendar variables
<b>AsFactor</b>	Set to TRUE if you want factor type columns returned; otherwise integer type columns will be returned
<b>TimeUnits</b>	Supply a character vector of time units for creating calendar variables. Options include: "second", "minute", "hour", "wday", "mday", "yday", "week", "isoweek", "month", "quarter", "year"

**Value**

Returns your data.table with the added calendar variables at the end

**Author(s)**

Adrian Antico

**See Also**

Other Feature Engineering: [AutoDataPartition](#), [AutoTransformationCreate](#), [AutoTransformationScore](#), [AutoWord2VecModeler](#), [DT\\_GDL\\_Feature\\_Engineering](#), [DummifyDT](#), [GDL\\_Feature\\_Engineering](#), [ModelDataPrep](#), [Partial\\_DT\\_GDL\\_Feature\\_Engineering2](#), [Partial\\_DT\\_GDL\\_Feature\\_Engineering](#), [Scoring\\_GDL\\_Feature\\_Engineering](#)

**Examples**

```
data <- data.table::data.table(Date = "2018-01-01 00:00:00")
data <- CreateCalendarVariables(data,
  DateCols = "Date",
  AsFactor = FALSE,
  TimeUnits = c("wday", "month", "year"))
```

---

DT\_GDL\_Feature\_Engineering

*An Automated Feature Engineering Function Using data.table  
frollmean*

---

## Description

Builds autoregressive and moving average from target columns and distributed lags and distributed moving average for independent features distributed across time. On top of that, you can also create time between instances along with their associated lags and moving averages. This function works for data with groups and without groups.

## Usage

```
DT_GDL_Feature_Engineering(data, lags = c(seq(1, 50, 1)),
  periods = c(seq(5, 95, 5)), statsNames = c("MA"),
  targets = c("qty"), groupingVars = c("Group1", "Group2"),
  sortDateName = c("date"), timeDiffTarget = c("TimeDiffName"),
  timeAgg = c("days"), WindowingLag = 0, Type = c("Lag"),
  Timer = TRUE, SimpleImpute = TRUE)
```

## Arguments

<code>data</code>	A data.table you want to run the function on
<code>lags</code>	A numeric vector of the specific lags you want to have generated. You must include 1 if WindowingLag = 1.
<code>periods</code>	A numeric vector of the specific rolling statistics window sizes you want to utilize in the calculations.
<code>statsNames</code>	A character vector of the corresponding names to create for the rollings stats variables.
<code>targets</code>	A character vector of the column names for the reference column in which you will build your lags and rolling stats
<code>groupingVars</code>	A character vector of categorical variable names you will build your lags and rolling stats by
<code>sortDateName</code>	The column name of your date column used to sort events over time
<code>timeDiffTarget</code>	Specify a desired name for features created for time between events. Set to NULL if you don't want time between events features created.
<code>timeAgg</code>	List the time aggregation level for the time between events features, such as "hour", "day", "week", "month", "quarter", or "year"
<code>WindowingLag</code>	Set to 0 to build rolling stats off of target columns directly or set to 1 to build the rolling stats off of the lag-1 target
<code>Type</code>	List either "Lag" if you want features built on historical values or "Lead" if you want features built on future values
<code>Timer</code>	Set to TRUE if you percentage complete tracker printout
<code>SimpleImpute</code>	Set to TRUE for factor level imputation of "0" and numeric imputation of -1

**Value**

data.table of original data plus created lags, rolling stats, and time between event lags and rolling stats

**Author(s)**

Adrian Antico

**See Also**

Other Feature Engineering: [AutoDataPartition](#), [AutoTransformationCreate](#), [AutoTransformationScore](#), [AutoWord2VecModeler](#), [CreateCalendarVariables](#), [DummifyDT](#), [GDL\\_Feature\\_Engineering](#), [ModelDataPrep](#), [Partial\\_DT\\_GDL\\_Feature\\_Engineering2](#), [Partial\\_DT\\_GDL\\_Feature\\_Engineering](#), [Scoring\\_GDL\\_Feature\\_Engineering](#)

**Examples**

```
N = 25116
data <- data.table::data.table(DateTime = as.Date(Sys.time()),
                               Target = stats::filter(rnorm(N,
                                                           mean = 50,
                                                           sd = 20),
                                                       filter=rep(1,10),
                                                       circular=TRUE))

data[, temp := seq(1:N)][, DateTime := DateTime - temp][, temp := NULL]
data <- data[order(DateTime)]
data <- DT_GDL_Feature_Engineering(data,
                                  lags           = c(seq(1,5,1)),
                                  periods        = c(3,5,10,15,20,25),
                                  statsNames     = c("MA"),
                                  targets        = c("Target"),
                                  groupingVars   = NULL,
                                  sortDateName  = "DateTime",
                                  timeDiffTarget = c("Time_Gap"),
                                  timeAgg       = c("days"),
                                  WindowingLag  = 1,
                                  Type           = "Lag",
                                  Timer          = TRUE,
                                  SimpleImpute  = TRUE)
```

---

DummifyDT

*DummifyDT creates dummy variables for the selected columns.*


---

**Description**

DummifyDT creates dummy variables for the selected columns. Either one-hot encoding, N+1 columns for N levels, or N columns for N levels.

**Usage**

```
DummifyDT(data, cols, KeepFactorCols = FALSE, OneHot = FALSE,
           SaveFactorLevels = FALSE, SavePath = NULL,
           ImportFactorLevels = FALSE, FactorLevelsList = NULL,
           ClustScore = FALSE, ReturnFactorLevels = FALSE)
```

**Arguments**

<code>data</code>	The data set to run the micro auc on
<code>cols</code>	A vector with the names of the columns you wish to dichotomize
<code>KeepFactorCols</code>	Set to TRUE to keep the original columns used in the dichotomization process
<code>OneHot</code>	Set to TRUE to run one hot encoding, FALSE to generate N columns for N levels
<code>SaveFactorLevels</code>	Set to TRUE to save unique levels of each factor column to file as a csv
<code>SavePath</code>	Provide a file path to save your factor levels. Use this for models that you have to create dummy variables for.
<code>ImportFactorLevels</code>	Instead of using the data you provide, import the factor levels csv to ensure you build out all of the columns you trained with in modeling.
<code>FactorLevelsList</code>	Supply a list of factor variable levels
<code>ClustScore</code>	This is for scoring AutoKMeans. Set to FALSE for all other applications.
<code>ReturnFactorLevels</code>	If you want a named list of all the factor levels returned, set this to TRUE. Doing so will cause the function to return a list with the source data.table and the list of factor variables' levels

**Value**

Either a data table with new dummy variables columns and optionally removes base columns (if `ReturnFactorLevels` is FALSE), otherwise a list with the data.table and a list of the factor levels.

**Author(s)**

Adrian Antico

**See Also**

Other Feature Engineering: [AutoDataPartition](#), [AutoTransformationCreate](#), [AutoTransformationScore](#), [AutoWord2VecModeler](#), [CreateCalendarVariables](#), [DT\\_GDL\\_Feature\\_Engineering](#), [GDL\\_Feature\\_Engineering](#), [ModelDataPrep](#), [Partial\\_DT\\_GDL\\_Feature\\_Engineering2](#), [Partial\\_DT\\_GDL\\_Feature\\_Engineering](#), [Scoring\\_GDL\\_Feature\\_Engineering](#)

**Examples**

```
test <- data.table::data.table(Value = runif(100000),
                               FactorCol = sample(x = c(letters,
                                                         LETTERS,
                                                         paste0(letters,letters),
                                                         paste0(LETTERS,LETTERS),
                                                         paste0(letters,LETTERS),
                                                         paste0(LETTERS,letters)),
                               size = 100000,
                               replace = TRUE))

test <- DummifyDT(data = test,
                  cols,
```

```

        KeepFactorCols = FALSE,
        OneHot = FALSE,
        SaveFactorLevels = FALSE,
        SavePath = NULL,
        ImportFactorLevels = FALSE,
        FactorLevelsList = NULL,
        ClustScore = FALSE,
        ReturnFactorLevels = FALSE)
ncol(test)
test[, sum(FactorCol_gg)]

```

---

EvalPlot	<i>EvalPlot automatically builds calibration plots for model evaluation</i>
----------	---

---

## Description

This function automatically builds calibration plots and calibration boxplots for model evaluation using regression, quantile regression, and binary and multinomial classification

## Usage

```

EvalPlot(data, PredictionColName = c("PredictedValues"),
        TargetColName = c("ActualValues"), GraphType = c("calibration"),
        PercentileBucket = 0.05, aggrfun = function(x) mean(x, na.rm = TRUE))

```

## Arguments

data	Data containing predicted values and actual values for comparison
PredictionColName	String representation of column name with predicted values from model
TargetColName	String representation of column name with target values from model
GraphType	Calibration or boxplot - calibration aggregated data based on summary statistic; boxplot shows variation
PercentileBucket	Number of buckets to partition the space on (0,1) for evaluation
aggrfun	The statistics function used in aggregation, listed as a function

## Value

Calibration plot or boxplot

## Author(s)

Adrian Antico

## See Also

Other Model Evaluation and Interpretation: [ParDepCalPlots](#), [RedYellowGreen](#), [threshOptim](#)

## Examples

```
Correl <- 0.85
data <- data.table::data.table(Target = runif(100))
data[, x1 := qnorm(Target)]
data[, x2 := runif(100)]
data[, Independent_Variable1 := log(pnorm(Correl * x1 +
                                         sqrt(1-Correl^2) * qnorm(x2)))]

data[, Predict := (pnorm(Correl * x1 +
                        sqrt(1-Correl^2) * qnorm(x2)))]

EvalPlot(data,
          PredictionColName = "Predict",
          TargetColName = "Target",
          GraphType = "calibration",
          PercentileBucket = 0.05,
          aggrfun = function(x) quantile(x, probs = 0.50, na.rm = TRUE))
```

---

GDL\_Feature\_Engineering

*An Automated Feature Engineering Function*

---

## Description

Builds autoregressive and rolling stats from target columns and distributed lags and distributed rolling stats for independent features distributed across time. On top of that, you can also create time between instances along with their associated lags and rolling stats. This function works for data with groups and without groups.

## Usage

```
GDL_Feature_Engineering(data, lags = seq(1, 5, 1), periods = seq(1, 5,
1), statsFUNs = "mean", statsNames = "mean", targets = NULL,
groupingVars = NULL, sortDateName = NULL, timeDiffTarget = NULL,
timeAgg = NULL, WindowingLag = 1, Type = "Lag", Timer = TRUE,
SkipCols = NULL, SimpleImpute = TRUE)
```

## Arguments

<b>data</b>	A data.table you want to run the function on
<b>lags</b>	A numeric vector of the specific lags you want to have generated. You must include 1 if WindowingLag = 1.
<b>periods</b>	A numeric vector of the specific rolling statistics window sizes you want to utilize in the calculations.
<b>statsFUNs</b>	Vector that holds functions for your rolling stats, such as function(x) mean(x), function(x) sd(x), or function(x) quantile(x)
<b>statsNames</b>	A character vector of the corresponding names to create for the rollings stats variables.
<b>targets</b>	A character vector of the column names for the reference column in which you will build your lags and rolling stats
<b>groupingVars</b>	A character vector of categorical variable names you will build your lags and rolling stats by

<code>sortDateName</code>	The column name of your date column used to sort events over time
<code>timeDiffTarget</code>	Specify a desired name for features created for time between events. Set to NULL if you don't want time between events features created.
<code>timeAgg</code>	List the time aggregation level for the time between events features, such as "hour", "day", "week", "month", "quarter", or "year"
<code>WindowingLag</code>	Set to 0 to build rolling stats off of target columns directly or set to 1 to build the rolling stats off of the lag-1 target
<code>Type</code>	List either "Lag" if you want features built on historical values or "Lead" if you want features built on future values
<code>Timer</code>	Set to TRUE if you percentage complete tracker printout
<code>SkipCols</code>	Defaults to NULL; otherwise supply a character vector of the names of columns to skip
<code>SimpleImpute</code>	Set to TRUE for factor level imputation of "0" and numeric imputation of -1

**Value**

data.table of original data plus created lags, rolling stats, and time between event lags and rolling stats

**Author(s)**

Adrian Antico

**See Also**

Other Feature Engineering: [AutoDataPartition](#), [AutoTransformationCreate](#), [AutoTransformationScore](#), [AutoWord2VecModeler](#), [CreateCalendarVariables](#), [DT\\_GDL\\_Feature\\_Engineering](#), [DummifyDT](#), [ModelDataPrep](#), [Partial\\_DT\\_GDL\\_Feature\\_Engineering2](#), [Partial\\_DT\\_GDL\\_Feature\\_Engineering](#), [Scoring\\_GDL\\_Feature\\_Engineering](#)

**Examples**

```

N = 25116
data <- data.table::data.table(DateTime = as.Date(Sys.time()),
  Target = stats::filter(rnorm(N,
    mean = 50,
    sd = 20),
    filter=rep(1,10),
    circular=TRUE))
data[, temp := seq(1:N)][, DateTime := DateTime - temp][, temp := NULL]
data <- data[order(DateTime)]
data <- GDL_Feature_Engineering(data,
  lags          = c(seq(1,1,1)),
  periods       = c(3),
  statsFUNs     = "mean",
  statsNames    = "mean",
  targets       = c("Target"),
  groupingVars  = NULL,
  sortDateName  = "DateTime",
  timeDiffTarget = NULL,
  timeAgg       = "days",
  WindowingLag  = 1,

```

```

Type          = "Lag",
Timer         = TRUE,
SkipCols      = FALSE,
SimpleImpute  = TRUE)

```

---

GenTSAnomVars	<i>GenTSAnomVars is an automated z-score anomaly detection via GLM-like procedure</i>
---------------	---

---

## Description

GenTSAnomVars is an automated z-score anomaly detection via GLM-like procedure. Data is z-scaled and grouped by factors and time periods to determine which points are above and below the control limits in a cumulative time fashion. Then a cumulative rate is created as the final variable. Set KeepAllCols to FALSE to utilize the intermediate features to create rolling stats from them. The anomalies are separated into those that are extreme on the positive end versus those that are on the negative end.

## Usage

```

GenTSAnomVars(data, ValueCol = "Value", GroupVar1 = "SKU",
  GroupVar2 = NULL, DateVar = "DATE", HighThreshold = 1.96,
  LowThreshold = -1.96, KeepAllCols = FALSE, IsDataScaled = TRUE)

```

## Arguments

data	the source residuals data.table
ValueCol	the numeric column to run anomaly detection over
GroupVar1	this is a group by variable
GroupVar2	this is another group by variable
DateVar	this is a time variable for grouping
HighThreshold	this is the threshold on the high end
LowThreshold	this is the threshold on the low end
KeepAllCols	set to TRUE to remove the intermediate features
IsDataScaled	set to TRUE if you already scaled your data

## Value

The original data.table with the added columns merged in. When KeepAllCols is set to FALSE, you will get back two columns: AnomHighRate and AnomLowRate - these are the cumulative anomaly rates over time for when you get anomalies from above the thresholds (e.g. 1.96) and below the thresholds.

## Author(s)

Adrian Antico

## See Also

Other Unsupervised Learning: [AutoKMeans](#), [ResidualOutliers](#)



## Examples

```
data <- data.table::data.table(DateTime = as.Date(Sys.time()),
  Target = stats::filter(rnorm(10000,
    mean = 50,
    sd = 20),
    filter=rep(1,10),
    circular=TRUE))
data[, temp := seq(1:10000)][, DateTime := DateTime - temp][, temp := NULL]
data <- data[order(DateTime)]
x <- data.table::as.data.table(sde::GBM(N=10000)*1000)
data[, predicted := x[-1,]]
stuff <- GenTSAnomVars(data,
  ValueCol = "Target",
  GroupVar1 = NULL,
  GroupVar2 = NULL,
  DateVar = "DateTime",
  HighThreshold = 1.96,
  LowThreshold = -1.96,
  KeepAllCols = TRUE,
  IsDataScaled = FALSE)
```

---

ModelDataPrep

*Final Data Preparation Function*


---

## Description

This function replaces inf values with NA, converts characters to factors, and imputes with constants

## Usage

```
ModelDataPrep(data, Impute = TRUE, CharToFactor = TRUE,
  RemoveDates = FALSE, MissFactor = "0", MissNum = -1,
  IgnoreCols = NULL)
```

## Arguments

<code>data</code>	This is your source data you'd like to modify
<code>Impute</code>	Defaults to TRUE which tells the function to impute the data
<code>CharToFactor</code>	Defaults to TRUE which tells the function to convert characters to factors
<code>RemoveDates</code>	Defaults to FALSE. Set to TRUE to remove date columns from your data.table
<code>MissFactor</code>	Supply the value to impute missing factor levels
<code>MissNum</code>	Supply the value to impute missing numeric values
<code>IgnoreCols</code>	Supply column numbers for columns you want the function to ignore

## Value

Returns the original data table with corrected values

**Author(s)**

Adrian Antico

**See Also**

Other Feature Engineering: [AutoDataPartition](#), [AutoTransformationCreate](#), [AutoTransformationScore](#), [AutoWord2VecModeler](#), [CreateCalendarVariables](#), [DT\\_GDL\\_Feature\\_Engineering](#), [DummifyDT](#), [GDL\\_Feature\\_Engineering](#), [Partial\\_DT\\_GDL\\_Feature\\_Engineering2](#), [Partial\\_DT\\_GDL\\_Feature\\_Engineering](#), [Scoring\\_GDL\\_Feature\\_Engineering](#)

**Examples**

```
data <- data.table::data.table(Value = runif(100000),
                              FactorCol = as.character(sample(x = c(letters,
                                                                    LETTERS,
                                                                    paste0(letters, letters),
                                                                    paste0(LETTERS, LETTERS),
                                                                    paste0(letters, LETTERS),
                                                                    paste0(LETTERS, letters)),
                                                                    size = 100000,
                                                                    replace = TRUE)))

data <- ModelDataPrep(data,
                      Impute = TRUE,
                      CharToFactor = TRUE,
                      MissFactor = "0",
                      MissNum    = -1)
```

---

multiplot

---

*Multiplot is a function for combining multiple plots*


---

**Description**

Sick of copying this one into your code? Well, not anymore.

**Usage**

```
multiplot(..., plotlist = NULL, cols = 2, layout = NULL)
```

**Arguments**

...	Passthrough arguments
plotlist	This is the list of your charts
cols	This is the number of columns in your multiplot
layout	Leave NULL

**Value**

Multiple ggplots on a single image

**Author(s)**

Adrian Antico

**See Also**

Other Misc: [AutoH2OTextPrepScoring](#), [ChartTheme](#), [PrintObjectsSize](#), [RemixTheme](#), [SimpleCap](#), [percRank](#), [tempDatesFun](#), [tokenizeH2O](#)

**Examples**

```
Correl <- 0.85
data <- data.table::data.table(Target = runif(100))
data[, x1 := qnorm(Target)]
data[, x2 := runif(100)]
data[, Independent_Variable1 := log(pnorm(Correl * x1 +
                                          sqrt(1-Correl^2) * qnorm(x2)))]

data[, Predict := (pnorm(Correl * x1 +
                        sqrt(1-Correl^2) * qnorm(x2)))]
p1 <- RemixAutoML::ParDepCalPlots(data,
                                PredictionColName = "Predict",
                                TargetColName = "Target",
                                IndepVar = "Independent_Variable1",
                                GraphType = "calibration",
                                PercentileBucket = 0.20,
                                FactLevels = 10,
                                Function = function(x) mean(x, na.rm = TRUE))
p2 <- RemixAutoML::ParDepCalPlots(data,
                                PredictionColName = "Predict",
                                TargetColName = "Target",
                                IndepVar = "Independent_Variable1",
                                GraphType = "boxplot",
                                PercentileBucket = 0.20,
                                FactLevels = 10,
                                Function = function(x) mean(x, na.rm = TRUE))
RemixAutoML::multiplot(plotlist = list(p1,p2), cols = 2)
```

---

ParDepCalPlots

---

*ParDepCalPlots automatically builds partial dependence calibration plots for model evaluation*


---

**Description**

This function automatically builds partial dependence calibration plots and partial dependence calibration boxplots for model evaluation using regression, quantile regression, and binary and multinomial classification

**Usage**

```
ParDepCalPlots(data, PredictionColName = c("PredictedValues"),
               TargetColName = c("ActualValues"),
               IndepVar = c("Independent_Variable_Name"),
               GraphType = c("calibration"), PercentileBucket = 0.05,
               FactLevels = 10, Function = function(x) base::mean(x, na.rm = TRUE))
```

**Arguments**

<code>data</code>	Data containing predicted values and actual values for comparison
<code>PredictionColName</code>	Predicted values column names
<code>TargetColName</code>	Target value column names
<code>IndepVar</code>	Independent variable column names
<code>GraphType</code>	calibration or boxplot - calibration aggregated data based on summary statistic; boxplot shows variation
<code>PercentileBucket</code>	Number of buckets to partition the space on (0,1) for evaluation
<code>FactLevels</code>	The number of levels to show on the chart (1. Levels are chosen based on frequency; 2. all other levels grouped and labeled as "Other")
<code>Function</code>	Supply the function you wish to use for aggregation.

**Value**

Partial dependence calibration plot or boxplot

**Author(s)**

Adrian Antico

**See Also**

Other Model Evaluation and Interpretation: [EvalPlot](#), [RedYellowGreen](#), [threshOptim](#)

**Examples**

```
Correl <- 0.85
data <- data.table::data.table(Target = runif(100))
data[, x1 := qnorm(Target)]
data[, x2 := runif(100)]
data[, Independent_Variable1 := log(pnorm(Correl * x1 +
                                          sqrt(1-Correl^2) * qnorm(x2)))]
data[, Predict := (pnorm(Correl * x1 +
                        sqrt(1-Correl^2) * qnorm(x2)))]
p1 <- RemixAutoML::ParDepCalPlots(data,
                                PredictionColName = "Predict",
                                TargetColName = "Target",
                                IndepVar = "Independent_Variable1",
                                GraphType = "calibration",
                                PercentileBucket = 0.20,
                                FactLevels = 10,
                                Function = function(x) mean(x, na.rm = TRUE))

p1
```

---

**Partial\_DT\_GDL\_Feature\_Engineering**

*A version of the DT\_GDL function for creating the GDL features for a subset of records in your data*

---

**Description**

A version of the DT\_GDL function for creating the GDL features for a subset of records in your data. Create the same lags, moving averages, and the like off of time between records. 100

For scoring models in production that have  $i$  1 grouping variables and for when you need  $i$  1 record (or records per grouping variables) returned. This function is for generating lags and moving averages (along with lags and moving averages off of time between records), for a partial set of records in your data set, typical new records that become available for model scoring. Column names and ordering will be identical to the output from the corresponding DT\_GDL\_Feature\_Engineering() function, which most likely was used to create features for model training.

**Usage**

```
Partial_DT_GDL_Feature_Engineering(data, lags = c(seq(1, 5, 1)),
  periods = c(3, 5, 10, 15, 20, 25), statsNames = c("MA"),
  targets = c("Target"), groupingVars = NULL,
  sortDateName = c("DateTime"), timeDiffTarget = c("Time_Gap"),
  timeAgg = "days", WindowingLag = 1, Type = "Lag", Timer = TRUE,
  SimpleImpute = TRUE, AscRowByGroup = "temp", RecordsKeep = 1,
  AscRowRemove = TRUE)
```

```
Partial_DT_GDL_Feature_Engineering(data, lags = c(seq(1, 5, 1)),
  periods = c(3, 5, 10, 15, 20, 25), statsNames = c("MA"),
  targets = c("Target"), groupingVars = NULL,
  sortDateName = c("DateTime"), timeDiffTarget = c("Time_Gap"),
  timeAgg = "days", WindowingLag = 1, Type = "Lag", Timer = TRUE,
  SimpleImpute = TRUE, AscRowByGroup = "temp", RecordsKeep = 1,
  AscRowRemove = TRUE)
```

**Arguments**

<b>data</b>	A data.table you want to run the function on
<b>lags</b>	A numeric vector of the specific lags you want to have generated. You must include 1 if WindowingLag = 1.
<b>periods</b>	A numeric vector of the specific rolling statistics window sizes you want to utilize in the calculations.
<b>statsNames</b>	A character vector of the corresponding names to create for the rollings stats variables.
<b>targets</b>	A character vector of the column names for the reference column in which you will build your lags and rolling stats
<b>groupingVars</b>	A character vector of categorical variable names you will build your lags and rolling stats by

sortDateName	The column name of your date column used to sort events over time
timeDiffTarget	Specify a desired name for features created for time between events. Set to NULL if you don't want time between events features created.
timeAgg	List the time aggregation level for the time between events features, such as "hour", "day", "week", "month", "quarter", or "year"
WindowingLag	Set to 0 to build rolling stats off of target columns directly or set to 1 to build the rolling stats off of the lag-1 target
Type	List either "Lag" if you want features built on historical values or "Lead" if you want features built on future values
Timer	Set to TRUE if you percentage complete tracker printout
SimpleImpute	Set to TRUE for factor level imputation of "0" and numeric imputation of -1
AscRowByGroup	Required to have a column with a Row Number by group (if grouping) with 1 being the record for scoring (typically the most current in time)
RecordsKeep	List the number of records to retain (1 for last record, 2 for last 2 records, etc.)
AscRowRemove	Set to TRUE to remove the AscRowByGroup column upon returning data.
statsFUNs	Vector of functions for your rolling windows, such as mean, sd, min, max, quantile
SkipCols	Defaults to NULL; otherwise supply a character vector of the names of columns to skip
data	A data.table you want to run the function on
lags	A numeric vector of the specific lags you want to have generated. You must include 1 if WindowingLag = 1.
periods	A numeric vector of the specific rolling statistics window sizes you want to utilize in the calculations.
statsFUNs	Vector of functions for your rolling windows, such as mean, sd, min, max, quantile
statsNames	A character vector of the corresponding names to create for the rollings stats variables.
targets	A character vector of the column names for the reference column in which you will build your lags and rolling stats
groupingVars	A character vector of categorical variable names you will build your lags and rolling stats by
sortDateName	The column name of your date column used to sort events over time
timeDiffTarget	Specify a desired name for features created for time between events. Set to NULL if you don't want time between events features created.
timeAgg	List the time aggregation level for the time between events features, such as "hour", "day", "week", "month", "quarter", or "year"
WindowingLag	Set to 0 to build rolling stats off of target columns directly or set to 1 to build the rolling stats off of the lag-1 target
Type	List either "Lag" if you want features built on historical values or "Lead" if you want features built on future values
Timer	Set to TRUE if you percentage complete tracker printout

SkipCols	Defaults to NULL; otherwise supply a character vector of the names of columns to skip
SimpleImpute	Set to TRUE for factor level imputation of "0" and numeric imputation of -1
AscRowByGroup	Required to have a column with a Row Number by group (if grouping) with the smallest numbers being the records for scoring (typically the most current in time).
RecordsKeep	List the number of records to retain (1 for last record, 2 for last 2 records, etc.)

**Value**

data.table of original data plus created lags, rolling stats, and time between event lags and rolling stats

data.table of original data plus created lags, rolling stats, and time between event lags and rolling stats

**Author(s)**

Adrian Antico

Adrian Antico

**See Also**

Other Feature Engineering: [AutoDataPartition](#), [AutoTransformationCreate](#), [AutoTransformationScore](#), [AutoWord2VecModeler](#), [CreateCalendarVariables](#), [DT\\_GDL\\_Feature\\_Engineering](#), [DummifyDT](#), [GDL\\_Feature\\_Engineering](#), [ModelDataPrep](#), [Partial\\_DT\\_GDL\\_Feature\\_Engineering2](#), [Scoring\\_GDL\\_Feature\\_Engi](#)

Other Feature Engineering: [AutoDataPartition](#), [AutoTransformationCreate](#), [AutoTransformationScore](#), [AutoWord2VecModeler](#), [CreateCalendarVariables](#), [DT\\_GDL\\_Feature\\_Engineering](#), [DummifyDT](#), [GDL\\_Feature\\_Engineering](#), [ModelDataPrep](#), [Partial\\_DT\\_GDL\\_Feature\\_Engineering2](#), [Scoring\\_GDL\\_Feature\\_Engi](#)

**Examples**

```

N = 25116
data <- data.table::data.table(DateTime = as.Date(Sys.time()),
  Target = stats::filter(rnorm(N,
    mean = 50,
    sd = 20),
    filter=rep(1,10),
    circular=TRUE))
data[, temp := seq(1:N)][, DateTime := DateTime - temp]
data <- data[order(DateTime)]
data <- Partial_DT_GDL_Feature_Engineering(data,
  lags          = c(1:5),
  periods       = c(seq(10,50,10)),
  statsNames    = "mean",
  targets       = c("Target"),
  groupingVars  = NULL,
  sortDateName  = "DateTime",
  timeDiffTarget = c("Time_Gap"),
  timeAgg       = "days",
  WindowingLag  = 1,
  Type          = "Lag",
  Timer         = TRUE,

```

```

SkipCols      = FALSE,
SimpleImpute  = TRUE,
AscRowByGroup = "temp",
RecordsKeep   = 1)

N = 25116
data <- data.table::data.table(DateTime = as.Date(Sys.time()),
  Target = stats::filter(rnorm(N,
    mean = 50,
    sd = 20),
    filter=rep(1,10),
    circular=TRUE))
data[, temp := seq(1:N)][, DateTime := DateTime - temp]
data <- data[order(DateTime)]
data <- Partial_DT_GDL_Feature_Engineering(data,
  lags      = c(1:5),
  periods   = c(seq(10,50,10)),
  statsNames = "mean",
  targets    = c("Target"),
  groupingVars = NULL,
  sortDateName = "DateTime",
  timeDiffTarget = c("Time_Gap"),
  timeAgg     = "days",
  WindowingLag = 1,
  Type        = "Lag",
  Timer       = TRUE,
  SimpleImpute = TRUE,
  AscRowByGroup = "temp",
  RecordsKeep  = 1,
  AscRowRemove = TRUE)

```

---

### Partial\_DT\_GDL\_Feature\_Engineering2

*A version of the DT\_GDL function for creating the GDL features for a new set of records*

---

## Description

For scoring models in production that have  $i \geq 1$  grouping variables and for when you need  $j \geq 1$  record (or records per grouping variables) returned. This function is for generating lags and moving averages (along with lags and moving averages off of time between records), for a partial set of records in your data set, typical new records that become available for model scoring. Column names and ordering will be identical to the output from the corresponding DT\_GDL\_Feature\_Engineering() function, which most likely was used to create features for model training.

## Usage

```

Partial_DT_GDL_Feature_Engineering2(data, lags = c(seq(1, 5, 1)),
  periods = c(3, 5, 10, 15, 20, 25), statsNames = c("MA"),
  targets = c("Target"), groupingVars = NULL,
  sortDateName = c("DateTime"), timeDiffTarget = c("Time_Gap"),
  timeAgg = "days", WindowingLag = 1, Type = "Lag", Timer = TRUE,
  SimpleImpute = TRUE, AscRowByGroup = "temp", RecordsKeep = 1,
  AscRowRemove = TRUE)

```



**Arguments**

<b>data</b>	A data.table you want to run the function on
<b>lags</b>	A numeric vector of the specific lags you want to have generated. You must include 1 if WindowingLag = 1.
<b>periods</b>	A numeric vector of the specific rolling statistics window sizes you want to utilize in the calculations.
<b>statsNames</b>	A character vector of the corresponding names to create for the rollings stats variables.
<b>targets</b>	A character vector of the column names for the reference column in which you will build your lags and rolling stats
<b>groupingVars</b>	A character vector of categorical variable names you will build your lags and rolling stats by
<b>sortDateName</b>	The column name of your date column used to sort events over time
<b>timeDiffTarget</b>	Specify a desired name for features created for time between events. Set to NULL if you don't want time between events features created.
<b>timeAgg</b>	List the time aggregation level for the time between events features, such as "hour", "day", "week", "month", "quarter", or "year"
<b>WindowingLag</b>	Set to 0 to build rolling stats off of target columns directly or set to 1 to build the rolling stats off of the lag-1 target
<b>Type</b>	List either "Lag" if you want features built on historical values or "Lead" if you want features built on future values
<b>Timer</b>	Set to TRUE if you percentage complete tracker printout
<b>SimpleImpute</b>	Set to TRUE for factor level imputation of "0" and numeric imputation of -1
<b>AscRowByGroup</b>	Required to have a column with a Row Number by group (if grouping) with the smallest numbers being the records for scoring (typically the most current in time).
<b>RecordsKeep</b>	List the number of records to retain (1 for last record, 2 for last 2 records, etc.)
<b>AscRowRemove</b>	Set to TRUE to remove the AscRowByGroup column upon returning data.
<b>statsFUNs</b>	Vector of functions for your rolling windows, such as mean, sd, min, max, quantile
<b>SkipCols</b>	Defaults to NULL; otherwise supply a character vector of the names of columns to skip

**Value**

data.table of original data plus created lags, rolling stats, and time between event lags and rolling stats

**Author(s)**

Adrian Antico

See Also

Other Feature Engineering: [AutoDataPartition](#), [AutoTransformationCreate](#), [AutoTransformationScore](#), [AutoWord2VecModeler](#), [CreateCalendarVariables](#), [DT\\_GDL\\_Feature\\_Engineering](#), [DummifyDT](#), [GDL\\_Feature\\_Engineering](#), [ModelDataPrep](#), [Partial\\_DT\\_GDL\\_Feature\\_Engineering](#), [Scoring\\_GDL\\_Feature\\_Engineering](#)

Examples

```
N = 25116
data <- data.table::data.table(DateTime = as.Date(Sys.time()),
  Target = stats::filter(rnorm(N,
    mean = 50,
    sd = 20),
    filter=rep(1,10),
    circular=TRUE))
data[, temp := seq(1:N)][, DateTime := DateTime - temp]
data <- data[order(DateTime)]
data <- Partial_DT_GDL_Feature_Engineering2(data,
  lags           = c(1:5),
  periods        = c(seq(10,50,10)),
  statsNames     = "mean",
  targets        = c("Target"),
  groupingVars   = NULL,
  sortDateName   = "DateTime",
  timeDiffTarget = c("Time_Gap"),
  timeAgg        = "days",
  WindowingLag   = 1,
  Type           = "Lag",
  Timer          = TRUE,
  SimpleImpute   = TRUE,
  AscRowByGroup  = "temp",
  RecordsKeep    = 1,
  AscRowRemove   = TRUE)
```

---

percRank	<i>Percentile rank function</i>
----------	---------------------------------

---

Description

This function computes percentile ranks for each row in your data like Excel’s PERCENT.RANK

Usage

```
percRank(x)
```

Arguments

x                      X is your variable of interest

Value

vector of percentile ranks

**Author(s)**

Adrian Antico

**See Also**

Other Misc: [AutoH20TextPrepScoring](#), [ChartTheme](#), [PrintObjectsSize](#), [RemixTheme](#), [SimpleCap](#), [multiplot](#), [tempDatesFun](#), [tokenizeH20](#)

**Examples**

```
data <- data.table::data.table(A = runif(100))
data[, Rank := percRank(A)]
data <- data.table::data.table(A = runif(100))
data[, Percentile := percRank(A)]
```

---

PrintObjectsSize

*PrintObjectsSize prints out the top N objects and their associated sizes, sorted by size*

---

**Description**

PrintObjectsSize prints out the top N objects and their associated sizes, sorted by size

**Usage**

```
PrintObjectsSize(N = 10)
```

**Arguments**

N                      The number of objects to display

**Value**

A print to your console of the sizes of the objects in your environment

**Author(s)**

Adrian Antico

**See Also**

Other Misc: [AutoH20TextPrepScoring](#), [ChartTheme](#), [RemixTheme](#), [SimpleCap](#), [multiplot](#), [percRank](#), [tempDatesFun](#), [tokenizeH20](#)

**Examples**

```
PrintObjectsSize(N = 10)
```

---

ProblematicFeatures	<i>ProblematicFeatures identifies problematic features for machine learning</i>
---------------------	---

---

## Description

ProblematicFeatures identifies problematic features for machine learning and outputs a data.table of the feature names in the first column and the metrics they failed to pass in the columns.

## Usage

```
ProblematicFeatures(data, ColumnNumbers = c(1:ncol(data)),
  NearZeroVarThresh = 0.05, CharUniqThresh = 0.5, NA_Rate = 0.2,
  Zero_Rate = 0.2, HighSkewThresh = 10)
```

## Arguments

data	The data.table with the columns you wish to have analyzed
ColumnNumbers	A vector with the column numbers you wish to analyze
NearZeroVarThresh	Set to NULL to not run NearZeroVar(). Checks to see if the percentage of values in your numeric columns that are not constant are greater than the value you set here. If not, the feature is collected and returned with the percentage unique value.
CharUniqThresh	Set to NULL to not run CharUniqthresh(). Checks to see if the percentage of unique levels / groups in your categorical feature is greater than the value you supply. If it is, the feature name is returned with the percentage unique value.
NA_Rate	Set to NULL to not run NA_Rate(). Checks to see if the percentage of NA's in your features is greater than the value you supply. If it is, the feature name is returned with the percentage of NA values.
Zero_Rate	Set to NULL to not run Zero_Rate(). Checks to see if the percentage of zero's in your features is greater than the value you supply. If it is, the feature name is returned with the percentage of zero values.
HighSkewThresh	Set to NULL to not run HighSkew(). Checks for numeric columns whose ratio of the sum of the top 5th percentile of values to the bottom 95th percentile of values is greater than the value you supply. If true, the column name and value is returned.

## Value

data table with new dummy variables columns and optionally removes base columns

## Author(s)

Adrian Antico

## See Also

Other EDA: [AutoWordFreq](#), [ProblematicRecords](#)

## Examples

```
test <- data.table::data.table(RandomNum = runif(1000))
test[, NearZeroVarEx := ifelse(runif(1000) > 0.99, runif(1), 1)]
test[, CharUniqueEx := as.factor(ifelse(RandomNum < 0.95, sample(letters, size = 1), "FFF"))]
test[, NA_RateEx := ifelse(RandomNum < 0.95, NA, "A")]
test[, ZeroRateEx := ifelse(RandomNum < 0.95, 0, runif(1))]
test[, HighSkewThreshEx := ifelse(RandomNum > 0.96, 100000, 1)]
ProblematicFeatures(test,
  ColumnNumbers = 2:ncol(test),
  NearZeroVarThresh = 0.05,
  CharUniqThresh = 0.50,
  NA_Rate = 0.20,
  Zero_Rate = 0.20,
  HighSkewThresh = 10)
```

---

ProblematicRecords	<i>ProblematicRecords identifies problematic records for further investigation</i>
--------------------	--

---

## Description

ProblematicRecords identifies problematic records for further investigation and data.table with 3 additional columns at the beginning of the data.table: PredictedOutlier (0 = no outlier, 1 = outlier), predict (raw H2O predicted value from Isolation Forest), and mean.length (mean length of number of splits)

## Usage

```
ProblematicRecords(data, ColumnNumbers = NULL, Threshold = 0.975,
  MaxMem = "28G", NThreads = -1, NTrees = 100,
  SampleRate = (sqrt(5) - 1)/2)
```

## Arguments

data	The data.table with the columns you wish to have analyzed
ColumnNumbers	A vector with the column numbers you wish to analyze
Threshold	Quantile value to find the cutoff value for classifying outliers
MaxMem	Specify the amount of memory to allocate to H2O. E.g. "28G"
NThreads	Specify the number of threads (E.g. cores * 2)
NTrees	Specify the number of decision trees to build
SampleRate	Specify the row sample rate per tree

## Value

A data.table

## Author(s)

Adrian Antico

## See Also

Other EDA: [AutoWordFreq](#), [ProblematicFeatures](#)

## Examples

```
Correl <- 0.85
N <- 10000
data <- data.table::data.table(Target = runif(N))
data[, x1 := qnorm(Target)]
data[, x2 := runif(N)]
data[, Independent_Variable1 := log(pnorm(Correl * x1 +
                                         sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable2 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable3 := exp(pnorm(Correl * x1 +
                                         sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable4 := exp(exp(pnorm(Correl * x1 +
                                             sqrt(1-Correl^2) * qnorm(x2))))]
data[, Independent_Variable5 := sqrt(pnorm(Correl * x1 +
                                           sqrt(1-Correl^2) * qnorm(x2)))]
data[, Independent_Variable6 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.10]
data[, Independent_Variable7 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.25]
data[, Independent_Variable8 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^0.75]
data[, Independent_Variable9 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^2]
data[, Independent_Variable10 := (pnorm(Correl * x1 +
                                       sqrt(1-Correl^2) * qnorm(x2)))^4]
data[, Target := as.factor(
  ifelse(Independent_Variable2 < 0.20, "A",
    ifelse(Independent_Variable2 < 0.40, "B",
      ifelse(Independent_Variable2 < 0.6, "C",
        ifelse(Independent_Variable2 < 0.8, "D", "E"))))]
data[, Independent_Variable11 := as.factor(
  ifelse(Independent_Variable2 < 0.15, "A",
    ifelse(Independent_Variable2 < 0.45, "B",
      ifelse(Independent_Variable2 < 0.65, "C",
        ifelse(Independent_Variable2 < 0.85, "D", "E"))))]
data[, ':= ' (x1 = NULL, x2 = NULL)]
Outliers <- ProblematicRecords(data,
                               ColumnNumbers = NULL,
                               Threshold = 0.95,
                               MaxMem = "28G",
                               NThreads = -1)
```

**Description**

This function will find the optimal thresholds for applying the main label and for finding the optimal range for doing nothing when you can quantify the cost of doing nothing

**Usage**

```
RedYellowGreen(data, PredictColNumber = 2, ActualColNumber = 1,
  TruePositiveCost = 0, TrueNegativeCost = 0,
  FalsePositiveCost = -10, FalseNegativeCost = -50, MidTierCost = -2,
  Cores = 8, Precision = 0.01, Boundaries = c(0.05, 0.75))
```

**Arguments**

<b>data</b>	data is the data table with your predicted and actual values from a classification model
<b>PredictColNumber</b>	The column number where the prediction variable is located (in binary form)
<b>ActualColNumber</b>	The column number where the target variable is located
<b>TruePositiveCost</b>	This is the utility for generating a true positive prediction
<b>TrueNegativeCost</b>	This is the utility for generating a true negative prediction
<b>FalsePositiveCost</b>	This is the cost of generating a false positive prediction
<b>FalseNegativeCost</b>	This is the cost of generating a false negative prediction
<b>MidTierCost</b>	This is the cost of doing nothing (or whatever it means to not classify in your case)
<b>Cores</b>	Number of cores on your machine
<b>Precision</b>	Set the decimal number to increment by between 0 and 1
<b>Boundaries</b>	Supply a vector of two values c(lower bound, upper bound) where the first value is the smallest threshold you want to test and the second value is the largest value you want to test. Note, if your results are at the boundaries you supplied, you should extend the boundary that was reached until the values is within both revised boundaries.

**Value**

A data table with all evaluated strategies, parameters, and utilities, along with a 3d scatterplot of the results

**Author(s)**

Adrian Antico

**See Also**

Other Model Evaluation and Interpretation: [EvalPlot](#), [ParDepCalPlots](#), [threshOptim](#)

## Examples

```
data <- data.table::data.table(Target = runif(10))
data[, x1 := qnorm(Target)]
data[, x2 := runif(10)]
data[, Predict := log(pnorm(0.85 * x1 +
                           sqrt(1-0.85^2) * qnorm(x2)))]
data[, ':= ' (x1 = NULL, x2 = NULL)]
data <- RedYellowGreen(data,
                        PredictColNumber = 2,
                        ActualColNumber = 1,
                        TruePositiveCost = 0,
                        TrueNegativeCost = 0,
                        FalsePositiveCost = -1,
                        FalseNegativeCost = -2,
                        MidTierCost = -0.5,
                        Precision = 0.5,
                        Cores = 1,
                        Boundaries = c(0.05,0.75))
```

---

RemixTheme

*RemixTheme function is a ggplot theme generator for ggplots*


---

## Description

This function adds the Remix Theme to ggplots

## Usage

```
RemixTheme()
```

## Value

An object to pass along to ggplot objects following the "+" sign

## Author(s)

Douglas Pestana

## See Also

Other Misc: [AutoH20TextPrepScoring](#), [ChartTheme](#), [PrintObjectsSize](#), [SimpleCap](#), [multiplot](#), [percRank](#), [tempDatesFun](#), [tokenizeH20](#)

## Examples

```
data <- data.table::data.table(DateTime = as.Date(Sys.time()),
                                Target = stats::filter(rnorm(1000,
                                                            mean = 50,
                                                            sd = 20),
                                                         filter=rep(1,10),
                                                         circular=TRUE))
data[, temp := seq(1:1000)][, DateTime := DateTime - temp][, temp := NULL]
data <- data[order(DateTime)]
p <- ggplot2::ggplot(data, ggplot2::aes(x = DateTime, y = Target)) + ggplot2::geom_line()
p <- p + RemixTheme()
```



---

ResidualOutliers	<i>ResidualOutliers is an automated time series outlier detection function</i>
------------------	--

---

## Description

ResidualOutliers is an automated time series outlier detection function that utilizes tsoutliers and auto.arima. It looks for five types of outliers: "AO" Additive outlier - a singular extreme outlier that surrounding values aren't affected by; "IO" Innovational outlier - Initial outlier with subsequent anomalous values; "LS" Level shift - An initial outlier with subsequent observations being shifted by some constant on average; "TC" Transient change - initial outlier with lingering effects that dissapate exponentially over time; "SLS" Seasonal level shift - similar to level shift but on a seasonal scale.

## Usage

```
ResidualOutliers(data, DateColName = "DateTime",
  TargetColName = "Target", PredictedColName = NULL,
  TimeUnit = "day", maxN = 5, tstat = 2)
```

## Arguments

<code>data</code>	the source residuals data.table
<code>DateColName</code>	The name of your data column to use in reference to the target variable
<code>TargetColName</code>	The name of your target variable column
<code>PredictedColName</code>	The name of your predicted value column. If you supply this, you will run anomaly detection of the difference between the target variable and your predicted value. If you leave PredictedColName NULL then you will run anomaly detection over the target variable.
<code>TimeUnit</code>	The time unit of your date column: hour, day, week, month, quarter, year
<code>maxN</code>	the largest lag or moving average (seasonal too) values for the arima fit
<code>tstat</code>	the t-stat value for tsoutliers

## Value

A named list containing FullData = original data.table with outliers data and ARIMA\_MODEL = the arima model.

## Author(s)

Adrian Antico

## See Also

Other Unsupervised Learning: [AutoKMeans](#), [GenTSAnomVars](#)

## Examples

```
data <- data.table::data.table(DateTime = as.Date(Sys.time()),
                               Target = as.numeric(stats::filter(rnorm(1000,
                                                                    mean = 50,
                                                                    sd = 20),
                                                                    filter=rep(1,10),
                                                                    circular=TRUE)))

data[, temp := seq(1:1000)][, DateTime := DateTime - temp][, temp := NULL]
data <- data[order(DateTime)]
data[, Predicted := as.numeric(stats::filter(rnorm(1000,
                                                    mean = 50,
                                                    sd = 20),
                                                    filter=rep(1,10),
                                                    circular=TRUE)))]

stuff <- ResidualOutliers(data = data,
                          DateColName = "DateTime",
                          TargetColName = "Target",
                          PredictedColName = NULL,
                          TimeUnit = "day",
                          maxN = 5,
                          tstat = 4)

data      <- stuff[[1]]
model     <- stuff[[2]]
outliers  <- data[type != "<NA>"]
```

---

Scoring\_GDL\_Feature\_Engineering

*An Automated Scoring Feature Engineering Function*

---

## Description

For scoring purposes of a single record when 1 or 0 groups were utilized in the DT\_GDL\_Feature\_Engineering() function. This function runs internally inside the CARMA functions but might have use outside of it.

## Usage

```
Scoring_GDL_Feature_Engineering(data, lags = c(seq(1, 5, 1)),
                                periods = c(3, 5, 10, 15, 20, 25), statsNames = c("MA"),
                                targets = c("Target"), groupingVars = NULL,
                                sortDateName = c("DateTime"), timeDiffTarget = c("Time_Gap"),
                                timeAgg = "days", WindowingLag = 1, Type = "Lag", Timer = TRUE,
                                SimpleImpute = TRUE, AscRowByGroup = "temp", RecordsKeep = 1)
```

## Arguments

<b>data</b>	A data.table you want to run the function on
<b>lags</b>	A numeric vector of the specific lags you want to have generated. You must include 1 if WindowingLag = 1.
<b>periods</b>	A numeric vector of the specific rolling statistics window sizes you want to utilize in the calculations.

statsNames	A character vector of the corresponding names to create for the rollings stats variables.
targets	A character vector of the column names for the reference column in which you will build your lags and rolling stats
groupingVars	A character vector of categorical variable names you will build your lags and rolling stats by
sortDateName	The column name of your date column used to sort events over time
timeDiffTarget	Specify a desired name for features created for time between events. Set to NULL if you don't want time between events features created.
timeAgg	List the time aggregation level for the time between events features, such as "hour", "day", "week", "month", "quarter", or "year"
WindowingLag	Set to 0 to build rolling stats off of target columns directly or set to 1 to build the rolling stats off of the lag-1 target
Type	List either "Lag" if you want features built on historical values or "Lead" if you want features built on future values
Timer	Set to TRUE if you percentage complete tracker printout
SimpleImpute	Set to TRUE for factor level imputation of "0" and numeric imputation of -1
AscRowByGroup	Required to have a column with a Row Number by group (if grouping) with 1 being the record for scoring (typically the most current in time)
RecordsKeep	Keep set to 1. Any larger value and the results will not be what you intend. I'll remove it eventually. If you want more than 1, look up <code>Partial_DT_GDL_Feature_Engineering()</code> .

**Value**

data.table of original data plus created lags, rolling stats, and time between event lags and rolling stats

**Author(s)**

Adrian Antico

**See Also**

Other Feature Engineering: [AutoDataPartition](#), [AutoTransformationCreate](#), [AutoTransformationScore](#), [AutoWord2VecModeler](#), [CreateCalendarVariables](#), [DT\\_GDL\\_Feature\\_Engineering](#), [DummifyDT](#), [GDL\\_Feature\\_Engineering](#), [ModelDataPrep](#), [Partial\\_DT\\_GDL\\_Feature\\_Engineering2](#), [Partial\\_DT\\_GDL\\_Feature\\_Engineering](#)

**Examples**

```

N = 25116
data1 <- data.table::data.table(DateTime = as.Date(Sys.time()),
                                Target = stats::filter(rnorm(N,
                                                            mean = 50,
                                                            sd = 20),
                                                         filter=rep(1,10),
                                                         circular=TRUE))

data1[, temp := seq(1:N)][, DateTime := DateTime - temp]
data1 <- data1[order(DateTime)]
data1 <- Scoring_GDL_Feature_Engineering(data1,
```

```
lags          = c(seq(1,5,1)),
periods       = c(3,5,10,15,20,25),
statsNames    = c("MA"),
targets       = c("Target"),
groupingVars   = NULL,
sortDateName  = c("DateTime"),
timeDiffTarget = c("Time_Gap"),
timeAgg       = "days",
WindowingLag  = 1,
Type          = "Lag",
Timer         = TRUE,
SimpleImpute  = TRUE,
AscRowByGroup = "temp",
RecordsKeep   = 1)
```

SimpleCap

*SimpleCap function is for capitalizing the first letter of words***Description**

SimpleCap function is for capitalizing the first letter of words (need I say more?)

**Usage**

```
SimpleCap(x)
```

**Arguments**

x                      Column of interest

**Value**

An object to pass along to ggplot objects following the "+" sign

**Author(s)**

Adrian Antico

**See Also**

Other Misc: [AutoH20TextPrepScoring](#), [ChartTheme](#), [PrintObjectsSize](#), [RemixTheme](#), [multiplot](#), [percRank](#), [tempDatesFun](#), [tokenizeH20](#)

**Examples**

```
x <- "adrian"
x <- SimpleCap(x)
```

---

tempDatesFun	<i>tempDatesFun Convert Excel datetime char columns to Date columns</i>
--------------	---

---

### Description

tempDatesFun takes the Excel datetime column, which imports as character, and converts it into a date type

### Usage

```
tempDatesFun(x)
```

### Arguments

x                      The column of interest

### Value

An object to pass along to ggplot objects following the "+" sign

### Author(s)

Adrian Antico

### See Also

Other Misc: [AutoH20TextPrepScoring](#), [ChartTheme](#), [PrintObjectsSize](#), [RemixTheme](#), [SimpleCap](#), [multiplot](#), [percRank](#), [tokenizeH20](#)

### Examples

```
Cdata <- data.table::data.table(DAY_DATE = "2018-01-01 8:53")
Cdata[, DAY_DATE := tempDatesFun(DAY_DATE)]
```

---

threshOptim	<i>Utility maximizing thresholds for binary classification</i>
-------------	--

---

### Description

This function will return the utility maximizing threshold for future predictions along with the data generated to estimate the threshold

### Usage

```
threshOptim(data, actTar = "target", predTar = "p1", tpProfit = 0,
  tnProfit = 0, fpProfit = -1, fnProfit = -2)
```

**Arguments**

<code>data</code>	data is the data table you are building the modeling on
<code>actTar</code>	The column name where the actual target variable is located (in binary form)
<code>predTar</code>	The column name where the predicted values are located
<code>tpProfit</code>	This is the utility for generating a true positive prediction
<code>tnProfit</code>	This is the utility for generating a true negative prediction
<code>fpProfit</code>	This is the cost of generating a false positive prediction
<code>fnProfit</code>	This is the cost of generating a false negative prediction

**Value**

Optimal threshold and corresponding utilities for the range of thresholds tested

**Author(s)**

Adrian Antico

**See Also**

Other Model Evaluation and Interpretation: [EvalPlot](#), [ParDepCalPlots](#), [RedYellowGreen](#)

**Examples**

```
data <- data.table::data.table(Target = runif(10))
data[, x1 := qnorm(Target)]
data[, x2 := runif(10)]
data[, Predict := log(pnorm(0.85 * x1 +
                           sqrt(1-0.85^2) * qnorm(x2)))]
data[, ':= ' (x1 = NULL, x2 = NULL)]
data <- threshOptim(data = data,
                    actTar = "Target",
                    predTar = "Predict",
                    tpProfit = 0,
                    tnProfit = 0,
                    fpProfit = -1,
                    fnProfit = -2)
optimalThreshold <- data$Thresholds
allResults <- data$EvaluationTable
```

---

tokenizeH2O

*For NLP work*

---

**Description**

This function tokenizes text data

**Usage**

```
tokenizeH2O(data)
```

**Arguments**

data                      The text data

**Author(s)**

Adrian Antico

**See Also**

Other Misc: [AutoH20TextPrepScoring](#), [ChartTheme](#), [PrintObjectsSize](#), [RemixTheme](#), [SimpleCap](#), [multiplot](#), [percRank](#), [tempDatesFun](#)

**Examples**

```
data <- tokenizeH2O(data = data[["StringColumn"]])
```

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