

# Package ‘RemixAutoML’

May 13, 2019

**Title** Remix Automated Machine Learning

**Version** 1.2.1

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**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 and AutoCatBoostRegression 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 AdditionalRepositories 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 ( $\geq$  3.5.0)

**Imports** base,  
catboost,  
caTools,  
data.table,  
doParallel,  
foreach,  
forecast,  
ggplot2,  
grid,  
h2o,  
xgboost,  
itertools,  
lubridate,  
methods,  
monreg,  
parallel,  
pROC,

RColorBrewer,  
 recommenderlab,  
 ROCR,  
 scatterplot3d,  
 stats,  
 stringr,  
 tm,  
 tsoutliers,  
 utils,  
 wordcloud,  
 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 ( $\geq$  7.0)

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## AutoCatBoostClassifier

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

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## Description

AutoCatBoostClassifier 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, 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, TestData = NULL, TargetColumnName = NULL,
  FeatureColNames = NULL, CatFeatures = NULL, TrainSplitRatio = 0.8,
  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)
```

**Arguments**

<code>data</code>	This is your data set for training and testing your model
<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. Also, not zero-indexed.
<code>CatFeatures</code>	A vector of column numbers of your categorical features, not zero indexed.
<code>TrainSplitRatio</code>	A decimal between 0.01 and 0.99 that tells the function how much data to keep for training and validation.
<code>task_type</code>	"GPU" 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. "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"
<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 "f", "auc", "tpr", "fnr", "fpr", "tnr", "prbe", "f", "odc"
<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

**Value**

Saves to file and returned in list: `_ModelID_VariableImportance.csv`, `_ModelID_` (the model), `_ModelID_ValidationData.csv`, `_ModelID_ROC_Plot.png`, `_ModelID_EvaluationPlot.png`, `_ModelID_EvaluationMetrics.csv`, `_ModelID_ParDepPlots.R` a named list of features with partial dependence calibration plots, `_ModelID_ParDepBoxPlots.R`, `_ModelID_GridCollect`, and `_ModelID_GridList`



```

FeatureColNames = c(2:11),
CatFeatures = NULL,
MaxModelsInGrid = 3,
TrainSplitRatio = 0.80,
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)

```

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### 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, 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, TestData = NULL, TargetColumnName = NULL,
  FeatureColNames = NULL, CatFeatures = NULL, TrainSplitRatio = 0.8,
  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)

```

## Arguments

<code>data</code>	This is your data set for training and testing your model
<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.

CatFeatures	A vector of column numbers of your categorical features, not zero indexed.
TrainSplitRatio	A decimal between 0.01 and 0.99 that tells the function how much data to keep for training and validation.
task_type	"GPU" Set to "GPU" to utilize your GPU for training. Default is "CPU".
eval_metric	This is the metric used inside catboost to measure performance on validation data during a grid-tune. "MultiClass" or "MultiClassOneVsAll"
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	This is the metric used to find the threshold "auc", "accuracy"
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

### Value

Saves to file and returned in list: `_ModelID_VariableImportance.csv`, `_ModelID_` (the model), `_ModelID_ValidationData.csv`, `_ModelID_EvaluationMetrics.csv`, `_ModelID_GridCollect`, and `_ModelID_GridList`

### Author(s)

Adrian Antico

### See Also

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostRegression](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFClassifier](#), [AutoH2oDRFMultiClass](#), [AutoH2oDRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoRecommenderScoring](#), [AutoRecommender](#), [AutoTS](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#)

### 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' := (x1 = NULL, x2 = NULL)]
TestModel <- AutoCatBoostMultiClass(data,
  TestData = NULL,
  TargetColumnName = "Target",
  FeatureColNames = c(2:11),
  CatFeatures = NULL,
  MaxModelsInGrid = 1,
  TrainSplitRatio = 0.80,
  task_type = "GPU",
  eval_metric = "MultiClass",
  grid_eval_metric = "Accuracy",
  Trees = 50,
  GridTune = FALSE,
  model_path = NULL,
  ModelID = "ModelTest",
  ReturnModelObjects = TRUE,
  SaveModelObjects = FALSE)

```

---

## 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, TestData = NULL, TargetColumnName = NULL,
  FeatureColNames = NULL, CatFeatures = NULL, TrainSplitRatio = 0.8,
  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)
```

**Arguments**

<b>data</b>	This is your data set for training and testing your model
<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>CatFeatures</b>	A vector of column numbers of your categorical features.
<b>TrainSplitRatio</b>	A decimal between 0.01 and 0.99 that tells the function how much data to keep for training and validation.
<b>task_type</b>	= "GPU" 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 c('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)

**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

**Value**

Saves to file: `_ModelID_VariableImportance.csv`, `_ModelID_`, `_ModelID_ValidationData.csv`, `_ModelID_EvaluationPlot.png`, `_ModelID_EvaluationBoxPlot.png`, `_ModelID_EvaluationMetrics.csv`, `_ModelID_ParDepPlots.R` a named list of features with partial dependence calibration plots, `_ModelID_ParDepBoxPlots.R`, `_ModelID_GridCollect`, and `_ModelID_catboostgrid`

**Author(s)**

Adrian Antico

**See Also**

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFClassifier](#), [AutoH2oDRFMultiClass](#), [AutoH2oDRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoRecommenderScoring](#), [AutoRecommender](#), [AutoTS](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#)

**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,
                                TestData = NULL,
                                TargetColumnName = "Target",
                                FeatureColNames = c(2:11),
                                CatFeatures = c(12),
                                MaxModelsInGrid = 1,
                                TrainSplitRatio = 0.80,
                                task_type = "GPU",
                                eval_metric = "RMSE",
                                grid_eval_metric = "r2",
                                Trees = 50,
                                GridTune = FALSE,
                                model_path = NULL,
                                ModelID = "ModelTest",
                                NumOfParDepPlots = 3,
                                ReturnModelObjects = TRUE,
                                SaveModelObjects = FALSE)

```

---

AutoH2oDRFClassifier     *AutoH2oDRFClassifier is an automated H2O modeling framework with grid-tuning and model evaluation*

---

## Description

AutoH2oDRFClassifier 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

```

AutoH2oDRFClassifier(data, TestData = NULL, TargetColumnName = NULL,
  FeatureColNames = NULL, eval_metric = "RMSE",
  TrainSplitRatio = 0.8, 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>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.

TargetColumnName	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.
FeatureColNames	Either supply the feature column names OR the column number where the target is located (but not mixed types)
eval_metric	This is the metric used to identify best grid tuned model. Choose from "AUC" or "logloss"
TrainSplitRatio	A decimal between 0.01 and 0.99 that tells the function how much data to keep for training and validation.
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.
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
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: `_ModelID_VariableImportance.csv`, `_ModelID_ValidationData.csv`, `_ModelID_EvaluationPlot.png`, `_ModelID_EvaluationBoxPlot.png`, `_ModelID_EvaluationMetrics.csv`, `_ModelID_ParDepPlots.R` a named list of features with partial dependence calibration plots

## Author(s)

Adrian Antico

## See Also

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoCatBoostRegression](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFMultiClass](#), [AutoH2oDRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoRecommenderScoring](#), [AutoRecommender](#), [AutoTS](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#)

## 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 <- AutoH2oDRFClassifier(data,
  TestData = NULL,
  TargetColumnName = "Target",
  FeatureColNames = 2:ncol(data),
  Alpha = NULL,
  eval_metric = "auc",
  TrainSplitRatio = 0.80,
  Trees = 50,
  GridTune = FALSE,
  MaxMem = "32G",
  MaxModelsInGrid = 10,
  model_path = NULL,
  ModelID = "FirstModel",
  NumOfParDepPlots = 3,
  ReturnModelObjects = TRUE,
  SaveModelObjects = FALSE,
  IfSaveModel = "mojo")

```

## Description

AutoH2oDRFMultiClass 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 metrics, confusion matrix, and variable importance.

## Usage

```
AutoH2oDRFMultiClass(data, TestData = NULL, TargetColumnName = NULL,
  FeatureColNames = NULL, eval_metric = "RMSE",
  TrainSplitRatio = 0.8, 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>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>TrainSplitRatio</b>	A decimal between 0.01 and 0.99 that tells the function how much data to keep for training and validation.
<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)

**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: `_ModelID_VariableImportance.csv`, `_ModelID_`, `_ModelID_ValidationData.csv`, `_ModelID_EvaluationPlot.png`, `_ModelID_EvaluationBoxPlot.png`, `_ModelID_EvaluationMetrics.csv`, `_ModelID_ParDepPlots.R` a named list of features with partial dependence calibration plots

**Author(s)**

Adrian Antico

**See Also**

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoCatBoostRegression](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFClassifier](#), [AutoH2oDRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoRecommenderScoring](#), [AutoRecommender](#), [AutoTS](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#)

**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 <- AutoH2oDRFMultiClass(data,
                                   TestData = NULL,
                                   TargetColumnName = "Target",
                                   FeatureColNames = 2:ncol(data),
                                   Alpha = NULL,
                                   eval_metric = "RMSE",
                                   TrainSplitRatio = 0.80,
                                   Trees = 50,
                                   GridTune = FALSE,
                                   MaxMem = "32G",
                                   MaxModelsInGrid = 10,
                                   model_path = NULL,
                                   ModelID = "FirstModel",
                                   NumOfParDepPlots = 3,
                                   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, TestData = NULL, TargetColumnName = NULL,
                      FeatureColNames = NULL, eval_metric = "RMSE",
                      TrainSplitRatio = 0.8, 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>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 "MSE", "RMSE", "MAE", "RMSLE"
<b>TrainSplitRatio</b>	A decimal between 0.01 and 0.99 that tells the function how much data to keep for training and validation.
<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. 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>IfSaveModel</b>	Set to "mojo" to save a mojo file, otherwise "standard" to save a regular H2O model object

**Value**

Saves to file: `_ModelID.VariableImportance.csv`, `_ModelID_`, `_ModelID.ValidationData.csv`, `_ModelID.EvaluationPlot.png`, `_ModelID.EvaluationBoxPlot.png`, `_ModelID.EvaluationMetrics.csv`, `_ModelID.ParDepPlots.R` a named list of features with partial dependence calibration plots, `_ModelID.ParDepBoxPlots.R`

**Author(s)**

Adrian Antico

**See Also**

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoCatBoostRegression](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFClassifier](#), [AutoH2oDRFMultiClass](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoRecommenderScoring](#), [AutoRecommender](#), [AutoTS](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#)

## 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,
  TestData = NULL,
  TargetColumnName = "Target",
  FeatureColNames = 2:ncol(data),
  Alpha = NULL,
  eval_metric = "RMSE",
  TrainSplitRatio = 0.80,
  Trees = 50,
  GridTune = FALSE,
  MaxMem = "32G",
  MaxModelsInGrid = 10,
  model_path = NULL,
  ModelID = "FirstModel",
  NumOfParDepPlots = 3,
  ReturnModelObjects = TRUE,
  SaveModelObjects = FALSE,
  IfSaveModel = "mojo")

```

## Description

AutoH2oGBMClassifier 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

```
AutoH2oGBMClassifier(data, TestData = NULL, TargetColumnName = NULL,
  FeatureColNames = NULL, eval_metric = "RMSE",
  TrainSplitRatio = 0.8, 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>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>TrainSplitRatio</b>	A decimal between 0.01 and 0.99 that tells the function how much data to keep for training and validation.
<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.

**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: `_ModelID_VariableImportance.csv`, `_ModelID_`, `_ModelID_ValidationData.csv`, `_ModelID_EvaluationPlot.png`, `_ModelID_EvaluationBoxPlot.png`, `_ModelID_EvaluationMetrics.csv`, `_ModelID_ParDepPlots.R` a named list of features with partial dependence calibration plots

**Author(s)**

Adrian Antico

**See Also**

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoCatBoostRegression](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFCClassifier](#), [AutoH2oDRFMultiClass](#), [AutoH2oDRFRegression](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoRecommenderScoring](#), [AutoRecommender](#), [AutoTS](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#)

**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 <- AutoH2oGBMClassifier(data,
                                TestData = NULL,
                                TargetColumnName = "Target",
                                FeatureColNames = 2:ncol(data),
                                Alpha = NULL,
                                eval_metric = "RMSE",
                                TrainSplitRatio = 0.80,
                                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, 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, TestData = NULL, TargetColumnName = NULL,
                      FeatureColNames = NULL, eval_metric = "RMSE",
                      TrainSplitRatio = 0.8, Trees = 50, GridTune = FALSE,
                      MaxMem = "32G", MaxModelsInGrid = 2, model_path = NULL,
                      ModelID = "FirstModel", NumOfParDepPlots = 3,
                      ReturnModelObjects = TRUE, SaveModelObjects = FALSE,
                      IfSaveModel = "mojo")

```

## Arguments

data	This is your data set for training and testing your model
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).
FeatureColNames	Either supply the feature column names OR the column number where the target is located (but not mixed types)
eval_metric	This is the metric used to identify best grid tuned model. Choose from "logloss", "r2", "RMSE", "MSE"
TrainSplitRatio	A decimal between 0.01 and 0.99 that tells the function how much data to keep for training and validation.
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.
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
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: `_ModelID_VariableImportance.csv`, `_ModelID_`, `_ModelID_ValidationData.csv`, `_ModelID_EvaluationPlot.png`, `_ModelID_EvaluationBoxPlot.png`, `_ModelID_EvaluationMetrics.csv`, `_ModelID_ParDepPlots.R` a named list of features with partial dependence calibration plots

### Author(s)

Adrian Antico

### See Also

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoCatBoostRegression](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFClassifier](#), [AutoH2oDRFMultiClass](#), [AutoH2oDRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoRecommenderScoring](#), [AutoRecommender](#), [AutoTS](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#)

## 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 <- AutoH2oGBMMultiClass(data,
  TestData = NULL,
  TargetColumnName = "Target",
  FeatureColNames = 2:ncol(data),
  Alpha = NULL,
  eval_metric = "RMSE",
  TrainSplitRatio = 0.80,
  Trees = 50,
  GridTune = FALSE,
  MaxMem = "32G",
  MaxModelsInGrid = 10,
  model_path = NULL,
  ModelID = "FirstModel",
  NumOfParDepPlots = 3,
  ReturnModelObjects = TRUE,
  SaveModelObjects = FALSE,
  IfSaveModel = "mojo")

```

## 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, TestData = NULL, TargetColumnName = NULL,
  FeatureColNames = NULL, Alpha = NULL, Distribution = "poisson",
  eval_metric = "RMSE", TrainSplitRatio = 0.8, 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>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>Alpha</b>	This is the quantile value you want to use for quantile regression. Must be a decimal between 0 and 1.
<b>Distribution</b>	Choose from "gaussian", "poisson", "gamma", "tweedie", "laplace", "quantile", "huber"
<b>eval_metric</b>	This is the metric used to identify best grid tuned model. Choose from "MSE", "RMSE", "MAE", "RMSLE"
<b>TrainSplitRatio</b>	A decimal between 0.01 and 0.99 that tells the function how much data to keep for training and validation.
<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



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

**Value**

Saves to file: `_ModelID_VariableImportance.csv`, `_ModelID_ValidationData.csv`, `_ModelID_EvaluationPlot.png`, `_ModelID_EvaluationBoxPlot.png`, `_ModelID_EvaluationMetrics.csv`, `_ModelID_ParDepPlots.R` a named list of features with partial dependence calibration plots, `_ModelID_ParDepBoxPlots.R`

**Author(s)**

Adrian Antico

**See Also**

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoCatBoostRegression](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFClassifier](#), [AutoH2oDRFMultiClass](#), [AutoH2oDRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoNLS](#), [AutoRecommenderScoring](#), [AutoRecommender](#), [AutoTS](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#)

**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 <- AutoH2oGBMRegression(data,
  TestData = NULL,
  TargetColumnName = "Target",
  FeatureColNames = 2:ncol(data),
  Alpha = NULL,
  Distribution = "poisson",
  eval_metric = "RMSE",
  TrainSplitRatio = 0.80,
  Trees = 50,
  GridTune = FALSE,
  MaxMem = "32G",
  MaxModelsInGrid = 10,
  model_path = NULL,
  ModelID = "FirstModel",
  NumOfParDepPlots = 3,
  ReturnModelObjects = TRUE,
  SaveModelObjects = FALSE,
  IfSaveModel = "mojo")

```

## 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

<code>model_path</code>	Directory path for where you want your models saved
<code>MaxRuntimeSeconds</code>	Number of seconds of run time for grid tuning
<code>MaxModels</code>	Number of models you'd like to have returned
<code>TrainData</code>	Set to NULL or supply a <code>data.table</code> for training data
<code>TestData</code>	Set to NULL or supply a <code>data.table</code> for validation data
<code>SaveToFile</code>	Set to TRUE to save models and output to <code>model_path</code>
<code>ReturnObjects</code>	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 let's 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`, `AutoH2OScoring`, `AutoH2oDRFCClassifier`, `AutoH2oDRFMultiClass`, `AutoH2oDRFRegression`, `AutoH2oGBMClassifier`, `AutoH2oGBMMultiClass`, `AutoH2oGBMRegression`, `AutoNLS`, `AutoRecommenderScoring`, `AutoRecommender`, `AutoTS`, `AutoXGBoostClassifier`, `AutoXGBoostMultiClass`, `AutoXGBoostRegression`

## Examples

```
# Classification 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 > 0.5,1,0))]
Construct <- data.table::data.table(Targets = rep("target",3),
                                   Distribution = c("bernoulli",
                                                    "bernoulli",
                                                    "bernoulli"),
                                   Loss = c("AUC", "AUC", "CrossEntropy"),
                                   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),
```



```

                                "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(NA,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)

# 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 +

```

[illegible]



```

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))

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)

```

---

AutoH2OScoring

*AutoH2OScoring is the complement of AutoH2OModeler.*


---

## Description

AutoH2OScoring is the complement of AutoH2OModeler. 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), NThreads = 6, MaxMem = "28G",
  JavaOptions = "-Xmx1g -XX:ReservedCodeCacheSize=256m",
  SaveToFile = FALSE, FilePath = NULL, H2OShutdown = rep(FALSE, 3))
```

## Arguments

Features	This is a data.table of features for scoring.
GridTuneRow	Numeric. The row numbers of grid_tuned_paths, KMeansModelFile, or StoreFile 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.



```

ParDep          = rep(NA,3),
PD_Data         = rep("A11",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","mojo","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)

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, MaxMem, NThreads)
```

**Arguments**

data	The text data
string	The name of the string column to prepare
MaxMem	Amount of memory you want to let H2O utilize
NThreads	The number of threads you want to let H2O utilize

**Author(s)**

Adrian Antico

**See Also**

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

**Examples**

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

---

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 = 4, MaxMem = "14G", 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**

data	is the source time series data.table
nthreads	set based on number of threads your machine has available
MaxMem	set based on the amount of memory your machine has available
SaveModels	Set to "standard", "mojo", or NULL (default)
PathFile	Set to folder where you will keep the models
GridTuneGLRM	If you want to grid tune the glrm model, set to TRUE, FALSE otherwise
GridTuneKMeans	If you want to grid tuen 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]]
```

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](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFClassifier](#), [AutoH2oDRFMultiClass](#), [AutoH2oDRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoRecommenderScoring](#), [AutoRecommender](#), [AutoTS](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#)

**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

This function returns the winning model that you pass onto AutoRecommenderScoring

## Usage

```
AutoRecommender(data, Partition = "Split", KFold = 2, Ratio = 0.75,
  RatingType = "TopN", RatingsKeep = 20,
  SkipModels = "AssociationRules", ModelMetric = "TPR")
```

## Arguments

<b>data</b>	This is your BinaryRatingsMatrix. See function RecomDataCreate
<b>Partition</b>	Choose from "split", "cross-validation", "bootstrap". See evaluation-Scheme in recommenderlab for details.
<b>KFold</b>	Choose 2 for traditional train and test. Choose greater than 2 for the number of cross validations
<b>Ratio</b>	The ratio for train and test. E.g. 0.75 for 75 percent data allocated to training
<b>RatingType</b>	Choose from "TopN", "ratings", "ratingMatrix"
<b>RatingsKeep</b>	The total ratings you wish to return. Default is 20.
<b>SkipModels</b>	AssociationRules runs the slowest and may crash your system. Choose from: "AssociationRules", "ItemBasedCF", "UserBasedCF", "PopularItems", "RandomItems"
<b>ModelMetric</b>	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 Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoCatBoostRegression](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2ODRFClassifier](#), [AutoH2ODRFMultiClass](#), [AutoH2ODRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoRecommenderScoring](#), [AutoTS](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#)

## Examples

```
WinningModel <- AutoRecommender(RatingsMatrix,
  Partition = "Split",
  KFold = 2,
  Ratio = 0.75,
  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")
```

**Arguments**

data	The binary ratings matrix from RecomDataCreate()
WinningModel	The winning model returned from AutoRecommender()
EntityColName	Typically your customer ID
ProductColName	Something like "StockCode"

**Value**

Returns the prediction data

**Author(s)**

Adrian Antico and Douglas Pestana

**See Also**

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

**Examples**

```
# F(G(Z(x))): AutoRecommenderScoring(AutoRecommender(RecomDataCreate(TransactionData)))
Results <- AutoRecommenderScoring(
  data = RecomDataCreate(
    data,
    EntityColName = "CustomerID",
    ProductColName = "StockCode",
    MetricColName = "TotalSales"),
  WinningModel = AutoRecommender(
    RecomDataCreate(
      data,
      EntityColName = "CustomerID",
      ProductColName = "StockCode",
      MetricColName = "TotalSales"),
```

```

Partition = "Split",
KFolds = 2,
Ratio = 0.75,
RatingType = "TopN",
RatingsKeep = 20,
SkipModels = "AssociationRules",
ModelMetric = "TPR"),
EntityColName = "CustomerID",
ProductColName = "StockCode")

```

AutoTS

*AutoTS is an automated time series modeling function*

## 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, TimeUnit = "day", Lags = 25,
       SLags = 2, NumCores = 4, SkipModels = NULL, StepWise = TRUE,
       TSClean = TRUE, ModelFreq = TRUE, PrintUpdates = FALSE)

```

## Arguments

<b>data</b>	is the source time series data as a data.table - or a data structure that can be converted to a data.table
<b>TargetName</b>	is the name of the target variable in your data.table
<b>DateName</b>	is the name of the date column in your data.table
<b>FCPeriods</b>	is the number of periods into the future you wish to forecast
<b>HoldOutPeriods</b>	is the number of periods to use for validation testing
<b>TimeUnit</b>	is the level of aggregation your dataset comes in
<b>Lags</b>	is the number of lags you wish to test in various models (same as moving averages)
<b>SLags</b>	is the number of seasonal lags you wish to test in various models (same as moving averages)
<b>NumCores</b>	is the number of cores available on your computer
<b>SkipModels</b>	Don't run specified models - e.g. exclude all models "DSHW" "ARFIMA" "ARIMA" "ETS" "NNET" "TBATS" "TSLM"
<b>StepWise</b>	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.

TSClean	Set to TRUE to have missing values interpolated and outliers replaced with interpolated values: creates separate models for a larger comparison set
ModelFreq	Set to TRUE to run a separate version of all models where the time series frequency is chosen algorithmically
PrintUpdates	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.

## Author(s)

Adrian Antico and Douglas Pestana

## See Also

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

## 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,
        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, 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 = 6, 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: [DT\\_GDL\\_Feature\\_Engineering](#), [DummifyDT](#), [FAST\\_GDL\\_Feature\\_Engineering](#), [GDL\\_Feature\\_Engineering](#), [ModelDataPrep](#), [ProblematicFeatures](#), [Scoring\\_GDL\\_Feature\\_Engineering](#)

**Examples**

```
data <- AutoWord2VecModeler(data,
  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 = 6,
  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)

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

**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,
Dwigt, Dwigt, Dwigt, Dwigt, Dwigt, Dwigt, Dwigt, Dwigt, Dwigt, Dwigt,
Schrute, 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"))
```

---

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, 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**

```
AutoXGBoostClassifier(data, TestData = NULL, TargetColumnName = NULL,
  FeatureColNames = NULL, CatFeatures = NULL, TrainSplitRatio = 0.8,
  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)
```

**Arguments**

<b>data</b>	This is your data set for training and testing your model
<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>CatFeatures</b>	Supply a vector of character names of the columns. Internally, the function will convert these columns to dummy variables.
<b>TrainSplitRatio</b>	A decimal between 0.01 and 0.99 that tells the function how much data to keep for training and validation.
<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)
<b>SaveModelObjects</b>	Set to TRUE to return all modeling objects to your environment



**Value**

Saves to file: `_ModelID_VariableImportance.csv`, `_ModelID_`, `_ModelID_ValidationData.csv`, `_ModelID_EvaluationPlot.png`, `_ModelID_EvaluationBoxPlot.png`, `_ModelID_EvaluationMetrics.csv`, `_ModelID_ParDepPlots.R` a named list of features with partial dependence calibration plots

**Author(s)**

Adrian Antico

**See Also**

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoCatBoostRegression](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFClassifier](#), [AutoH2oDRFMultiClass](#), [AutoH2oDRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoRecommenderScoring](#), [AutoRecommender](#), [AutoTS](#), [AutoXGBoostMultiClass](#), [AutoXGBoostRegression](#)

**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)]
data <- data.table::rbindlist(list(data,data))
test <- data[5001:10000]
data <- data[1:5000]
TestModel <- AutoXGBoostClassifier(data,
                                   TestData = test,
```

```

TargetColumnName = 1,
FeatureColNames = 2:12,
CatFeatures = 12,
TrainSplitRatio = 0.80,
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)

```

---

**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, 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

```

AutoXGBoostMultiClass(data, TestData = NULL, TargetColumnName = NULL,
  FeatureColNames = NULL, CatFeatures = NULL, TrainSplitRatio = 0.8,
  eval_metric = "merror", Trees = 50, GridTune = FALSE,
  grid_eval_metric = "merror", TreeMethod = "hist",
  MaxModelsInGrid = 10, NThreads = 8, model_path = NULL,
  ModelID = "FirstModel", NumOfParDepPlots = 3, Verbose = 0,
  ReturnModelObjects = TRUE, SaveModelObjects = FALSE)

```

## Arguments

<b>data</b>	This is your data set for training and testing your model
<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). 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)
CatFeatures	Supply a vector of character names of the columns. Internally, the function will convert these columns to dummy variables.
TrainSplitRatio	A decimal between 0.01 and 0.99 that tells the function how much data to keep for training and validation.
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.
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
NumOfParDepPlots	Tell the function the number of partial dependence calibration plots you want to create.
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

**Value**

Saves to file: `_ModelID_VariableImportance.csv`, `_ModelID_`, `_ModelID_ValidationData.csv`, `_ModelID_EvaluationMetrics.csv`

**Author(s)**

Adrian Antico

**See Also**

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoCatBoostRegression](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFClassifier](#), [AutoH2oDRFMultiClass](#), [AutoH2oDRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoRecommenderScoring](#), [AutoRecommender](#), [AutoTS](#), [AutoXGBoostClassifier](#), [AutoXGBoostRegression](#)

## 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[, ':= ' (x1 = NULL, x2 = NULL)]
data[, Target := ifelse(Target > 0.5, 1, 0)]
data <- data.table::rbindlist(list(data,data))
test <- data[5001:10000]
data <- data[1:5000]
TestModel <- AutoXGBoostMultiClass(data,
  TestData = test,
  TargetColumnName = 1,
  FeatureColNames = 2:12,
  CatFeatures = 12,
  TrainSplitRatio = 0.80,
  eval_metric = "merror",
  Trees = 50,
  GridTune = TRUE,
  grid_eval_metric = "merror",
  MaxModelsInGrid = 10,
  NThreads = 8,
  TreeMethod = "hist",
  model_path = getwd(),
  ModelID = "FirstModel",
  NumOfParDepPlots = 3,
  ReturnModelObjects = TRUE,
  SaveModelObjects = FALSE)

```

---

**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, TestData = NULL, TargetColumnName = NULL,
  FeatureColNames = NULL, CatFeatures = NULL, TrainSplitRatio = 0.8,
  eval_metric = "RMSE", Trees = 50, GridTune = FALSE,
  grid_eval_metric = "mae", TreeMethod = "hist",
  MaxModelsInGrid = 10, NThreads = 8, model_path = NULL,
  ModelID = "FirstModel", NumOfParDepPlots = 3, Verbose = 0,
  ReturnModelObjects = TRUE, SaveModelObjects = FALSE)
```

## Arguments

<b>data</b>	This is your data set for training and testing your model
<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>CatFeatures</b>	Supply a vector of character names of the columns. Internally, the function will convert these columns to dummy variables.
<b>TrainSplitRatio</b>	A decimal between 0.01 and 0.99 that tells the function how much data to keep for training and validation.
<b>eval_metric</b>	This is the metric used to identify best grid tuned model. Choose from "r2", "RMSE", "MSE", "MAE"
<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>	Choose from "poisson", "mae", "mape", "mse", "msle", "kl", "cs", "r2"
<b>TreeMethod</b>	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
NumOfParDepPlots	Tell the function the number of partial dependence calibration plots you want to create.
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

### Value

Saves to file: `_ModelID_VariableImportance.csv`, `_ModelID_`, `_ModelID_ValidationData.csv`, `_ModelID_EvaluationPlot.png`, `_ModelID_EvaluationBoxPlot.png`, `_ModelID_EvaluationMetrics.csv`, `_ModelID_ParDepPlots.R` a named list of features with partial dependence calibration plots

### Author(s)

Adrian Antico

### See Also

Other Supervised Learning: [AutoCatBoostClassifier](#), [AutoCatBoostMultiClass](#), [AutoCatBoostRegression](#), [AutoH2OModeler](#), [AutoH2OScoring](#), [AutoH2oDRFClassifier](#), [AutoH2oDRFMultiClass](#), [AutoH2oDRFRegression](#), [AutoH2oGBMClassifier](#), [AutoH2oGBMMultiClass](#), [AutoH2oGBMRegression](#), [AutoNLS](#), [AutoRecommenderScoring](#), [AutoRecommender](#), [AutoTS](#), [AutoXGBoostClassifier](#), [AutoXGBoostMultiClass](#)

### 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 <- data.table::rbindlist(list(data,data))
test <- data[5001:10000]
data <- data[1:5000]
TestModel <- AutoXGBoostRegression(data,
  TestData = test,
  TargetColumnName = 1,
  FeatureColNames = 2:12,
  CatFeatures = 12,
  TrainSplitRatio = 0.80,
  eval_metric = "RMSE",
  Trees = 50,
  GridTune = TRUE,
  grid_eval_metric = "mae",
  MaxModelsInGrid = 10,
  NThreads = 8,
  TreeMethod = "hist",
  model_path = getwd(),
  ModelID = "FirstModel",
  NumOfParDepPlots = 3,
  ReturnModelObjects = TRUE,
  SaveModelObjects = FALSE)

```

---

ChartTheme

---

*ChartTheme function is a ggplot theme generator for ggplots*


---

## 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: [AutoH2OTextPrepScoring](#), [AutoWordFreq](#), [PrintObjectsSize](#), [RecomDataCreate](#), [RemixTheme](#), [SimpleCap](#), [multiplot](#), [percRank](#), [tempDatesFun](#), [tokenizeH2O](#)

**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)
```

---

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, 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>statsNames</b>	A character vector of the corresponding names to create for the rollings stats variables.







```

                                size = 100000,
                                replace = TRUE))
test <- DummifyDT(data = test,
                  cols = "FactorCol",
                  KeepFactorCols = FALSE)
ncol(test)
test[, sum(FactorCol_gg)]

```

---

EvalPlot

*EvalPlot automatically builds calibration plots for model evaluation*


---

## 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) base::mean(x, na.rm =
TRUE))

```

## Arguments

<code>data</code>	Data containing predicted values and actual values for comparison
<code>PredictionColName</code>	String representation of column name with predicted values from model
<code>TargetColName</code>	String representation of column name with target values from model
<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>aggrfun</code>	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))
```

---

FAST\_GDL\_Feature\_Engineering

*An Fast Automated Feature Engineering Function*

---

## Description

For models with target variables within the realm of the current time frame but not too far back in time, this function creates 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

```
FAST_GDL_Feature_Engineering(data, lags = c(1:5), periods = c(seq(10,
  50, 10)), statsFUNs = c("mean", "median", "sd", "quantile85",
  "quantile95"), statsNames = c("mean", "median", "sd", "quantile85",
  "quantile95"), targets = c("Target"),
  groupingVars = c("GroupVariable"), sortDateName = c("DateTime"),
  timeDiffTarget = NULL, timeAgg = c("hours"), WindowingLag = 1,
  Type = c("Lag"), Timer = FALSE, SkipCols = FALSE,
  SimpleImpute = TRUE, AscRowByGroup = c("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.
<b>statsFUNs</b>	Vector of functions for your rolling windows, such as mean, sd, min, max, quantile
<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>SkipCols</b>	Defaults to NULL; otherwise supply a character vector of the names of columns to skip
<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 1 being the record 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.)

**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: [AutoWord2VecModeler](#), [DT\\_GDL\\_Feature\\_Engineering](#), [DummifyDT](#), [GDL\\_Feature\\_Engineering](#), [ModelDataPrep](#), [ProblematicFeatures](#), [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 <- FAST_GDL_Feature_Engineering(data,
  lags = c(1:5),
  periods = c(seq(10,50,10)),

```

```

statsFUNs      = c("mean",
                   "median",
                   "sd",
                   "quantile85",
                   "quantile95"),
statsNames     = c("mean",
                   "median",
                   "sd",
                   "quantile85",
                   "quantile95"),
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")

```

---

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 = c(seq(1, 5, 1)), periods = c(3, 5,
10, 15, 20, 25), statsFUNs = c(function(x) quantile(x, probs = 0.1,
na.rm = TRUE), function(x) quantile(x, probs = 0.9, na.rm = TRUE),
function(x) base::mean(x, na.rm = TRUE), function(x) sd(x, na.rm = TRUE),
function(x) quantile(x, probs = 0.25, na.rm = TRUE), function(x)
quantile(x, probs = 0.75, na.rm = TRUE)), statsNames = c("q10", "q90",
"mean", "sd", "q25", "q75"), targets = c("qty"),
groupingVars = c("Group1", "Group2"), sortDateName = c("date"),
timeDiffTarget = c("TimeDiffName"), timeAgg = c("days"),
WindowingLag = 0, Type = c("Lag"), Timer = TRUE, SkipCols = NULL,
SimpleImpute = TRUE)

```

## Arguments

<code>data</code>	A <code>data.table</code> 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 <code>WindowingLag = 1</code> .

<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 <code>function(x) mean(x)</code> , <code>function(x) sd(x)</code> , or <code>function(x) quantile(x)</code>
<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>SkipCols</b>	Defaults to NULL; otherwise supply a character vector of the names of columns to skip
<b>SimpleImpute</b>	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: [AutoWord2VecModeler](#), [DT\\_GDL\\_Feature\\_Engineering](#), [DummifyDT](#), [FAST\\_GDL\\_Feature\\_Engineering](#), [ModelDataPrep](#), [ProblematicFeatures](#), [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     = c(function(x) quantile(x, probs = 0.20, na.rm = TRUE)),
statsNames    = c("q20"),
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,
  MissFactor = "0", MissNum = -1)
```

**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>MissFactor</code>	Supply the value to impute missing factor levels
<code>MissNum</code>	Supply the value to impute missing numeric values

**Value**

Returns the original data table with corrected values

**Author(s)**

Adrian Antico

**See Also**

Other Feature Engineering: [AutoWord2VecModeler](#), [DT\\_GDL\\_Feature\\_Engineering](#), [DummifyDT](#), [FAST\\_GDL\\_Feature\\_Engineering](#), [GDL\\_Feature\\_Engineering](#), [ProblematicFeatures](#), [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	<i>Multiplot is a function for combining multiple plots</i>
-----------	---

---

**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](#), [AutoWordFreq](#), [ChartTheme](#), [PrintObjectsSize](#), [RecomDataCreate](#), [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
```

---

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](#), [AutoWordFreq](#), [ChartTheme](#), [PrintObjectsSize](#), [RecomDataCreate](#), [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	<i>PrintObjectsSize prints out the top N objects and their associated sizes, sorted by size</i>
------------------	---

---

**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](#), [AutoWordFreq](#), [ChartTheme](#), [RecomDataCreate](#), [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**

<b>data</b>	The data.table with the columns you wish to have analyzed
<b>ColumnNumbers</b>	A vector with the column numbers you wish to analyze
<b>NearZeroVarThresh</b>	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.
<b>CharUniqThresh</b>	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.
<b>NA_Rate</b>	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 Feature Engineering: [AutoWord2VecModeler](#), [DT\\_GDL\\_Feature\\_Engineering](#), [DummifyDT](#), [FAST\\_GDL\\_Feature\\_Engineering](#), [GDL\\_Feature\\_Engineering](#), [ModelDataPrep](#), [Scoring\\_GDL\\_Feature\\_Engineering](#)

### 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)
```

---

RecomDataCreate

*Convert transactional data.table to a binary ratings matrix*

---

### Description

Convert transactional data.table to a binary ratings matrix

### Usage

```
RecomDataCreate(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 Misc: [AutoH20TextPrepScoring](#), [AutoWordFreq](#), [ChartTheme](#), [PrintObjectsSize](#), [RemixTheme](#), [SimpleCap](#), [multiplot](#), [percRank](#), [tempDatesFun](#), [tokenizeH20](#)

**Examples**

```
data <- data.table::data.table(CustomerID = c(1,1,2,2,3,3),
                               StockCode = c("A","B","A","A","B","A"),
                               TotalSales = c(2,3,4,5,1,2))
RatingsMatrix <- RecomDataCreate(data,
                                  EntityColName = "CustomerID",
                                  ProductColName = "StockCode",
                                  MetricColName = "TotalSales",
                                  ReturnMatrix = TRUE)
```

---

RedYellowGreen

*RedYellowGreen is for determining the optimal thresholds for binary classification when do-nothing is an option*

---

**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**

<code>data</code>	data is the data table with your predicted and actual values from a classification model
<code>PredictColNumber</code>	The column number where the prediction variable is located (in binary form)
<code>ActualColNumber</code>	The column number where the target variable is located
<code>TruePositiveCost</code>	This is the utility for generating a true positive prediction
<code>TrueNegativeCost</code>	This is the utility for generating a true negative prediction
<code>FalsePositiveCost</code>	This is the cost of generating a false positive prediction
<code>FalseNegativeCost</code>	This is the cost of generating a false negative prediction
<code>MidTierCost</code>	This is the cost of doing nothing (or whatever it means to not classify in your case)
<code>Cores</code>	Number of cores on your machine
<code>Precision</code>	Set the decimal number to increment by between 0 and 1
<code>Boundaries</code>	Supply a vector of two values <code>c(lower bound, upper bound)</code> 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 extent 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](#), [AutoWordFreq](#), [ChartTheme](#), [PrintObjectsSize](#), [RecomDataCreate](#), [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

data	the source residuals data.table
DateColName	The name of your data column to use in reference to the target variable
TargetColName	The name of your target variable column
PredictedColName	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.
TimeUnit	The time unit of your date column: hour, day, week, month, quarter, year
maxN	the largest lag or moving average (seasonal too) values for the arima fit
tstat	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 (brings back a single row by group), this function creates 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

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

## Arguments

<code>data</code>	A <code>data.table</code> 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 <code>WindowingLag = 1</code> .



```

                                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),
                                statsFUNs    = c(function(x) mean(x,na.rm = TRUE)),
                                statsNames   = c("MA"),
                                targets      = c("Target"),
                                groupingVars = NULL,
                                sortDateName = c("DateTime"),
                                timeDiffTarget = c("Time_Gap"),
                                timeAgg      = "days",
                                WindowingLag = 1,
                                Type         = "Lag",
                                Timer        = TRUE,
                                SkipCols     = FALSE,
                                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](#), [AutoWordFreq](#), [ChartTheme](#), [PrintObjectsSize](#), [RecomDataCreate](#), [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](#), [AutoWordFreq](#), [ChartTheme](#), [PrintObjectsSize](#), [RecomDataCreate](#), [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: [AutoH2OTextPrepScoring](#), [AutoWordFreq](#), [ChartTheme](#), [PrintObjectsSize](#), [RecomDataCreate](#), [RemixTheme](#), [SimpleCap](#), [multiplot](#), [percRank](#), [tempDatesFun](#)

**Examples**

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

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