ARESLab

Adaptive Regression Splines toolbox for Matlab/Octave

ver. 1.8.2

Gints Jekabsons

Institute of Applied Computer Systems
Riga Technical University
Meza 1/3, LV-1048, Riga, Latvia
E-mail: gints.jekabsons@rtu.lv
URL: http://www.cs.rtu.lv/jekabsons/

Reference manual

June, 2015

CONTENTS

1. INTRODUCTION

What is ARESLab

ARESLab is a Matlab/Octave toolbox for building piecewise-linear and piecewise-cubic regression models using the Multivariate Adaptive Regression Splines technique (also known as MARS). (The term "MARS" is a registered trademark and thus not used in the name of the toolbox.) The original author of MARS technique is Jerome Friedman (Friedman 1991, Friedman 1993).

With this toolbox you can build MARS models (here referred to as ARES models), test them on a separate test set or using k-fold Cross-Validation, use the models for prediction, print their equations, perform ANOVA decomposition, as well as plot the models.

This reference manual provides overview of the functions available in the ARESLab.

ARESLab can be downloaded at http://www.cs.rtu.lv/jekabsons/.

The toolbox code is licensed under the GNU GPL ver. 3 or any later version. Some parts of `aresbuild` and `createList` functions were initially derived from ENTOOL toolbox (Merkwirth & Wichard 2003, Norgaard 2000) which also falls under the GPL licence.

Details

The ARESLab toolbox is written entirely in Matlab/Octave. I implement the main functionality of the MARS technique for regression as close to the description in the Friedman's original paper (Friedman 1991) as possible. While implementing the knot placement part (see remarks about `minSpan` and `endSpan` in Section 2.2), I also took a look at the source code of the R Earth package (Milborrow 2015) and implemented it very similarly to Earth version 4.3.0.

One major difference is that the model building is not accelerated using the "fast least-squares update technique" (Friedman 1991). This difference however affects only the speed of the algorithm execution, not predictive performance of built models.

The absence of the acceleration means that the code might be slow for large data sets (however, see description of `aresparams` on how to make the process faster by using the "Fast MARS" algorithm and/or setting more conservative values for algorithm parameters). An alternative is to use the R Earth package which is faster and in some aspects more sophisticated however currently lacks the ability to create piecewise-cubic models. Yet another open source alternative is to use py-earth for Python (Rudy 2015).

ARESLab does not automatically handle missing data or categorical input variables.

Feedback

For any feedback on the toolbox including bug reports feel free to contact me via the email address given on the title page of this reference manual.

Citing the ARESLab toolbox

Please give a reference to the webpage in any publication describing research performed using the toolbox e.g., like this:

Jekabsons G., ARESLab: Adaptive Regression Splines toolbox for Matlab/Octave, 2015, available at http://www.cs.rtu.lv/jekabsons/

2. AVAILABLE FUNCTIONS

ARESLab toolbox provides the following list of functions:

- `aresbuild` – builds an ARES model;
- `aresparams` – creates a structure of ARES configuration parameter values for further use with `aresbuild`, `arescv`, and `arescvc` functions;
- `arespredict` – makes predictions using an ARES model;
- `arestest` – tests an ARES model on a test data set;
- `arescv` – tests ARES performance using k-fold Cross-Validation;
- `arescvc` – finds the "best" value for penalty *c* of the Generalized Cross-Validation criterion from a set of candidate values using k-fold Cross-Validation;
- `aresplot` – plots surface of an ARES model;
- `areseq` – prints an ARES model;
- `aresanova` – performs ANOVA decomposition;
- `aresanovareduce` – reduces an ARES model according to ANOVA decomposition.

2.1. Function `aresbuild`

Purpose:

Builds a regression model using the Multivariate Adaptive Regression Splines technique.

Call:

```
[model, time] = aresbuild(Xtr, Ytr, trainParams, weights, keepX, modelOld, verbose)
```

All the arguments, except the first two, of this function are optional. Empty values are also accepted (the corresponding default values will be used).

Input:

`Xtr, Ytr` : Training data cases (`Xtr`(*i*,:), `Ytr`(*i*)), *i* = 1,...,*n*. Note that it is recommended to pre-scale `Xtr` values to [0,1] (Friedman 1991) and to standardize `Ytr` values (Milborrow 2015). This is because widely different locations and scales for the input variables can cause instabilities that could affect the quality of the final model. The MARS technique is (except for numerics) invariant to the locations and scales of the input variables. It is therefore reasonable to perform a transformation that causes resulting locations and scales to be most favourable from the point of view of numeric stability (Friedman 1991).

`trainParams` : A structure of training parameters for the algorithm. If not provided, default values will be used (see function `aresparams` for details).

`weights` : A vector of data case weights; if supplied, the algorithm calculates the sum of squared errors multiplying the squared residuals by the supplied weights. The length of `weights` vector must be the same as the number of data cases. The weights must be nonnegative.

`keepX` : Set this to true to retain `model.X` matrix (see description of `model.X`). (default value = `false`)

`modelOld` : If here an already built ARES model is provided, no forward phase will be done. Instead this model will be taken directly to the backward phase and pruned. This is useful for fast selection of the "best" penalty

trainParams.c value using Cross-Validation in arescvc function or for fast tuning of other parameters of backward phase (cubic, maxFinalFuncs).

verbose : Set to false for no verbose. (default value = true)

Output:

model : The built ARES model – a structure with the following elements:

- coefs : Coefficient vector of the regression model (for the intercept term and each basis function).
- knotdims : Cell array of indexes of used input variables for each knot in each basis function.
- knotsites : Cell array of knot sites for each knot and used input variable in each basis function.
- knotdirs : Cell array of directions (-1 or 1) of the hinge functions for each used input variable in each basis function.
- parents : Vector of indexes of direct parents for each basis function (0 if there is no direct parent or it is the intercept term).
- trainParams : A structure of training parameters for the algorithm.
- MSE : Mean Squared Error of the model in the training data set.
- GCV : Generalized Cross-Validation (GCV) of the model in the training data set. The GCV is calculated using trainParams.c argument (for details on GCV calculation, see Friedman 1991). The value may also be Inf if model's effective number of parameters (see Eq. 1) is larger than or equal to *n*.
- t1 : For piecewise-cubic models only. Matrix of knot sites for the additional side knots on the left of the central knot.
- t2 : For piecewise-cubic models only. Matrix of knot sites for the additional side knots on the right of the central knot.
- minX : Vector of minimums for input variables (used for t1 and t2 placements as well as for model plotting).
- maxX : Vector of maximums for input variables (used for t1 and t2 placements as well as for model plotting).
- endSpan : The used value of endSpan.
- X : Matrix of values of basis functions applied to Xtr. Each column corresponds to a selected basis function. Each row corresponds to a row in Xtr. Multiplying X by coefs gives ARES prediction for Ytr. This variable is available only if parameter keepX is true.

time : Algorithm execution time (in seconds)

Remarks:

The model building algorithm builds a model in two phases: forward selection and backward deletion. In the forward phase the algorithm starts with a model consisting of just the intercept term and iteratively adds reflected pairs of basis functions giving the largest reduction of training error. The forward phase is executed until one of the following conditions is met:

1) reached maximum number of basis functions (trainParams.maxFuncs);
2) the difference between err and newErr is smaller than trainParams.threshold, where newErr is calculated by dividing sum of squared residuals by the variance of Ytr and err is the newErr value from the previous iteration;
3) the newErr is smaller than trainParams.threshold;
4) the number of model's coefficients (i.e., the number of all the basis functions including the intercept term) in the next iteration is expected to be equal to or larger than *n*.

At the end of the forward phase we have a large model which typically overfits the data, and so a backward deletion phase is engaged. In the backward phase the model is simplified by deleting one least important basis function (according to GCV) at a time until the model again has only the intercept term. At the end of the backward phase, from those "best" models of each size one model of lowest GCV value is selected and outputted as the final one.

GCV for a model is calculated as follows (Hastie et al. 2009, Milborrow 2015):

$$GCV = MSE_{train} \Big/ \left(1 - \frac{enp}{n}\right)^2, \quad (1)$$

where MSE_{train} is Mean Squared Error of the evaluated model in the training data, n is the number of data cases in the training data, and enp is the effective number of parameters:

$$enp = k + c \times (k - 1)/2, \quad (2)$$

where k is the number of basis functions in the model (including the intercept term) and c is `trainParams.c`. Note that $(k-1)/2$ is the number of hinge function knots, so the formula penalizes not only the number of model's basis functions but also the number of knots. Also note that in ARESLab in the situation when $enp \geq n$ the GCV value will be equal to `Inf` (the model is considered infinitely bad).

After the pruning, the largest possible final model has $k = \text{int}((n + c/2)/(1 + c/2))$ basis functions for `maxInteractions` > 1 and $k = \text{int}((n + c/3)/(1 + c/3))$ basis functions for `maxInteractions` = 1. In the forward phase the models may also get larger than this however for such models GCV = `Inf` as then $enp \geq n$.

2.2. Function `aresparams`

Purpose:

Creates a structure of ARES configuration parameter values for further use with `aresbuild`, `arescv`, or `arescvc` functions.

Call:

```
trainParams = aresparams(maxFuncs, c, cubic, cubicFastLevel, selfInteractions, maxInteractions, threshold, prune, fastK, fastBeta, fastH, useMinSpan, useEndSpan, maxFinalFuncs, endSpanAdjust, newVarPenalty)
```

All the arguments of this function are optional. Empty values are also accepted (the corresponding default values will be used).

Parameters `c`, `prune`, and `maxFinalFuncs` are used in backward phase. Parameters `cubic` and `cubicFastLevel` can be used used in both phases. All other parameters are used in forward phase only.

For most applications, it can be expected that the most attention should be paid to the following parameters: `maxFuncs`, `c`, `cubic`, `maxInteractions` as well as `fastK`, `fastBeta`, and `fastH` if Fast MARS algorithm is to be used.

Input:

`maxFuncs` : The maximal number of basis functions included in the model in the forward model building phase (before pruning in the backward phase). Includes the intercept term. The recommended value for this parameter is about two times the expected number of basis functions in the final model (Friedman 1991). Note that the algorithm may also not reach the number. This can happen when the number of coefficients in the model exceeds the number of data cases or because of the threshold parameter. The default value for this parameter is -1 in which case maxFuncs is calculated automatically using formula $\min(200, \max(20, 2d)) + 1$, where d is the

number of input variables (Milborrow 2015). This is fairly arbitrary but can be useful for first experiments.

`c` : Generalized Cross-Validation (GCV) penalty per knot. Theory suggests values in the range of about 2 to 4. Larger values will lead to fewer knots being placed (i.e., final models will be simpler). A value of 0 penalizes only terms, not knots (can be useful e.g., with lots of data and low noise). The recommended (and default) value is 3 (Friedman 1991). Note that if `maxInteractions` = 1 (additive modelling) then function `aresbuild` will automatically recalculate *c* so that the actually used value is 2*c* / 3; this is recommended for additive modelling (Friedman 1991).

`cubic` : Whether to use piecewise-cubic (`true`) or piecewise-linear (`false`) type of modelling (Friedman 1991). In general, it is expected that the piecewise-cubic modelling will give slightly better predictive performance for smoother and less noisy data. (default value = `true`)

`cubicFastLevel` : ARESLab implements three levels of piecewise-cubic modelling. In level 0 cubic modelling for each candidate model is done in both phases of the technique (slow). In level 1 cubic modelling is done only in the backward phase (much faster). In level 2 cubic modelling is done after both phases only for the final model (fastest). The default and recommended level is 2. Theoretically, levels 0 and 1 may bring extra accuracy in the modelling process, however the results can actually also be worse. It is expected that the two much slower levels will mostly be not worth the waiting.

`selfInteractions` : This is experimental feature. The maximum degree of self interactions for any input variable. In ARESLab, it can be larger than 1 only for piecewise-linear modelling. Usually the self interactions are not allowed. (default value = 1, no self interactions)

`maxInteractions` : The maximum degree of interactions between input variables. Set to 1 for additive modelling (i.e., no interaction terms). For maximal interactivity between the variables, set the parameter to *d* ‖ `selfInteractions`, where *d* is the number of input variables – this way the modelling procedure will have the most freedom building a complex model. Typically only a low degree of interaction is allowed, but higher degrees can be used when the data warrants it. (default value = 1)

`threshold` : One of the stopping criteria for the forward phase. The larger the value of `threshold` the potentially simpler models are generated (see remarks section of `aresparams` and `aresbuild` for details). Default value = 1e-4. For noise-free data the value may be lowered.

`prune` : Whether to perform the model pruning (the backward phase). (default value = `true`)

`fastK` : Parameter (integer) for Fast MARS algorithm (Friedman 1993, Section 3.0). Maximum number of parent basis functions considered at each step of the forward phase. Typical values for `fastK` are 20, 10, 5 (default value = `Inf`, i.e., no Fast MARS). With lower `fastK` values model building is faster at the expense of some accuracy. Good starting values for fast exploratory work are `fastK` = 20, `fastBeta` = 1, `fastH` = 5 (Friedman 1993). Friedman in his paper concluded that while changing the values of `fastK` and `fastH` can have big effect on training computation times, predictive performance is largely unaffected over a wide range of their values (Friedman 1993).

`fastBeta` : Artificial ageing factor for Fast MARS algorithm (Friedman 1993, Section 3.1). Typical value for `fastBeta` is 1 (default value = 0, i.e., no artificial ageing). The parameter is ignored if `fastK` = `Inf`.

`fastH` : Parameter (integer) for Fast MARS algorithm (Friedman 1993, Section 4.0). Number of iterations till next full optimization over all input variables for each parent basis function. Higher values make the search faster. Typical values for `fastH` are 1, 5, 10 (default value = 1, i.e., full optimization in every iteration). Computational reduction associated with increasing `fastH` is most pronounced for data sets with many input variables and when large `fastK` is used. There seems to be little gain in increasing `fastH` beyond the value `fastH` = 5 (Friedman 1993). The parameter is ignored if `fastK` = `Inf`.

`useMinSpan` : In order to lower the local variance of the estimates, a minimum span is imposed that makes the technique resistant to runs of positive or negative error values between knots (by jumping over a `minSpan` number of data cases each time the next potential knot placement is requested) (Friedman 1991). `useMinSpan` allows to disable (set to 0 or 1) the protection so that all *x* values are considered for knot placement in each dimension (except, see `useEndSpan`). Disabling `minSpan` may allow creating a model which is more responsive to local variations in the data however this can also lead to an overfitted model. Setting the `useMinSpan` to > 1, enables to manually tune the value. (default and recommended value = -1 which corresponds to the automatic mode)

`useEndSpan` : In order to lower the local variance of the estimates near the ends of data intervals, a minimum span is imposed that makes the technique resistant to runs of positive or negative error values between extreme knot locations and the corresponding ends of data intervals (by not allowing to place a knot too near to the end of data interval) (Friedman 1991). `useEndSpan` allows to disable (set to 0) the protection so that all the data cases are considered for knot placement in each dimension (except, see `useMinSpan`). Disabling `endSpan` may allow creating a model which is more responsive to local variations in the data however this can also lead to a model that is overfitted near the edges of the data. Setting the `useMinSpan` to > 1, enables to manually tune the value. (default and recommended value = -1 which corresponds to the automatic mode)

`maxFinalFuncs` : Maximum number of basis functions (including the intercept term) in the pruned model. Use this (rather than the `maxFuncs` parameter) to enforce an upper bound on the final model size. (default value = `Inf`).

`endSpanAdjust` : For basis functions with variable interactions, `endSpan` gets multiplied by this value. This reduces probability of getting overfitted interaction terms supported by just a few cases on the boundaries of data intervals. Still, at least one knot will always be allowed in the middle, even if `endSpanAdjust` would prohibit it. Useful values range from 1 to 10. (default value = 1, i.e., no adjustment)

`newVarPenalty` : Penalty for adding a new variable to a model in the forward phase. This is the gamma parameter of Eq. 74 in the original paper (Friedman 1991). The higher is the penalty, the more reluctant will be the forward phase to add a new variable to the model – it will rather try to use variables already in the model. This can be useful when some of the variables are highly collinear. As a result, the final model may be easier to interpret although usually the built models also will have worse predictive performance. Useful non-zero values typically range from 0.01 to 0.2 (Milborrow 2015). (default value = 0, i.e., no penalty)

Output:

`trainParams` : A structure of training parameters for `aresbuild` function containing the provided values of the parameters (or default ones, if not provided).

Remarks:

The knot placement in `aresbuild` is implemented very similarly to R Earth package version 4.3.0 (Milborrow 2015) with calculations of `minSpan` and `endSpan` values using formulas given in Eq. 45 and Eq. 43 of the Friedman's original paper (Friedman 1991) with *alpha* = 0.05. Note that for a fixed dimensionality of the data, the `endSpan` value always stays the same but the value of the `minSpan` is recalculated for each individual parent basis function that is used for generation of new basis functions. The knots are placed symmetrically so that there are approximately equal number of skipped cases at each end of data intervals.

If more speed is required, try to use the Fast MARS algorithm by setting `fastK` parameter to something other than `Inf`. Good starting values for fast exploratory work are `fastK` = 20, `fastBeta` = 1, `fastH` = 5 (Friedman 1993). For more information, see descriptions of the mentioned parameters and the Friedman's paper.

Alternatively for more speed you can try some of the following options:

1) decreasing `maxFuncs` (less iterations in the forward phase);
2) increasing `cubicFastLevel` (see the description of this parameter);
3) decreasing `selfInteractions` (less candidate models in the forward phase);
4) decreasing `maxInteractions` (less candidate models in the forward phase);
5) increasing `threshold` (may result in less iterations in the forward phase);
6) manually increasing `useMinSpan` and/or `useEndSpan` (less candidate models in the forward phase).

Note that decreasing number of iterations or candidate models in the forward phase may also result in worse models.

2.3. Function `arespredict`

Purpose:

Predicts output values for the given query points using an ARES model.

Call:

```
Yq = arespredict(model, Xq)
```

Input:

`model` : ARES model

`Xq` : Inputs of query data points `Xq`(*i*,:), *i* = 1,...,*nq*

Output:

`Yq` : Predicted outputs of the query data points `Yq`(*i*), *i* = 1,...,*nq*

2.4. Function `arestest`

Purpose:

Tests an ARES model on a test data set (`Xtst`, `Ytst`).

Call:

```
[MSE, RMSE, RRMSE, R2] = arestest(model, Xtst, Ytst)
```

Input:

`model` : ARES model
`Xtst, Ytst` : Test data cases (`Xtst`(*i*,:), `Ytst`(*i*)), *i* = 1,...,*ntst*

Output:

`MSE` : Mean Squared Error
`RMSE` : Root Mean Squared Error
`RRMSE` : Relative Root Mean Squared Error
`R2` : Coefficient of Determination

2.5. Function `arescv`

Purpose:

Tests ARES performance using k-fold Cross-Validation.

Call:

```
[avgMSE, avgRMSE, avgRRMSE, avgR2, avgTime] = arescv(X, Y, trainParams, weights, k, shuffle, cvc_cTry, cvc_k, verbose)
```

All the arguments, except the first two, of this function are optional. Empty values are also accepted (the corresponding default values will be used).

Input:

`X, Y` : Data cases (`X`(*i*,:), `Y`(*i*)), *i* = 1,...,*n*
`trainParams` : See function `aresbuild`.
`weights` : See function `aresbuild`.
`k` : Value of *k* for k-fold Cross-Validation. The typical values are 5 or 10. For Leave-One-Out Cross-Validation set *k* equal to *n*. (default value = 10)
`shuffle` : Whether to shuffle the order of the data cases before performing Cross-Validation. Note that the random seed value can be controlled externally before calling `arescv`. (default value = `true`)
`cvc_cTry, cvc_k` : `cTry` and k values for `arescvc` function. Supply these values if you want to perform another Cross-Validation loop for finding the "best" penalty *c* value in each iteration of the outer Cross-Validation loop. (default values = [], meaning that a fixed *c* is used)
`verbose` : Set to `false` for no verbose. (default value = `true`)

Output:

`avgMSE` : Average Mean Squared Error
`avgRMSE` : Average Root Mean Squared Error
`avgRRMSE` : Average Relative Root Mean Squared Error
`avgR2` : Average Coefficient of Determination
`avgTime` : Average execution time

2.6. Function `arescvc`

Purpose:

Finds the "best" value for penalty *c* of the Generalized Cross-Validation criterion from a set of candidate values using k-fold Cross-Validation.

Call:

```
cBest = arescvc(X, Y, trainParams, cTry, weights, k, shuffle, verbose)
```

All the arguments, except the first three, of this function are optional. Empty values are also accepted (the corresponding default values will be used).

Input:

`X, Y`	: Data cases (`X`(*i*,:), `Y`(*i*)), *i* = 1,...,*n*
`trainParams`	: See function `aresbuild`.
`cTry`	: A set of candidate values for *c*. (default value = `1:5`)
`weights`	: See function `aresbuild`.
`k`	: Value of *k* for k-fold Cross-Validation. The typical values are 5 or 10. For Leave-One-Out Cross-Validation set *k* equal to *n*. (default value = 10)
`shuffle`	: Whether to shuffle the order of the data cases before performing Cross-Validation. Note that the random seed value can be controlled externally before calling `arescvc`. (default value = `true`)
`verbose`	: Set to `false` for no verbose. (default value = `true`)

Output:

`cBest`	: The "best" value for penalty *c*.

Remarks:

This function finds the "best" penalty *c* value using Cross-Validation in a clever way using function `aresbuild`. In each CV iteration the forward phase in `aresbuild` is done only once while the backward phase is done separately for each `cTry` value. The results will be the same as if each time a full model building process would be performed because in the forward phase the GCV criterion is not used.

2.7. Function `aresplot`

Purpose:

Plots surface of an ARES model.

Call:

```
aresplot(model, minX, maxX, vals, gridSize)
```

All the arguments, except the first one, of this function are optional. Empty values are also accepted (the corresponding default values will be used).

Input:

`model`	: ARES model
`minX, maxX`	: User defined minimum and maximum values for each input variable (this is the same type of data as in `model.minX` and `model.maxX`). If not supplied, the `model.minX` and `model.maxX` values will be used.
`vals`	: Only used when the number of input variables is larger than 2. This is a vector of fixed values for all the input variables except the two varied in the plot. The two varied variables are identified in `vals` using `NaN` values. By default the two first variables will be varied and all the other will be fixed at (*max* – *min*) / 2.
`gridSize`	: Grid size. (default value = 50)

2.8. Function `areseq`

Purpose:

Prints ARES model.

Call:

```
eq = areseq(model, precision, varNames)
```

Piecewise-cubic spline equations are very complex as they involve smoothing. For easier interpretation use piecewise-linear models.

Input:

`model`	: ARES model
`precision`	: Optional. Number of digits in the model coefficients and knot sites. (default value = 15)
`varNames`	: Optional. A cell array of variable names to show instead of the generic ones. This is used for piecewise-linear models only.

Output:

`eq`	: A cell array of equations for individual basis functions and the main model.

2.9. Function `aresanova`

Purpose:

Performs ANOVA decomposition (see sections 3.5 and 4.3 of the original paper by Jerome Friedman (Friedman 1991) for details) of the given ARES model and reports the results.

Call:

```
aresanova(model, Xtr, Ytr)
```

Input:

`model`	: ARES model
`Xtr, Ytr`	: Training data cases (`Xtr`(i,:), `Ytr`(i)), $i = 1,...,n$.

Remarks:

To understand the table outputted by the function, here is an excerpt from the original paper by Jerome Friedman (Friedman 1991) Section 4.3:

> *"The ANOVA decomposition is summarized by one row for each ANOVA function. The columns represent summary quantities for each one. The first column lists the function number. The second gives the standard deviation of the function. This gives one indication of its (relative) importance to the overall model and can be interpreted in a manner similar to a standardized regression coefficient in a linear model. The third column provides another indication of the importance of the corresponding ANOVA function, by listing the GCV, score for a model with all of the basis functions corresponding to that particular ANOVA function removed. This can be used to judge whether this ANOVA function is making an important contribution to the model, or whether it just slightly helps to improve the global GCV score. The fourth column gives the number of basis functions comprising the ANOVA function while the fifth column provides an estimate of the additional number of linear degrees-of-freedom used by including it. The last column gives the particular predictor variables associated with the ANOVA function."*

2.10. Function `aresanovareduce`

Purpose:

Deletes all the basis functions from an ARES model in which at least one used variable is not in the given list of allowed variables. This can be used to perform ANOVA decomposition as well as for investigation of individual and joint contributions of variables in the model, i.e., the reduced model can be plotted using function aresplot to visualize the contributions.

Call:

```
[modelReduced, usedBasis] = aresanovareduce(model, vars, exact)
```

Input:

`model`	: ARES model
`vars`	: A vector of indexes for input variables to stay in the model. The size of the vector should be between 1 and *d*, where *d* is the total number of input variables.
`exact`	: Set this to true to get a model with only those basis functions where the exact combination of variables is present (default value = `false`). This is used from function `aresanova`.

Output:

`modelReduced`	: The reduced model
`usedBasis`	: The list of original indexes for the used basis functions

3. EXAMPLES OF USAGE

3.1. Ten-dimensional function with noise

We start by creating a data set using a ten-dimensional function with Gaussian noise. The data consists of 200 cases randomly uniformly distributed in a ten-dimensional unit hypercube.

```
clear
X = rand([200,10]);
Y = 10.*sin(pi.*X(:,1).*X(:,2)) + 20.*(X(:,3)-0.5).^2 + ...
    10.*X(:,4) + 5.*X(:,5) + randn(200,1)*0.5;
```

We define the maximal number of basis functions to be 21 (including the intercept term), and limit maximum interaction level to 2 (only pairwise products of basis functions will be allowed), leaving all the other parameters to their defaults. The model will be of piecewise-cubic type as it is the default.

```
params = aresparams(21, [], [], [], [], 2);
```

If we want to find a better value for parameter `c` than the default, we run `arescvc`. Let's try all values from 1 to 8.

```
cBest = arescvc(X, Y, params, 1:8)

c = 1  MSE = 0.36969
c = 2  MSE = 0.36567
c = 3  MSE = 0.36427
c = 4  MSE = 0.36427
c = 5  MSE = 0.37936
c = 6  MSE = 0.39912
c = 7  MSE = 0.40659
c = 8  MSE = 0.42365
cBest = 3
```

In this case the “best” value for `c` turned out to be the same as the default. So we won't reconfigure.

ARES model is built by calling `aresbuild`.

```
model = aresbuild(X, Y, params)
```

As the model building process ends, we can examine the data structure of the final model. It has 16 basis functions including the intercept term.

```
model =
          coefs: [16x1 double]
       knotdims: {15x1 cell}
      knotsites: {15x1 cell}
       knotdirs: {15x1 cell}
        parents: [15x1 double]
    trainParams: [1x1 struct]
            MSE: 0.3234
            GCV: 0.4960
             t1: [15x10 double]
             t2: [15x10 double]
           minX: [1x10 double]
```

```
          maxX: [1x10 double]
        endSpan: 10
```

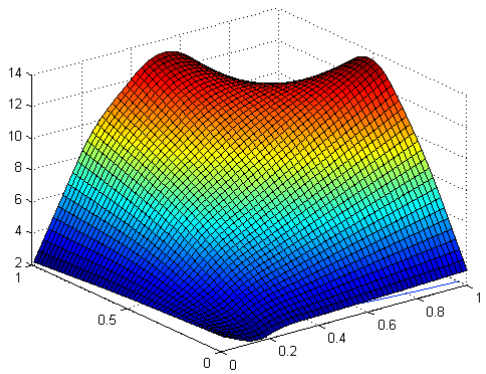
Now we can perform ANOVA decomposition.

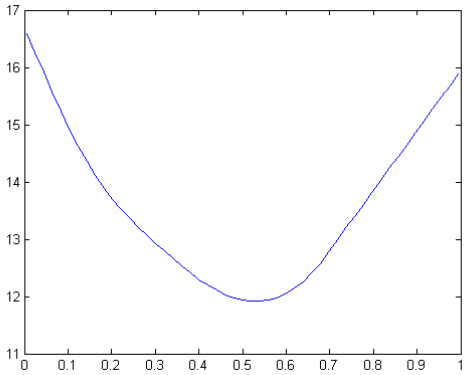
```
aresanova(model, X, Y)

Type: piecewise-cubic
GCV: 0.496
Total number of basis functions: 16
Total effective number of parameters: 38.5
ANOVA decomposition:
Func.     STD    GCV          #basis       #params      variable(s)
1         4.675 52.925       2             5.0         1
2         2.621 10.342       2             5.0         2
3         1.281 59.738       2             5.0         3
4         2.956 17.665       2             5.0         4
5         1.445  8.879       1             2.5         5
6         3.732 37.938       5            12.5         1 2
7         0.289  0.617       1             2.5         3 5
```

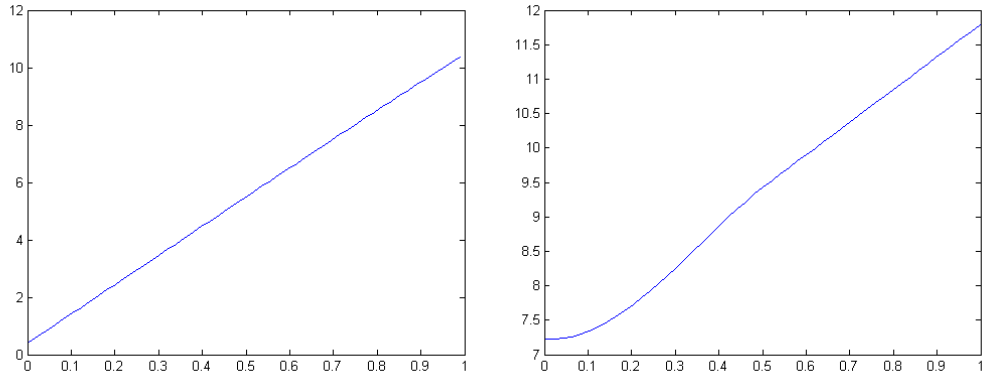
We can see that the last ANOVA function gives very small contribution and maybe should be deleted (it corresponds to one basis function which uses the 3rd and the 5th input variable).

Let's plot pair-wise (for variables x_1 and x_2) and individual (for variables x_3, x_4, and x_5 contributions of variables.

```
modelReduced = aresanovareduce(model, [1 2])
aresplot(modelReduced)

for i = 3 : 5
    modelReduced = aresanovareduce(model, i);
    Xtmp = zeros(51,length(model.minX));
    Xtmp(:,i) = [model.minX(i):((model.maxX(i)-model.minX(i))/50):model.maxX(i)]';
    figure
    plot(Xtmp(:,i), arespredict(modelReduced, Xtmp));
end
```







Now let's evaluate performance of this ARES configuration on the data using 10-fold Cross-Validation.

```
rand('state',0);
[avgMSE, avgRMSE, avgRRMSE, avgR2] = arescv(X, Y, params)

avgMSE = 0.4427
avgRMSE = 0.6543
avgRRMSE = 0.1339
avgR2 = 0.9818
```

Now let's try piecewise-linear modelling.

```
params = aresparams(21, [], false, [], [], 2);
model = aresbuild(X, Y, params)

model =
          coefs: [16x1 double]
       knotdims: {15x1 cell}
      knotsites: {15x1 cell}
       knotdirs: {15x1 cell}
        parents: [15x1 double]
    trainParams: [1x1 struct]
            MSE: 0.3802
            GCV: 0.5830
           minX: [1x10 double]
           maxX: [1x10 double]
        endSpan: 10

rand('state',0);
[avgMSE, avgRMSE, avgRRMSE, avgR2] = arescv(X, Y, params)

avgMSE = 0.5254
avgRMSE = 0.7165
avgRRMSE = 0.1469
avgR2 = 0.9783
```

Finally we output the equation of the piecewise-linear model with all its basis functions.

```
areseq(model, 5);

BF1 = max(0, x4 -0.66938)
BF2 = max(0, 0.66938 -x4)
BF3 = max(0, x2 -0.57231)
BF4 = max(0, 0.57231 -x2)
BF5 = max(0, x1 -0.23961)
BF6 = max(0, 0.23961 -x1)
BF7 = max(0, x5 -0.036179)
BF8 = max(0, 0.553 -x3)
```

```
    BF9 = BF3 * max(0, x1 -0.5981)
    BF10 = BF3 * max(0, 0.5981 -x1)
    BF11 = BF4 * max(0, x1 -0.14539)
    BF12 = BF4 * max(0, 0.14539 -x1)
    BF13 = max(0, x3 -0.16723)
    BF14 = BF7 * max(0, x3 -0.8273)
    BF15 = BF5 * max(0, x2 -0.25707)
    y =  7.6853 +9.8771*BF1 -10.312*BF2 +15.52*BF3 -6.1597*BF4 +15.121*BF5 -20.444*BF6 +4.9892*BF7
+15.757*BF8 -65.272*BF9 -28.664*BF10 -25.491*BF11 +53.099*BF12 +9.7016*BF13 +18.575*BF14 -18.91*BF15
```

3.2. Noise-free two-dimensional function

We start by creating training and test data using a two-dimensional noise-free function. The training data consists of 121 cases distributed in a regular 11×11 grid. The test data has 10000 cases distributed randomly.

```
clear
[tmpX1,tmpX2] = meshgrid(-1:0.2:1, -1:0.2:1);
X(:,1) = reshape(tmpX1, numel(tmpX1), 1);
X(:,2) = reshape(tmpX2, numel(tmpX2), 1);
clear tmpX1; clear tmpX2;
Y = sin(0.83.*pi.*X(:,1)) .* cos(1.25.*pi.*X(:,2));

Xt = rand([10000,2]);
Yt = sin(0.83.*pi.*Xt(:,1)) .* cos(1.25.*pi.*Xt(:,2));
```

Such noise-free functions can be approximated very precisely. We define the maximal number of basis functions to be 121, no penalty for knots, and maximum interaction level equal to 2 (the number of input variables), leaving all the other parameters to their defaults. The model will be of piecewise-cubic type as it is the default.

```
params = aresparams(121, 0, [], [], [], 2);
```

ARES model is built by calling `aresbuild`.

```
model = aresbuild(X, Y, params)
```

As the model building process ends, we can examine the data structure of the new model. The model has 44 basis functions including the intercept term.

```
model =
          coefs: [44x1 double]
       knotdims: {43x1 cell}
      knotsites: {43x1 cell}
       knotdirs: {43x1 cell}
        parents: [43x1 double]
    trainParams: [1x1 struct]
            MSE: 1.6901e-004
            GCV: 4.1734e-004
             t1: [43x2 double]
             t2: [43x2 double]
           minX: [-1 -1]
           maxX: [1 1]
        endSpan: 8
```

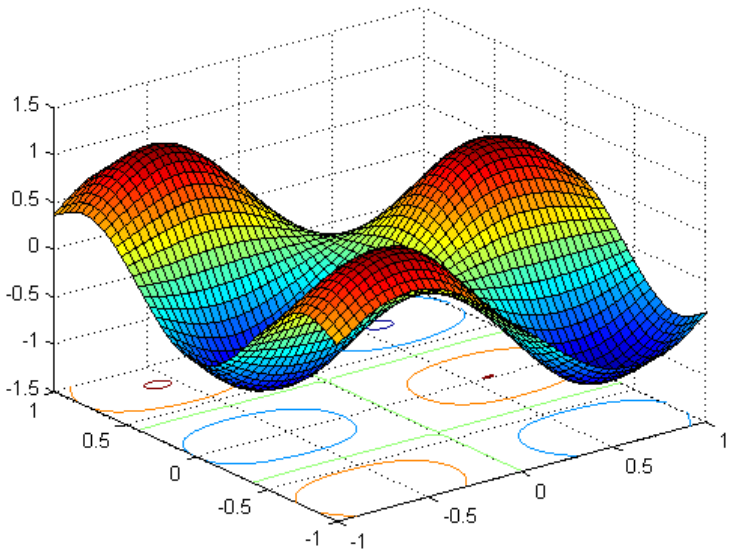
We test the model using test data.

```
[MSE, RMSE, RRMSE, R2] = arestest(model, Xt, Yt)
```

```
MSE = 1.9909e-004
RMSE = 0.0141
RRMSE = 0.0244
R2 = 0.9994
```

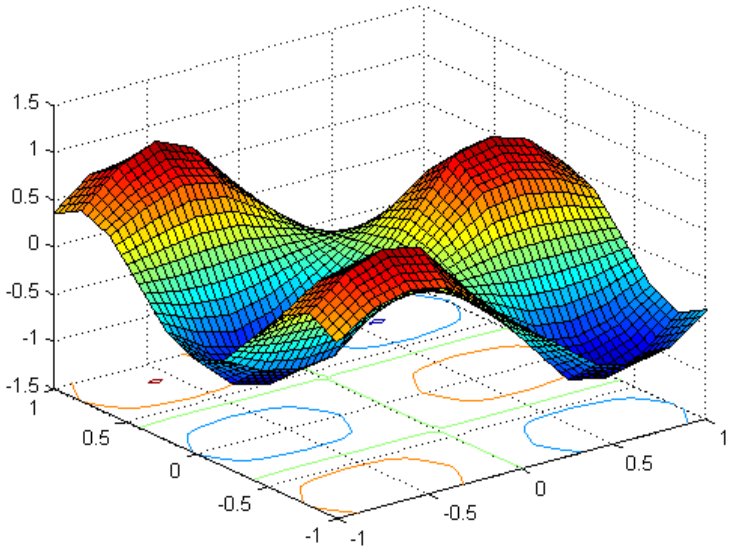
Plot the surface of the model.

```
aresplot(model);
```



Let's try doing the same but instead of piecewise-cubic modelling we will use piecewise-linear.

```
params = aresparams(121, 0, false, [], [], 2);
model = aresbuild(X, Y, params);
[MSE, RMSE, RRMSE, R2] = arestest(model, Xt, Yt)

MSE = 0.0023
RMSE = 0.0480
RRMSE = 0.0829
R2 = 0.9931

aresplot(model);
```



4. REFERENCES

1. Friedman J.H. Multivariate Adaptive Regression Splines (with discussion), The Annals of Statistics, Vol. 19, No. 1, 1991, pp. 1-141
2. Friedman J.H. Fast MARS, Department of Statistics, Stanford University, Tech. Report LCS110, 1993
3. Hastie T., Tibshirani R., Friedman J. The elements of statistical learning: Data mining, inference and prediction, 2nd edition, Springer, 2009
4. Merkwirth C. and Wichard J. A Matlab toolbox for ensemble modelling, 2003, available at http://www.j-wichard.de
5. Milborrow S., Earth: Multivariate Adaptive Regression Spline Models (derived from code by T. Hastie and R. Tibshriani), 2015, R package available at http://cran.r-project.org/src/contrib/Descriptions/earth.html
6. Norgaard M. Neural network based system identification toolbox, ver. 2, Tech. Report. 00-E-891, Department of Automation, Technical University of Denmark, 2000
7. Rudy J., py-earth: A Python implementation of Jerome Friedman's Multivariate Adaptive Regression Splines, 2015, available at https://github.com/jcrudy/py-earth