

# Package ‘GaSP’

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**Type** Package

**Title** Train and Apply a Gaussian Stochastic Process Model

**Version** 1.0.1

## Description

Train a Gaussian stochastic process model of an unknown function, possibly observed with error, via maximum likelihood or MAP estimation, run model diagnostics, and make predictions, following Sacks, J., Welch, W.J., Mitchell, T.J., and Wynn, H.P. (1989) "Design and Analysis of Computer Experiments", *Statistical Science*, <[doi:10.1214/ss/1177012413](https://doi.org/10.1214/ss/1177012413)>. Perform sensitivity analysis and visualize low-order effects, following Schonlau, M. and Welch, W.J. (2006), "Screening the Input Variables to a Computer Model Via Analysis of Variance and Visualization", <[doi:10.1007/0-387-28014-6\\_14](https://doi.org/10.1007/0-387-28014-6_14)>.

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borehole	<i>Data for the borehole function</i>
----------	---------------------------------------

---

## Description

Training and test data for the borehole function; see source for background.

## Usage

```
borehole
```

## Format

A list with the following four data frames:

**x** 8-dimensional input for 40 training runs.

**y** Output (the flow) for the 40 training runs in **x**.

**x\_pred** 8-dimensional input for 1000 test runs at which to predict **y**.

**y\_true** Output for the 1000 runs in **x\_pred**.

## Source

<https://www.sfu.ca/~ssurjano/borehole.html>

---

CrossValidate	<i>Cross-validated predictions for a GaSPModel object.</i>
---------------	--

---

**Description**

Compute leave-one-out cross-validated predictions for a GaSPModel object.

**Usage**

```
CrossValidate(GaSP_model)
```

**Arguments**

GaSP\_model      Object of class [GaSPModel](#).

**Value**

A data frame with two columns: the cross-validated predictions Pred and their standard errors SE.

**Note**

[RMSE](#) computes the root mean squared error of the predictions. [PlotPredictions](#) and [PlotResiduals](#) plot the predictions or their residuals; [PlotStdResiduals](#) and [PlotQQ](#) plot the standardized residuals.

**Examples**

```
borehole_cv <- CrossValidate(borehole_fit)
```

---

DescribeX	<i>Describe the input variables.</i>
-----------	--------------------------------------

---

**Description**

Describe the input variables to set up integration or summation ranges for Visualize.

**Usage**

```
DescribeX(  
  x_names,  
  x_min,  
  x_max,  
  support = NULL,  
  num_levels = NULL,  
  distribution = NULL  
)
```

**Arguments**

<code>x_names</code>	A vector of character strings containing the names of the input variables.
<code>x_min, x_max</code>	Vectors of the same length as <code>x_names</code> containing the minima and maxima, respectively, of the input variables.
<code>support</code>	Optional vector of character strings of the same length as <code>x_names</code> . Valid strings for a variable are: "Continuous" (continuous between the input's <code>x_min</code> and <code>x_max</code> ); "Fixed" (the input's <code>x_min</code> must equal its <code>x_max</code> ); and "Grid" (which requires the next argument).
<code>num_levels</code>	An optional vector of integers for the number of levels of each input; must be present if the <code>support</code> argument includes "Grid". An input's number of levels is 0 if it is "Continuous", 1 if it is "Fixed", or > 1 if it is "Grid" to define an equally spaced grid inclusive of the input's <code>x_min</code> and <code>x_max</code> .
<code>distribution</code>	An optional vector of character strings of the same length as <code>x_names</code> to define the weight distributions of the input variables. Valid strings are "Uniform" or "Normal" (ignored for "Fixed" inputs).

**Value**

A data frame with the following columns: `Variable` (containing `x_names`), `Min` (containing `x_min`), and `Max` (containing `x_max`), plus the optional columns `Support` (from `support`), `NumberLevels` (from `num_levels`), and `Distribution` (from `distribution`).

**Note**

Does not check against [GaSPModel](#) and all characters are CASE SENSITIVE.

**Examples**

```
borehole_x_names <- colnames(borehole$x)
borehole_min <- c(0.05, 100.00, 63070.00, 990.00, 63.10, 700.00, 1120.00, 9855.00)
borehole_max <- c(0.15, 50000.00, 115600.00, 1110.00, 116.00, 820.00, 1680.00, 12045.00)
borehole_x_desc <- DescribeX(borehole_x_names, borehole_min, borehole_max)
```

---

 Fit

*Fit a GaSP model.*


---

**Description**

Fit (train) a GaSP model.

**Usage**

```
Fit(
  x,
  y,
  reg_model,
  sp_model = NULL,
  cor_family = c("PowerExponential", "Matern"),
  cor_par = data.frame(0),
  random_error = c(FALSE, TRUE),
  sp_var = -1,
  error_var = -1,
  nugget = 1e-09,
  tries = 10,
  seed = 500,
  fit_objective = c("Likelihood", "Posterior"),
  theta_standardized_min = 0,
  theta_standardized_max = .Machine$double.xmax,
  alpha_min = 0,
  alpha_max = 1,
  derivatives_min = 0,
  derivatives_max = 3,
  log_obj_tol = 1e-05,
  log_obj_diff = 0,
  lambda_prior = 0.1,
  model_comparison = c("Objective", "CV")
)
```

**Arguments**

<code>x</code>	A data frame containing the input (explanatory variable) training data.
<code>y</code>	A vector or a data frame with one column containing the output (response) training data.
<code>reg_model</code>	The regression model, specified as a formula, but note the left-hand side of the formula is unused; see example.
<code>sp_model</code>	An optional stochastic process model, specified as a formula, but note the left-hand side of the formula and the intercept are unused. The default NULL uses all column names in <code>x</code> .
<code>cor_family</code>	A character string specifying the (product, anisotropic) correlation-function family: "PowerExponential" for the power-exponential family or "Matern" for the Matern family.
<code>cor_par</code>	An optional data frame containing the correlation parameters with one row per <code>sp_model</code> term and two columns set up as described in <a href="#">GaSPModel</a> Details; only used to start the first objective optimization (see Details).
<code>random_error</code>	A boolean for the presence or not of a random (measurement, white-noise) error term.

sp_var, error_var	Starting values of the stochastic process and error variances for the first try to optimize the objective (see Details); valid (i.e., nonnegative) values will only be used if random_error = TRUE. The invalid default value of -1 indicates that a starting value will be chosen by Fit.
nugget	For numerical stability the proportion of the total variance due to random error is fixed at this value (random_error = FALSE) or bounded below by it (random_error = TRUE).
tries	Number of optimizations of the objective from different random starting points.
seed	The random-number seed to generate starting points.
fit_objective	The objective that Fit attempts to optimize: "Likelihood" (maximum likelihood estimation) or "Posterior" (Bayesian maximum a posteriori estimation).
theta_standardized_min, theta_standardized_max	The minimum and maximum of the standardized $\theta$ parameter (see Details).
alpha_min, alpha_max	The minimum and maximum of the $\alpha$ parameter of power-exponential.
derivatives_min, derivatives_max	The minimum and maximum of the $\delta$ parameter of Matern.
log_obj_tol	An absolute tolerance for terminating the optimization of the log of the objective.
log_obj_diff	The critical value for the change in the log objective for informal tests during optimization of correlation parameters. No testing is done with the default of 0; a larger critical value such as 2 may give a more parsimonious model.
lambda_prior	The rate parameter of an exponential prior for each $\theta$ parameter; used only if fit_objective = "Posterior".
model_comparison	The criterion used to select from multiple solutions when tries > 1: the objective function ("Objective") or leave-one-out cross validation ("CV").

## Details

Fit numerically optimizes the profile objective function with respect to the correlation parameters; the mean and overall variance parameters are estimated in closed form given the correlation parameters.

A cor\_par data frame supplied by the user is the starting point for the first optimization try. If random\_error = TRUE, then sp\_var / (sp\_var + error\_var) is another correlation parameter to be optimized; sp\_var and error\_var values supplied by the user will initialize this parameter for the first try.

Set random\_error = TRUE to estimate the variance of the random (measurement, white-noise) error; a small nugget error variance is for numerical stability.

For term  $j$  in the stochastic-process model, the estimate of  $\theta_j$  is constrained between theta\_standardized\_min /  $r_j^2$  and theta\_standardized\_max /  $r_j^2$ , where  $r_j$  is the range of term  $j$ . Note that Fit returns unscaled estimates relating to the original, unscaled inputs.

**Value**

A GaSPModel object, which is a list with the following components:

x	The data frame containing the input training data.
y	The training output data, now as a vector.
reg_model	The regression model, now in the form of a data frame.
sp_model	The stochastic process model, now in the form of a data frame.
cor_family	The correlation family.
cor_par	A data frame for the estimated correlation parameters.
random_error	The boolean for the presence or not of a random error term.
sp_var	The estimated stochastic process variance.
error_var	The estimated random error variance.
beta	A data frame holding the estimated regression-model parameters.
objective	The maximum value found for the objective function: the log likelihood (fit_objective = "Likelihood") or the log posterior (fit_objective = "Posterior").
cond_num	The condition number.
CVRMSE	The leave-one-out cross-validation root mean squared error.

**References**

Sacks, J., Welch, W.J., Mitchell, T.J., and Wynn, H.P. (1989) "Design and Analysis of Computer Experiments", *Statistical Science*, 4, pp. 409-423, doi:10.1214/ss/1177012413.

**Examples**

```
x <- borehole$x
y <- borehole$y
borehole_fit <- Fit(
  reg_model = ~1, x = x, y = y, cor_family = "Matern",
  random_error = FALSE, nugget = 0, fit_objective = "Posterior"
)
```

---

GaSPModel

---

*Create a GaSPModel object.*


---

**Description**

Return a template for a GaSPModel object.

**Usage**

```

GaSPModel(
  x,
  y,
  reg_model,
  sp_model = NULL,
  cor_family = c("PowerExponential", "Matern"),
  cor_par,
  random_error = c(FALSE, TRUE),
  sp_var,
  error_var = 0
)

```

**Arguments**

x	A data frame containing the input (explanatory variable) training data.
y	A vector or a data frame with one column containing the output (response) training data.
reg_model	The regression model, specified as a formula, but note the left-hand side of the formula is unused; see example.
sp_model	An optional stochastic process model, specified as a formula, but note the left-hand side of the formula and the intercept are unused. The default NULL uses all column names in x.
cor_family	A character string specifying the (product, anisotropic) correlation-function family: "PowerExponential" for the power-exponential family or "Matern" for the Matern family.
cor_par	A data frame containing the correlation parameters with one row per sp_model term and two columns (see Details).
random_error	A boolean for the presence or not of a random (measurement, white-noise) error term.
sp_var	The stochastic process variance.
error_var	The random error variance, with default 0.

**Details**

The data frame `cor_par` contains one row for each term in the stochastic process model. There are two columns. The first is named `Theta`, and the second is either `Alpha` (power-exponential) or `Derivatives` (Matern). Let  $h_j$  be a distance between points for term  $j$  in the stochastic-process model. For power-exponential, the contribution to the product correlation from term  $j$  depends on a distance-scale parameter  $\theta_j$  from the `Theta` column and a smoothness parameter  $\alpha_j$  from the `Alpha` column; the contribution is  $\exp(-\theta_j h_j^{2-\alpha_j})$ . For example,  $\alpha_j = 0$  gives the squared-exponential (Gaussian) correlation. The contribution to the product correlation for Matern also depends on  $\theta_j$ , and the second parameter is the number of derivatives  $\delta_j = 0, 1, 2, 3$  from the `Derivatives` column. The contribution is  $\exp(-\theta_j h_j)$  for  $\delta_j = 0$  (the exponential correlation),  $\exp(-\theta_j h_j)(\theta_j h_j + 1)$  for  $\delta_j = 1$ ,  $\exp(-\theta_j h_j)((\theta_j h_j)^2/3 + \theta_j h_j + 1)$  for  $\delta_j = 2$ , and  $\exp(-\theta_j h_j^2)$  for  $\delta_j = 3$  (the squared-exponential correlation). Note that  $\delta_j = 3$  codes for a limiting infinite number of derivatives. This

is not the usual parameterization of the Matern, but it is consistent with power-exponential for the exponential and squared-exponential special cases common to both.

A value should be given to `error_var` if the model has a random-error term (`random_error = TRUE`), and a small "nugget" such as  $10^{-9}$  may be needed for improved numerical conditioning.

### Value

A `GaSPModel` object, which is a list with the following components:

<code>x</code>	The data frame containing the input training data.
<code>y</code>	The training output data, now as a vector.
<code>reg_model</code>	The regression model, now in the form of a data frame.
<code>sp_model</code>	The stochastic process model, now in the form of a data frame.
<code>cor_family</code>	The correlation family.
<code>cor_par</code>	The data frame containing the correlation parameters.
<code>random_error</code>	The boolean for the presence or not of a random error term.
<code>sp_var</code>	The stochastic process variance.
<code>error_var</code>	The random error variance.
<code>beta</code>	A placeholder for a data frame to hold the regression-model parameters.
<code>objective</code>	A placeholder for the maximum fit objective.
<code>cond_num</code>	A placeholder for the condition number.
<code>CVRMSE</code>	A placeholder for the model's cross-validated root mean squared error.

### Note

This function does not execute `Fit` and is intended for `CrossValidate`, `Predict` and `Visualize` with models trained otherwise by the user. Placeholders do not need to be specified to execute these further functions, as they are always recomputed as needed.

### References

Sacks, J., Welch, W.J., Mitchell, T.J., and Wynn, H.P. (1989) "Design and Analysis of Computer Experiments", *Statistical Science*, 4, pp. 409-423, doi:10.1214/ss/1177012413.

### Examples

```
x <- borehole$x
y <- borehole$y
theta <- c(
  5.767699e+01, 0.000000e+00, 0.000000e+00, 1.433571e-06,
  0.000000e+00, 2.366557e-06, 1.695619e-07, 2.454376e-09
)
alpha <- c(
  1.110223e-16, 0.000000e+00, 0.000000e+00, 0.000000e+00,
  0.000000e+00, 0.000000e+00, 2.494862e-03, 0.000000e+00
)
cor_par <- data.frame(Theta = theta, Alpha = alpha)
```

```

rownames(cor_par) <- colnames(borehole$x)
sp_var <- 38783.7
borehole_gasp <- GaSPModel(
  x = borehole$x, y = borehole$y,
  reg_model = ~1, cor_family = "PowerExponential",
  cor_par = cor_par, random_error = FALSE,
  sp_var = sp_var
)

```

---

PlotAll	<i>Execute</i> PlotPredictions, PlotResiduals, PlotStdResiduals, PlotMainEffects, and PlotJointEffects.
---------	---

---

### Description

Execute PlotPredictions, PlotResiduals and PlotStdResiduals (all applied to CV only), PlotMainEffects, and PlotJointEffects.

### Usage

```

PlotAll(
  GaSP_model,
  cross_validation,
  visualization,
  y_name = "y",
  y_units = "",
  x_units = NULL,
  se_plot = TRUE,
  y_values = NULL,
  se_values = NULL,
  pch = 1
)

```

### Arguments

GaSP_model	Object of class <a href="#">GaSPModel</a> , the entire model will be verified but only x and y will be used.
cross_validation	A data frame returned by <a href="#">CrossValidate</a> .
visualization	A list object returned by <a href="#">Visualize</a> .
y_name	An optional character string containing the output variable name (for labels).
y_units	An optional character string containing the units of the output variable (for labels).
x_units	An optional vector of character strings containing the units of the input variables (for labels).
se_plot	An optional boolean indicating whether to make standard-error contour plots.

y_values	An optional vector of contour values for the estimated joint effects.
se_values	An optional vector of contour values for the standard errors.
pch	Plotting symbol for plot; default is open circle.

**Value**

No return value, generates plots.

**Examples**

```
PlotAll(borehole_fit, borehole_cv, borehole_vis)
```

---

PlotJointEffects	<i>Plot the estimated joint effects.</i>
------------------	--

---

**Description**

Plot the estimated joint effects.

**Usage**

```
PlotJointEffects(
  joint_effect,
  anova_percent,
  x_units = NULL,
  y_name = "y",
  y_units = "",
  se_plot = TRUE,
  y_values = NULL,
  se_values = NULL
)
```

**Arguments**

joint_effect	A data frame from <a href="#">Visualize</a> with plotting coordinates for the estimated joint effects.
anova_percent	A data frame from <a href="#">Visualize</a> of ANOVA percentages.
x_units	An optional vector of character strings containing the units of the input variables (for labels).
y_name	An optional character string containing the output variable name (for labels).
y_units	An optional character string containing the units of the output variable (for labels).
se_plot	An optional boolean indicating whether to make standard-error contour plots.
y_values	An optional vector of contour values for the estimated joint effects.
se_values	An optional vector of contour values for the standard errors.

**Details**

Plots are sent to the active device.

**Value**

No return value, generates plots.

**Examples**

```
PlotJointEffects(borehole_vis$joint_effect, borehole_vis$anova_percent)
```

---

PlotMainEffects	<i>Plot the estimated main effects.</i>
-----------------	---

---

**Description**

Plot the estimated main effects.

**Usage**

```
PlotMainEffects(
  main_effect,
  anova_percent,
  x_units = NULL,
  y_name = "y",
  y_units = ""
)
```

**Arguments**

main_effect	A data frame from <a href="#">Visualize</a> with plotting coordinates for the estimated main effects.
anova_percent	A data frame from <a href="#">Visualize</a> of ANOVA percentages.
x_units	An optional vector of character strings containing the units of the input variables (for labels).
y_name	An optional character string containing the output variable name (for labels).
y_units	An optional character string containing the units of the output variable (for labels).

**Details**

Plots are sent to the active device. Each plot shows an estimated main effect (red solid line) and pointwise approximate 95% confidence limits (green dashed line).

**Value**

No return value, generates plots.

**Examples**

```
PlotMainEffects(borehole_vis$main_effect, borehole_vis$anova_percent)
```

---

PlotPredictions      *Plot true versus predicted output.*

---

**Description**

Plot true versus predicted output (response) made by Predict or CrossValidate.

**Usage**

```
PlotPredictions(  
  y_pred,  
  y,  
  y_name = "y",  
  y_units = "",  
  title = c("Predict", "CrossValidate"),  
  pch = 1  
)
```

**Arguments**

y_pred	A data frame of predicted output values made by Predict or CrossValidate.
y	A vector of length equal to the number of rows in y_pred containing the true output values.
y_name	An optional character string containing the output variable name (for labels).
y_units	An optional character string containing the units of the output variable (for labels).
title	A character string for the name of the function generating the predictions (for an appropriate title): "Predict" from <a href="#">Predict</a> or "CrossValidate" from <a href="#">CrossValidate</a> ; "" for no title.
pch	Plotting symbol for plot; default is open circle.

**Value**

No return value, generates plots.

**Examples**

```
PlotPredictions(borehole_cv, y, title = "CrossValidate")
PlotPredictions(borehole_pred$y_pred, borehole$y_true, title = "Predict")
```

---

PlotQQ	<i>Normal Q-Q plot.</i>
--------	-------------------------

---

**Description**

Normal Q-Q plot of the standardized residuals of predictions from Predict or CrossValidate.

**Usage**

```
PlotQQ(y_pred, y, y_name = "y")
```

**Arguments**

y_pred	A data frame of predicted output values made by Predict or CrossValidate.
y	A vector of length equal to the number of rows in y_pred containing the true output values.
y_name	An optional character string containing the output variable name (for labels).

**Value**

No return value, generates plots.

**Examples**

```
PlotQQ(borehole_cv, y)
```

---

PlotResiduals	<i>Plot residuals versus each input variable.</i>
---------------	---

---

**Description**

Plot residuals versus each input variable.

**Usage**

```
PlotResiduals(
  x,
  y_pred,
  y,
  x_units = NULL,
  y_name = "y",
  y_units = "",
  pch = 1
)
```

**Arguments**

x	A data frame with number of rows equal to the number of rows in y_pred containing the input (explanatory) variables.
y_pred	A data frame of predicted output values made by Predict or CrossValidate.
y	A vector of length equal to the number of rows in y_pred containing the true output values.
x_units	An optional vector of character strings containing the units of the input variables in x (for labels).
y_name	An optional character string containing the output variable name (for labels).
y_units	An optional character string containing the units of the output variable (for labels).
pch	Plotting symbol for plot; default is open circle.

**Value**

No return value, generates plots.

**Examples**

```
PlotResiduals(x, borehole_cv, y)
```

---

PlotStdResiduals	<i>Plot standardized residuals versus predictions.</i>
------------------	--

---

**Description**

Plot standardized residuals versus predictions made by Predict or CrossValidate.

**Usage**

```
PlotStdResiduals(
  y_pred,
  y,
  y_name = "y",
  y_units = "",
  title = c("Predict", "CrossValidate"),
  pch = 1
)
```

**Arguments**

<code>y_pred</code>	A data frame of predicted output values made by <code>Predict</code> or <code>CrossValidate</code> .
<code>y</code>	A vector of length equal to the number of rows in <code>y_pred</code> containing the true output values.
<code>y_name</code>	An optional character string containing the output variable name (for labels).
<code>y_units</code>	An optional character string containing the units of the output variable (for labels).
<code>title</code>	A character string for the name of the function generating the predictions (for an appropriate title): "Predict" from <code>Predict</code> or "CrossValidate" from <code>CrossValidate</code> ; "" for no title.
<code>pch</code>	Plotting symbol for plot; default is open circle.

**Value**

No return value, generates plots.

**Examples**

```
PlotStdResiduals(borehole_cv, y, title = "CrossValidate")
```

---

Predict

*Predict from a GaSPModel object.*

---

**Description**

Predict from a `GaSPModel` object.

**Usage**

```
Predict(GaSP_model, x_pred, generate_coefficients = c(FALSE, TRUE))
```

**Arguments**

<code>GaSP_model</code>	Object of class <a href="#">GaSPModel</a> .
<code>x_pred</code>	A data frame containing the values of the input variables at which to predict the output.
<code>generate_coefficients</code>	A boolean indicating whether coefficients for further external predictions are generated.

**Value**

A list with the following elements:

<code>y_pred</code>	A data frame with two columns: the predictions <code>Pred</code> and their standard errors <code>SE</code> .
<code>pred_coeffs</code>	A vector of coefficients for further predictions; NULL if <code>generate_coefficients</code> is FALSE.

**Note**

The vector of prediction coefficients in `pred_coeffs` can be used as follows. Let  $c$  denote the coefficients and let  $r$  denote a vector with element  $i$  containing the correlation between the output at a given new point and the output at training point  $i$ . Then the prediction for the output at the new point is the dot product of  $c$  and  $r$ .

[RMSE](#) computes the root mean squared error of the predictions. [PlotPredictions](#) and [PlotResiduals](#) plot the predictions or their residuals; [PlotStdResiduals](#) and [PlotQQ](#) plot the standardized residuals.

**Examples**

```
borehole_pred <- Predict(
  GaSP_model = borehole_fit,
  x_pred = borehole$x_pred,
  generate_coefficients = TRUE
)
```

---

 RMSE

*Calculate the root mean squared error (RMSE) of prediction*

---

**Description**

Calculate the root mean squared error (RMSE) of prediction

**Usage**

```
RMSE(y_pred, y_true, normalized = FALSE)
```

**Arguments**

<code>y_pred</code>	A vector of predicted output values.
<code>y_true</code>	A vector of true output values.
<code>normalized</code>	An optional boolean: if TRUE, the RMSE is normalized by dividing it by the standard deviation of <code>y_true</code> .

**Value**

The RMSE or normalized RMSE.

**Examples**

```
RMSE(borehole_pred$y_pred$Pred, borehole$y_true)
```

```
RMSE(borehole_cv$Pred, y)
```

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 Visualize

*Visualize a GaSPModel object.*


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**Description**

Carry out a functional analysis of variance (ANOVA) of a `GaSPModel` object and generate plotting coordinates for its estimated main and 2-input joint effects.

**Usage**

```
Visualize(GaSP_model, x_description, main_percent = 0, interaction_percent = 0)
```

**Arguments**

<code>GaSP_model</code>	Object of class <code>GaSPModel</code> .
<code>x_description</code>	A data frame describing the input variables. See <a href="#">DescribeX</a> .
<code>main_percent</code>	An optional minimum percentage of variation explained by an input's main effect to return the effect's plotting coordinates; the default of zero gives plotting coordinates for all inputs.
<code>interaction_percent</code>	An optional minimum percentage of variation explained by the interaction effect of a pair of inputs to return the plotting coordinates for their joint effect (main effects plus interaction effect); the default of zero gives plotting coordinates for all pairs of inputs.

**Details**

If there are many inputs, to avoid excessive plotting of many trivial joint effects set `interaction_percent = 1` say.

**Value**

A list with the following elements:

<code>anova_percent</code>	A data frame containing the ANOVA percentages for all main effects and 2-input interaction effects.
<code>main_effect</code>	A data frame with plotting coordinates for the estimated main effects.
<code>joint_effect</code>	A data frame with plotting coordinates for the estimated 2-input joint effects.
<code>total_percent</code>	Total percentage of the prediction variation accounted for by all main effects and 2-input interaction effects.
<code>average</code>	Overall average of the prediction function, averaged with respect to all inputs.
<code>SE_average</code>	Standard error of the overall average.

**References**

Schonlau, M. and Welch, W.J. (2006), "Screening the Input Variables to a Computer Model Via Analysis of Variance and Visualization", in *Screening: Methods for Experimentation in Industry, Drug Discovery, and Genetics*, Dean. A. and Lewis, S., eds., pp. 308-327, Springer, New York, doi:10.1007/0-387-28014-6\_14.

**Examples**

```
borehole_vis <- Visualize(borehole_fit, borehole_x_desc)
```

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