

Package ‘janusplot’

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Type Package

Title Asymmetric Smoothed-Association Matrices via GAM Fits

Version 0.1.0

Description Render a pairwise, asymmetric smoothed-association matrix of continuous variables. Each cell shows the fitted spline from an 'mgcv' generalised additive model, with the upper triangle displaying 'gam(x_j ~ s(x_i))' and the lower triangle 'gam(x_i ~ s(x_j))'. Unlike Pearson's correlation matrix, the visualisation is intentionally asymmetric, revealing heteroscedasticity, leverage, and directional non-linearity that a single scalar correlation hides. An asymmetry index and a 24-category shape taxonomy quantify the directional difference and qualitative form of each fitted smooth.

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URL <https://github.com/max578/janusplot>,
<https://max578.github.io/janusplot/>

BugReports <https://github.com/max578/janusplot/issues>

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Contents

janusplot	2
janusplot_data	8
janusplot_shape_cutoffs	10
janusplot_shape_hierarchy	11
janusplot_shape_metrics	13
janusplot_shape_sensitivity	14
janusplot_shape_sensitivity_plot	16
janusplot_shape_sensitivity_shapes	17
janusplot_shape_sensitivity_summary	18
shape_sensitivity_demo	18
Index	20

janusplot	<i>Asymmetric smoothed-association matrix</i>
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Description

[Experimental]

Render a pairwise, asymmetric matrix of smoothed associations between numeric variables. Each cell $[i, j]$ where $i \neq j$ shows the fitted spline from `mgcv::gam()`:

- Upper triangle ($i < j$): $\text{gam}(x_j \sim s(x_i) + \langle \text{adjust} \rangle)$.
- Lower triangle ($i > j$): $\text{gam}(x_i \sim s(x_j) + \langle \text{adjust} \rangle)$.
- Diagonal: blank panel when labels live on the border (default), or a variable-name label when `labels = "diagonal"`.

The two triangles intentionally differ — the asymmetry reveals heteroscedasticity, leverage, and directional non-linearity that a single scalar correlation hides.

Usage

```

janusplot(
  data,
  vars = NULL,
  adjust = NULL,
  method = "REML",
  k = -1L,
  bs = "tp",
  order = c("original", "hclust", "alphabetical"),
  show_data = TRUE,
  show_ci = TRUE,
  display = c("fit", "d1", "d2"),
  derivative_ci = c("none", "pointwise", "simultaneous"),
  derivative_ci_nsim = 1000L,
  n_grid = NULL,
  colour_by = c("pearson", "spearman", "kendall", "edf", "deviance_gap", "none"),
  fill_by = NULL,
  palette = NULL,
  annotations = c("edf", "A"),
  shape_cutoffs = janusplot_shape_cutoffs(),
  show_shape_legend = TRUE,
  glyph_style = c("ascii", "unicode"),
  labels = c("border", "diagonal", "none"),
  diagonal = c("auto", "blank", "name", "density"),
  label_srt = 45,
  label_cex = 1,
  signif_glyph = TRUE,
  show_asymmetry = NULL,
  na_action = c("pairwise", "complete"),
  parallel = FALSE,
  with_data = FALSE,
  text_scale_diag = 1,
  text_scale_off_diag = 1,
  show_glossary = TRUE,
  glossary_scale = 1,
  ...
)

```

Arguments

<code>data</code>	A data frame with numeric columns to include.
<code>vars</code>	Character vector of column names to use. <code>NULL</code> (default) uses all numeric columns in data. Non-numeric columns trigger an error listing offenders.
<code>adjust</code>	A one-sided formula RHS giving additional covariates and/or random effects to include in every pairwise GAM. For example, <code>adjust = ~ s(age) + s(site, bs = "re")</code> fits <code>gam(y ~ s(x) + s(age) + s(site, bs = "re"))</code> for each pair. Default <code>NULL</code> fits unadjusted pairwise smooths.

method	Smoothing-parameter estimation method passed to <code>mgcv::gam()</code> . Default "REML" per mgcv recommendation.
k	Integer, or named list mapping variable names to integers. Basis dimension for <code>s()</code> . Default -1L (mgcv's automatic choice).
bs	Basis type for <code>s()</code> . Default "tp" (thin plate).
order	One of "original" (default), "hclust" (reorder by hierarchical clustering of Pearson correlations), or "alphabetical".
show_data	Logical. If TRUE (default), overlay raw data points (low alpha) behind each spline. Only applies when <code>display = "fit"</code> ; derivative panels never overlay raw data.
show_ci	Logical. If TRUE (default), overlay the 95% confidence envelope from <code>predict(gam, se.fit = TRUE)</code> on the fit panel (i.e. when <code>display = "fit"</code>). CI rendering on derivative panels is controlled separately by <code>derivative_ci</code> .
display	<p>One of "fit" (default), "d1", or "d2". Selects which single quantity is rendered in every off-diagonal cell of the matrix.</p> <ul style="list-style-type: none"> • "fit" — the fitted smooth $\hat{f}(x)$; default, behaviour identical to the pre-derivative release. • "d1" — the first derivative $\hat{f}'(x)$ of the fitted smooth. Zero crossings localise turning points of \hat{f}. • "d2" — the second derivative $\hat{f}''(x)$. Zero crossings localise inflection points of \hat{f}. <p>A single matrix shows a single quantity by design: stacked multi-panel cells crowd the matrix at any realistic variable count. To compare fit against derivative, render two or three <code>janusplot()</code> calls side-by-side; each call keeps its own <code>with_data = TRUE</code> summary table tagged with the <code>display</code> column.</p> <p>Orders $k \geq 3$ are not exposed — higher-order derivatives of penalised regression splines amplify noise and rarely carry usable signal at realistic sample sizes. See <code>vignette("janusplot")</code> for the theoretical justification and applied use-cases.</p>
derivative_ci	<p>One of "none" (default), "pointwise", or "simultaneous". Controls whether — and how — a 95% confidence ribbon is drawn underneath the derivative curve when <code>display %in% c("d1", "d2")</code>. Ignored when <code>display = "fit"</code>.</p> <ul style="list-style-type: none"> • "none" — no ribbon. The curve and the zero reference line are all you see. Default, because pointwise ribbons overshoot nominal coverage as a joint region and can invite over-reading of local features. • "pointwise" — 95% pointwise ribbon from $\sqrt{\text{diag}(DV_p D^T)}$ (Wood 2017 §7.2.4). Valid marginally; not a simultaneous statement. • "simultaneous" — 95% simultaneous band via the Monte Carlo construction of Ruppert, Wand & Carroll (2003) popularised for GAMs by Simpson (2018, <i>Frontiers Ecol. Evol.</i> 6:149): draw B samples $\tilde{\beta} \sim \mathcal{N}(\hat{\beta}, V_p)$, compute $\max_x D_i(\tilde{\beta} - \hat{\beta}) /se_i$, and use the $(1 - \alpha)$ quantile as a critical multiplier on the pointwise SE. Valid for feature localisation ("where is $\hat{f}'(x)$ significantly non-zero").
derivative_ci_nsim	<p>Integer. Number of Monte Carlo samples used when <code>derivative_ci = "simultaneous"</code>. Default 1000L — a compromise between coverage accuracy (Simpson 2018 uses</p>

	10000) and CPU budget across every pair in a medium-sized matrix. Ignored for any other derivative_ci.
n_grid	Integer or NULL. Number of equally-spaced points used to evaluate each fitted smooth (and its derivatives). Default NULL resolves to 100 when display = "fit" and 200 otherwise, because finite-difference second derivatives visibly degrade below ~ 150 points on moderate-k smooths. Supplying n_grid directly overrides both defaults. Larger grids shift the numerical shape-metric values (M , C , turning / inflection counts) slightly because they are computed on this same grid. Shapes and asymmetry are the primary reading; M , C and the counts are secondary diagnostics and the grid-induced drift is tolerable.
colour_by	One of "pearson" (default), "spearman", "kendall", "edf", "deviance_gap", or "none". Encodes the per-cell fill colour by the chosen scalar. Correlation choices use a diverging palette with limits $c(-1, 1)$ and a shared corr colour-bar title; "edf" and "deviance_gap" use a sequential palette labelled by the metric.
fill_by	Deprecated alias for colour_by. If supplied emits a single soft deprecation warning and is forwarded to colour_by.
palette	Character. Colour palette for the cell fill scale. Defaults to "RdBu" when colour_by is a correlation and "viridis" otherwise. Sequential choices: "viridis", "magma", "inferno", "plasma", "cividis", "mako", "rocket", "turbo" (not CB-safe), "YlOrRd", "YlGnBu", "Blues", "Greens". Diverging choices: "RdYlBu", "RdBu", "PuOr", "Spectral" (not CB-safe). Passing a sequential palette while colour_by is a correlation silently upgrades to the default diverging palette.
annotations	Character vector, a subset of $c("edf", "A", "shape", "code")$. Controls which corner annotations appear on each off-diagonal cell: <ul style="list-style-type: none"> • "code" — 2-letter ASCII shape code, top-left corner. • "A" and "edf" — asymmetry index and effective degrees of freedom, stacked bottom-left. • "shape" — shape glyph (Unicode or ASCII per glyph_style), bottom-right corner. <p>Default $c("edf", "A")$. "code" and "shape" occupy distinct corners so both can be requested together. See janusplot_shape_hierarchy() for the full code list.</p>
shape_cutoffs	Named list of classification thresholds used to map the continuous shape indices into discrete shape_category labels; see janusplot_shape_cutoffs() .
show_shape_legend	Logical. If TRUE (default), attach a standing shape-types legend plate below the matrix that illustrates every category in the taxonomy as a canonical thumbnail spline. Independent of annotations.
glyph_style	One of "ascii" (default) or "unicode". Controls how cell shape glyphs render when "shape" is included in annotations. Default is "ascii" for maximum portability across typesetting pipelines; switch to "unicode" only when the target font is known to cover the curve glyph set.
labels	One of "border" (default), "diagonal", or "none". Controls where variable names are rendered:

	<ul style="list-style-type: none"> • "border" — names along the top (rotated per <code>label_srt</code>) and left margins of the matrix; diagonal cells are left blank. Mirrors <code>corrplot</code>'s <code>tl.pos = "lt"</code> convention. • "diagonal" — names centred on the diagonal cells (the pre-0.1 layout). • "none" — labels suppressed entirely; diagonal cells blank.
<code>diagonal</code>	<p>One of "auto" (default), "blank", "name", or "density". Controls what is rendered in the diagonal cells of the matrix.</p> <ul style="list-style-type: none"> • "auto" — preserves the historical behaviour: variable name when <code>labels = "diagonal"</code>, blank otherwise. • "blank" — empty bordered panel (uniform grid reading). • "name" — variable name centred in the cell, bold. • "density" — kernel density of the variable filled in translucent grey, with a rug of raw values along the bottom edge. Mirrors the <code>GGally::ggpairs</code> convention; surfaces tail weight, bimodality, and support clipping that the pairwise smooths alone cannot reveal. Variable names should come from the border (<code>labels = "border"</code>, the default) when this mode is on.
<code>label_srt</code>	Numeric. Rotation (degrees) of top labels when <code>labels = "border"</code> . Default 45; set to 0 for horizontal or 90 for vertical. Ignored when <code>labels != "border"</code> .
<code>label_cex</code>	Positive numeric multiplier on the border-label font size. Default 1. Ignored when <code>labels = "none"</code> .
<code>signif_glyph</code>	Logical. If TRUE (default), annotate cells with <code>· / * / **</code> reflecting the smooth's F-test p-value.
<code>show_asymmetry</code>	Deprecated. Use annotations instead (" <code>A</code> " <code>%in%</code> annotations). When supplied, a soft deprecation warning fires and the argument is merged into annotations.
<code>na_action</code>	One of "pairwise" (default; per-cell complete observations) or "complete" (listwise; all cells use the same rows).
<code>parallel</code>	Logical. If TRUE, use <code>future.apply::future_mapply()</code> to fit pairs in parallel. Requires the <code>future.apply</code> package and a user-configured <code>future::plan()</code> . Default FALSE.
<code>with_data</code>	Logical. If TRUE, return a two-element list <code>list(plot, data)</code> where <code>data</code> is a flat per-cell summary (one row per off-diagonal cell) of everything the plot displays. The <code>data</code> element is always a plain <code>data.frame</code> (base R — no <code>data.table</code> dependency). Default FALSE — in which case only the <code>ggplot</code> is returned.
<code>text_scale_diag</code>	Positive numeric multiplier applied to the diagonal variable-name labels. Default 1. Diagonal labels additionally auto-shrink for long variable names (<code>nchar(var) > 10</code>) so they fit the cell regardless of this value.
<code>text_scale_off_diag</code>	Positive numeric multiplier applied to all off-diagonal annotations (n / EDF read-outs, significance glyphs, asymmetry-index labels). Default 1. Use <code>< 1</code> when cells are small and the annotations crowd the fit line; use <code>> 1</code> for presentation plots.
<code>show_glossary</code>	Logical. If TRUE (default), attach a multi-line caption below the matrix describing the on-plot abbreviations (n, EDF, A, fill encoding, significance glyphs). Only keys actually displayed are listed.

`glossary_scale` Positive numeric multiplier on the glossary caption font size. Default 1.
 ... Additional arguments passed to `mgcv::gam()`.

Value

If `with_data = FALSE` (default), a `ggplot2::ggplot` object (via `patchwork::wrap_plots()`) carrying a top-of-matrix title that names the displayed quantity ("Direct fit", "First derivative f' ", or "Second derivative f'' "). If `with_data = TRUE`, a list with two elements: `plot` (the `ggplot`) and `data` (a tidy table with columns `var_x`, `var_y`, `position`, `n_used`, `edf`, `pvalue`, `signif`, `dev_exp`, `asymmetry_index`, `cor_pearson`, `cor_spearman`, `cor_kendall`, `tie_ratio`, `monotonicity_index`, `convexity_index`, `n_turning_points`, `n_inflections`, `flat_range_ratio`, `shape_category`, `colour_value`, `display`, one row per off-diagonal cell). The `display` column tags which quantity the call rendered, so separate calls for `fit` / `d1` / `d2` yield comparable, stackable tables. Derivative *curves* themselves (grid of x , fitted $\hat{f}^{(k)}$, SE) live on `janusplot_data()` — see there.

See Also

[janusplot_data\(\)](#) for the raw per-cell fits + metrics.

Other smooth-associations: [janusplot_data\(\)](#)

Examples

```
# Minimal runnable example – 3 variables, 6 asymmetric pairwise GAM fits.
janusplot(mtcars[, c("mpg", "hp", "wt")])
```

```
# Heteroscedastic DGP: Pearson r is ~ 0.9 but the inverse fit is
# clearly non-linear, yielding asymmetry index > 0.5.
set.seed(2026L)
n <- 200L
x1 <- stats::runif(n, 0, 10)
x2 <- x1 + stats::rnorm(n, sd = 0.2 * x1)
janusplot(data.frame(x1 = x1, x2 = x2, x3 = stats::rnorm(n)))
```

```
# A single matrix renders a single quantity. To compare the fit
# against its derivatives, render three calls and place them
# side-by-side; each call's title makes the quantity explicit.
set.seed(2026L)
xs <- stats::runif(300L, -3, 3)
df <- data.frame(
  x = xs,
  y1 = sin(xs) + stats::rnorm(300L, sd = 0.3),
  y2 = xs^2 + stats::rnorm(300L, sd = 0.6)
)
janusplot(df, display = "fit")
janusplot(df, display = "d1")
janusplot(df, display = "d2")
```

```
# Simultaneous CI bands on a derivative panel, per Simpson (2018).
janusplot(df, display = "d1", derivative_ci = "simultaneous")
```

janusplot_data	<i>Raw GAM fits and per-cell metrics for a smoothed-association matrix</i>
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Description

[Experimental]

Companion to `janusplot()` returning the raw list of GAM fits plus per-cell metrics (EDF, F-test p-value, deviance explained, asymmetry index, pairwise correlations, shape descriptors) without constructing the `ggplot`. Useful for custom rendering or downstream analysis.

Usage

```
janusplot_data(
  data,
  vars = NULL,
  adjust = NULL,
  method = "REML",
  k = -1L,
  bs = "tp",
  na_action = c("pairwise", "complete"),
  parallel = FALSE,
  keep_fits = FALSE,
  derivatives = integer(),
  derivative_ci = c("pointwise", "none", "simultaneous"),
  derivative_ci_nsim = 1000L,
  n_grid = NULL,
  shape_cutoffs = janusplot_shape_cutoffs(),
  ...
)
```

Arguments

<code>data</code>	A data frame with numeric columns to include.
<code>vars</code>	Character vector of column names to use. <code>NULL</code> (default) uses all numeric columns in <code>data</code> . Non-numeric columns trigger an error listing offenders.
<code>adjust</code>	A one-sided formula RHS giving additional covariates and/or random effects to include in every pairwise GAM. For example, <code>adjust = ~ s(age) + s(site, bs = "re")</code> fits <code>gam(y ~ s(x) + s(age) + s(site, bs = "re"))</code> for each pair. Default <code>NULL</code> fits unadjusted pairwise smooths.
<code>method</code>	Smoothing-parameter estimation method passed to <code>mgcv::gam()</code> . Default "REML" per <code>mgcv</code> recommendation.
<code>k</code>	Integer, or named list mapping variable names to integers. Basis dimension for <code>s()</code> . Default <code>-1L</code> (<code>mgcv</code> 's automatic choice).

bs	Basis type for <code>s()</code> . Default "tp" (thin plate).
na_action	One of "pairwise" (default; per-cell complete observations) or "complete" (listwise; all cells use the same rows).
parallel	Logical. If TRUE, use <code>future.apply::future_mapply()</code> to fit pairs in parallel. Requires the <code>future.apply</code> package and a user-configured <code>future::plan()</code> . Default FALSE.
keep_fits	Logical. If TRUE, retain full <code>mgcv::gam()</code> model objects in the return (large memory footprint for <code>k</code> above ~15). Default FALSE — retains summary metrics and prediction grids only.
derivatives	Integer vector of derivative orders to compute on every pair (subset of 1:2). Default <code>integer()</code> — no derivatives. Unlike <code>janusplot()</code> , the data companion can return multiple orders from a single call for programmatic analysis; pass <code>c(1L, 2L)</code> to surface both.
derivative_ci	One of "none" (default), "pointwise", or "simultaneous". Controls whether — and how — a 95% confidence ribbon is drawn underneath the derivative curve when <code>display %in% c("d1", "d2")</code> . Ignored when <code>display = "fit"</code> . <ul style="list-style-type: none"> "none" — no ribbon. The curve and the zero reference line are all you see. Default, because pointwise ribbons overshoot nominal coverage as a joint region and can invite over-reading of local features. "pointwise" — 95% pointwise ribbon from $\sqrt{\text{diag}(DV_p D^T)}$ (Wood 2017 §7.2.4). Valid marginally; not a simultaneous statement. "simultaneous" — 95% simultaneous band via the Monte Carlo construction of Ruppert, Wand & Carroll (2003) popularised for GAMs by Simpson (2018, <i>Frontiers Ecol. Evol.</i> 6:149): draw B samples $\tilde{\beta} \sim \mathcal{N}(\hat{\beta}, V_p)$, compute $\max_x D_i(\tilde{\beta} - \hat{\beta}) /se_i$, and use the $(1 - \alpha)$ quantile as a critical multiplier on the pointwise SE. Valid for feature localisation ("where is $\hat{f}'(x)$ significantly non-zero").
derivative_ci_nsim	Integer. Number of Monte Carlo samples used when <code>derivative_ci = "simultaneous"</code> . Default 1000L — a compromise between coverage accuracy (Simpson 2018 uses 10000) and CPU budget across every pair in a medium-sized matrix. Ignored for any other <code>derivative_ci</code> .
n_grid	Integer or NULL. Number of equally-spaced points used to evaluate each fitted smooth (and its derivatives). Default NULL resolves to 100 when <code>display = "fit"</code> and 200 otherwise, because finite-difference second derivatives visibly degrade below ~ 150 points on moderate- <code>k</code> smooths. Supplying <code>n_grid</code> directly overrides both defaults. Larger grids shift the numerical shape-metric values (M , C , turning / inflection counts) slightly because they are computed on this same grid. Shapes and asymmetry are the primary reading; M , C and the counts are secondary diagnostics and the grid-induced drift is tolerable.
shape_cutoffs	Named list of classification thresholds used to map the continuous shape indices (<code>monotonicity_index</code> , <code>convexity_index</code>) into discrete <code>shape_category</code> labels. Defaults from <code>janusplot_shape_cutoffs()</code> .
...	Additional arguments passed to <code>mgcv::gam()</code> .

Value

A list with components:

`vars` Character vector of variables used, in plotted order.

`pairs` List of per-pair results. Each element has `i`, `j`, `var_i`, `var_j`, `fit_yx`, `fit_xy` (NULL if `keep_fits = FALSE`), `pred_yx`, `pred_xy` (data frames with `x`, `fit`, `se`, `lo`, `hi`), `edf_yx`, `edf_xy`, `pvalue_yx`, `pvalue_xy`, `dev_exp_yx`, `dev_exp_xy`, `n_used`, `asymmetry_index`, plus Pearson / Spearman / Kendall correlations (`cor_pearson`, `cor_spearman`, `cor_kendall`), the maximum tie ratio across `x` and `y` (`tie_ratio`), and per-direction shape descriptors (`monotonicity_index_yx`, `convexity_index_yx`, `monotonicity_index_xy`, `convexity_index_xy`, `n_turning_yx`, `n_inflect_yx`, `n_turning_xy`, `n_inflect_xy`, `shape_yx`, `shape_xy`). When `derivatives` is non-empty, each pair additionally carries `deriv_yx` and `deriv_xy`, each a named list keyed by order ("1", "2") whose entries are data frames with columns `x`, `fit`, `se`, `lo`, `hi`, `ci_type` matching the schema of `pred_yx` / `pred_xy`. The `ci_type` column records whether the `lo` / `hi` columns are "pointwise" (default), "simultaneous" (Ruppert–Wand–Carroll / Simpson 2018 critical-multiplier bands), or "none". When `derivative_ci = "simultaneous"`, each derivative frame also carries a "crit_multiplier" attribute giving the MC-derived critical multiplier used. See [janusplot_shape_metrics\(\)](#) for the definition of the monotonicity and convexity indices.

`call` Match call.

See Also

[janusplot\(\)](#) for the ggplot front-end, [janusplot_shape_metrics\(\)](#) for the shape-metric primitives.

Other smooth-associations: [janusplot\(\)](#)

Examples

```
# Per-pair fits + metrics on a small mtcars slice
out <- janusplot_data(mtcars[, c("mpg", "hp", "wt")])
out$pairs[[1L]]$asymmetry_index
out$pairs[[1L]]$cor_spearman
out$pairs[[1L]]$shape_yx
```

janusplot_shape_cutoffs

Default cutoff thresholds for shape_category classification

Description**[Experimental]**

Returns the named list of thresholds used to map the continuous monotonicity (M) and convexity (C) indices (plus inflection counts) into a discrete `shape_category`. Expose so callers can override individual thresholds or pass a fully custom list to [janusplot\(\)](#) / [janusplot_shape_metrics\(\)](#).

Usage

```
janusplot_shape_cutoffs(...)
```

Arguments

... Optional named overrides to merge into the defaults.

Value

A named list with numeric thresholds:

mono_strong |M| threshold for a strictly monotone smooth (default 0.9).
 mono_mod |M| threshold for a curved-but-monotone smooth (default 0.5).
 mono_nonmono |M| below this is considered non-monotone (default 0.3).
 mono_s |M| threshold for labelling an S-shape (default 0.5).
 curv_low |C| below this is considered near-linear curvature (default 0.2).
 curv_mod |C| threshold for a clearly curved monotone (default 0.5).
 curv_strong |C| threshold for a U-shape / inverted-U shape (default 0.5).
 flat range(fit) / sd(y) below this is called flat (default 0.05).

See Also

Other shape-metrics: [janusplot_shape_hierarchy\(\)](#), [janusplot_shape_metrics\(\)](#)

Examples

```
janusplot_shape_cutoffs()
janusplot_shape_cutoffs(curv_mod = 0.6, flat = 0.02)
```

```
janusplot_shape_hierarchy
```

Shape-category taxonomy table

Description**[Experimental]**

Return the full janusplot shape taxonomy as a data frame with four hierarchy columns plus presentation fields. The taxonomy is the single source of truth consumed by the classifier, the cell renderer, the legend plate, and the `janusplot_data()` output.

Hierarchy columns (finest → coarsest):

category 24-way fine label (linear_up, skewed_peak, bimodal, ...). Computed per cell by [janusplot\(\)](#).

code Unique two-letter ASCII shorthand (safe on any font or typesetting pipeline) — e.g. lu for linear_up.

archetype Seven-family grouping: monotone_linear, monotone_curved, unimodal, wave, multimodal, chaotic, degenerate.

monotonic Three-way coarse classification: monotone / non_monotone / degenerate.

linear Binary: linear / non_linear / degenerate.

The broader tiers (linear/non-linear, monotone/non-monotone) are textbook calculus; the archetype layer maps cleanly to shape-constrained regression vocabulary (Pya & Wood 2015; Meyer 2008) and to dose-response shape categories (Calabrese 2008; Calabrese & Baldwin 2001). The (T, I) dispatch underlying each fine category is a coarsened Morse-theoretic critical-point classification (Milnor 1963).

Usage

```
janusplot_shape_hierarchy()
```

Value

A data frame with 24 rows and columns category, code, archetype, monotonic, linear, glyph, ascii, label, gloss.

References

Calabrese, E. J. (2008). Hormesis: why it is important to toxicology and toxicologists. *Environmental Toxicology and Chemistry*, **27**(7), 1451–1474.

Meyer, M. C. (2008). Inference using shape-restricted regression splines. *Annals of Applied Statistics*, **2**(3), 1013–1033.

Milnor, J. (1963). *Morse Theory*. Princeton University Press.

Pya, N., & Wood, S. N. (2015). Shape constrained additive models. *Statistics and Computing*, **25**(3), 543–559.

See Also

Other shape-metrics: [janusplot_shape_cutoffs\(\)](#), [janusplot_shape_metrics\(\)](#)

Examples

```
tax <- janusplot_shape_hierarchy()
head(tax[, c("category", "code", "archetype", "monotonic", "linear")])
# Count how many categories live in each archetype
table(tax$archetype)
```

 janusplot_shape_metrics

Shape metrics for a fitted univariate smooth

Description

[Experimental]

Compute the continuous monotonicity and convexity indices, inflection and turning-point counts, and rule-based shape category for a fitted univariate smooth. Works on either a per-pair fit object returned from the janusplot internal machinery or a freshly fitted `mgcv::gam()` with a single `s()` term.

Both indices are bounded in $[-1, 1]$ and weighted by the empirical density of the predictor:

- `monotonicity_index` (paper symbol M). Let f be the fitted smooth evaluated on a dense grid of `n_grid` equally-spaced points across the predictor range, f' its numerical first derivative, and w the empirical density of the predictor on the same grid with $\text{sum}(w) = 1$. Then $\text{monotonicity_index} = \text{sum}(w * f') / \text{sum}(w * |f'|)$ in $[-1, 1]$. +1 is strictly increasing, -1 strictly decreasing, 0 non-monotone.
- `convexity_index` (paper symbol C). With f'' the numerical second derivative on the same grid, $\text{convexity_index} = \text{sum}(w * f'') / \text{sum}(w * |f''|)$ in $[-1, 1]$. +1 is globally convex (bowl-up), -1 globally concave (bowl-down), 0 inflection-dominated (S-curve, sine, flat).

Both indices are scale-invariant (replacing $y \rightarrow a*y + b$ leaves them unchanged) and density-weighted so they describe the smooth *where the data actually live*, not extrapolated tails.

Usage

```
janusplot_shape_metrics(
  fit,
  x_name = NULL,
  newdata = NULL,
  n_grid = 200L,
  cutoffs = janusplot_shape_cutoffs()
)
```

Arguments

<code>fit</code>	Either a list returned by a janusplot pair-fit helper (must contain <code>pred</code> and <code>raw</code>), or a fitted <code>mgcv::gam()</code> with a single <code>s(x)</code> term.
<code>x_name</code>	Character. Column name of the predictor when <code>fit</code> is a <code>mgcv::gam()</code> object. Ignored for pair-fit lists.
<code>newdata</code>	Optional data frame supplying the raw predictor values used for density weighting when <code>fit</code> is a <code>mgcv::gam()</code> object. If <code>NULL</code> , the model frame is used.
<code>n_grid</code>	Integer. Prediction grid length when <code>fit</code> is a <code>mgcv::gam()</code> object. Default 200L.
<code>cutoffs</code>	Named list of classification thresholds; see <code>janusplot_shape_cutoffs()</code> . Default uses package defaults.

Value

A named list with components:

`monotonicity_index` M in $[-1, 1]$. See Description.

`convexity_index` C in $[-1, 1]$. See Description.

`n_turning_points` Integer count of lobe-mass-weighted sign changes of f' . Equals the number of interior extrema.

`n_inflections` Integer count of lobe-mass-weighted sign changes of f'' .

`flat_range_ratio` $\text{range}(f) / \text{sd}(y)$ — small values indicate a degenerate flat smooth.

`shape_category` One of 24 labels from `janusplot_shape_hierarchy()` dispatched on `(n_turning_points, n_inflect)` with `(monotonicity_index, convexity_index)` disambiguation for the monotone case.

See Also

[janusplot_shape_cutoffs\(\)](#), [janusplot\(\)](#), [janusplot_data\(\)](#).

Other shape-metrics: [janusplot_shape_cutoffs\(\)](#), [janusplot_shape_hierarchy\(\)](#)

Examples

```
# On a fitted gam
set.seed(2026L)
n <- 200L
x <- stats::runif(n, 0, 10)
y <- log1p(x) + stats::rnorm(n, sd = 0.3)
d <- data.frame(x = x, y = y)
fit <- mgcv::gam(y ~ s(x), data = d, method = "REML")
janusplot_shape_metrics(fit, x_name = "x", newdata = d)
```

janusplot_shape_sensitivity

Shape-recognition sensitivity study

Description**[Experimental]**

Run a full-factorial sensitivity sweep for the `janusplot` 24-category shape classifier. For each combination of ground-truth shape, sample size n , noise level σ , and replicate, the sweep:

1. Generates n points from the noiseless canonical curve on $[0, 1] + \text{Gaussian noise with SD} = \sigma$ (fraction of the y -range, so signal-to-noise is comparable across shapes).
2. Fits `mgcv::gam(y ~ s(x), method = "REML")`.
3. Runs `janusplot_shape_metrics()` to classify the fitted smooth.
4. Records correctness at both the fine (24-category) and archetype (7-family) levels.

The function is the package-native implementation of `simulation/scripts/scenario_4_shape_recognition.R`. A small precomputed dataset is shipped as [shape_sensitivity_demo](#) for downstream examples without requiring users to re-run the sweep.

Usage

```
janusplot_shape_sensitivity(
  shapes = NULL,
  n_grid = c(50L, 100L, 200L, 500L),
  sigma_grid = c(0.02, 0.05, 0.1, 0.2, 0.4),
  n_rep = 200L,
  cutoffs = janusplot_shape_cutoffs(),
  parallel = FALSE,
  seed = 2026L,
  verbose = interactive()
)
```

Arguments

shapes	Character vector of ground-truth names from janusplot_shape_sensitivity_shapes() . Default NULL → all 14.
n_grid	Integer vector of sample sizes. Default <code>c(50L, 100L, 200L, 500L)</code> .
sigma_grid	Numeric vector of noise levels (fraction of the y-range). Default <code>c(0.02, 0.05, 0.10, 0.20, 0.40)</code> .
n_rep	Integer. Replicates per cell. Default 200L.
cutoffs	Named list of classification thresholds; see janusplot_shape_cutoffs() .
parallel	Logical. If TRUE and <code>future.apply</code> is installed, dispatch replicates in parallel. The caller is responsible for configuring <code>future::plan()</code> (e.g. <code>future::plan(future::multisession)</code>).
seed	Integer. Base seed — each fit uses <code>seed + row_index</code> so results are reproducible and cell-permutation-invariant.
verbose	Logical. Print progress messages to the console. Default is <code>interactive()</code> .

Value

A data frame with one row per fit. Columns:

truth	Ground-truth shape name.
n	Sample size for this fit.
sigma	Noise level for this fit.
seed	RNG seed used.
predicted	Classifier output at the fine (24-category) level.
correct	Logical — does <code>predicted == truth</code> ?
archetype_truth	Expected archetype for truth.
archetype_pred	Archetype of predicted.
archetype_correct	Logical — archetype-level correctness.
monotonicity_index	Monotonicity index M (see janusplot_shape_metrics()).
convexity_index	Convexity index C (see janusplot_shape_metrics()).
n_turn, n_inflect	Recovered turning-point and inflection counts.
error	"gam_fit_failed" when <code>mgcv::gam()</code> errored; NA otherwise.

See Also

[janusplot_shape_sensitivity_summary\(\)](#), [janusplot_shape_sensitivity_plot\(\)](#), [janusplot_shape_sensitivity_shape_sensitivity_demo](#).

Other shape-sensitivity: [janusplot_shape_sensitivity_plot\(\)](#), [janusplot_shape_sensitivity_shapes\(\)](#), [janusplot_shape_sensitivity_summary\(\)](#)

Examples

```
# Tiny-run smoke test (< 2 seconds): 3 shapes x 2 n x 2 sigma x 5 reps.
res <- janusplot_shape_sensitivity(
  shapes      = c("linear_up", "u_shape", "wave"),
  n_grid      = c(100L, 200L),
  sigma_grid  = c(0.05, 0.20),
  n_rep       = 5L,
  verbose     = FALSE
)
head(res)
janusplot_shape_sensitivity_summary(res, level = "archetype")
```

janusplot_shape_sensitivity_plot

Visualise a shape-sensitivity sweep

Description**[Experimental]**

Produce one of four diagnostic plots from the raw data frame returned by [janusplot_shape_sensitivity\(\)](#):

"confusion_fine" 24 x (lshapes) confusion matrix at the fine category level — rows = ground truth, columns = predicted, cells coloured by P(pred | truth).

"confusion_archetype" 7 x 7 confusion matrix at the archetype level.

"accuracy_grid" per-shape heatmap of archetype-level accuracy across the (n, sigma) design.

"recovery_curves" accuracy as a function of sigma, one line per sample size, faceted by shape.

Usage

```
janusplot_shape_sensitivity_plot(
  results,
  type = c("confusion_fine", "confusion_archetype", "accuracy_grid", "recovery_curves")
)
```

Arguments

results Data frame from [janusplot_shape_sensitivity\(\)](#) or the precomputed [shape_sensitivity_demo](#).

type One of "confusion_fine", "confusion_archetype", "accuracy_grid", or "recovery_curves".

Value

A [ggplot2::ggplot](#) object.

See Also

Other shape-sensitivity: [janusplot_shape_sensitivity\(\)](#), [janusplot_shape_sensitivity_shapes\(\)](#), [janusplot_shape_sensitivity_summary\(\)](#)

Examples

```
data("shape_sensitivity_demo", package = "janusplot")
janusplot_shape_sensitivity_plot(shape_sensitivity_demo,
                                "recovery_curves")
```

`janusplot_shape_sensitivity_shapes`

Canonical ground-truth shapes for the sensitivity study

Description**[Experimental]**

Return the names of every canonical ground-truth shape that [janusplot_shape_sensitivity\(\)](#) can simulate from. Fourteen shapes spanning five archetypes (`monotone_linear`, `monotone_curved`, `unimodal`, `wave`, `multimodal`). The chaotic and degenerate archetypes are out of scope (no realistic deterministic generator).

Usage

```
janusplot_shape_sensitivity_shapes()
```

Value

Character vector of length 14 — the generator names.

See Also

[janusplot_shape_sensitivity\(\)](#), [janusplot_shape_hierarchy\(\)](#).

Other shape-sensitivity: [janusplot_shape_sensitivity\(\)](#), [janusplot_shape_sensitivity_plot\(\)](#), [janusplot_shape_sensitivity_summary\(\)](#)

Examples

```
janusplot_shape_sensitivity_shapes()
```

janusplot_shape_sensitivity_summary

Summarise a shape-sensitivity sweep

Description

[Experimental]

Aggregate the raw output of `janusplot_shape_sensitivity()` into a per-cell mean-accuracy table at either the fine (24-category) or archetype (7-family) level.

Usage

```
janusplot_shape_sensitivity_summary(results, level = c("fine", "archetype"))
```

Arguments

results	Data frame returned by <code>janusplot_shape_sensitivity()</code> .
level	One of "fine" (default) or "archetype".

Value

A data frame with columns truth, n, sigma, accuracy.

See Also

Other shape-sensitivity: `janusplot_shape_sensitivity()`, `janusplot_shape_sensitivity_plot()`, `janusplot_shape_sensitivity_shapes()`

Examples

```
data("shape_sensitivity_demo", package = "janusplot")
head(janusplot_shape_sensitivity_summary(shape_sensitivity_demo,
                                       level = "archetype"))
```

shape_sensitivity_demo

Precomputed shape-recognition sensitivity results (demo)

Description

Raw output from a small-footprint invocation of `janusplot_shape_sensitivity()`. Shipped so users can explore the sensitivity API and regenerate every figure in the `shape-recognition-sensitivity` vignette without having to re-run the sweep themselves. Regenerated via `data-raw/shape_sensitivity_demo.R`.

Design:

- **Shapes** (6, one per non-degenerate archetype): `linear_up`, `concave_up`, `u_shape`, `inverted_u`, `wave`, `bimodal`.
- **Sample sizes** (3): `c(100, 200, 500)`.
- **Noise levels** (4): `c(0.05, 0.10, 0.20, 0.40)` fraction of y-range.
- **Replicates**: 30.
- **Total fits**: 2160.
- **Seed**: 2026.

Usage

```
shape_sensitivity_demo
```

Format

A data frame with 2160 rows and 14 columns — see the "Value" section of `janusplot_shape_sensitivity()` for the column schema.

See Also

[janusplot_shape_sensitivity\(\)](#), [janusplot_shape_sensitivity_plot\(\)](#), [janusplot_shape_sensitivity_summary\(\)](#)

Examples

```
data("shape_sensitivity_demo", package = "janusplot")
head(shape_sensitivity_demo)
janusplot_shape_sensitivity_plot(shape_sensitivity_demo,
                                "recovery_curves")
```

Index

- * **datasets**
 - shape_sensitivity_demo, 18
 - * **shape-metrics**
 - janusplot_shape_cutoffs, 10
 - janusplot_shape_hierarchy, 11
 - janusplot_shape_metrics, 13
 - * **shape-sensitivity**
 - janusplot_shape_sensitivity, 14
 - janusplot_shape_sensitivity_plot, 16
 - janusplot_shape_sensitivity_shapes, 17
 - janusplot_shape_sensitivity_summary, 18
 - * **smooth-associations**
 - janusplot, 2
 - janusplot_data, 8
- ggplot2::ggplot, 7, 17
- janusplot, 2, 10
- janusplot(), 8, 10, 11, 14
- janusplot_data, 7, 8
- janusplot_data(), 7, 14
- janusplot_shape_cutoffs, 10, 12, 14
- janusplot_shape_cutoffs(), 5, 9, 13–15
- janusplot_shape_hierarchy, 11, 11, 14
- janusplot_shape_hierarchy(), 5, 14, 17
- janusplot_shape_metrics, 11, 12, 13
- janusplot_shape_metrics(), 10, 14, 15
- janusplot_shape_sensitivity, 14, 17, 18
- janusplot_shape_sensitivity(), 16–19
- janusplot_shape_sensitivity_plot, 16, 16, 17, 18
- janusplot_shape_sensitivity_plot(), 16, 19
- janusplot_shape_sensitivity_shapes, 16, 17, 17, 18
- janusplot_shape_sensitivity_shapes(), 15, 16
- janusplot_shape_sensitivity_summary, 16, 17, 18
- janusplot_shape_sensitivity_summary(), 16, 19
- mgcv::gam(), 2, 4, 7–9, 13
- patchwork::wrap_plots(), 7
- shape_sensitivity_demo, 14, 16, 18