

Package ‘polyCub’

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Title Cubature over Polygonal Domains

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Description Numerical integration of continuously differentiable functions $f(x,y)$ over simple closed polygonal domains. The following cubature methods are implemented: product Gauss cubature (Sommariva and Vianello, 2007, <[doi:10.1007/s10543-007-0131-2](https://doi.org/10.1007/s10543-007-0131-2)>), the simple two-dimensional midpoint rule (wrapping 'spatstat.geom' functions), and adaptive cubature for radially symmetric functions via line integrate() along the polygon boundary (Meyer and Held, 2014, <[doi:10.1214/14-AOAS743](https://doi.org/10.1214/14-AOAS743)>, Supplement B). For simple integration along the axes, the 'cubature' package is more appropriate.

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URL <https://github.com/bastistician/polyCub>

BugReports <https://github.com/bastistician/polyCub/issues>

Note Building the package requires R \geq 4.6.0 for \bibshow{ } et al.

Depends R (\geq 3.4.0), methods

Imports grDevices, graphics, stats, sp (\geq 1.0-11)

Suggests spatstat.geom, lattice, mvtnorm, statmod, sf, cubature, litedown (\geq 0.9), microbenchmark

VignetteBuilder litedown

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polyCub-package	<i>Cubature over Polygonal Domains</i>
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Description

The R package **polyCub** implements *cubature* (numerical integration) over *polygonal* domains. It solves the problem of integrating a continuously differentiable function $f(x, y)$ over simple closed polygons.

Details

polyCub provides the following cubature methods:

polyCub.SV: General-purpose *product Gauss cubature* (Sommariva and Vianello 2007)

polyCub.midpoint: Simple *two-dimensional midpoint rule* based on `as.im.function` from **spat-stat.geom** (Baddeley, Rubak, and Turner 2015)

polyCub.iso: Adaptive cubature for *radially symmetric functions* via line `integrate()` along the polygon boundary (Meyer and Held 2014, Supplement B, Section 2.4)

A brief description and benchmark experiment of the above cubature methods can be found in the vignette("polyCub").

There is also **polyCub.exact.Gauss**, intended to accurately (but slowly) integrate the *bivariate Gaussian density*; however, this implementation is disabled as of **polyCub** 0.9.0: it needs a reliable implementation of polygon triangulation.

Meyer (2010, Section 3.2) discusses and compares some of these methods.

Note

To cite package **polyCub** in publications, please use `citation("polyCub")`:

Meyer S (2019). “polyCub: An R package for Integration over Polygons.” *Journal of Open Source Software*, **4**(34), 1056. ISSN 2475-9066. doi:10.21105/joss.01056.

Author(s)

Sebastian Meyer

References

Baddeley A, Rubak E, Turner R (2015). *Spatial Point Patterns: Methodology and Applications with R*. Chapman and Hall/CRC Press, London. ISBN 9781482210200.

Meyer S (2010). *Spatio-Temporal Infectious Disease Epidemiology based on Point Processes*. Master’s thesis, Department of Statistics, LMU, Munich, Germany. <https://epub.uni-muenchen.de/11703/>.

Meyer S, Held L (2014). “Power-law models for infectious disease spread.” *Annals of Applied Statistics*, **8**(3), 1612–1639. doi:10.1214/14AOAS743.

Sommariva A, Vianello M (2007). “Product Gauss cubature over polygons based on Green’s integration formula.” *BIT Numerical Mathematics*, **47**(2), 441–453. doi:10.1007/s1054300701312.

See Also

`vignette("polyCub")`

For the special case of a rectangular domain along the axes (e.g., a bounding box), the **cubature** package is more appropriate.

checkintrfr

Check the Integral of $r f_r(r)$

Description

This function is auxiliary to **polyCub.iso** for the cubature of a radially symmetric function $f(x, y) = f_r(\|(x, y) - \mu\|)$, with μ being the center of isotropy, over a polygonal domain. The (analytical) integral of $r f_r(r)$ from 0 to R , `intrfr`, is checked against an **integrate**-based approximation for various values (rs) of the upper bound R . A warning is issued if inconsistencies are found.

Usage

```
checkintrfr(intrfr, f, ..., center, control = list(), rs = numeric(0L),
  tolerance = control$rel.tol)
```

Arguments

intrfr	a function(R, \dots), which implements the (analytical) antiderivative of $r f_r(r)$ from 0 to R . The first argument must be vectorized but not necessarily named R . If <code>intrfr</code> is missing, it will be approximated numerically via <pre>integrate(function(r, ...) r * f(cbind(x0 + r, y0), ...), 0, R, ..., control = control)</pre> where $c(x_0, y_0)$ is the center of isotropy. Note that f will <i>not</i> be checked for isotropy.
f	a two-dimensional real-valued function. As its first argument it must take a coordinate matrix, i.e., a numeric matrix with two columns, and it must return a numeric vector of length the number of coordinates.
...	further arguments for <code>f</code> or <code>intrfr</code> .
center	numeric vector of length 2, the center of isotropy.
control	list of arguments passed to <code>integrate</code> , the quadrature rule used for the line integral along the polygon boundary.
rs	numeric vector of upper bounds for which to check the validity of <code>intrfr</code> . If it has length 0 (default), no checks are performed.
tolerance	of <code>all.equal.numeric</code> when comparing <code>intrfr</code> results with numerical integration. Defaults to the relative tolerance used for <code>integrate</code> .

Value

The `intrfr` function, invisibly. If only `f` was given, an `integrate`-based approximation of `intrfr` is returned.

Examples

```
f_const <- function (coords) rep(1, nrow(coords))
intrfr_const <- function (R) R^2/2 # = \int_0^R r f_r(r) dr
checkintrfr(intrfr_const, f = f_const, center = c(0,0), rs = 1:10) # OK
checkintrfr(function(R) R, f = f_const, center = c(0,0), rs = 1:10) # warns
```

circleCub.Gauss

Integration of the Isotropic Gaussian Density over Circular Domains

Description

This function calculates the integral of the bivariate, isotropic Gaussian density (i.e., $\Sigma = \text{sd}^2 \cdot \text{diag}(2)$) over a circular domain via the cumulative distribution function `pchisq` of the (non-central) Chi-Squared distribution (Abramowitz and Stegun 1972, Formula 26.3.24).

Usage

```
circleCub.Gauss(center, r, mean, sd)
```

Arguments

center	numeric vector of length 2 (center of the circle).
r	numeric (radius of the circle). Several radii may be supplied.
mean	numeric vector of length 2 (mean of the bivariate Gaussian density).
sd	numeric (common standard deviation of the isotropic Gaussian density in both dimensions).

Value

The integral value (one for each supplied radius).

Note

The non-centrality parameter of the evaluated chi-squared distribution equals the squared distance between the mean and the center. If this becomes too large, the result becomes inaccurate, see [pchisq](#).

References

Abramowitz M, Stegun IA (1972). *Handbook of Mathematical Functions with Formulas, Graphs, and Mathematical Tables*. Dover Publications, New York. ISBN 0486612724.

Examples

```
circleCub.Gauss(center=c(1,2), r=3, mean=c(4,5), sd=6)

## compare with cubature over a polygonal approximation of a circle
d2norm <- function(s, mean, sd)
  dnorm(s[,1], mean=mean[1], sd=sd) * dnorm(s[,2], mean=mean[2], sd=sd)
if (requireNamespace("spatstat.geom")) { # for the disc()
  npoly <- 32 # increase this for a closer match
  disc.poly <- spatstat.geom::disc(radius=3, centre=c(1,2), npoly=npoly)
  polyCub.iso(disc.poly, d2norm, mean=c(4,5), sd=6, center=c(4,5))
}
```

coerce-gpc-methods *Conversion between polygonal "owin" and "gpc.poly"*

Description

Package **polyCub** implements converters between the classes *"owin"* of package **spatstat.geom** and *"gpc.poly"* of package **gpclib**.

Usage

```

owin2gpc(object)

gpc2owin(object, ...)

as.owin.gpc.poly(W, ...)

```

Arguments

object	an object of class "gpc.poly" or "owin", respectively.
...	further arguments passed to owin .
W	an object of class "gpc.poly".

Value

The converted polygon of class "gpc.poly" or "owin", respectively. If package **gpclib** is not available, `owin2gpc` will just return the `pts` slot of the "gpc.poly" (no formal class) with a warning.

Note

The converter `owin2gpc` requires the package **gpclib** for the formal class definition of a "gpc.poly". It will produce vertices ordered according to the **sp** convention, i.e. clockwise for normal boundaries and anticlockwise for holes, where, however, the first vertex is *not* repeated!

Author(s)

Sebastian Meyer

See Also

[xymatrix](#)

Examples

```

## use example polygons from
example(plotpolyf, ask = FALSE)
letterR # a simple "xymatrix"

letterR.owin <- spatstat.geom::owin(poly = letterR)
letterR.gpc_from_owin <- owin2gpc(letterR.owin)
## warns if "gpclib" is unavailable

if (is(letterR.gpc_from_owin, "gpc.poly")) {
  letterR.xymatrix_from_gpc <- xymatrix(letterR.gpc_from_owin)
  stopifnot(all.equal(letterR, lapply(letterR.xymatrix_from_gpc, `[`, 1:2)))
  letterR.owin_from_gpc <- gpc2owin(letterR.gpc_from_owin)
  stopifnot(all.equal(letterR.owin, letterR.owin_from_gpc))
}

```

coerce-sp-methods *Coerce "SpatialPolygons" to "owin"*

Description

Package **polyCub** implements coerce-methods (`as(object, Class)`) to convert "[SpatialPolygons](#)" (or "[Polygon](#)" or "[Polygons](#)") of package **sp** to "[owin](#)" of package **spatstat.geom**. They are also available as `as.owin.*` functions to support `polyCub.midpoint`.

Usage

```
as.owin.SpatialPolygons(W, ...)
```

```
as.owin.Polygons(W, ...)
```

```
as.owin.Polygon(W, ...)
```

Arguments

`W` an object of class "[SpatialPolygons](#)", "[Polygons](#)", or "[Polygon](#)".
`...` further arguments passed to [owin](#).

Value

The polygon(s) as an "[owin](#)" object.

Author(s)

Sebastian Meyer

See Also

[xylist](#)

Examples

```
diamond <- list(x = c(1,2,1,0), y = c(1,2,3,2)) # anti-clockwise
diamond.sp <- sp::Polygon(lapply(diamond, rev)) # clockwise

diamond.Ps <- sp::Polygons(list(diamond.sp), ID = "my diamond")
diamond.SpPs <- sp::SpatialPolygons(list(diamond.Ps))

if (require("spatstat.geom")) {
  diamond.owin <- owin(poly = diamond)
  diamond.owin_from_Polygon <- as.owin(diamond.sp)
  stopifnot(all.equal(diamond.owin, diamond.owin_from_Polygon))
  ## also for "Polygons" and "SpatialPolygons", using S3 or S4 methods:
  stopifnot(identical(diamond.owin, as.owin(diamond.Ps)))
  stopifnot(identical(diamond.owin, as(diamond.SpPs, "owin")))
```

```
}

```

plotpolyf

Plot Polygonal Domain on Image of Bivariate Function

Description

Produces a combined plot of a polygonal domain and an image of a bivariate function, using either [lattice::levelplot](#) or [image](#).

Usage

```
plotpolyf(polyregion, f, ..., npixel = 100, cuts = 15,
  col = rev(heat.colors(cuts + 1)), lwd = 3, xlim = NULL, ylim = NULL,
  use.lattice = TRUE, print.args = list())
```

Arguments

polyregion	a polygonal domain. The following classes are supported: "owin" from package spatstat.geom , "gpc.poly" from gpclib , "SpatialPolygons", "Polygons", and "Polygon" from package sp , as well as "(MULTI)POLYGON" from package sf . (For these classes, polyCub knows how to get an xylist .)
f	a two-dimensional real-valued function. As its first argument it must take a coordinate matrix, i.e., a numeric matrix with two columns, and it must return a numeric vector of length the number of coordinates.
...	further arguments for f.
npixel	numeric vector of length 1 or 2 setting the number of pixels in each dimension.
cuts	number of cut points in the z dimension. The range of function values will be divided into cuts+1 levels.
col	color vector used for the function levels.
lwd	line width of the polygon edges.
xlim, ylim	numeric vectors of length 2 setting the axis limits. NULL means using the bounding box of polyregion.
use.lattice	logical indicating if lattice graphics (levelplot) should be used.
print.args	a list of arguments passed to print.trellis for plotting the produced "trellis" object (if use.lattice = TRUE). The print step is omitted if print.args is not a list.

Author(s)

Sebastian Meyer

Examples

```

### a polygonal domain (a simplified version of spatstat.data::letterR$bdry)
letterR <- list(
  list(x = c(2.7, 3, 3.3, 3.9, 3.7, 3.4, 3.8, 3.7, 3.4, 2, 2, 2.7),
       y = c(1.7, 1.6, 0.7, 0.7, 1.3, 1.8, 2.2, 2.9, 3.3, 3.3, 0.7, 0.7)),
  list(x = c(2.6, 2.6, 3, 3.2, 3),
       y = c(2.2, 2.7, 2.7, 2.5, 2.2))
)

### f: isotropic exponential decay
fr <- function(r, rate = 1) dexp(r, rate = rate)
fcenter <- c(2,3)
f <- function (s, rate = 1) fr(sqrt(rowSums(t(t(s)-fcenter)^2)), rate = rate)

### plot
plotpolyf(letterR, f, use.lattice = FALSE)
plotpolyf(letterR, f, use.lattice = TRUE)

```

polyCub

*Wrapper Function for the Various Cubature Methods***Description**

The wrapper function `polyCub` can be used to call specific cubature methods via its `method` argument. It calls the `polyCub.SV` function by default, which implements general-purpose product Gauss cubature. The desired cubature function should usually be called directly.

Usage

```
polyCub(polyregion, f, method = c("SV", "midpoint", "iso", "exact.Gauss"),
  ..., plot = FALSE)
```

Arguments

<code>polyregion</code>	a polygonal domain. The following classes are supported: <code>"owin"</code> from package <code>spatstat.geom</code> , <code>"gpc.poly"</code> from <code>gpclib</code> , <code>"SpatialPolygons"</code> , <code>"Polygons"</code> , and <code>"Polygon"</code> from package <code>sp</code> , as well as <code>"(MULTI)POLYGON"</code> from package <code>sf</code> . (For these classes, <code>polyCub</code> knows how to get an <code>xylist</code> .)
<code>f</code>	a two-dimensional real-valued function to be integrated over <code>polyregion</code> . As its first argument it must take a coordinate matrix, i.e., a numeric matrix with two columns, and it must return a numeric vector of length the number of coordinates. For the <code>"exact.Gauss"</code> method, <code>f</code> is ignored since it is specific to the bivariate normal density.
<code>method</code>	choose one of the implemented cubature methods (partial argument matching is applied), see <code>help("polyCub-package")</code> for an overview. Defaults to using product Gauss cubature implemented in <code>polyCub.SV</code> .

... arguments of `f` or of the specific method.
`plot` logical indicating if an illustrative plot of the numerical integration should be produced.

Value

The approximated integral of `f` over `polyregion`.

See Also

Details and examples in the vignette("polyCub") and on the method-specific help pages.

Other cubature methods: `polyCub.SV()`, `polyCub.iso()`, `polyCub.midpoint()`

`polyCub.exact.Gauss` *Quasi-Exact Cubature of the Bivariate Normal Density (DEFUNCT)*

Description

This cubature method is **defunct** as of **polyCub** version 0.9.0. It relied on `tristrip()` from package **gpclib** for polygon triangulation, but that package did not have a FOSS license and was no longer maintained on a mainstream repository.

Contributions to resurrect this cubature method are welcome: an alternative implementation for constrained polygon triangulation is needed, see <https://github.com/bastistician/polyCub/issues/2>.

Usage

```
polyCub.exact.Gauss(polyregion, mean = c(0, 0), Sigma = diag(2),
  plot = FALSE)
```

Arguments

`polyregion` a "gpc.poly" polygon or something that can be coerced to this class, e.g., an "owin" polygon (via `owin2gpc`), or an "sfg" polygon (via `sfg2gpc`).

`mean, Sigma` mean and covariance matrix of the bivariate normal density to be integrated.

`plot` logical indicating if an illustrative plot of the numerical integration should be produced. Note that the `polyregion` will be transformed (shifted and scaled).

Details

The bivariate Gaussian density can be integrated based on a triangulation of the (transformed) polygonal domain, using formulae from the Abramowitz and Stegun (1972) handbook (Section 26.9, Example 9, pp. 956f.). This method is quite cumbersome because the A&S formula is only for triangles where one vertex is the origin (0,0). For each triangle we have to check in which of the 6 outer regions of the triangle the origin (0,0) lies and adapt the signs in the formula appropriately: $(AOB + BOC - AOC)$ or $(AOB - AOC - BOC)$ or $(AOB + AOC - BOC)$ or $(AOC + BOC - AOB)$ or However, the most time consuming step is the evaluation of `pmvnorm`.

Value

The integral of the bivariate normal density over polyregion. Two attributes are appended to the integral value:

nEval	number of triangles over which the standard bivariate normal density had to be integrated, i.e. number of calls to <code>pmvnorm</code> and <code>pnorm</code> , the former of which being the most time-consuming operation.
error	Approximate absolute integration error stemming from the error introduced by the nEval <code>pmvnorm</code> evaluations. For this reason, the cubature method is in fact only quasi-exact (as is the <code>pmvnorm</code> function).

References

Abramowitz M, Stegun IA (1972). *Handbook of Mathematical Functions with Formulas, Graphs, and Mathematical Tables*. Dover Publications, New York. ISBN 0486612724.

See Also

[circleCub.Gauss](#) for quasi-exact cubature of the isotropic Gaussian density over a circular domain.

polyCub.iso

Cubature of Isotropic Functions over Polygonal Domains

Description

`polyCub.iso` numerically integrates a radially symmetric function $f(x, y) = f_r(\|(x, y) - \mu\|)$, with μ being the center of isotropy, over a polygonal domain. It internally approximates a line integral along the polygon boundary using `integrate`. The integrand requires the antiderivative of $r f_r(r)$, which should be supplied as argument `intrfr` (`f` itself is only required if `check.intrfr=TRUE`). The two-dimensional integration problem thereby reduces to an efficient adaptive quadrature in one dimension. If `intrfr` is not available analytically, `polyCub.iso` can use a numerical approximation (meaning `integrate` within `integrate`), but the general-purpose cubature method [polyCub.SV](#) might be more efficient in this case. See Meyer and Held (2014, Supplement B, Section 2.4) for mathematical details.

`.polyCub.iso` is a “bare-bone” version of `polyCub.iso`.

Usage

```
polyCub.iso(polyregion, f, intrfr, ..., center, control = list(),
  check.intrfr = FALSE, plot = FALSE)
```

```
.polyCub.iso(polys, intrfr, ..., center, control = list(),
  .witherror = FALSE)
```

Arguments

polyregion	a polygonal domain. The following classes are supported: "owin" from package spatstat.geom , "gpc.poly" from gpclib , "SpatialPolygons", "Polygons", and "Polygon" from package sp , as well as "(MULTI)POLYGON" from package sf . (For these classes, polyCub knows how to get an xylist .)
f	a two-dimensional real-valued function. As its first argument it must take a coordinate matrix, i.e., a numeric matrix with two columns, and it must return a numeric vector of length the number of coordinates.
intrfr	a function(R, \dots), which implements the (analytical) antiderivative of $rf_r(r)$ from 0 to R . The first argument must be vectorized but not necessarily named R . If intrfr is missing, it will be approximated numerically via <pre>integrate(function(r, ...) r * f(cbind(x0 + r, y0), ...), 0, R, ..., control = control)</pre> <p>where $c(x_0, y_0)$ is the center of isotropy. Note that f will <i>not</i> be checked for isotropy.</p>
...	further arguments for f or intrfr .
center	numeric vector of length 2, the center of isotropy.
control	list of arguments passed to integrate , the quadrature rule used for the line integral along the polygon boundary.
check.intrfr	logical (or numeric vector) indicating if (for which r 's) the supplied intrfr function should be checked against a numerical approximation. This check requires f to be specified. If TRUE, the set of test r 's defaults to a seq of length 20 from 1 to the maximum absolute x or y coordinate of any edge of the polyregion.
plot	logical indicating if an image of the function should be plotted together with the polygonal domain, i.e., plotpolyf (polyregion, f, \dots).
polys	something like <code>owin\$bdry</code> , but see xylist .
.witherror	logical indicating if an upper bound for the absolute integration error should be attached as an attribute to the result?

Value

The approximate integral of the isotropic function f over polyregion.

If the **intrfr** function is provided (which is assumed to be exact), an upper bound for the absolute integration error is appended to the result as attribute "abs.error". It equals the sum of the absolute errors reported by all **integrate** calls (there is one for each edge of polyregion).

Author(s)

Sebastian Meyer

The basic mathematical formulation of this efficient integration for radially symmetric functions was ascertained with great support by Emil Hedevang (Dept. of Mathematics, Aarhus University, Denmark) during the Summer School on Topics in Space-Time Modeling and Inference (May 2013, Aalborg, Denmark).

References

Meyer S, Held L (2014). “Power-law models for infectious disease spread.” *Annals of Applied Statistics*, **8**(3), 1612–1639. doi:10.1214/14AOAS743.

See Also

system.file("include", "polyCubAPI.h", package = "polyCub") for a full C-implementation of this cubature method (for a *single* polygon). The corresponding C-routine polyCub_iso can be used by other R packages, notably **surveillance**, via ‘LinkingTo: polyCub’ (in the ‘DESCRIPTION’) and ‘#include <polyCubAPI.h>’ (in suitable ‘/src’ files). Note that the intrfr function must then also be supplied as a C-routine. An example can be found in the package tests.

Other cubature methods: [polyCub\(\)](#), [polyCub.SV\(\)](#), [polyCub.midpoint\(\)](#)

Examples

```
## we use the example polygon and f (exponential decay) from
example(plotpolyf)

## numerical approximation of 'intrfr' (not recommended)
(intISOnum <- polyCub.iso(letterR, f, center = fcenter))

## analytical 'intrfr'
## intrfr(R) = int_0^R r*f(r) dr, for f(r) = dexp(r), gives
intrfr <- function (R, rate = 1) pgamma(R, 2, rate) / rate
(intISOana <- polyCub.iso(letterR, f, intrfr = intrfr, center = fcenter,
                        check.intrfr = TRUE))
## f is only used to check 'intrfr' against a numerical approximation

stopifnot(all.equal(intISOana, intISOnum, check.attributes = FALSE))

### polygon area: f(r) = 1, f(x,y) = 1, center does not really matter

## intrfr(R) = int_0^R r*f(r) dr = int_0^R r dr = R^2/2
intrfr.const <- function (R) R^2/2
(area.ISO <- polyCub.iso(letterR, intrfr = intrfr.const, center = c(0,0)))

if (require("spatstat.geom")) { # check against area.owin()
  stopifnot(all.equal(area.owin(owin(poly = letterR)),
                    area.ISO, check.attributes = FALSE))
}
```

polyCub.midpoint

Two-Dimensional Midpoint Rule

Description

The surface is converted to a binary pixel image using the [as.im.function](#) method from package **spatstat.geom**. The integral under the surface is then approximated as the sum over (pixel area * f(pixel midpoint)).

Usage

```
polyCub.midpoint(polyregion, f, ..., eps = NULL, dimyx = NULL,
  plot = FALSE)
```

Arguments

polyregion	a polygonal integration domain. It can be any object coercible to the spatstat.geom class "owin" via a corresponding <code>as.owin</code> -method. Note that this includes polygons of the classes "gpc.poly" and "SpatialPolygons", because polyCub defines methods <code>as.owin.gpc.poly</code> and <code>as.owin.SpatialPolygons</code> , respectively. sf also registers suitable <code>as.owin</code> methods for its "(MULTI)POLYGON" classes.
f	a two-dimensional real-valued function. As its first argument it must take a coordinate matrix, i.e., a numeric matrix with two columns, and it must return a numeric vector of length the number of coordinates.
...	further arguments for f.
eps	width and height of the pixels (squares), see <code>as.mask</code> .
dimyx	number of subdivisions in each dimension, see <code>as.mask</code> .
plot	logical indicating if an illustrative plot of the numerical integration should be produced.

Value

The approximated value of the integral of f over polyregion.

See Also

Other cubature methods: `polyCub()`, `polyCub.SV()`, `polyCub.iso()`

Examples

```
## a function to integrate (here: isotropic zero-mean Gaussian density)
f <- function (s, sigma = 5)
  exp(-rowSums(s^2)/2/sigma^2) / (2*pi*sigma^2)

## a simple polygon as integration domain
hexagon <- list(
  list(x = c(7.33, 7.33, 3, -1.33, -1.33, 3),
    y = c(-0.5, 4.5, 7, 4.5, -0.5, -3))
)

if (require("spatstat.geom")) {
  hexagon.owin <- owin(poly = hexagon)

  show_midpoint <- function (eps)
  {
    plotpolyf(hexagon.owin, f, xlim = c(-8,8), ylim = c(-8,8),
      use.lattice = FALSE)
    ## add evaluation points to plot
  }
}
```

```

    with(as.mask(hexagon.owin, eps = eps),
         points(expand.grid(xcol, yrow), col = t(m), pch = 20))
    title(main = paste("2D midpoint rule with eps =", eps))
  }

  ## show nodes (eps = 0.5)
  show_midpoint(0.5)

  ## show pixel image (eps = 0.5)
  polyCub.midpoint(hexagon.owin, f, eps = 0.5, plot = TRUE)

  ## use a decreasing pixel size (increasing number of nodes)
  for (eps in c(5, 3, 1, 0.5, 0.3, 0.1))
    cat(sprintf("eps = %.1f: %.7f\n", eps,
               polyCub.midpoint(hexagon.owin, f, eps = eps)))
}

```

polyCub.SV

*Product Gauss Cubature over Polygonal Domains***Description**

Product Gauss cubature over polygons as proposed by Sommariva and Vianello (2007).

Usage

```
polyCub.SV(polyregion, f, ..., nGQ = 20, alpha = NULL, rotation = FALSE,
           engine = "C", plot = FALSE)
```

Arguments

polyregion	a polygonal domain. The following classes are supported: "owin" from package spatstat.geom , "gpc.poly" from gpclib , "SpatialPolygons", "Polygons", and "Polygon" from package sp , as well as "(MULTI)POLYGON" from package sf . (For these classes, polyCub knows how to get an xylist .)
f	a two-dimensional real-valued function to be integrated over polyregion (or NULL to only compute nodes and weights). As its first argument it must take a coordinate matrix, i.e., a numeric matrix with two columns, and it must return a numeric vector of length the number of coordinates.
...	further arguments for f.
nGQ	degree of the one-dimensional Gauss-Legendre quadrature rule (default: 20) as implemented in function gauss.quad of package statmod . Nodes and weights up to nGQ=60 are cached in polyCub , for larger degrees statmod is required.
alpha	base-line of the (rotated) polygon at $x = \alpha$ (see Sommariva and Vianello 2007, for an explication). If NULL (default), the midpoint of the x-range of each polygon is chosen if no rotation is performed, and otherwise the x -coordinate of the rotated point "P" (see rotation). If f has its maximum value at the origin (0,0), e.g., the bivariate Gaussian density with zero mean, alpha = 0 is a reasonable choice.

rotation	logical (default: FALSE) or a list of points "P" and "Q" describing the preferred direction. If TRUE, the polygon is rotated according to the vertices "P" and "Q", which are farthest apart (see Sommariva and Vianello 2007). For convex polygons, this rotation guarantees that all nodes fall inside the polygon.
engine	character string specifying the implementation to use. Up to polyCub version 0.4-3, the two-dimensional nodes and weights were computed by R functions and these are still available by setting <code>engine = "R"</code> . The new C-implementation is now the default (<code>engine = "C"</code>) and requires approximately 30% less computation time. The special setting <code>engine = "C+reduce"</code> will discard redundant nodes at (0,0) with zero weight resulting from edges on the base-line $x = \alpha$ or orthogonal to it. This extra cleaning is only worth its cost for computationally intensive functions <code>f</code> over polygons which really have some edges on the baseline or parallel to the x-axis. Note that the old R implementation does not have such unset zero nodes and weights.
plot	logical indicating if an illustrative plot of the numerical integration should be produced.

Value

The approximated value of the integral of `f` over polyregion.
In the case `f = NULL`, only the computed nodes and weights are returned in a list of length the number of polygons of polyregion, where each component is a list with nodes (a numeric matrix with two columns), weights (a numeric vector of length `nrow(nodes)`), the rotation `angle`, and `alpha`.

Author(s)

Sebastian Meyer

References

Sommariva A, Vianello M (2007). "Product Gauss cubature over polygons based on Green's integration formula." *BIT Numerical Mathematics*, **47**(2), 441–453. doi:10.1007/s1054300701312.
Their MATLAB implementation 'polygauss', on which this R implementation was based, is available (in revised versions) at https://sites.google.com/view/alvisesommarivaunipd/home-page/software/software_matlab under the GNU GPL (>=2) license.

See Also

Other cubature methods: `polyCub()`, `polyCub.iso()`, `polyCub.midpoint()`

Examples

```
## a function to integrate (here: isotropic zero-mean Gaussian density)
f <- function (s, sigma = 5)
  exp(-rowSums(s^2)/2/sigma^2) / (2*pi*sigma^2)

## a simple polygon as integration domain
hexagon <- list(
  list(x = c(7.33, 7.33, 3, -1.33, -1.33, 3),
```

```

        y = c(-0.5, 4.5, 7, 4.5, -0.5, -3))
    )

## image of the function and integration domain
plotpolyf(hexagon, f)

## use a degree of nGQ = 3 and show the corresponding nodes
polyCub.SV(hexagon, f, nGQ = 3, plot = TRUE)

## extract nodes and weights
nw <- polyCub.SV(hexagon, f = NULL, nGQ = 3)[[1]]
nrow(nw$nodes)

## manually apply the cubature rule
sum(nw$weights * f(nw$nodes))

## use an increasing number of nodes
for (nGQ in c(1:5, 10, 20, 60))
  cat(sprintf("nGQ = %2i: %.16f\n", nGQ,
             polyCub.SV(hexagon, f, nGQ = nGQ)))

## polyCub.SV() is the default method used by the polyCub() wrapper
polyCub(hexagon, f, nGQ = 3) # calls polyCub.SV()

### now using a simple *rectangular* integration domain

rectangle <- list(list(x = c(-1, 7, 7, -1), y = c(-3, -3, 7, 7)))
polyCub.SV(rectangle, f, plot = TRUE)

## effect of rotation given a very low nGQ
opar <- par(mfrow = c(1,3))
polyCub.SV(rectangle, f, nGQ = 4, rotation = FALSE, plot = TRUE)
  title(main = "without rotation (default)")
polyCub.SV(rectangle, f, nGQ = 4, rotation = TRUE, plot = TRUE)
  title(main = "standard rotation")
polyCub.SV(rectangle, f, nGQ = 4,
           rotation = list(P = c(0,0), Q = c(2,-3)), plot = TRUE)
  title(main = "custom rotation")
par(opar)

## comparison with the "cubature" package
if (requireNamespace("cubature")) {
  fc <- function(s, sigma = 5) # non-vectorized version of f
    exp(-sum(s^2)/2/sigma^2) / (2*pi*sigma^2)
  cubature::hcubature(fc, lowerLimit = c(-1, -3), upperLimit = c(7, 7))
}

```

Description

Package **polyCub** implements a converter from class "**(MULTI)POLYGON**" of package **sf** to "gpc.poly" of package **gpclib** such that `polyCub.exact.Gauss` can be used with simple feature polygons.

Usage

```
sfg2gpc(object)
```

Arguments

`object` a "POLYGON" or "MULTIPOLYGON" "sfg" object.

Value

The converted polygon of class "gpc.poly". If package **gpclib** is not available, `sfg2gpc` will just return the `pts` slot of the "gpc.poly" (no formal class) with a warning.

Note

Package **gpclib** is required for the formal class definition of a "gpc.poly".

Author(s)

Sebastian Meyer

See Also

[xylist](#)

Examples

```
## use example polygons from
example(plotpolyf, ask = FALSE)
letterR # a simple "xylist"

letterR.sfg <- sf::st_polygon(lapply(letterR, function(xy)
  rbind(cbind(xy$x, xy$y), c(xy$x[1], xy$y[1]))))
letterR.sfg
stopifnot(identical(letterR, xylist(letterR.sfg)))

## convert sf "POLYGON" to a "gpc.poly"
letterR.gpc_from_sfg <- sfg2gpc(letterR.sfg)
letterR.gpc_from_sfg
```

Description

Different packages concerned with spatial data use different polygon specifications, which sometimes becomes very confusing (see Details below). To be compatible with the various polygon classes, package **polyCub** uses an S3 class "xylist", which represents a polygonal domain (of potentially multiple polygons) by its core feature only: a list of lists of vertex coordinates (see the "Value" section below). The generic function xylist can deal with the following polygon classes:

- "owin" from package **spatstat.geom**
- "gpc.poly" from package **gpplib**
- "Polygons" from package **sp** (as well as "Polygon" and "SpatialPolygons")
- "(MULTI)POLYGON" from package **sf**

The (somehow useless) default xylist-method does not perform any transformation but only ensures that the polygons are not closed (first vertex not repeated).

Usage

```
xylist(object, ...)  
  
## S3 method for class 'owin'  
xylist(object, ...)  
  
## S3 method for class 'sfg'  
xylist(object, ...)  
  
## S3 method for class 'gpc.poly'  
xylist(object, ...)  
  
## S3 method for class 'SpatialPolygons'  
xylist(object, reverse = TRUE, ...)  
  
## S3 method for class 'Polygons'  
xylist(object, reverse = TRUE, ...)  
  
## S3 method for class 'Polygon'  
xylist(object, reverse = TRUE, ...)  
  
## Default S3 method:  
xylist(object, ...)
```

Arguments

object	an object of one of the supported spatial classes.
...	(unused) argument of the generic.
reverse	logical (TRUE) indicating if the vertex order of the sp classes should be reversed to get the xylist/owin convention.

Details

Polygon specifications differ with respect to:

- is the first vertex repeated?
- which ring direction represents holes?

Package overview:

spatstat.geom: "owin" does *not repeat* the first vertex, and anticlockwise = normal boundary, clockwise = hole. This convention is also used for the return value of xylist.

sp: Repeat first vertex at the end (closed), anticlockwise = hole, clockwise = normal boundary

sf: Repeat first vertex at the end (closed), clockwise = hole, anticlockwise = normal boundary; however, **sf** does not check the ring direction by default, so it cannot be relied upon.

gpclib: There seem to be no such conventions for polygons of class "gpc.poly".

Thus, for polygons from **sf** and **gpclib**, xylist needs to check the ring direction, which makes these two formats the least efficient for integration domains in **polyCub**.

Value

Applying xylist to a polygon object, one gets a simple list, where each component (polygon) is a list of "x" and "y" coordinates. These represent vertex coordinates following **spatstat.geom**'s "owin" convention (anticlockwise order for exterior boundaries, without repeating any vertex).

Author(s)

Sebastian Meyer

Examples

```
diamond <- list(x = c(1,2,1,0), y = c(1,2,3,2)) # anti-clockwise
diamond.sp <- sp::Polygon(lapply(diamond, rev)) # clockwise

diamond.Ps <- sp::Polygons(list(diamond.sp), ID = "my diamond")
diamond.SpPs <- sp::SpatialPolygons(list(diamond.Ps))

stopifnot(identical(xylist(diamond.sp), list(diamond)))
stopifnot(identical(xylist(diamond.Ps), list(diamond)))
stopifnot(identical(xylist(diamond.SpPs), list(diamond)))
```

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