Package ‘subplex’

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Title Unconstrained Optimization using the Subplex Algorithm
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Depends R(>= 2.5.1)

URL https://github.com/kingaa/subplex/

BugReports https://github.com/kingaa/subplex/issues/

Description The subplex algorithm for unconstrained optimization, developed by Tom Rowan <http://www.netlib.org/opt/subplex.tgz>.

Encoding UTF-8

RoxygenNote 7.1.2

Collate 'subplex-package.R' 'subplex.R'

NeedsCompilation yes

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subplex-package

Subplex unconstrained optimization algorithm

Description

The subplex package implements Tom Rowan’s subspace-searching simplex algorithm for unconstrained minimization of a function.

Details

Subplex is a subspace-searching simplex method for the unconstrained optimization of general multivariate functions. Like the Nelder-Mead simplex method it generalizes, the subplex method is well suited for optimizing noisy objective functions. The number of function evaluations required for convergence typically increases only linearly with the problem size, so for most applications the subplex method is much more efficient than the simplex method.

Subplex was written in FORTRAN by Tom Rowan (Oak Ridge National Laboratory). The FORTRAN source code is maintained on the netlib repository (netlib.org).

Author(s)

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References


See Also

subplex, optim

subplex

Minimization of a function by the subplex algorithm

Description

subplex minimizes a function.

Usage

subplex(par, fn, control = list(), hessian = FALSE, ...)
**subplex**

**Arguments**

- **par**: Initial guess of the parameters to be optimized over.
- **fn**: The function to be minimized. Its first argument must be the vector of parameters to be optimized over. It should return a scalar result.
- **control**: A list of control parameters, consisting of some or all of the following:
  - **reitol**: The relative optimization tolerance for `par`. This must be a positive number. The default value is `.Machine$double.eps`.
  - **maxit**: Maximum number of function evaluations to perform before giving up. The default value is 10000.
  - **parscale**: The scale and initial stepsizes for the components of `par`. This must either be a single scalar, in which case the same scale is used for all parameters, or a vector of length equal to the length of `par`. For `parscale` to be valid, it must not be too small relative to `par`: if `par + parscale = par` in any component, `parscale` is too small. The default value is 1.
  - **hessian**: If `hessian=TRUE`, the Hessian of the objective at the estimated optimum will be numerically computed.
  - ... Additional arguments to be passed to the function `fn`.

**Details**

The convergence codes are as follows:

- `-2`: invalid input
- `-1`: number of function evaluations needed exceeds `maxnfe`
- `0`: success: tolerance `tol` satisfied
- `1`: limit of machine precision reached

For more details, see the source code.

**Value**

`subplex` returns a list containing the following:

- **par**: Estimated parameters that minimize the function.
- **value**: Minimized value of the function.
- **count**: Number of function evaluations required.
- **convergence**: Convergence code (see Details).
- **message**: A character string giving a diagnostic message from the optimizer, or ’NULL’.
- **hessian**: Hessian matrix.

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References


See Also

optim

Examples

```r
rosen <- function (x) {  # Rosenbrock Banana function
  x1 <- x[1]
  x2 <- x[2]
  100*(x2-x1*x1)^2+(1-x1)^2
}
subplex(par=c(11,-33),fn=rosen)

rosen2 <- function (x) {
  X <- matrix(x,nrow=2)
  sum(apply(X,1,rosen))
}
subplex(par=c(-33,11,14,9,0,12),fn=rosen2,control=list(maxit=30000))
## compare with optim:
optim(par=c(-33,11,14,9,0,12),fn=rosen2,control=list(maxit=30000))

ripple <- function (x) {
  r <- sqrt(sum(x^2))
  1-exp(-r^2)*cos(10*r)^2
}
subplex(par=c(1),fn=ripple,hessian=TRUE)
subplex(par=c(0.1,3),fn=ripple,hessian=TRUE)
subplex(par=c(0.1,3,2),fn=ripple,hessian=TRUE)

rosen <- function (x, g = 0, h = 0) {  # Rosenbrock Banana function (using names)
  x1 <- x["a"]
  x2 <- x["b"]-h
  100*(x2-x1*x1)^2+(1-x1)^2+g
}
subplex(par=c(b=11,a=-33),fn=rosen,h=22,control=list(abstol=1e-9,parscale=5),hessian=TRUE)
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