



AHELP for CIAO 3.4

# simplex

Context: [sherpa](#)

*Jump to:* [Description](#) [Parameters](#) [Bugs](#) [See Also](#)

## Synopsis

A simplex optimization method.

## Syntax

```
simplex [iters] [eps] [alpha] [beta] [gamma]
```

## Description

The SIMPLEX method is a single-shot method which attempts to find the local fit-statistic minimum nearest to the starting point. Its principal advantage is that it can work well with complicated statistic surfaces (more so than LEVENBERG–MARQUARDT), while also working quickly (more so than POWELL). Its principal disadvantages are that it has a tendency to “get stuck” in regions with complicated topology before reaching the local fit-statistic minimum, and that there is no guarantee it will find the global fit-statistic minimum. Its tendency to stick means that the user may be best-served by repeating fits until the best-fit point does not change.

A simplex is geometrical form in N-dimensional in parameter space which has N + 1 vertices (e.g., in 3-D it is a tetrahedron). The fit statistic is evaluated for each vertex, and one or more points of the simplex are moved, so that the simplex moves towards the nearest local fit-statistic minimum. When a minimum is reached, the simplex may also contract itself, as an amoeba might; hence, the routine is also sometimes called “amoeba.” Convergence is reached when the simplex settles into a minimum and all the vertices are within some value eps of each other.

The eps parameter controls when the optimization will cease; for SIMPLEX, this will occur when

$$| S_i - S_{(i-1)} | < \text{eps} ,$$

where  $S_{(i-1)}$  and  $S_i$  are the observed statistic values for the (i-1)th and ith iteration, respectively.

## Parameters

name	type	def	min	max
<a href="#">iters</a>	integer	2000	1	10000
<a href="#">eps</a>	real	1.e-6	1.e-9	0.001
<a href="#">alpha</a>	real	1	0.1	2

<u>beta</u>	real	0.5	0.05	1
<u>gamma</u>	real	2	1.1	20

## Detailed Parameter Descriptions

**Parameter=iters** (integer default=2000 min=1 max=10000)

*Maximum number of iterations.*

**Parameter=eps** (real default=1.e-6 min=1.e-9 max=0.001)

*Criterion to stop fit.*

**Parameter=alpha** (real default=1 min=0.1 max=2)

*Algorithm convergence factor.*

**Parameter=beta** (real default=0.5 min=0.05 max=1)

*Algorithm convergence factor.*

**Parameter=gamma** (real default=2 min=1.1 max=20)

*Algorithm convergence factor.*

## Bugs

See the [Sherpa bug pages](#) online for an up-to-date listing of known bugs.

## See Also

*sherpa*

[get\\_method](#), [expr](#), [grid](#), [grid-powell](#), [levenberg-marquardt](#), [method](#), [monte-lm](#), [monte-powell](#), [montecarlo](#), [powell](#), [sigma-rejection](#), [simul-ann-1](#), [simul-ann-2](#), [simul-pow-1](#), [simul-pow-2](#), [usermethod](#)

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URL:  
<http://cxc.harvard.edu/ciao3.4/simplex.html>  
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