

# RiemannianRobustMestimator Usage Guide

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This guide explains how to use **RiemannianRobustMestimator** to fit robust location (mean/center) estimators for manifold-valued data using gradient-based optimization.

**Goal.** Clear usage for both non-autograd and autograd workflows, including built-in M-estimators, optimizer options, initialization, and custom losses.

## 1) What it does

Given samples on a manifold with a metric that supports *log* / *exp*, the estimator minimizes a robust objective of the form below (notation shown for reference):

$$L(\mu) = (1/\sum_i \omega_i) \sum_{i=1}^n \omega_i \rho(r_i(\mu)), \quad r_i(\mu) = \|\log_{\mu}(x_i)\| \\ \nabla_{\mu} L = -(1/\sum_i \omega_i) \sum_i \omega_i w(r_i) \log_{\mu}(x_i), \quad w(r) = \rho'(r)/r$$

## 2) Backends and prerequisites

### Supported methods.

- **default** and **adaptive** work on the NumPy backend and autodiff backends.
- **autograd** requires an autodiff backend (geomstats *autograd* or *pytorch*) and is not available on pure NumPy.

**Minimum metric operations.** Your space.metric should provide log, exp, and a compatible norm/distance.

## 3) Inputs and outputs

### Constructor inputs

Argument	Meaning
space	Equipped manifold (geomstats)
m_estimator	Built-in loss name or 'custom'
critical_value	Cutoff/scale parameter(s) for the loss (may be auto-filled)
method	Optimization method: 'default', 'adaptive', or 'autograd'
init_point_method	Initialization: 'first', 'mean-projection', 'midpoint'

### fit inputs

- **X**: array-like, shape [n\_samples, \*metric.shape]
- **weights** (optional): array-like, shape [n\_samples] (defaults to uniform)

## Output: estimate\_

Field	Meaning
x	Final estimate (fitted center)
losses	Loss values over iterations
bases	Base points over iterations
n_iter	Number of accepted iterations
time	Wall-clock time (seconds)

## 4) Quick start

### Non-autograd (NumPy-compatible): method='default' or 'adaptive'

```
from geomstats.geometry.hypersphere import Hypersphere
from geomstats.learning.riemannian_robust_m_estimator import RiemannianRobustMestimator

space = Hypersphere(3)
X = space.random_point(200)

est = RiemannianRobustMestimator(
    space=space,
    method="default",          # or "adaptive"
    m_estimator="huber",
    critical_value=None,       # auto default for ~95% ARE (if available)
    init_point_method="mean-projection",
)

est.set(init_step_size=0.5, max_iter=1024, epsilon=1e-7, verbose=False)
est.fit(X)

mu_hat = est.estimate_.x
```

### Autodiff: method='autograd' (requires autodiff backend)

```
est = RiemannianRobustMestimator(
    space=space,
    method="autograd",
    m_estimator="huber",
    critical_value=None,
    init_point_method="mean-projection",
)

est.set(init_step_size=0.2, max_iter=512, epsilon=1e-7)
est.fit(X)

mu_hat = est.estimate_.x
```

## 5) Built-in M-estimators and default critical\_value

Accepted names: default, huber, pseudo-huber, cauchy, biweight, fair, hampel, welsch, logistic, lorentzian, correntropy, custom. ('default' maps to 'huber').

If `critical_value` is `None` and a built-in estimator is used, the code fills a default value (intended to target ~95% ARE under a reference model).

Loss	Default critical_value
huber / default	1.345
pseudo-huber	1.345
cauchy	2.3849
biweight	4.6851
fair	1.3998
hampel	1.35 (base scale)
welsch	2.9846
logistic	1.205
lorentzian	2.678
correntropy	2.1105

## 6) Optimization methods

**default:** standard Riemannian gradient descent. Current non-autograd path expects (loss, gradient).

**adaptive:** step size adaptation (Levenberg–Marquardt-style). Current path expects (loss, gradient).

**autograd:** gradients via autodiff; loss only needs to return a scalar.

## 7) Initialization (`init_point_method`)

Value	Meaning
first	Use <code>X[0]</code>
mean-projection	Project the Euclidean mean onto the manifold
midpoint	Choose a 'midpoint' sample by sorting along the first coordinate

## 8) Custom loss functions

### A) Function-style custom loss (simple)

For `method='default'/'adaptive'`, your function should be able to return (loss, grad) when requested.

```
def my_loss(space, points, base, critical_value, weights=None, loss_and_grad=True):
    # return loss only if loss_and_grad is False
    # return (loss, grad) if loss_and_grad is True
    ...

est = RiemannianRobustMestimator(space=space, method="default", m_estimator="custom")
est.set_loss(my_loss)
est.set(init_step_size=0.5, max_iter=1024, epsilon=1e-7)
```

```
est.fit(X)
```

## B) Bound loss object (recommended for unified base-only calls)

A cleaner pattern is a loss object with `bind(...)` and `__call__(base, return_grad=...)`. After `bind`, you can make a base-only callable `loss_with_base(base)`. To fully unify, update the non-autograd optimizers to call this base-only function as well.

## 9) Optimizer hyperparameters (set)

Parameter	Meaning
<code>init_step_size</code>	Step size / learning rate
<code>max_iter</code>	Maximum iterations
<code>epsilon</code>	Stopping tolerance
<code>verbose</code>	Print progress periodically

```
est.set(init_step_size=0.2, max_iter=512, epsilon=1e-7, verbose=True)
```

## 10) Common pitfalls

- autograd method on NumPy backend is unsupported.
- non-autograd custom losses must match the expected signature and provide gradients.
- add a tiny epsilon in terms like  $c/r$  to avoid division by zero at  $r=0$ .
- some spaces/metrics (e.g., SPD) may need numerical safeguards in  $\exp/\log$ .