

# PhD Progress

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# 1 Introduction

This document is a log-book for all the work done during my PhD project. All code used can be found on github repository <https://github.com/marco-cucchi/L96gev>.

## 2 EVT and the Lorenz-96 model

Aim of this work is to find EVT parameters for observables of the Lorenz-96 (L96) model, and compare them with the bounds provided in [1].

### 2.1 Lorenz-96 model simulations

As a first step, a number of independent simulations of the L96 model are performed. The L96 model is defined as follows. For  $i = 1, \dots, N$ :

$$\frac{dx_i}{dt} = (x_{i+1} - x_{i-2})x_{i-1} - x_i + F \quad (1)$$

where it is assumed that  $x_{-1} = x_{N-1}$ ,  $x_0 = x_N$  and  $x_{N+1} = x_1$ . Here  $x_i$  is the state of the system on the  $i$ -th coordinate, and  $F$  is the forcing constant. For this set of simulations, values of  $F$  and  $N$  have been set to  $F = 8$  and  $N = 32$ .

Integration has been conducted using 4-th order Runge-Kutta scheme, with integration step  $dt = 10^{-2}$ . Initial conditions for each simulation have been set equal to

$$x_i^0 = 8 + \epsilon, \quad \epsilon \sim U([-0.05, +0.05]) \quad (2)$$

Different levels of spatial aggregation, defined as  $A = 32, 16, 8, 4, 2, 1$ , have been considered:

- For  $A = 32$  no aggregation is performed, and each value  $x_i$  is treated independently;
- For  $A = 1$  all original  $N$   $x_i$  values are spatially averaged into one single value  $\bar{x}$  for each time-step;
- More in general, for  $A = K$  the  $N$  spatial coordinates indicated by the index  $i$  are divided into  $K$  non-overlapping clusters  $c_j$  fixed in time, and corresponding  $x_i$  values belonging to the same cluster are averaged at each time-step.

As observable, the local energy of the system for different levels of aggregation

$$E_j = \frac{1}{2}x_j^2, \quad x_j = \begin{cases} x_i, & A = 32 \\ \bar{x} = \frac{1}{N} \sum_{i=1}^N x_i & A = 1 \\ \frac{1}{\#c_j} \sum_{i \in c_j, K} x_i & A = K \end{cases} \quad (3)$$

is considered.

In order to extract information on the statistics of extremes, very long simulations have to be performed. In order to find a good compromise between this requirement and the limited amount of disk space, a similar procedure to the one adopted in [2] has been followed: instead of keeping all values of each simulation, only block-maxima are retained, with block size  $\Delta t = 0.5$ . It is important to highlight that block-maxima are computed *after* aggregation (spatial average).

Following this procedure, for each simulation (initial condition) 6 different files are obtained, each corresponding to one particular aggregation level  $A$ : each of these files, then, contain  $A$  time-series of block-maxima, one for each of the  $A$  clusters.

Script: c003e11.

## 2.2 Statistics of Extreme Events

Parameters defining GEV distribution are estimated using three different approaches:

- Direct fit using *block-maxima* approach;
- Direct fit using *POT* approach (still not described here);
- Method of *moments* described in [1].

Finally, estimates derived with these approaches are compared among them and with bounds related to attractor's dimensions described in [2].

### 2.2.1 Block-maxima approach

In this approach, each time-series for each different cluster of each simulation is fitted against GEV family of distribution separately. More specifically, for each cluster time-series belonging to a different simulation the following procedure is carried out:

1. Percentiles' orders  $p$  of interest are fixed (e.g. 0.99, 0.995, ...), and corresponding percentiles (thresholds)  $T_p$  are computed;<sup>1</sup>
2. Time-series is divided in  $n$  blocks, where  $n = \text{length}(\text{time-series})(1 - p)$ ;
3. Compute maxima for each block;
4. Fit GEVD family to the block-maxima series.

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<sup>1</sup>This could be something to think upon; in this way I have (slightly?) different percentiles for different time-series in the same simulation and for different simulations. Is this right? The underlying assumption in this procedure should be that all time-series belonging to all simulations should come from the same distribution. So shouldn't the percentiles be the same for all of them?

The fit is performed with the R function `gevFit` from the package `fExtremes`, using MLE approach. As a result estimations of shape parameter  $\xi$ , location parameter  $\mu$  and scale parameter  $\sigma$  are returned, with respective uncertainties as computed via MLE.

Location parameter  $\mu$  is actually assigned the value  $T_p$ ; the *absolute maximum* from each time-series is also kept; *modified scale parameter*  $\sigma^*$  is computed as

$$\sigma^* = \sigma - \xi T_p \quad (4)$$

Error on  $\sigma^*$  is estimated via propagation of error.

As explained in [3], in order to find a valid threshold value  $T_0$  for excess to follow generalized Pareto distribution (and, consequently, GEV distribution), it is a good practice to plot  $\xi$  and  $\sigma^*$  against  $T_p$  and look for the value where both start to be approximately constant: that value is  $T_0$ .

Once parameters have been estimated for all clusters in a simulation, a single estimation of each parameter is saved as the average among all estimates.<sup>2</sup> Furthermore, the following parameters are estimated for each simulation:

- *scale parameter*  $\sigma$  is computed with inverse of equation 4, and relative error is computed via propagation of errors;<sup>3</sup>
- *upper end-point* is computed as<sup>4</sup>

$$u\hat{e}p = \hat{\mu} - \frac{\hat{\sigma}}{\hat{\xi}} \quad (5)$$

and the relative error is computed via propagation of errors.

Finally, for each different aggregation, ensemble averages of *shape* and *modified scale* among all simulations are computed.

### 2.2.2 Method of Moments

Following theory described in [1], we want to estimate *shape* and *scale* parameters using the following equations (Par 8.2.6 in [1]):

$$\xi_A^T = \frac{1}{2} \left( 1 - \frac{(\langle \tilde{A}_1^T \rangle)^2}{\langle \tilde{A}_0^T \rangle \langle \tilde{A}_2^T \rangle - (\langle \tilde{A}_1^T \rangle)^2} \right) \quad (6)$$

$$\sigma_A^T = \frac{1}{2} \frac{\langle \tilde{A}_1^T \rangle \langle \tilde{A}_2^T \rangle}{\langle \tilde{A}_2^T \rangle \langle \tilde{A}_0^T \rangle - \langle \tilde{A}_1^T \rangle^2} \quad (7)$$

---

<sup>2</sup>The *absolute maximum* is also averaged, and this could be an error. The average of the *location parameter*  $\mu$  is also a little disturbing, but this could be solved following reasoning in footnote ???. Error computation should be checked.

<sup>3</sup>This sounds very stupid, since  $\sigma$  was originally estimated (but not saved) via MLE fit to GEVD.

<sup>4</sup>Find reference

where  $A(x)$  is an observable of the system,  $T$  is a threshold value and

$$\langle \tilde{A}_n^T \rangle = \int \mu(dx) \Theta(A(x) - T) (A(x) - T)^n, \quad (8)$$

being  $\Theta$  the Heaviside distribution. This results are exact in the limit for  $T \rightarrow A_{max}$ .

In order to perform this computation, the following procedure has been adopted. First, for each cluster time-series belonging to a different simulation:

1. Percentiles' orders  $p$  of interest are fixed, and corresponding percentile (thresholds)  $T_p$  are computed (footnote 1);
2.  $\langle \tilde{A}_n^{T_p} \rangle$  for  $n = 0, 1, 2$  are computed, using temporal average in place of ensemble average (assuming ergodicity).<sup>5</sup>

Once moments have been estimated for all clusters in a simulation, a single estimation of each moment is saved as the average among all estimates, and relative standard deviations are computed.

Using these estimates, *shape* parameter is computed through equation 6 and estimation of uncertainty is computed via propagation of error. Finally, for each different aggregation, ensemble averages among all simulations are computed.

### 2.2.3 Bounds to the *shape* parameter from the attractor's dimensions

We want to verify relation (8.2.15) in [1], which states that

$$(d_s + d_u + d_n) / 2 \leq \delta \leq d_s + (d_u + d_n) / 2, \quad (9)$$

where

- $d_u$  is equal to the number of positive Lyapunov exponents of the system [4];
- $d_n$  is equal to the number of zero Lyapunov exponents of the system, and in particular it is 1 for Axiom A systems<sup>6</sup>;
- $d_s = n + \sum_{k=1}^n \lambda_k / |\lambda_{n+1}| - d_u - d_n$  [2], with  $\lambda_k$  denoting the Lyapunov exponents of the system, in a descending order, and  $n$  is such that  $\sum_{k=1}^n \lambda_k$  is positive and  $\sum_{k=1}^{n+1} \lambda_k$  is negative;
- $\xi = -1/\delta$ ;
- $\sigma = (A_{max} - T) / \delta$ , with  $A_{max}$  and  $T$  denoting the maximum observed value of the observable<sup>7</sup> and the threshold value.

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<sup>5</sup>No standard deviation has been computed at this stage!

<sup>6</sup>We are taking this for true in our system

<sup>7</sup>or the *upper end point*?

Lyapunov exponents have been computed using Benettin algorithm with QR decomposition. Bounds have been computed and averaged over 50 iterations (simulations).

Script:

- Lyapunov exponents computation: 136afd4
- Average bounds computation:

## 2.3 Statistics of Extreme Events: Corrections and Results

In this section results of the analyses reported in Sec. 2.2 are described, after issues highlighted in the footnotes 1,2,3.

Script:

- quantiles computation: 6880d20

### 2.3.1 Block-maxima approach

The following corrections have been applied:

- Percentiles are computed once, concatenating the first 80 simulations of the first clusters for each aggregation;
- Shape, scale and location parameters from fit procedures are saved for each cluster in each simulation. Averages and computation of derived parameters come after;

Results are shown in Fig. 1 and 2.

Script:

- fit: 240d6d0
- parameters derivation and plots: 97925ef

### 2.3.2 Method of Moments

Results are shown in Fig. 3 and 4.

Script:

- moments computation: 240d6d0
- parameters derivation and plots: 97925ef

## 3 LRT and the Lorenz-96 model

We aim to apply Ruelle response theory [5][6] to predict the response of different observables to the action of both constant and time-dependent forcings to our Lorenz-96 model. Among these observables, we will focus our efforts on ones describing statistics of extreme events.

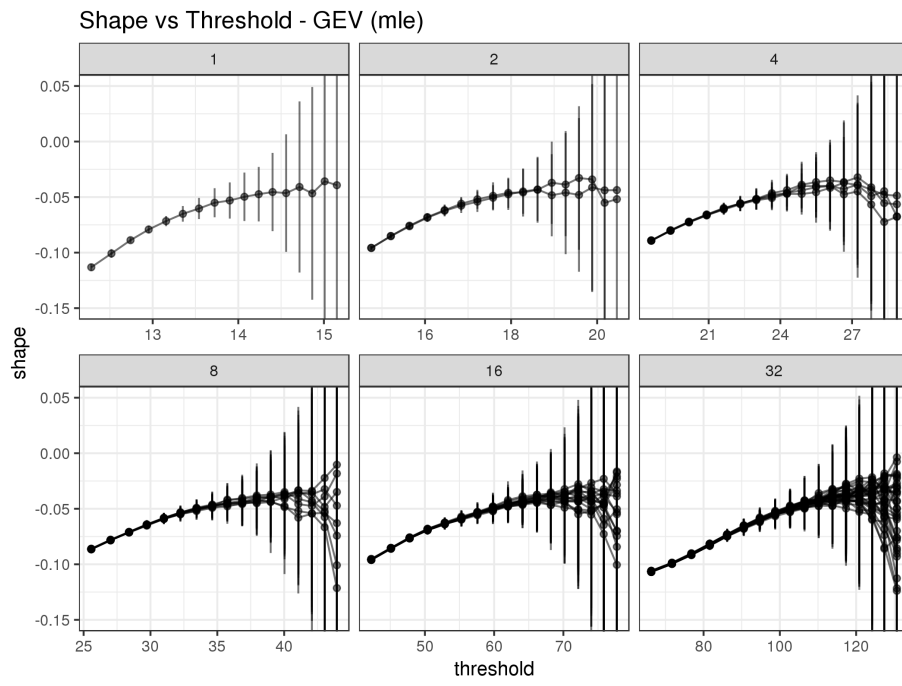


Figure 1: Ensemble average of *shape* parameter over 92 simulations. Each cluster is treated separately.

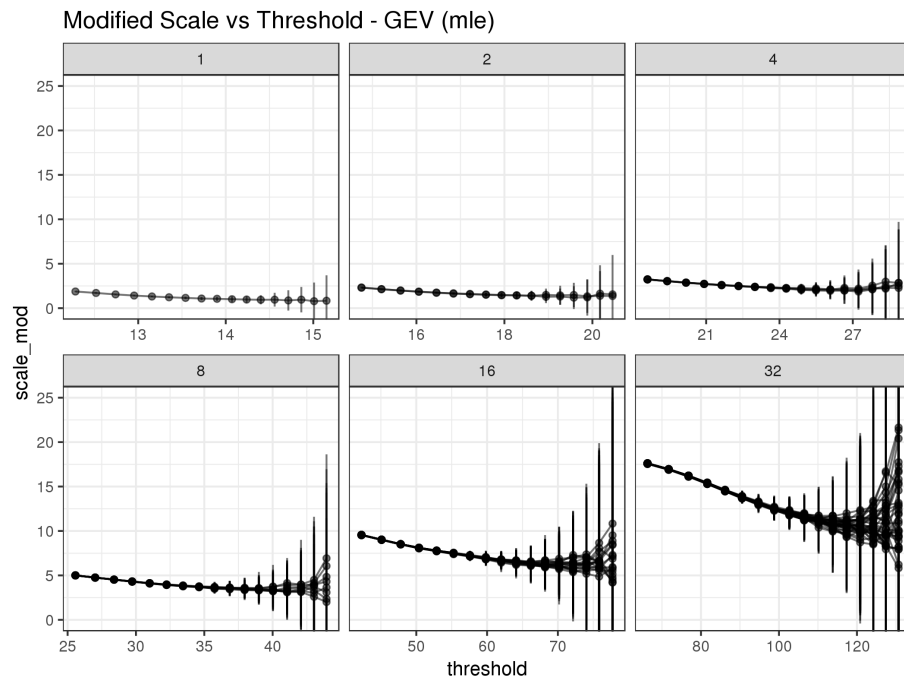


Figure 2: Ensemble average of *modified scale* parameter over 92 simulations. Each cluster is treated separately.



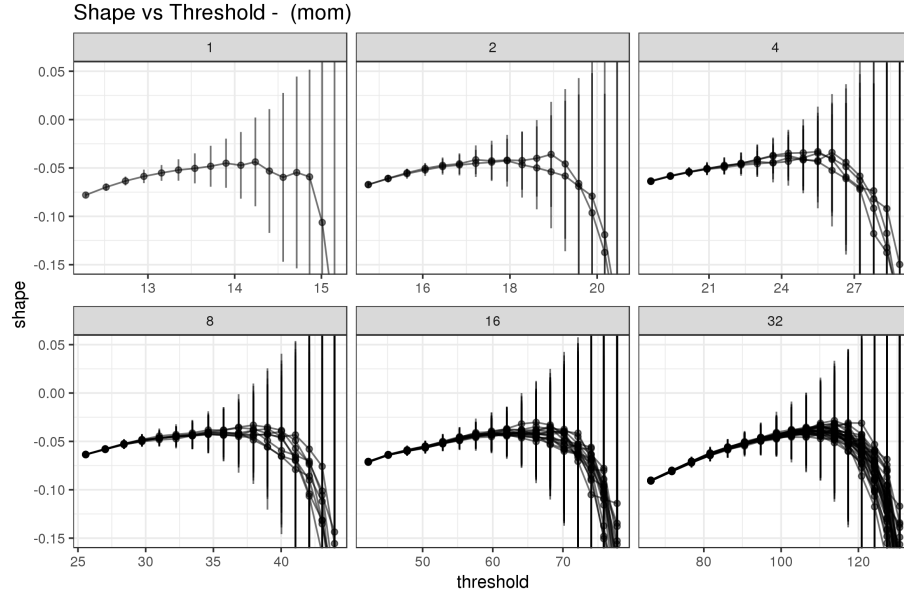


Figure 3: Ensemble average of *shape* parameter over 92 simulations. Each cluster is treated separately.

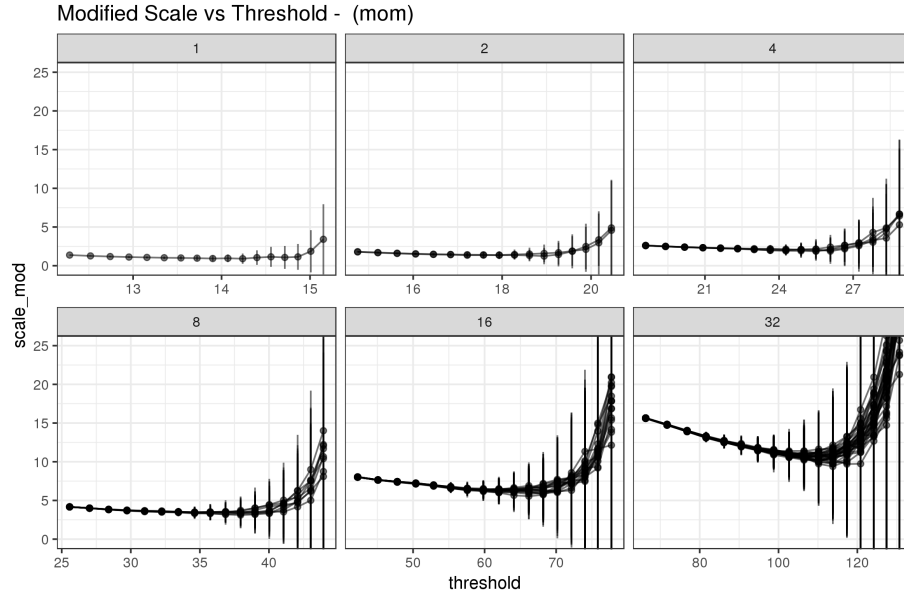


Figure 4: Ensemble average of *modified scale* parameter over 92 simulations. Each cluster is treated separately.

### 3.1 Background

Given a *nonautonomous dissipative dynamical system* in the form

$$\dot{x} = F(x) + \epsilon g(x)f(t) \quad (10)$$

and a scalar observable  $\Psi(x)$ , Ruelle's response theory [5] asserts that its mean  $\langle \Psi \rangle = \int \mu_t(dx) \Psi(x)$  can be decomposed as

$$\langle \Psi \rangle(t) = \sum_{j=1}^{\infty} \epsilon^j \langle \Psi \rangle^{(j)} + \langle \Psi \rangle_0 \quad (11)$$

where the  $\langle \Psi \rangle$  can be expressed as multiple convolution integrals involving the pertinent Green's functions [7]. In the linear (first-order) approximation, we can thus express the response to the forcing  $f(t)$  as

$$\Delta \langle \Psi \rangle(t) = \langle \Psi \rangle^{(1)}(t) = G_{\Psi}^{(1)}(t) * f(t) = \int_{-\infty}^{\infty} d\tau G_{\Psi}^{(1)}(\tau) f(t - \tau) \quad (12)$$

where the Green's function has been established by Ruelle to take the form of

$$G_{\Psi}^{(1)}(t) = \int dx \Psi(x) (\exp [tL_f] [L_g \bar{\mu}]) (x) \quad (13)$$

where  $\bar{\mu}(dx)$  is the natural invariant measure/probability distribution of the autonomous system ( $f = 0$ ), and operators are defined as  $L_f \mu = -\text{div}(f\mu)$  and  $L_g \mu = -\text{div}(g\mu)$ , in the notation of [8].

## References

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