

# The yaev package: Yet Another Extreme Value package?

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## 1 Probability functions of Extreme-Value distributions

The probability functions for the Generalised Pareto (GP) and Generalised Extreme Value (GEV) distributions Coles (2001) depend smoothly on the parameters: they are infinitely differentiable functions of the parameters. However, these functions are not *analytic* functions of the parameters and a singularity exists for all of them when the shape parameter, say  $\xi$ , is zero. In practice, the functions are given with different formulas depending on whether  $\xi$  is zero or not; the formulas for  $\xi = 0$  relate to the exponential and Gumbel distributions and correspond to the limit for  $\xi \rightarrow 0$  of the functions given by the formulas for  $\xi \neq 0$ . As an example, consider the quantile function of the Generalised Pareto distribution with shape  $\xi$  and unit scale

$$q(p) = \begin{cases} [(1-p)^{-\xi} - 1]/\xi & \xi \neq 0 \\ -\log(1-p) & \xi = 0, \end{cases} \quad 0 < p < 1. \quad (1)$$

It can be shown that for  $\xi \approx 0$  whatever be  $p$

$$q \approx -\log(1-p), \quad \frac{\partial q}{\partial \xi} \approx \frac{1}{2} \log^2(1-p), \quad \frac{\partial^2 q}{\partial \xi^2} \approx -\frac{1}{3} \log^2(1-p).$$

It is quite easy to obtain expressions for the derivatives w.r.t.  $\xi$  using the definition (1). We can even rely on the symbolic differentiation method `D` available in R which, as opposed to me and many other humans, never makes any mistake when differentiating.

```

qEx <- function(p, xi) ((1 - p)^(-xi) - 1) / xi
dqEx <- D(expression(((1 - p)^(-xi) - 1) / xi), name = "xi")
d2qEx <- D(dqEx, name = "xi")
p <- 0.99; pBar <- 1 - p
for (xi in c(1e-4, 1e-6, 1e-8)) {
  r <- rbind("ord 0" = c("lim" = -log(pBar), "der" = qEx(p = p, xi = xi)),
            "ord 1" = c("lim" = log(pBar)^2 / 2, "der" = eval(dqEx, list(p = p, xi = xi))),
            "ord 2" = c("lim" = -log(pBar)^3 / 3, "der" = eval(d2qEx, list(x = p, xi = xi))))
  cat("xi = ", xi, "\n")
  print(r)
}

## xi = 1e-04
##          lim          der
## ord 0  4.60517  4.606231
## ord 1 10.60380 10.607052
## ord 2 32.55486 32.566137
## xi = 1e-06
##          lim          der
## ord 0  4.60517  4.605181
## ord 1 10.60380 10.603811
## ord 2 32.55486 68.213867
## xi = 1e-08
##          lim          der
## ord 0  4.60517  4.605170e+00
## ord 1 10.60380  1.085129e+01
## ord 2 32.55486 -4.949912e+07

```

We see that the formula for the function works fine. However, the formula for the 2-nd order derivative can be completely wrong when  $\xi$  is about  $1e-6$  and the formula for the 1-st order derivative can also be wrong when  $\xi$  is about  $1e-8$ . The reason is that the formulas for the derivatives involve difference and/or fractions or small quantities since  $\xi$  or  $\xi^2$  comes at the denominator. As a general rule, the derivatives with higher order are more difficult to evaluate numerically, since they involve more complex expressions. Note that using a shape  $\xi$  with  $|\xi| \leq 1e-6$  is quite common in EVA because the values of  $\xi$  used in practice are often quite small, and moreover very small values of  $\xi$  are often used in the initialisation of the Maximum-Likelihood (ML) optimisation.

Although not yet widespread, the use of the exact formulas for the derivatives w.r.t. the parameters can be of great help in the optimisation tasks required in EVA. These tasks of course involve the ML estimation, but also profile-likelihood inference for models with covariates. Differential equations methods can be also used to derive confidence intervals. Note that the use of formulas for the derivatives is called *symbolic* differentiation and differs from *automatic* differentiation as increasingly available.

Our strategy consists in fixing a small  $\epsilon > 0$  and use the formulas for  $\xi \neq 0$  only when  $|\xi| > \epsilon$ . When  $|\xi| \leq \epsilon$ , we use a few terms from the Taylor series at  $\xi = 0$  e.g.,

$$q(p; \xi) \approx q(p; 0) + \partial_{\xi} q(p; \xi)|_{\xi=0} \times \xi + \frac{1}{2} \partial_{\xi, \xi}^2 q(p; \xi)|_{\xi=0} \times \xi^2 + o(\xi^2).$$

In order to maintain the consistency between the derivatives, it seems good to use the same  $\epsilon$  for all the derivatives and use consistent Taylor approximations, so use the order 1 for the derivative  $\partial_{\xi} q$  and the order order 0 for  $\partial_{\xi, \xi}^2 q$  or for a crossed derivative such as  $\partial_{\sigma, \xi}^2 q$ . Since the 2-nd order derivatives can be required, we must take a value for  $\epsilon$  which is not too small:  $1e-4$  or  $1e-5$ , not

much smaller. Note that  $\epsilon$  gives the level of error for the 2-nd order derivative; since the error on the function is  $O(\xi^3)$ , using  $\epsilon = 1e-4$  leads to an error of order  $1e-12$  on the function, which seems acceptable in practice. This kind of approximation is used in some codes of the **revdbayes** package by Paul Northrop, see the code if the **dgev** and **pgev** functions on GitHub repos.

## 2 Deriving the formulas

The reports provided with **yaev**: **GEV.pdf**, **GPD2.pdf**, **PoisGP2PP.pdf** and **PP2PoisGP.pdf** give the exact expressions for the first-order and the second-order derivatives of the probability functions w.r.t. the parameters and also provides workable approximations for the case  $\xi \approx 0$ . We used the Maxima Computer Algebra System along with the **maxiplot** package for L<sup>A</sup>T<sub>E</sub>X.

- The **raw expressions given by Maxima are reported in green**. The expressions can be regarded as exact, not being influenced by manual computations. However these formulas are usually difficult to use in a compiled code and require some manual transformation for this aim.
- The **simplified expressions are reported in red**. These expressions are derived by us from the raw expressions; they are influenced by manual computations hence could in principle contain errors, although they have been carefully checked. These formulas are used to write the compiled code. They often use auxiliary variables that are shared across several formulas.

## 3 Testing the derivatives

The **yaev** package comes with a series of tests in the format of the **testthat** package. The **numDeriv** package is used to compute the derivatives by numeric differentiation up to the order 2; these derivatives are compared to those provided by the formulas.

A quite difficult task when checking derivatives is to give a threshold used to decide if the difference between the numeric derivative and the symbolic derivative, say the “error”, is acceptable or not. This error has two sources: one is the numerical evaluation of the symbolic derivative or of its approximation for  $\xi \approx 0$  (see example above), and the other is the approximation used in numeric differentiation where the limit defining the derivative is replaced by a finite difference. We should use small values of  $\xi$  with  $|\xi| < \epsilon$  to check that the approximation for small  $\xi$  is correct, although we can only test the approximation at the first order by doing so. Then, with a good choice of  $\epsilon$  the error should be mainly due to the numeric differentiation. But when the true derivative is small, the relative error may be large (think of a true derivative which is exactly zero). On the other hand, when a derivative is large in absolute value, the absolute error may also be quite large. Mind that the derivatives can in practice be very small or very large, and also that a gradient vector or a Hessian matrix often contain values that are not of the same order of magnitude.

We check that either the *absolute* error or the *relative* error is small. The idea is that none of these two things can come by chance, and if one holds, the formula used must be good even if the other criterion suggests an opposed conclusion. The test is made *elementwise*, meaning that the relative error is computed for each element of a gradient vector or Hessian matrix ignoring the other elements.

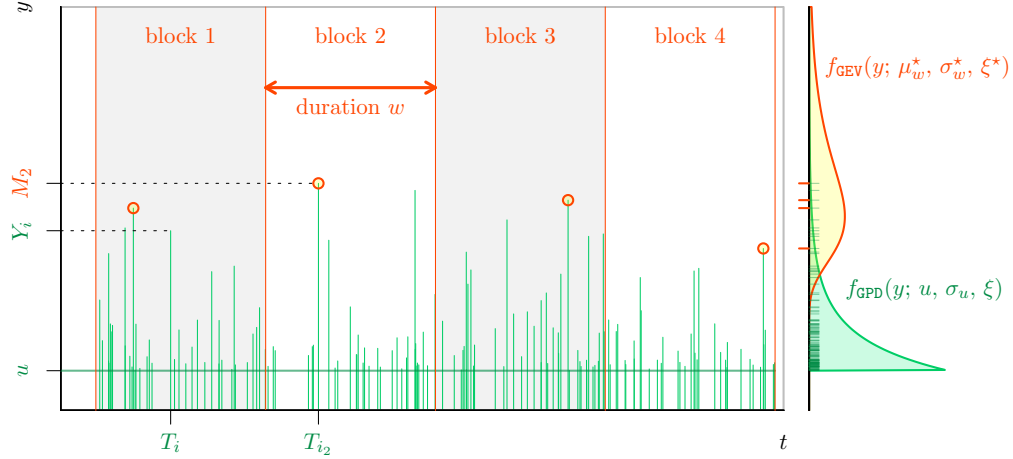


Figure 1: Block maxima by aggregation of the Poisson-GP.

## 4 Parameterisations for Peaks Over Threshold models

Peaks Over Threshold (POT) models are very popular in EVA. These models relate to marked Poisson Process: at each time  $T_i$  in a sequence or random times  $T_1 < T_2 < \dots$  we observe a random variable  $Y_i$  called the *mark*. In applications, the time  $T_i$  is typically for the occurrence a storm and the mark  $Y_i$  can be an amount of precipitation or a sea levels. The following two frameworks are commonly used, see Northrop et al. (2016) for more details.

- The *Poisson-GP* framework involves a given threshold  $u$  having the same physical dimension as the marks  $Y_i$ . It assumes that the  $T_i$  form an homogeneous Poisson Process with rate  $\lambda_u$ , and that  $Y_i$  are i.i.d. with a Generalised Pareto distribution  $\text{GPD}(u, \sigma_u, \xi)$  or equivalently that the excesses over the threshold  $Y_i - u$  follow the two-parameter GP distribution with scale  $\sigma > 0$  and shape  $\xi$ . The marks  $Y_i$  are further assumed to be independent of the event process  $\{T_i\}_i$ . The parameters form the vector  $\theta_u = [\lambda_u, \sigma_u, \xi]$ .
- An alternative framework is often named *Point Process* (PP) or *Non-Homogeneous Poisson Process* (NHPP). It involves a reference duration  $w > 0$ , usually chosen to be one year, and a vector of parameters  $\theta^* := [\mu_w^*, \sigma_w^*, \xi^*]$ . The random observations  $[T_i, Y_i]$  are given by a Poisson process on the  $(t, y)$ -plane with intensity

$$\gamma_w^*(t, y) := \frac{1}{w} \times \frac{1}{\sigma_w^*} \left[ 1 + \xi^* \frac{y - \mu_w^*}{\sigma_w^*} \right]^{-1/\xi^* - 1} 1_{\mathcal{S}_{\theta_w^*}}(t, y),$$

where  $\mathcal{S}_{\theta_w^*}$  is the domain of the plane

$$\mathcal{S}_{\theta_w^*} := \{[t, y] : \sigma_w^* + \xi^*[y - \mu_w^*] > 0\}.$$

This is a half-plane for  $\xi^* \neq 0$  and the whole plane for  $\xi^* = 0$ .

If we take the observation region for the PP model as being the product of the time interval  $(0, t^\dagger)$  by the  $y$ -interval  $(u, \infty)$  where the threshold  $u$  is such that  $\sigma_w^* + \xi^*[y - \mu_w^*] > 0$ , we get the

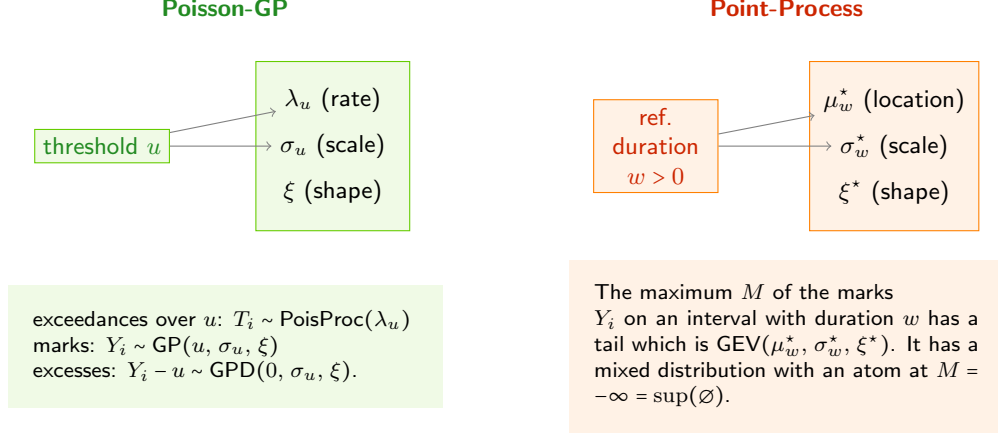


Figure 2: Two parameterisations for POT models.

same model as the Poisson-GP on  $(0, t^\dagger)$ , up to a re-parameterisation. As an interesting feature of the PP parameter  $\theta_w^*$ , it does not depend on the threshold  $u$ . It relates to the distribution of the maximum  $M$  of the marks  $Y_i$  corresponding to a time interval with duration  $w$ , see Figure 4. This distribution is indeed essentially the  $\text{GEV}(\mu_w^*, \sigma_w^*, \xi^*)$ , up to an atom corresponding to the possibility that no mark is observed during the time interval. We can then define  $M := -\infty$  since this is arguably the maximum of the empty set, and the probability of the corresponding event is  $\exp\{-\lambda_u w\}$ . We may speak of  $\text{GEV}(\mu_w^*, \sigma_w^*, \xi^*)$  as the *GEV reference distribution* in relation to  $w$ , although this is not exactly the distribution of a maximum  $M$  over the reference duration.

For a given threshold  $u$  and a given reference duration  $w > 0$ , the one-to-one correspondence between the vectors  $\theta_u$  and  $\theta_w^*$  is given by

$$\begin{cases} \mu_w^* &= u + \frac{(\lambda_u w)^\xi - 1}{\xi} \sigma_u, \\ \sigma_w^* &= (\lambda_u w)^\xi \sigma_u, \\ \xi^* &= \xi, \end{cases} \quad (2)$$

the fraction  $[(\lambda_u w)^\xi - 1]/\xi$  of the first equation being to be replaced for  $\xi = 0$  by its limit  $\log(\lambda_u w)$ . The reciprocal transformation is

$$\begin{cases} \sigma_u &= \sigma_w^* + \xi^* [u - \mu_w^*], \\ \lambda_u &= w^{-1} [\sigma_u / \sigma_w^*]^{-1/\xi^*}, \\ \xi &= \xi^*, \end{cases} \quad (3)$$

where the second equation becomes  $\lambda_u = w^{-1}$  for  $\xi^* = 0$ .

When a vector  $\mathbf{x}$  of covariates can be used, two different kinds of so-called *non-stationary* POT models can be obtained by relating or “linking” the three parameters to the covariates, either parametrically or non-parametrically. For instance the Poisson-GP scale or its logarithm can be specified as having the parametric form  $\mathbf{x}^\top \boldsymbol{\beta}^\sigma$  where  $\boldsymbol{\beta}^\sigma$  is a vector of parameters. Different forms of models arise from the two frameworks. Moreover, parametric Poisson-GP models relate to a specific threshold since the same form of link can not persist when the threshold is changed. Remind also that the threshold should in general depend on the covariates. Anyway, for non-stationary models

it is often useful to transform one of the two parameterisations into the other. If a Poisson-GP model is used, we may be interested in the GEV reference distribution conditional on a given value  $\mathbf{x}$ . When instead a PP model is used, it may be useful to investigate the relation of the implied rate  $\lambda_u$  with the covariates, and possibly to compare the POT estimate with a non-parametric estimate.

The **yaev** package provides the two transformations (2) and (3) required for all these tasks – along with their Jacobian, under the names **poisGP2PP** and **PP2poisGP**. As for the probability functions, the singularity for  $\xi = 0$  or  $\xi^* = 0$  is coped with by using a second-order Taylor approximation. As often required when coping with non-stationary POT models, the functions **poisGP2PP** and **PP2poisGP** are vectorised w.r.t. their arguments including **threshold** for the **PP2poisGP** transformation. Each element  $i$  in the provided vector arguments (such as **lambda**) correspond to a value  $\mathbf{x}_i$  of the covariates, and most often the threshold is then also chosen as depending on the covariates, hence used via a vector with  $n$  elements  $u_i = u(\mathbf{x}_i)$ . The condition

$$\sigma_i^* + \xi_i^*[u_i - \mu_i^*] > 0, \quad i = 1, \dots, n$$

should then hold.

## 5 EV distributions from other R packages

The EV distributions are implemented in many R packages. A variety of strategies regarding the problem  $\xi = 0$  can be found. We now describe these strategies and provide for each of them a “code” (shown as framed: NT, ...) that is used in Table 1, columns  $\xi = 0$ . Each strategy is briefly discussed.

1. NT Use only the formula for  $\xi \neq 0$  i.e., “do nothing”. In practice, an optimisation or sampling algorithm will never come to the case  $\xi = 0$  *exactly* and this can only happen when the user gives this value e.g., as an initial value. There will be some numerical problems when  $\xi$  is very small, say  $\xi = 1\text{e-}14$  or less. These problems are not so crucial for the usual probability functions: we get some wiggling when plotting the curves and zooming. Mind however that the random generation **rgev** or **rgpd** will produce silly results with a very small  $\xi$  if they use **qgev** or **qgpd**.
2. 0.0 Test the exact equality  $\xi = 0$ , and if this is true, *switch to the exponential/Gumbel formula*. This helps only when the user gives **xi** = 0.0, but we are essentially doing the same thing as in NT.
3.  $\epsilon/\text{S}$  Test the equality  $|\xi| \leq \epsilon$  where  $\epsilon > 0$  is very small, and if this is true, *switch to the exponential/Gumbel formula*. So this produces a (very small) discontinuity. E.g., **Renext** uses  $\epsilon \approx 2\text{e-}14$ .
4.  $\epsilon/\text{AI}$  Test the equality  $|\xi| \leq \epsilon$  where  $\epsilon > 0$  is very small, and if this is true, use a dedicated *approximation or interpolation*. Several methods can be used including Taylor approximations. The discontinuity should then be undetectable. Mind the probability functions although not being *analytic* functions, are infinitely differentiable  $C^\infty$  w.r.t. the parameters.

Note that some of the cited packages are quite old: **evir** (Pfaff and McNeil, 2018), **evd** (Stephenson, 2002), **ismev** (Stephenson and Heffernan, 2018), **Renext** (Deville and IRSN, 2022), **POT** (Ribatet and Dutang, 2022) and **SpatialExtremes** Ribatet (2022). The packages **revdbayes** (Northrop,

Package	GEV					GPD				
	Lang.	Vec. $\theta$	Grad.	Hess.	$\xi = 0$	Lang.	Vec. $\theta$	Grad.	Hess.	$\xi = 0$
<b>evir</b>	R	no	no	no	NT	R	no	no	no	NT
<b>evd</b>	R	no	no	no	0.0	R	no	no	no	0.0
<b>ismev</b>	R*	no	no	no	NT	R*	no	no	no	NT
<b>Renext</b>						R	no	yes	yes	$\epsilon/S$
<b>POT</b>						R*	yes	no	no	0.0
<b>SpatialExtremes</b>	R	yes	no	no	0.0	R	yes	no	no	0.0
<b>revdbayes</b>	R	yes	no	no	$\epsilon/AI$	R	yes	no	no	$\epsilon/AI$
<b>mev</b>	R	yes	yes*	yes*	NT	R	yes	yes*	yes*	NT
<b>yaev</b>	C	yes	yes	yes	$\epsilon/AI$	C	yes	yes	yes	$\epsilon/AI$

Table 1: Features of some CRAN packages. *Lang.*: the implementation language, *Vec.  $\theta$* : vectorised w.r.t. the parameters. The columns *Grad.* and the *Hess.* indicate if the gradient and Hessian are provided, and the columns  $\epsilon = 0$  indicate the strategy used to cope with a zero or small shape, as described in the text. A star  $\star$  means that the functions are not exported.

2022) and **mev** (Belzile et al., 2022) are more recent. See the CRAN Task View on Extreme Value Analysis (Dutang, 2022) for an extended list of packages devoted to EVA. Also it is worth mentioning that the **extRemes** package (Gilleland and Katz, 2016) optionally uses the exact gradient of the log-likelihood for models with GEV and GP margins but the derivatives are coded (in R) only for internal use in optimisation tasks.

## Acknowledgments

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## References

- Belzile, L., Wadworth, J. L., Northrop, P. J., Grimshaw, S. D., Jin, Z., A., S. M., Owen, A. B., and Huser, R. (2022). **mev**: *Modelling Extreme Values*. R package version 1.14.
- Coles, S. (2001). *An Introduction to Statistical Modelling of Extreme Values*. Springer.
- Deville, Y. and IRSN (2022). **Renext**: *Renewal Method for Extreme Values Extrapolation*. R package version 3.1-3.
- Dutang, C. (2022). CRAN Task View: Extreme Value Analysis. Version 2022-08-04.
- Gilleland, E. and Katz, R. W. (2016). **extRemes 2.0**: An Extreme Value Analysis Package in R. *Journal of Statistical Software*, 72(8):1–39.
- Northrop, P. J. (2022). **revdbayes**: *Ratio-of-Uniforms Sampling for Bayesian Extreme Value Analysis*. R package version 1.4.9.

- Northrop, P. J., Jonathan, P., and Randell, D. (2016). Threshold Modelling of Non-Stationary Extremes. In Dey, D. and Yan, J., editors, *Extreme Value Modeling and Risk Analysis*. Chapman & Hall.
- Pfaff, B. and McNeil, A. (2018). *evir: Extreme Values in R*. R package version 1.7-4.
- Ribatet, M. (2022). *SpatialExtremes: Modelling Spatial Extremes*. R package version 2.1-0.
- Ribatet, M. and Dutang, C. (2022). *POT: Generalized Pareto Distribution and Peaks Over Threshold*. R package version 1.1-10.
- Stephenson, A. G. (2002). **evd**: Extreme value distributions. *R News*, 2(2):0.
- Stephenson, A. G. and Heffernan, J. E. (2018). *ismev: An Introduction to Statistical Modeling of Extreme Values*. R package version 1.42.