# Scalable BFAST: R package optimizations and arraybased data management

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### **Topics**

### Part I: BFAST R package optimizations

Part II: Scalable EO data management with SciDB

Part III: Hands-on with SciDB, Landsat, and BFAST

- 1. SciDB installation (with Docker)
- 2. Data ingestion
- 3. Analysis (practical part)

### Online Material

Slides, reports, and tutorial available at:

- https://github.com/appelmar/scalbf-wur/
- https://appelmar.github.io/scalbf-wur/

### Overview

- Change detection and monitoring are computationally intensive operations:
  - Example from bfastSpatial package with the TURA dataset [1,2]
     (148 x 143 pixels and 166 images, approx. 4x4 km, 7 MB) takes around 20 mins on this computer
- → Change detection / monitoring on national scale?
- → Optimize R package and enable scalable distributed processing of large areas

[1] DeVries, B., Verbesselt, J., Kooistra, L., & Herold, M. (2015). Robust monitoring of small-scale forest disturbances in a tropical montane forest using Landsat time series. Remote Sensing of Environment, 161, 107-121.

## BFAST R package

### bfast()

#### Main input:

- time series of response variable
- seasonal model
- type of fluctuation process
- maximum number of breaks / minimum segment size between breaks

#### Main output:

 time, number, and magnitude of changes

### bfastmonitor()

#### Main input:

- time series of response variable (and optional regressors)
- start of the monitoring period
- model formula
- type of monitoring process: OLS-CUSUM, OLS-MOSUM, RE, ME
- model for identifying a stable history
   (Reverse CUSUM, Bai and Perron, all, ...)

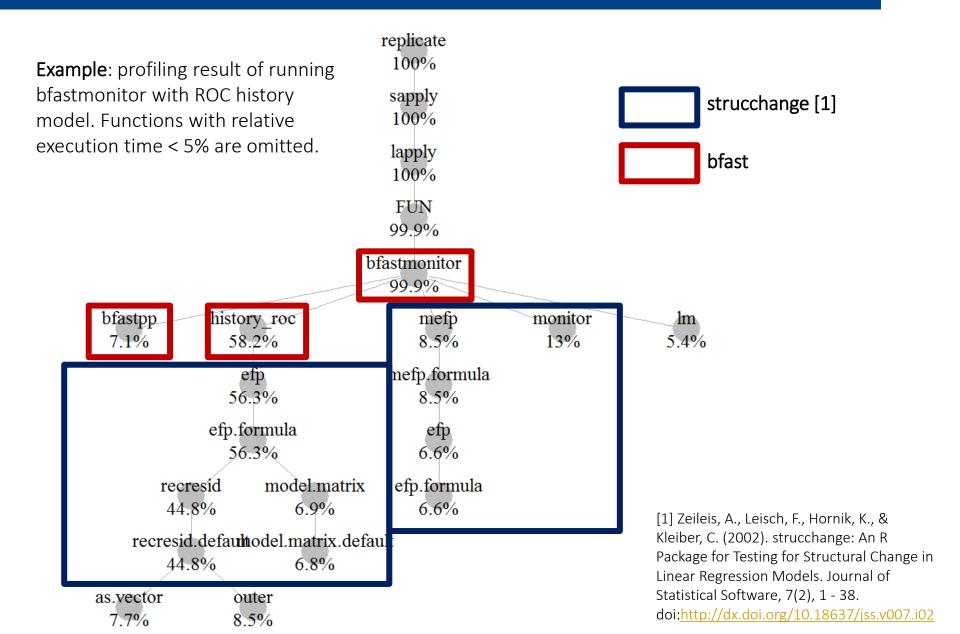
#### Main output:

change magnitude, change date (if any)

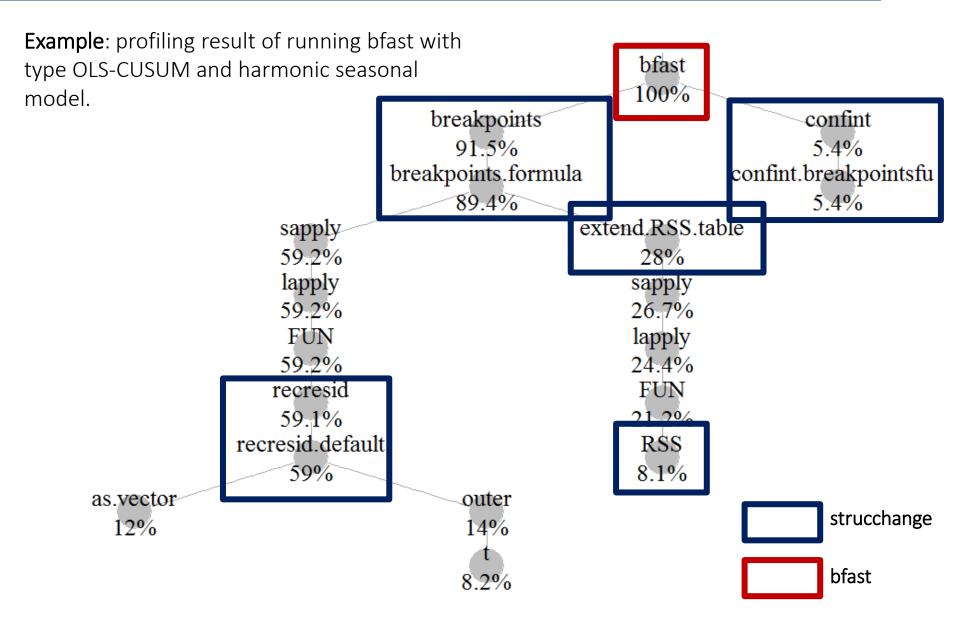
### Identifying computational bottlenecks with profiling

- Profiling = executing R expressions and regularly checking which function is currently being evaluated (e.g. every 5 ms)
- Counting how often the execution is within a specific function allows to estimates how much of the overall time that function takes
- For each sample, the full call stack is available
- Profiling in R: Rprof, lineprof, profr

## Profiling: example for bfastmonitor()



## Profiling: example bfast()



### Package optimizations (overview)

 Move operations to C++ with Rcpp and RcppArmadillo [1,2] if possible

```
> recresid, extend_rss_table, efp_process_me, efp_process_re
```

- Use design matrix and response vector instead of data.frame and formula and lm.fit() instead of lm()
  - bfastpp, bfastmonitor, bfast
  - breakpoints, efp, mefp, monitor,
- bfastts: avoid as.ts(zoo(...))

[1] Conrad Sanderson and Ryan Curtin. *Armadillo: a template-based C++ library for linear algebra*. Journal of Open Source Software, Vol. 1, pp. 26, 2016.

### Moving computations to C++: recursive residuals

#### Recursive residuals in C++ with RcppArmadillo

#### Recursive residuals in R

```
arma::vec sc cpp recresid arma(const mat& X, const vec& y, unsigned int
                                                                                recresid.default <- function(x, y, start = ncol(x) + 1, end = nrow(x),
start, unsigned int end, cons
                                         _const double& rcond min)
                                                                                 tol = sqrt(.Machine$double.eps)/ncol(x), ...)
 if(!(start > X.r
 if(!(end >= star
                                                                                            nd data dimensions
                    vec sc cpp recresid arma (const mat&
 --start;
                                                                                            tart > ncol(x) & start <= nrow(x))
 --end;
                                                                                            nd >= start & end <= nrow(x))
                                 X, const vec& y,
 int n=end;
 int q=start-1;
 int k = X.n cols,
 vec rval = vec(n-q, fill::zeros);
                                                                                           betar <- betar + X1 %*% xr *
                cur coef += X1 * trans(xr) *
 mat X1
                                                                                                 rval[r-q-1] * sqrt(fr)
 colvec
                     rval(r-q-2) * sqrt(fr);
                                                                                 rval[1] \leftarrow (y)
                                                                                                               betar)/sqrt(fr)
 mat xr = X.row(q+1)
                                                                                 check <- TRUE
 vec fr = (1 + xr *
                                                                                  if((q+1) < n
 rval(0) = as scala
                               xr * cur coef)/sqrt(fr));
 bool check = true
                                                                                   for(r in
 if((q+1) < n)
                                                                                                (!is.na(fm$coefficients))
   for (int r=q+2;
                                                                                              - (X1 %*% outer(xr, xr) %*% X1)/fr
                   ans(xr) * xr * X1)/as_scalar(fr);
                                                                                     betar <- betar + X1 %*% xr * rval[r-q-1] * sqrt(fr)
     cur coef += 1 \times 1 \times trans(xr) \times rval(r-q-2) \times sqrt(fr);
                                                                                     if(check) {
     if (check) {
                                                                                       y1 < - y[1:(r-1)]
       cur y = y.subvec(0, r-1);
                                                                                       fm \leftarrow lm.fit(x[1:(r-1), drop = FALSE], y1)
       cur X = X.submat(0, 0, r-1, k-1);
                                                                                       nona <- nona & all(!is.na(betar)) & all(!is.na(fm$coefficients))
                                                                                       if (nona && isTRUE(all.equal(as.vector(fm$coefficients), as.vector(betar),
       solve(cur coef full, cur X, cur y, solve opts::no approx);
                                                                               tol = tol))) {
       inv sympd(X1, trans(cur X) * cur X);
                                                                                         check <- FALSE
       bool nona = is finite( cur coef) && is finite( cur coef full );
       if(nona && approx equal(cur coef full,cur coef, "absdiff", tol)) {
                                                                                       X1 < - .Xinv0(fm)
         check = false;
                                                                                       betar <- .coef0(fm)
       cur coef = cur coef full;
                                                                                     xr <- as.vector(x[r,])</pre>
                                                                                     fr <- as.vector((1 + (t(xr) %*% X1 %*% xr)))
     xr = X.row(r); // This a a row
                                                                                     rval[r-q] \leftarrow (y[r] - sum(xr * betar, na.rm = TRUE))/sqrt(fr)
     fr = (1 + xr * X1 * trans(xr));
     rval(r-q-1) = as scalar((y(r) - xr * cur coef)/sqrt(fr));
                                                                                 return(rval)
 return rval;
```

### Moving computations to C++: recursive residuals

- fits linear models in the first few iterations, until recursive model parameter updates are stable
- for ill-conditioned systems (e.g. dummy variables), R's lm.fit is used, otherwise Armadillo's solve
- speedup varies with length of time series, number of explanatory variables, and stability of the system, mostly around 15-20 on this machine
- changes have effect on performance of both bfast and bfastmonitor (unless history = "all")

## bfastpp

```
bfastpp(data, order = 3,
  lag = NULL, slag = NULL, na.action = na.omit,
  stl = c("none", "trend", "seasonal", "both"),
  formula = NULL)
```

Derives variables that occur in the formula only

- Instead of a data.frame, modified output is a list with elements
  - y: response vector
  - X: design matrix
  - t: vector of dates (same as time (data))

### Avoiding formulas and data.frames

### S3 methods for matrix input:

• breakpoints, efp, mefp, monitor



```
breakpoints.matrix(\mathbf{X}, \mathbf{y}, h = 0.15, breaks = NULL, hpc = c("none", "foreach"), ...)
```

#### Matrix methods

- avoid model.frame() and model.matrix()
- use lm.fit() instead of lm()

### bfastts modifications

 bfastts() not directly called from bfast() or bfastmonitor() but often used before (e.g. in bfmSpatial())

 takes up to 30% of computation time of bfastSpatial example with the TURA dataset

modification does not use zoo

around 2 -3 times faster

## Using package modifications

Install strucchange and bfast from github (Rtools needed for Windows)

```
library(devtools)
install_github("appelmar/strucchange")
install_github("appelmar/bfast")
```

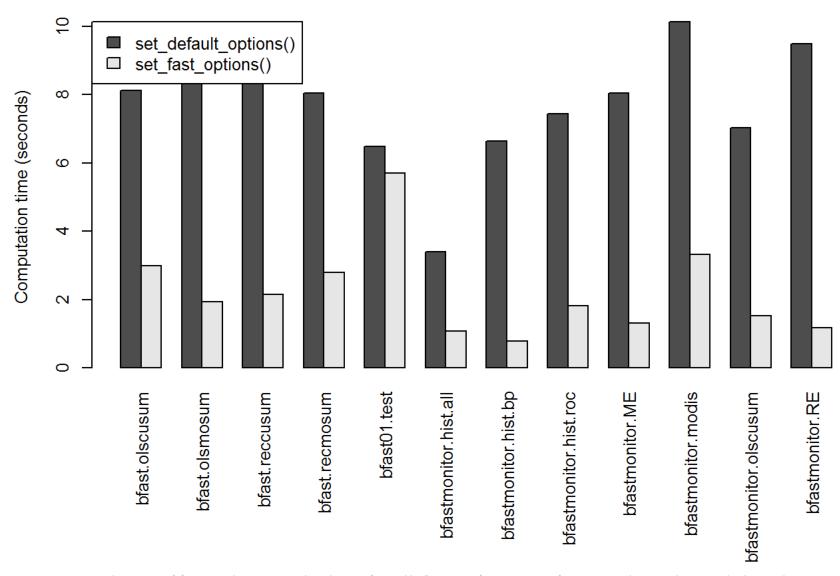
2. Load the package and enable modifications

```
library(bfast)
set_fast_options() # use modifications
...
set_default_options() # use default implementation
...
```

# Package Options

Package option	Description
strucchange.use_armadillo	Defines whether or not C++ functions should be used if available
strucchange.armadillo_rcond_min	For ill-conditioned systems in the recursive residual computation, the minimum reciprocal conditioning number to use armadillo solve instead of column pivoting QR from R
bfast.prefer_matrix_methods	Defines whether or not bfastpp generates a data frame or a matrix
bfast.use_bfastts_modifications	Defines whether or not bfastts modifications should be used

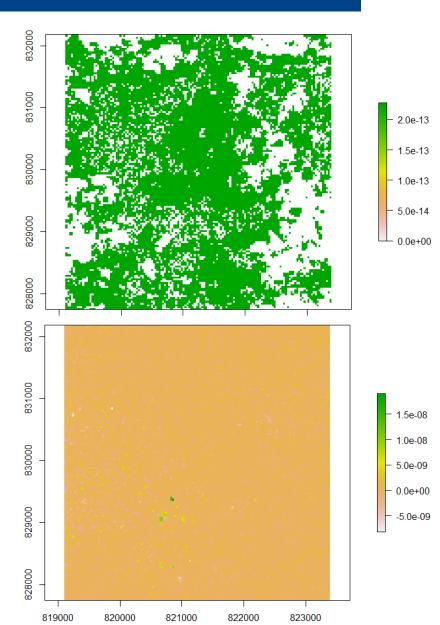
## Results: speedup



see <a href="https://appelmar.github.io/scalbf-wur/reports/report.benchmark.html">https://appelmar.github.io/scalbf-wur/reports/report.benchmark.html</a>

### Example: bfastSpatial with tura dataset

```
library(bfastSpatial)
data(tura)
set default options()
system.time(bfm.tura.new <-</pre>
    bfmSpatial(tura, start=c(2009, 1),
    history = "ROC")) \# 1318.52 s
set fast options()
system.time(bfm.tura.new <-</pre>
    bfmSpatial(tura, start=c(2009, 1),
    history = "ROC")) \# 288.75 s
plot(bfm.tura.new$breakpoint -
     bfm.tura.reference$breakpoint)
plot(bfm.tura.new$magnitude -
     bfm.tura.reference$magnitude)
```



### Package reports

The package comes with some R Markdown reports to test, benchmark, and profile the modifications.

```
library(rmarkdown)

render(system.file("reports/report.test.Rmd",
    package = "bfast"), output_file = "report.test.html")

render(system.file("reports/report.benchmark.Rmd",
    package = "bfast"), output_file = "report.benchmark.html")

render(system.file("reports/report.profiling.Rmd",
    package = "bfast"), output_file = "report.profiling.html")
```

#### Examples:

- https://appelmar.github.io/scalbf-wur/reports/report.test.html
- https://appelmar.github.io/scalbf-wur/reports/report.benchmark.html
- https://appelmar.github.io/scalbf-wur/reports/report.profiling.html

### Summary and conclusions

- speedup of bfastmonitor and bfast varies between 2 and 10, depending on computations
- There is no single computational bottleneck, probably difficult to employ e.g. GPUs
- Strongest speedup in bfastmonitor with (type="RE", Bai and Perron history (history = "BP")
- Optimizations also work with bfastSpatial even with parallelization
- bfast and strucchange now need to be compiled and linked to Rcpp and RcppArmadillo respectively

# Thank you

Questions?