Using LIKELTD

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NULL

1 Code usage

Computing a likehood model takes place in three-parts: (i) creation of the hypothesis, (ii) creation of the model itself from the hypothesis, (iii) maximizing the likehood function over the set of nuisance parameters.

1.1 Creating a hypothesis

A hypothesis consists of all the parameters are used to describe the model, e. g. the known profiles to include, the number of unprofiled contributors, and whether to include dropin. A proper description of the hypothesis should be independent of other information, including neither information about the maximization of the likelihood nor any information which needs to be computed.

```
require(likeLTD)
require(DEoptim)
# Case we are going to be looking at.
caseName = 'hammer'
datapath = file.path(system.file("extdata", package="likeLTD"),
                     caseName)
args = list(
  databaseFile = NULL,
             = file.path(datapath, 'hammer-CSP.csv'),
  cspFile
               = file.path(datapath, 'hammer-reference.csv'),
 refFile
               = 0,
 nUnknowns
  doDropin
               = TRUE,
               = "EA1",
  ethnic
  adj
               = 1.0,
  fst
               = 0.02,
 relatedness = c(0, 0)/4
# Create hypothesis for defence and prosecution.
defenceHyp = do.call(defence.hypothesis, args)
prosecuHyp = do.call(prosecution.hypothesis, args)
```

Two methods are provided to facilitate the creation of a hypothesis from a common minimal set of input parameters, defence.hypothesis for the defence

and prosecution.hypothesis for the prosecution. These two methods read the allele database, the known profiles, and the crime-scene profile from file. They also automate sensible default decisions about the input, (determining which known profiles need be subject to dropout; in the defence hypothesis the queried individual is replaced by an unprofiled contributor; in the prosecution case, relatedness is set to zero.) These default choices may be further modified by the user at this point. These methods return lists with all the input needed for model execution.

```
> defenceModel <- create.likelihood(defenceHyp)
> prosecuModel <- create.likelihood(prosecuHyp)</pre>
```

create.likelihood returns a method which takes as arguments the nuisance parameters and computes the full weight of evidence, e. g. the product of the likelihoods and penalties associated with each locus.

The function above returns a scalar which represents the weight of evidence for the given values of the nuisance parameters. One could then use defence-Model to perform an optimisation or to create a plot with respect to various arguments. For instance, the following leads to Fig. 1:

```
require(ggplot2)
>
   require(scales)
>
    # Function that winnows down to a single value
    scalarWoE <- function(x) {</pre>
       defenceModel(locusAdjustment=list(D3=0.983, vWA=1.010,
+
                                          D16=1.028, D2=1.072,
                                          D8=1.020, D21=0.930,
                                          D18=0.850, D19=0.932,
                                          TH01=1.041, FGA=0.916),
                    dropout=c(0.5072, 1e-8),
                     degradation=c(10^-2.27, 10^-2.74, 10^-2.47),
                    rcont=c(x, 1e-8, 1),
                     dropin=1.0216,
                    power=-4.4462)
   }
   x = 0:30/30 * 3e0
   data = data.frame(x=x, y=sapply(x, scalarWoE))
   plots <- ggplot(data, aes(x=x, y=y))</pre>
                geom_line()
                xlab("Relative contribution of Victim 1") +
```

```
+ ylab("Weight of Evidence") +
+ scale_y_log10(
+ labels=trans_format("log10", math_format(10^.x)))
> print(plots)
```

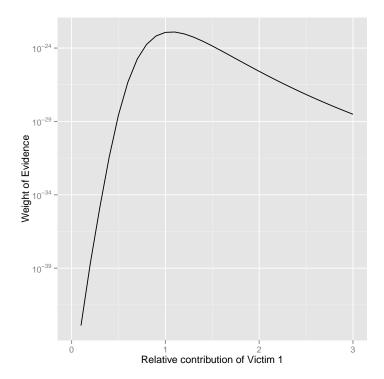


Figure 1: Logarithmic plot of the weight of evidence versus the relative contribution from "Victim 1". The likelihood is for the Hammer case, with one unprofiled contributor ("X"), and including dropin. The fixed parameters are given in Tab. 1.

1.2 Maximizing the likehood

Once we have a likelihood method, it is possible to use the stats package to maximize it. However, the likelihood method takes several arguments (rcont, degradation, etc), whereas DEoptim::DEoptim expects a methods which takes only a vector as argument. Hence, we need to transform our method into the form the optimisation method expects:

```
> skeleton = initial.arguments(defenceHyp)
> vector.model <- function(x) {
+ args <- relist(x, skeleton)
+ args[["degradation"]] = 10^args[["degradation"]]
+ result <- do.call(defenceModel, args)
+ log10(result)</pre>
```

```
+ }
> # Call vector.model with vector argument.
> arguments = skeleton
> arguments[["degradation"]] = log10(arguments[["degradation"]])
> vector.model( as.vector(unlist(arguments)) )
```

The new method vectorModel achieves three objectives: (i) it recreates the list of arguments for defenceModel, (ii) it transforms the degradation parameter from an exponential form, (iii) it takes the logarithm of the weight of evidence. The last two points make optimisation somewhat easier. We can now apply the optimisation methods on vectorModel.

```
require(stats)
   # define upper and lower bounds for constrained maximization
>
   nloci = ncol(defenceHyp$cspProfile)
>
   upper = list(locusAdjustment = rep(1.5, nloci),
                                = c(1-1e-3, 1-1e-3),
                 dropout
                 degradation
                                 = rep(-1e-3, 3),
                 rcont
                                 = rep(100, 2),
+
                 dropin
                                 = 1,
                            = -2 )[names(arguments)]
                 power
   lower = list(locusAdjustment = rep(0.5, nloci),
                 dropout
                                 = c(1e-3, 1e-3),
+
                                 = rep(-20, 3),
                 degradation
                 rcont
                                 = rep(1e-3, 2),
                 dropin
                                 = 1e-3,
                 power
                            = -6 )[names(arguments)]
>
   # perform maximization
>
   result <- DEoptim(fn = vector.model,
                    upper = unlist(upper),
                    lower = unlist(lower),
                    control = list(strategy=3, itermax=500)
+
   opti = relist(result$optim$bestmem, skeleton)
   cat(sprintf("Resulting Weight of Evidence: 10^%f\n",
                -result$optim$bestval))
[1] -44.7266
```

Resulting Weight of Evidence: 10°Inf

The above calculates the maximum of the likelihood using a differential evolution (DE) algorithm to to perform evolutionary global optimization. The particular flavor of DE algorithm used here allows the user to set upper and lower bounds for parameters. Upon convergence, it returns a list with the optimum and its location. Please see the <code>DEoptim</code> package for description.

The functionality of the above code can be achieved more succinctly through a convenience method provided by LikeLTD optimisation.params. It returns a list of adequate arguments for optim given a hypothesis:

```
> params = optimisation.params(defenceHyp, verbose=FALSE)
```

> params\$control\$itermax=50 # Less strict convergence, for demo purposes.

```
> results <- do.call(DEoptim, params)
```

> arguments <- relistArguments(results\$optim\$bestmem, defenceHyp)</pre>

Running the above yields the parameters in Tab. 1. The last line transforms the linear vector of arguments returned by DEoptim back into a more meaningful list, much as relist did earlier. However, it takes care of some specialized problems with the operation, and should be preferred over relist.

A method for ensuring proper convergence is provided by LIKELTD DEoptimLoop. This calls the external function DEoptim::DEoptim, but crucially compares the optimised result every 50 generations with the previous optimised result, and quits once the relative difference is below the given tolerance.

1.3 Testing

LIKELTDcomes a fairly extensive suite of tests. The tests can be run as part of the installation process, or using the following commands:

```
> library(svUnit)
> library(likeLTD)
> runTest( svSuite("package:likeLTD") )
> Log()
```

Although not shown here, this snippet will print results for each tests. Each should return "OK".

Iteration	: 1 bestvalit: 58.030	118 bestmemit:	1.414272	0.842718	0.868778	1.1701
Iteration	: 2 bestvalit: 57.6550	014 bestmemit:	1.414272	0.842718	0.957482	1.1203
Iteration	: 3 bestvalit: 57.6550	014 bestmemit:	1.414272	0.842718	0.957482	1.1203
Iteration	: 4 bestvalit: 57.6550	014 bestmemit:	1.414272	0.842718	0.957482	1.1203
Iteration	: 5 bestvalit: 57.6550	014 bestmemit:	1.414272	0.842718	0.957482	1.1203
Iteration	: 6 bestvalit: 57.6550	014 bestmemit:	1.414272	0.842718	0.957482	1.1203
Iteration	: 7 bestvalit: 57.6550	014 bestmemit:	1.414272	0.842718	0.957482	1.1203
Iteration	: 8 bestvalit: 56.4883	168 bestmemit:	1.071043	0.969552	0.661102	0.9482
Iteration	: 9 bestvalit: 56.4793	366 bestmemit:	1.069704	0.969552	0.661102	0.9482
Iteration	: 10 bestvalit: 56.479	9366 bestmemit:	1.069704	0.969552	0.661102	0.948
Iteration	: 11 bestvalit: 56.479	9366 bestmemit:	1.069704	0.969552	0.661102	0.948
Iteration	: 12 bestvalit: 56.479	9366 bestmemit:	1.069704	0.969552	0.661102	0.948
Iteration	: 13 bestvalit: 56.219	9777 bestmemit:	1.069704	0.969552	0.661102	0.948
Iteration	: 14 bestvalit: 56.219	9777 bestmemit:	1.069704	0.969552	0.661102	0.948
Iteration	: 15 bestvalit: 56.219	9777 bestmemit:	1.069704	0.969552	0.661102	0.948
Iteration	: 16 bestvalit: 54.760	918 bestmemit:	1.414272	0.842718	0.957482	1.120
	: 17 bestvalit: 54.760		1.414272	0.842718	0.957482	1.120
Iteration	: 18 bestvalit: 54.226	3147 bestmemit:	1.414272	0.842718	0.957482	1.120
Iteration	: 19 bestvalit: 54.226	3147 bestmemit:	1.414272	0.842718	0.957482	1.120
Iteration	: 20 bestvalit: 54.226	3137 bestmemit:	1.414272	0.842718	0.957482	1.120
Iteration	: 21 bestvalit: 48.38	6621 bestmemit:	0.817168	0.944136	1.008518	1.072
Iteration	: 22 bestvalit: 48.382	2249 bestmemit:	0.817168	0.944136	1.008518	1.072
Iteration	: 23 bestvalit: 48.382	2249 bestmemit:	0.817168	0.944136	1.008518	1.072
Iteration	: 24 bestvalit: 48.382	2249 bestmemit:	0.817168	0.944136	1.008518	1.072
Iteration	: 25 bestvalit: 47.21	1077 bestmemit:	0.817168	0.944136	1.008518	1.072
Iteration	: 26 bestvalit: 47.21	1077 bestmemit:	0.817168	0.944136	1.008518	1.072
Iteration	: 27 bestvalit: 47.21	1077 bestmemit:	0.817168	0.944136	1.008518	1.072
Iteration	: 28 bestvalit: 47.21	1077 bestmemit:	0.817168	0.944136	1.008518	1.072
Iteration	: 29 bestvalit: 47.153	3099 bestmemit:	0.817168	0.944136	1.008518	1.205
Iteration	: 30 bestvalit: 47.153	3099 bestmemit:	0.817168	0.944136	1.008518	1.205
Iteration	: 31 bestvalit: 46.808	3181 bestmemit:	0.817168	0.944136	1.008518	1.205
Iteration	: 32 bestvalit: 46.808	3181 bestmemit:	0.817168	0.944136	1.008518	1.205
	: 33 bestvalit: 43.28		0.863960	1.128802	0.935080	0.963
	: 34 bestvalit: 43.28		0.863960	1.128802	0.935080	0.963
Iteration	: 35 bestvalit: 43.247	7553 bestmemit:	0.863960	1.128802	0.935080	0.963
Iteration	: 36 bestvalit: 43.247	7553 bestmemit:	0.863960	1.128802	0.935080	0.963
Iteration	: 37 bestvalit: 43.247	7553 bestmemit:	0.863960	1.128802	0.935080	0.963
Iteration	: 38 bestvalit: 43.247	7553 bestmemit:	0.863960	1.128802	0.935080	0.963
	: 39 bestvalit: 43.247		0.863960	1.128802	0.935080	0.963
	: 40 bestvalit: 42.853		0.863960	1.037335	1.075849	0.963
Iteration	: 41 bestvalit: 42.853	3003 bestmemit:	0.863960	1.037335	1.075849	0.963
	: 42 bestvalit: 41.963		0.863960	1.037335	1.075849	0.963
	: 43 bestvalit: 41.963		0.863960	1.037335	1.075849	0.963
	: 44 bestvalit: 41.963		0.863960	1.037335	1.075849	0.963
Iteration	: 45 bestvalit: 41.963	3643 bestmemit:	0.863960	1.037335	1.075849	0.963
	: 46 bestvalit: 41.963		0.863960	1.037335	1.075849	0.963
Iteration	: 47 bestvalit: 41.848	3468 bestmemit:	0.863960	1.037335	1.075849	0.963
	: 48 bestvalit: 41.848		0.863960	1.037335	1.075849	0.963
Iteration	: 49 bestvalit: 41.848	3468 bestmemit:	0.863960	1.037335	1.075849	0.963
Iteration	: 50 bestvalit: 41.65	7689 bestmemit:	0.745327	0.947655	1.124808	0.994

	Victim 1	Victim 2	X
rcont	1.000	3.542 6	0.262
$\operatorname{degradation}$	$10^{5.68e-09}$	$10^{2.97e-13}$	$10^{1.95e-18}$

Locus Ajustments for each locus								
D3	vWA	D16	D2	D8				
0.745	0.948	1.125	0.995	0.961				
D21	D18	D19	TH01	FGA				
1.030	1.004	1.022	1.193	0.746				