

installation and help

<code>dependencies(prompt=)</code>	obtain supporting packages (with or without interactive prompting)
<code>msession\$setAuth(file=)</code>	enter key for private data (use <code>file</code> to read from disk)
<code>demo(package="matR"); demo2("demo")</code>	list demos; or step through demo named "demo"
<code>data(package="matR")</code>	list metagenome collections prepackaged
<code>?command; library(help="matR"); msession\$debug()</code>	get help and examples for <code>command</code> ; see index of help topics; create debug report

metagenome collections

<code>cc <- collection(x, ..., file=)</code>	create collection with sample IDs as in <code>x</code> or <code>file</code> , and data views specified by ...
<code>samples(cc); projects(cc)</code>	see samples in a collection, or projects appearing (fully or partially) in a collection
<code>views(cc)</code>	see data views included in a collection
<code>names(cc); names(cc) <- x</code>	see sample names; or name or rename samples
<code>groups(cc); groups(cc) <- x</code>	see sample grouping; or group samples of a collection
<code>cc[sub]; as(x, "collection")</code>	subset collection, retaining samples <code>sub</code> ; formally convert matrix <code>x</code> to collection

data views in a collection

<code>view.parameters; view.descriptions; view.defaults</code>	see parameters that define data views and meaning of valid values; or see default views
<code>cc\$raw</code>	extract data view named <code>raw</code> as an ordinary matrix
<code>cc[["raw", full=, plain=]]</code>	extract data view named <code>raw</code> with options
<code>views(cc); viewnames(cc) <- x</code>	see or rename data views of a collection
<code>x <- c(annot=, entry=, level=, source=, hit=)</code>	describe a data view by selecting values for view parameters
<code>cc\$new <- x</code>	add data view named <code>new</code> , as described by <code>x</code> , to an existing collection

metadata

<code>metadata(cc)</code>	all metadata of a collection
<code>metadata(cc)["str"]</code>	metadata fields matching specifier "str"
<code>metadata(cc)["str1", "str2", ..., bygroup=]</code>	fields matching any of given specifiers; NA inserted as needed when <code>bygroup=TRUE</code>
<code>mm <- metadata(x, file=)</code>	retrieve (only) metadata of samples specified by ID by <code>x</code> or <code>file</code>

analysis utilities (x is matrix)

<code>remove.singletons(x, lim.entry=, lim.row=)</code>	replace entries with zero and/or remove rows below a threshold
<code>normalize(x, method=)</code>	apply logarithmic transformation, scale, and center to [0,1]
<code>dist(x, y=, groups=, method=)</code>	compute distances among columns of <code>x</code> ; mean-pairwise among <code>groups</code> ; or from <code>y</code> to each
<code>sigtest(x, groups, test=, fdr.level=, qvalue=)</code>	apply significance <code>test</code> across rows to <code>groups</code> of columns of <code>x</code>
<code>randomize(x, n=, method=, seed=, FUN=)</code>	apply <code>FUN</code> to each of <code>n</code> permutations of <code>x</code> according to <code>method</code>

analysis routines (cc is collection)

<code>boxplot(cc, ..., view=)</code>	summarize distribution of annotations per sample; graphical parameters as in <i>boxplot</i>
<code>pco(cc, comp=, method=, ..., view=, rows=)</code>	compute and plot principal coordinates; graphical parameters as in <i>plot</i> and <i>points</i>
<code>heatmap(cc, ..., view=, rows=)</code>	compute and view heatmap dendrogram; graphical parameters as in <i>gplots::heatmap.2</i>
<code>parcoord(cc, groups=, test=, p.lim=, n.lim=, ..., view=, rows=)</code>	compute and plot <code>n.lim</code> parallel coordinates (or all with p-value less than <code>p.lim</code>)