



R Package **DoE.base** for Factorial Experiments

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Abstract

The R package **DoE.base** can be used for creating full factorial designs and general factorial experiments based on orthogonal arrays. Besides design creation, some analysis functionality is also available, particularly (augmented) half-normal effects plots. In addition to this specific functionality, the package provides convenience features for analyzing experimental designs and the infrastructure for a suite of further packages on designing and analyzing experiments. This infrastructure is available for use also by further design of experiments packages.

Keywords: design of experiments, DoE, factorial designs, **DoE.base**.

1. Introduction

Factorial experiments are very common in industrial experimentation. The most widely spread such experiments use 2-level factors only, but experiments with mixed level factors are also quite common, for example with the 18 run experimental plan proposed by Taguchi ([NIST/SEMATECH 2012](#)). The design and execution of such experiments is often done during everyday work without support from a statistical expert – thus it is important to have a software available that can be safely used by non-experts. At the same time, statisticians are often involved in the more important industrial experiments, and there are many facets to construction of such experiments for which a statistician very much appreciates support from a powerful software. The R package **DoE.base** ([Grömping 2018b](#)) targets both non-experts and statisticians. It is part of a larger package suite containing also the packages **FrF2**, **DoE.wrapper** and **RcmdrPlugin.DoE**, and a fifth supporting package **FrF2.catlg128** ([Grömping 2011b, 2013, 2017b, 2014b,c](#)). All these packages and all packages on which **DoE.base** depends ([Chasalow 2012; Venables 2013; Venables and Ripley 2002; Meyer, Zeileis, and Hornik 2006; Sarkar 2008](#)) are available from the Comprehensive R Archive Network (CRAN), which also holds the software R itself ([R Development Core Team 2018](#)). The graphical user interface (GUI) package **RcmdrPlugin.DoE**, which will not be described in this article, provides access to some functionality from the package suite. [Grömping \(2011b\)](#) gives a detailed

example-based tutorial for using it. Package **DoE.base** provides the infrastructure for the entire package suite, in particular the class ‘`design`’, functions for importing and exporting experimental designs, and simple analysis functions for printing, summarizing, plotting, and modeling design data.

Besides providing infrastructure, the main contribution of package **DoE.base** is to offer features for creating factorial designs: Potentially blocked full factorials (function `fac.design`) and catalog-based general factorials (function `oa.design`) are available. The two functions `fac.design` and `oa.design` have taken inspiration from the “white book” ([Chambers and Hastie 1984](#)), where these S functions are described that never made it into base R. The most advanced contributions of the package are the features around orthogonal arrays (function `oa.design`), which are subject to ongoing research. These rely heavily on a catalog of orthogonal arrays, most of which have been taken from [Kuhfeld \(2010\)](#). To the author’s knowledge, the package is the only place in R where non-regular orthogonal arrays other than Plackett-Burman designs are provided for experimentation. Non-regularity of an array has been discussed to be beneficial for screening experiments because it implies good projectivity properties (see, e.g., [Box and Tyssedal 1996](#); [Deng and Tang 1999](#); [Tang and Deng 1999](#)). This discussion has so far focused on 2-level designs, but should analogously apply to more general factorial designs.

There is another R package closely related to the design creation functionality of package **DoE.base**: The R package **planor** ([Kobilinsky, Bouvier, and Monod 2018](#)) can create regular fractional factorial designs in a general sense (see also [Kobilinsky, Monod, and Bailey 2017](#)). Package **DoE.base** is more general than package **planor** in that it also creates non-regular designs, can calculate various types of quality criteria, and does not require specification of a model but can optimize a design with respect to model robustness criteria. It is less general than **planor** in that it does not allow to specify a model and estimable effects, i.e., it treats all effects of the same order on an equal footing. Sections 4 and 5.3 will illustrate function `regular.design` from package **planor** as an alternative to functions from packages **DoE.base** and **FrF2**.

The remainder of this article is organized as follows: Section 2 briefly explains and exemplifies full factorial designs and orthogonal arrays and explains the basic principles of experimental design. Section 3 presents the mathematical background and terminology for general orthogonal arrays and quality criteria for them. Section 4 discusses creation of full factorial designs, in particular also with the possibility of blocking them. Section 5 provides insights into usage and inspection of the orthogonal arrays implemented in package **DoE.base**. Section 6 discusses design creation and analysis tools, using the example of an experimental design in biotechnology ([Vasilev, Schmitz, Grömping, Fischer, and Schillberg 2014](#)). Section 7 describes in more detail the half-normal plotting functionality provided by package **DoE.base**. Finally, a brief overview of further developments is provided.

2. Basics

2.1. Full factorial designs and designs based on orthogonal arrays

A factorial design is an experimental plan in which k “factors” are systematically varied. The j -th factor has l_j “levels”, $j = 1 \dots k$. If all factors have the same number of levels,

i.e., $l_1 = \dots = l_k$, the design is called a “fixed level” or “symmetric” design, otherwise it is called “mixed level” or “asymmetric”. A “full factorial” design contains (a multiple of) *all* factor level combinations, i.e., a multiple of $l_1 \cdot \dots \cdot l_k$ experimental runs. In a full factorial design, all coefficients for an adequately coded linear model with all main effects, two-factor interactions, ..., up to k -factor interactions are estimable. The number of estimable effects remains the same, regardless of the choice of adequate coding. Section 7 will discuss how the coding affects correlation between coefficient estimates.

Full factorial designs are often not feasible in the real world, if the number of factors or the numbers of factor levels are not very small. For example, a full factorial experiment with one 2-level factor and six 3-level factors requires 1458 runs. There are several possibilities for designs with fewer runs: D -optimal designs require the specification of a model to be estimated; they can be created with R packages **AlgDesign** or **DoE.wrapper** (Wheeler 2014; Grömping 2017b), but are not the topic of this article. Here, we consider experimental designs based on orthogonal arrays: These do not require specification of a model but assume that (i) all effects of the same degree (main effect = degree 1, two-factor interaction = degree 2, etc.) are equally important and (ii) that effects of lower degree are more important than those of higher degree. Orthogonal array designs are often used with the intention of estimating main effects only; they are particularly common for qualitative factors, although they can also be used for quantitative factors. For a design based on an orthogonal array, each factor is contained with the same number of times for each level, and each pair of factors is contained with the same number of times for each pair of levels. Genizi Taguchi provided various orthogonal arrays for engineering experimentation; one of the most-well-known ones is an 18-run array for up to one 2-level factor and up to seven 3-level factors. This array can for example be found in NIST/SEMATECH (2012), and is also contained in package **DoE.base**:

R> L18

```
A B C D E F G H
1 1 1 1 1 1 1 1
2 1 1 2 2 2 2 2
3 1 1 3 3 3 3 3
4 1 2 1 1 2 2 3 3
5 1 2 2 2 3 3 1 1
6 1 2 3 3 1 1 2 2
7 1 3 1 2 1 3 2 3
8 1 3 2 3 2 1 3 1
9 1 3 3 1 3 2 1 2
10 2 1 1 3 3 2 2 1
11 2 1 2 1 1 3 3 2
12 2 1 3 2 2 1 1 3
13 2 2 1 2 3 1 3 2
14 2 2 2 3 1 2 1 3
15 2 2 3 1 2 3 2 1
16 2 3 1 3 2 3 1 2
17 2 3 2 1 3 1 2 3
18 2 3 3 2 1 2 3 1
attr(,"origin")
```

```
[1] "Taguchi"
attr(,"class")
[1] "oa"      "matrix"

further attribute(s) (accessible with attr( res$value , attrname)):
[1] "comment"
```

This small array can already accommodate the above-mentioned experiment with one 2-level factor and six 3-level factors. Of course, as it is very much smaller than the 1458 runs for a full factorial, there is a substantial amount of confounding built into the array. If a small array like the L18 is to be used, two things are very important: picking the best possible columns for the design, and understanding the limitations of the resulting design. Package **DoE.base** can help with both. However, except perhaps for very preliminary investigations, it will usually be preferable to use less severely confounded designs. Package **DoE.base** can also help with optimizing the selection of an array and the column selection within the array. This will be demonstrated in Sections 5 and 6.

Orthogonal arrays may be regular or non-regular. In the regular case, it is possible to describe the array by a few defining relations, similar to the well-known way of doing so for regular fractional factorial 2-level designs: Starting from a full factorial design in some “generating” or “base” factors, additional factors are accommodated by assigning them to interactions between the base factors, which are consequently completely confounded with the new factors’ main effects. This is a little more complicated for factors with more than two levels, but the general principle remains the same. One complication arises from base level factors with non-prime numbers of levels; these can be decomposed into full factorials of factors with prime numbers of levels, so-called “pseudo factors”. The aforementioned catalog of arrays contains quite a few regular arrays. Regular orthogonal designs can also be created using function `regular.design` from package **planor**.

Non-regular orthogonal arrays cannot be described by defining relations. Some of the cataloged arrays in package **DoE.base** are non-regular. Section 5.1 provides details on the catalog and its usage. Note that the catalog is by no means complete; in particular, it is much more difficult to completely enumerate all orthogonal arrays than it is to enumerate all *regular* orthogonal arrays. Partially complete catalogs of orthogonal arrays are available, e.g., from the website by [Eendebak and Schoen \(2010\)](#) based on the algorithm described in [Schoen, Eendebak, and Nguyen \(2010\)](#). In many cases, the web site provides the best arrays only, or does not provide an array at all (in case of very large numbers of arrays). Where a number of arrays is shown, the complete set of arrays can in principle be obtained from Eric Schoen; however, with large numbers of arrays the complete catalogs are so large that it is not easily feasible to work with them.

2.2. Principles of experimental design

A very important principle of experimentation is replication: When comparing two different setups, one will usually not rely on a single instance of each setup, but will replicate each setup a specified number of times. This serves the purpose of making sure that differences are only interpreted if they are sufficiently larger than can be expected from experimental variation. Replication is quite different from repeating measurements only: For a proper replication, all experimental settings have to be redone for each replicate. Sometimes, with

very variable measurement devices, it may make sense not to include replications but to repeat the measurement process only. This is called “repeated measurements” and has to be treated quite differently from proper replication. Several design generation functions of packages **DoE.base** and **FrF2** offer the option `replications` for specifying the number of replications and `repeat.only` for indicating whether these are proper replications (default) or repeated measurements only.

One of the very useful aspects of factorial experiments is *implicit* replication: When experimenting with many factors, one can often expect higher order interactions to be irrelevant. If this is the case, the degrees of freedom that would have to be dedicated to higher order interactions can instead be used for estimating error variation (or for accommodating further experimental factors). Therefore, in factorial experimentation, one will encounter experiments without replicated runs.

A further important principle is blocking, which can be used to control for known influential factors that are not of interest in themselves, like batch-to-batch variation. For an orthogonal array design, one can simply include the block factor as an additional factor and thus has to find an array of the desired structure. A full factorial design can be blocked without increasing the number of runs, by allocating the degrees of freedom for the block factor to portions from interaction effects. This functionality is implemented in function `fac.design` (see Section 4).

Randomization means that the experimental runs are conducted in random order; it is a safeguard against bias from unknown influences. If the run order is completely randomized, all experimental runs can be treated as independent observations, and there is little risk of systematic bias from experimental order or unknown factors related to experimental order or time. In real life, there are sometimes so-called randomization restrictions; for example, experimental runs may be randomized within each block only. Function `fac.design` allows randomization within blocks, while designs created with function `oa.design` have to be re-randomized with function `rerandomize.design` for using one of the factors as a block factor.

Whenever proper replication is used, package **DoE.base** separately randomizes each replication as though it were a block; however, it does not include a block factor for the replications. Users who want to include a block factor for replications in the analysis can obtain such a factor using the function `getblock`. Users who want to change the randomization, i.e., randomize all replications together instead of in separate blocks, can use the function `rerandomize.design`. Using the `[` method for the class ‘`design`’, users can also reorder a design according to a randomization scheme that has been worked out outside of R. Of course, whenever the randomization involves non-trivial restrictions like randomizing in meaningful blocks, the analysis has to be conducted accordingly.

3. General orthogonal arrays

This section provides the mathematical background for general orthogonal arrays, as far as it can be helpful for using the orthogonal arrays available in R package **DoE.base**.

3.1. Terminology for orthogonal arrays

An array in the sense of this article is a rectangular table of numbers with n rows and k columns, like the L18 shown on p. 3. The rows correspond to experimental runs, the columns to experimental factors. In the catalogued arrays in **DoE.base**, the levels of an l -

level factor are denoted by the numbers $1 \dots l$. An array becomes an experimental design by allocating numbers to factor levels. The array is orthogonal, i.e., an OA, if for each pair of columns each combination of levels occurs equally often. If this is the case, main effects of all factors can be estimated separately from each other (provided, no higher order effects are in the model).

An OA is said to be of strength s , if each combination of levels occurs equally often for each subset of s columns. Thus, each OA is at least of strength 2. Strength of an OA is directly related to resolution of an array: Resolution, denoted by roman numerals, is always one higher than the strength, i.e., strength 2 arrays are of resolution III and so forth. For an array of resolution III, main effects are not aliased with main effects, but can be aliased with two-factor interactions (three factors involved); for an array of resolution IV, main effects are not aliased with two-factor interactions, but can be aliased with three-factor interactions, while two-factor interactions can be aliased with other two-factor interactions (four factors involved). This notion is well-known for regular fractional factorial 2-level designs, and is completely analogous for non-regular designs and for designs with factors at more than 2 levels or in mixed level situations. Note that a full factorial in k factors has strength k .

3.2. Generalized word length pattern and refinements

[Xu and Wu \(2001\)](#) introduced the generalized word length pattern (GWLP) for general orthogonal arrays. It is an extension of the well-known word length pattern (WLP) for regular fractional factorial 2-level designs: In the latter, one starts out with a set of base factors and allocates additional factors to interactions among these (the generating contrasts). Coding all main effects model matrix columns with “−1” (one level) and “+1” (the other level), this way of design generation causes products of model matrix columns to be either half “−1” and half “+1”, or constant columns. Factors, whose product of model matrix columns yields a constant column, form a “word” together. The word length pattern is a frequency table of word lengths. For regular fractional factorial 2-level designs, each group of c factors either does or does not form a word, i.e., contributes one or zero to the count for words of length c . This results in a word length pattern with only integer entries.

In general, partial aliasing is possible. Even if there are only 2-level factors, e.g., in a Plackett-Burman design ([Plackett and Burman 1946](#)), a set of factors can contribute a fraction of a word to the GWLP count for the respective word length. Consequently, GWLP entries need not be integers. For example, the GWLP of the L18 is

```
R> GWLP(L18)
```

0	1	2	3	4	5	6	7	8
1.0	0.0	0.0	28.0	52.5	52.5	70.0	33.0	6.0

The GWLP is denoted as $A_0, A_1, A_2, A_3, \dots, A_k$ with A_c the number of generalized words of length c . The entry “1” for A_0 is generic and does not indicate confounding. The GWLP for orthogonal arrays and designs based on them is usually presented starting with A_3 , since orthogonality implies absence of words of lengths one or two. The GWLP coincides with the WLP for regular fractional factorial 2-level designs; for details, consult [Xu and Wu \(2001\)](#) themselves or [Grömping \(2011a\)](#) for a more accessible account. The concepts of strength and resolution directly relate to the (G)WLP: The shortest word length with a non-zero count

is the resolution of the design, the longest word length with a zero count is the strength. Hence, the L18 has resolution III and strength 2. Note that it is not adequate to use the term “generalized resolution” here, because that term is already in use for a different concept that is also implemented in package **DoE.base** (see below).

The GWLP can be used for selecting a best design by comparing designs with respect to their so-called aberration: A design is better than another one, if it has higher resolution; in case of equal resolution, a design has smaller aberration, if its number of shortest words is smaller, in case of ties, successively considering longer words until a difference is encountered. If this principle is applied to a complete set of possible designs, the best design is said to have “generalized minimum aberration” (GMA; analogous to minimum aberration for regular fractional factorial 2-level designs). For orthogonal arrays, one seldom compares complete sets of possible designs; however, the website by [Eendebak and Schoen \(2010\)](#) provides GMA arrays for various scenarii.

The number A_c of generalized words of length c is the sum over contributions from all sets of c factors. For a set of R factors with $l_1 \dots l_R$ levels in a resolution R design (equivalent to $s + 1$ factors in a strength s design), [Grömping and Xu \(2014\)](#) have shown an upper bound for the contribution to the count A_R to be $\min((l_1 - 1), \dots, (l_R - 1))$, i.e., the upper bound for the number of generalized words in a set of R factors depends on the pattern of levels in the set and is given by the main effect degrees of freedom for the factor with the fewest levels (the analogous result for symmetric designs was shown earlier by [Xu, Cheng, and Wu 2004](#)). Furthermore, [Grömping and Xu \(2014\)](#) provided a statistical rationale for the contributions of sets of R factors to A_R in resolution R designs, i.e., the building blocks for the number of shortest words: Each R -set contribution can be seen as the sum of R^2 values from linear models explaining orthogonal contrast columns for any one of the R factors by a full model in the other $R - 1$ factors (provided the factor to be explained has orthogonally coded model matrix columns; otherwise, R^2 values have to be replaced by squared canonical correlations). Thus, the number of shortest words measures the extent of worst case confounding in a plausible way. It is therefore particularly instructive to study R factor sets in resolution R designs. Based on these, Table 1 and Figure 1 illustrate the meaning of words in an informal sense, using mosaic plots of contingency tables as proposed in [Grömping \(2014a\)](#). The main purpose of a mosaic plot in this context is the visualization of balance or the absence thereof; the most critical imbalances occur in case of empty cells, which are visualized by flat lines (present in situations (a) and (b)); the ideal balance occurs in a (replicated) full factorial design, for which a mosaic plot shows equally-sized rectangles only (situation (d)). The mosaic plots in Figure 1 show the first row factor and the column factor of Table 1 in the row and column subdivision (nine approximately square areas that represent 4 runs each) and subdivide this area proportionally to the frequencies of the second row factor of Table 1. Let us consider the top left square for all situations: In situation (a), all four runs belong to level “1” of factor C with both other levels represented by flat lines; in situation (b), all four runs belong to level “2” of factor E , with level “1” represented by a flat line, the allocation in situation (c) is $\frac{3}{4}$ of the runs for level “1” and $\frac{1}{4}$ for level “2” of factor E , and that of situation (d) is completely balanced with half of the runs for each of the two levels of factor D .

The severe imbalance shown in situation (a) of Table 1 and Figure 1 implies that the factor level combination of any pair of factors completely determines the level of the third factor. This is a resolution III regular array, for which each main effect is completely confounded by the two-factor interaction of the other two factors. This is reflected in the number of words,

			Situation (a)						Situation (b)		
			B						C		
A	C		1	2	3		B	E	1	2	3
1	1		4	0	0		1	1	0	4	2
	2		0	0	4		2	4	0	2	
	3		0	4	0						
2	1		0	0	4		2	1	2	0	4
	2		0	4	0		2	2	4	0	
	3		4	0	0						
3	1		0	4	0		3	1	4	2	0
	2		4	0	0		2	0	2	4	
	3		0	0	4						

			Situation (c)						Situation (d)		
			C						C		
B	E		1	2	3		B	D	1	2	3
1	1		3	2	1		1	1	2	2	2
	2		1	2	3		2	2	2	2	2
2	1		1	3	2		2	1	2	2	2
	2		3	1	2		2	2	2	2	2
3	1		2	1	3		3	1	2	2	2
	2		2	3	1		2	2	2	2	2

Table 1: Contingency tables for the 36-run designs of Figure 1.

which coincides with the aforementioned upper bound. The plot shows a triple from the array L36.2.16.3.4 from package **DoE.base**; an analogous plot can be produced from columns 2, 4 and 5 of the well-known Taguchi L18 shown on p. 3. For the other factor sets shown in Table 1 and Figure 1, the upper bound is one because of one degree of freedom only for the 2-level factor in the set. However, with a 2-level and two 3-level factors in an *orthogonal* array, this upper bound cannot be attained. The upper bound can only be attained if the smallest number of levels is a divisor of all the other numbers of levels, which is for example the case in symmetric designs. In summary, the four situations illustrate that more words are related to less balance.

Over and above the GWLP, package **DoE.base** allows to look at individual *R*-factor set confounding through mosaic plots (Grömping 2014a), like the ones shown in Figure 1, and provides an overview of the confounding in *R*-factor projections through projection frequency tables (functions P2.2, P3.3 or P4.4), average R^2 frequency tables (ARFT) and squared canonical correlation frequency tables (SCFT); the latter are based on the results by Grömping and Xu (2014) and detailed in Grömping (2017a). Sometimes, several designs that have GMA can be distinguished further by the more detailed criteria. Grömping and Xu (2014) also introduced a generalization of the generalized resolution *GR* as proposed by Deng and Tang (1999): *GR* refines the resolution *R* by indicating the distance from complete confounding. We have $R \leq GR < R + 1$; the larger *GR*, the less severe the worst case confounding in the design; if *GR* = *R*, there is at least one instance of complete confounding in an *R*-factor set. Furthermore, *GR_{ind}* is a stricter version of *GR* which already becomes equal to *R* if there is a triple of factors for which there is a coding such that at least one degree of freedom of

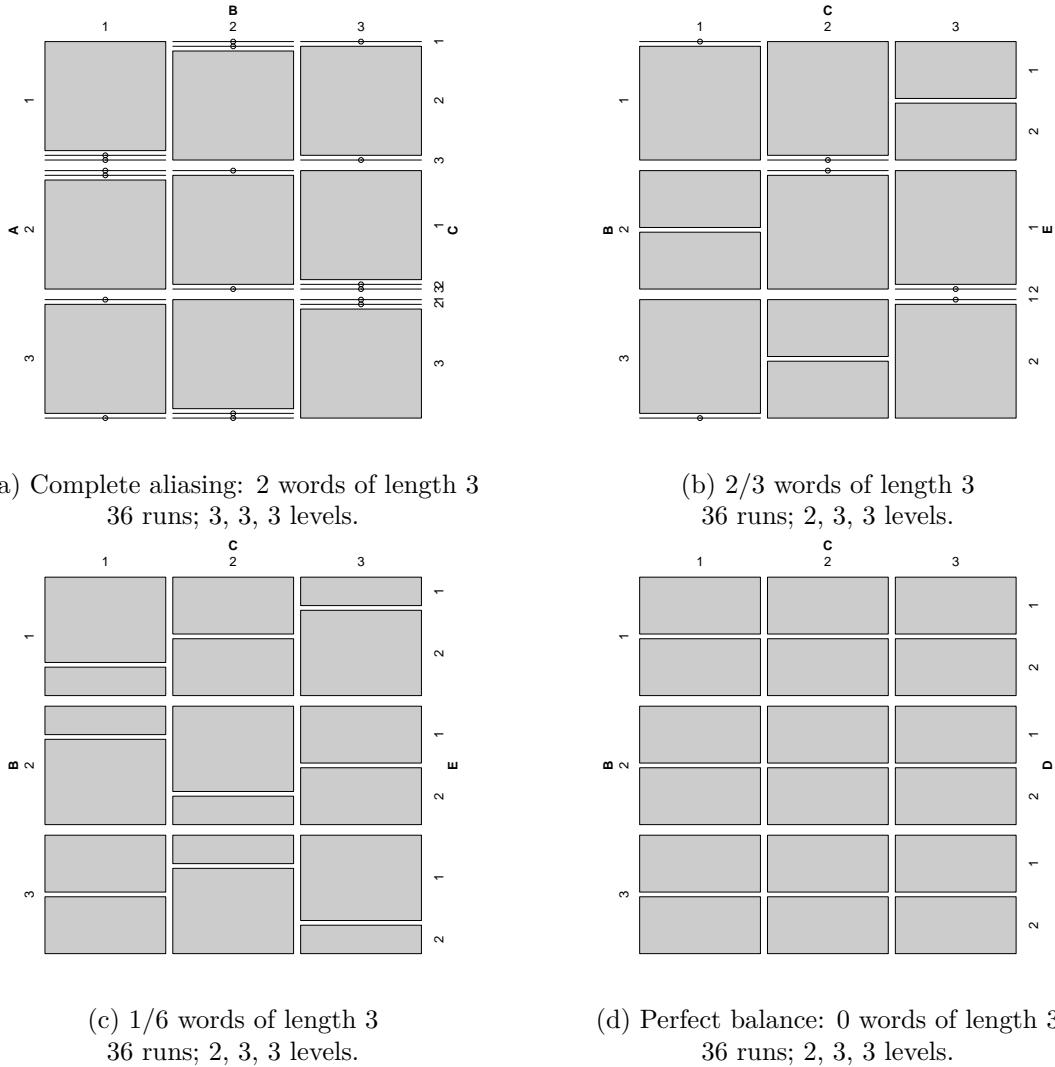


Figure 1: Mosaic plots of different degrees of confounding for triples of factors in 36-run designs.

at least one factor is completely confounded. Besides the overall GR and GR_{ind} , individual factor versions $GR_{\text{tot},i}$ and $GR_{\text{ind},i}$ capture the corresponding worst cases for R factor sets involving the i -th factor.

The GWLP can be obtained with function `GWLP` (or with the older function `lengths`, which is usually slower but performs better for designs with many runs); GR can be obtained with the (old and fast) function `GR` or – together with GR_{ind} , the individual $GR_{\text{tot},i}$ and $GR_{\text{ind},i}$, ARFT and SCFT – with function `GRind`.

4. Full factorial designs with function `fac.design`

Function `fac.design` creates full factorial designs. There is also a simple way in base R for creating all combinations of factor levels: Function `expand.grid` with subsequent random-

ization of the run order will do the job. The benefits of using function `fac.design` lie in the inclusion into the general framework, and in the possibility of automatically blocking designs. The blocking method makes use of the aforementioned pseudo factors: Whenever the number of levels of a factor is not prime, it can of course be factored into primes, e.g., 6 into 2 and 3; thus, a six-level factor can be obtained by the six different factor level combinations of a full factorial in a 2- and a 3-level factor; such component factors are called pseudo factors (e.g., F_1 and F_2 below).

Function `fac.design` uses a method by Collings (1984, 1989) for creating the block factor; in case of automatic blocking, the blocking pattern for several 2- or 3-level factors is taken from optimal blocked catalogs (internal objects `block.catlg` and `block.catlg3`); if a factor contains several pseudo factors with the same prime, its use in block generators ensures that different pseudo factors are used for different block generators involving the factor (where possible). However, the procedure does not ensure overall optimality of the blocking strategy. Neither does function `regular.design` from package `planor`; however, that function may be worth a try if the result from function `fac.design` is not satisfactory, and it can be used for situations outside `fac.design`'s scope for automatic blocking.

The following two code examples exemplify situations for which automated blocking works without or with confounding of two-factor interactions of experimental factors (with a warning in the latter case).

For a full factorial design with six factors with 2, 3, 3, 2, 2 and 6 levels (hence 432 runs), running in six blocks is possible without confounding blocks with two-factor interactions in experimental factors:

```
R> full.factorial.blocked6 <- fac.design(nlevels = c(2, 3, 3, 2, 2, 6),
+     blocks = 6)
R> summary(full.factorial.blocked6)
```

```
Call:
fac.design(nlevels = c(2, 3, 3, 2, 2, 6), blocks = 6)

Experimental design of type  full factorial.blocked
432  runs
blocked design with  6  blocks of size  72

Confounding of  2 -level pseudo-factors with blocks
(each row gives one independent confounded effect):
A B C D E F F
1 0 0 1 1 1 0
```

```
Confounding of  3 -level pseudo-factors with blocks
(each row gives one independent confounded effect):
A B C D E F F
0 1 1 0 0 0 1
```

```
Factor settings (scale ends):
```

	A	B	C	D	E	F
1	1	1	1	1	1	1
2	2	2	2	2	2	2
3	3	3				3
4						4
5						5
6						6

The summary indicates that the design confounds the block factor with the interactions $ADEF_1$ and BCF_2 , where F_1 and F_2 denote two different pseudo factors that make up the six-level factor F . This means, in particular, that there is no confounding of the block factor with two-factor interactions of experimental factors.

Blocking this full factorial in four or nine blocks is also possible, but confounds the block factor with a two-factor interaction among experimental factors, which is signaled by a warning message and is also visible from the summary:

```
R> full.factorial.blocked4 <- fac.design(nlevels = c(2, 3, 3, 2, 2, 6),
+   blocks = 4)
R> summary(full.factorial.blocked4)
```

Call:

```
fac.design(nlevels = c(2, 3, 3, 2, 2, 6), blocks = 4)
```

Experimental design of type full factorial.blocked
 432 runs
 blocked design with 4 blocks of size 108

Confounding of 2 -level pseudo-factors with blocks
 (each row gives one independent confounded effect):

	A	B	C	D	E	F
[1,]	1	0	0	0	1	1
[2,]	0	0	0	1	1	0
[3,]	1	0	0	1	0	0

```
Factor settings (scale ends):
```

	A	B	C	D	E	F
1	1	1	1	1	1	1
2	2	2	2	2	2	2
3	3	3				3
4						4
5						5
6						6

Function `fac.design` allows automatic blocking for the most frequent situations, where most prime level (pseudo) factors have two or three levels, and only single prime level (pseudo)

factors have more than three levels; there is also a limit on the number of 2- and 3-level factors (see the package manual).

We now consider an example, for which block generators have to be manually specified: For blocking a full factorial in one 2-level factor, three 5-level factors and one 10-level factor into 25 blocks, the prime 5 is needed twice for creating the block factor; thus, two block generators for the prime 5 need to be specified. These can be given as a matrix with two rows (one for each block generator) and a column for each prime factor (in the order of factors, and within each factor, in increasing order, i.e., 2, 5, 5, 5, 2, 5 for the present design). The following code yields the desired blocked full factorial for the above requirement.

```
R> BG <- rbind(c(0, 1, 1, 2, 0, 0), c(0, 0, 1, 1, 0, 1))
R> full.factorial.blocked25 <- fac.design(nlevels = c(2, 5, 5, 5, 10),
+     blocks = 25, block.gen = BG)
R> summary(full.factorial.blocked25)
```

Call:

```
fac.design(nlevels = c(2, 5, 5, 5, 10), blocks = 25, block.gen = BG)
```

Experimental design of type full factorial.blocked
 2500 runs
 blocked design with 25 blocks of size 100

Confounding of 5 -level pseudo-factors with blocks
 (each row gives one independent confounded effect):

	A	B	C	D	E1	E2
[1,]	0	1	1	2	0	0
[2,]	0	0	1	1	0	1
[3,]	0	1	2	3	0	1
[4,]	0	1	3	4	0	2
[5,]	0	1	4	0	0	3
[6,]	0	1	0	1	0	4

Factor settings (scale ends):

	A	B	C	D	E
1	1	1	1	1	1
2	2	2	2	2	2
3	3	3	3	3	3
4	4	4	4	4	4
5	5	5	5	5	5
6					6
7					7
8					8
9					9
10					10

The above design is reasonable and does not confound two-factor interactions of experimental

factors with the block factor. Function `fac.design` would throw an error, if the chosen block generator confounded a block effect with a main effect contrast of experimental factors or did not provide an appropriate number of blocks. Apart from these gross issues, responsibility for an appropriate choice of `block.gen` is completely with the user.

If the user is not able to come up with a satisfactory block structure, function `regular.design` from package `planor` can be used to create a design with the required properties (see code below). If that function takes a very long time for a reasonably-sized problem, there is in many cases no solution for the requested situation; it can, however, also mean that an existing solution is difficult to find for regular designs with factors that have non-prime numbers of levels.

```
R> library("planor")
R> planor.blocked25 <- regular.design(
+   factors = c("Block", "A", "B", "C", "D", "E"),
+   nlevels = c(25, 2, 5, 5, 5, 10), block = ~Block, nunits = 2500,
+   model = ~ (A + B + C + D + E) ^ 2 + Block,
+   estimate = ~ (A + B + C + D + E) ^ 2)

The search is closed: max.sol = 1 solution(s) found
The search is closed: max.sol = 1 solution(s) found
```

For the design `planor.blocked25`, the `estimate` option guarantees that blocks are not confounded with two-factor interactions. The block generator used by function `regular.design` is different from the one chosen in function `fac.design` above, but of the same quality with respect to the GMA criterion. This can be verified by applying function `GWLP` or function `lengths` to both designs (as mentioned before, `lengths` is faster for designs with many runs):

```
R> round(lengths(full.factorial.blocked25, with.blocks = TRUE), 2)
2 3 4 5
0 0 16 8

R> round(lengths(data2design(getDesign(planor.blocked25))), 2)
2 3 4 5
0 0 16 8
```

Neither function `fac.design` nor function `regular.design` from package `planor` guarantees optimality of the confounding structure, and neither is generally superior to the other.

5. Orthogonal arrays with package `DoE.base`

If more than two levels are required for some factors, but a full factorial cannot be afforded, a regular or non-regular orthogonal array can be used. For the creation of *regular* designs, also with mixed levels, function `regular.design` from package `planor` is very useful. As seen in the code of the previous section, the function can specify a model and separately a (sub)model

to be estimable; in this way, it was, e.g., possible to treat the block factor differently from the other factors. Note, however, that it is not possible to generate non-regular designs, and that no effort is made at a better design in terms of overall model robustness, whenever the requested estimability requirements are satisfied. Furthermore, creation of some designs takes a very long time, even for relatively small designs (e.g., 32 runs).

Package **DoE.base** pursues a different route: It contains the previously-mentioned catalog of orthogonal arrays; most of its arrays have been taken from [Kuhfeld \(2010\)](#). In addition, since version 0.28, it contains a small separate catalog `oacat3` of orthogonal arrays of strength at least 3 from other sources. Similarly to most experimental design software, the default approach of function `oa.design` is to use the smallest array possible and to take columns with the requested numbers of levels from left to right until the design has been filled. If the user does not influence the chosen columns with the `columns` option, a warning is issued to make the user aware of potential improvements. The following subsections discuss the catalogs `oacat` and `oacat3` and ways to select arrays from them, the optimization of column selection from a selected array, and ways to inspect experimental plans regarding their suitability for the experiment at hand.

5.1. The data frames `oacat` and `oacat3` and the function `show.oas`

The arrays available in package **DoE.base** are documented in the data frames `oacat` and `oacat3`; `oacat` contains strength 2 arrays, `oacat3` a selection of stronger arrays. Both data frames contain the following columns:

- `name` gives a structured array name, which indicates the number of runs and the frequency of factors with different numbers of levels; for example, the name `L18.2.1.3.7` indicates 18 runs with one 2-level factor and seven 3-level factors. `nruns` directly gives the number of runs. `n2` to `n72` give the number of factors with 2 to 72 levels. Thus, for `L18.2.1.3.7`, `nruns = 18`, `n2 = 1`, `n3 = 7`, and all other `nx` entries have the value 0.
- `lineage` contains a string variable which indicates how the array was constructed from so-called parent arrays. An empty string indicates that the array itself is a parent array. Parent arrays are stored in the package and are objects of class ‘`oa`’, which are matrices that usually contain an “`origin`” attribute and sometimes also a “`comment`” attribute. Arrays with a lineage entry are constructed from the parent arrays.
- Logical columns indicate a full factorial array (`ff`), an array that is R^2 regular according to the definition by [Grömping and Bailey \(2016\)](#) (`regular.strict` with only average R^2 values 0 and 1 for all factor sets), or an array that is CC regular according to the definition by [Grömping and Bailey \(2016\)](#), if their conjecture holds, i.e., if it suffices to check for squared canonical correlations 0 and 1 for all full resolution factor sets only (`regular`); for some large arrays, the necessary checks for CC regularity – even given the conjecture – have not been fully done, so that it is conceivable (but considered highly unlikely) that some of the designs indicated as regular may be non-regular according to a higher order non-regularity.
- Column `SCones` contains the number of squared canonical correlations in R factor sets that are one.
- Columns `GR` and `GRind` contain the GR and GR_{ind} values, respectively.

- Columns `maxAR` and `maxSC` contain the maximum average R^2 or squared canonical correlation, respectively.
- Column `dfe` provides the number of error degrees of freedom, if all columns of the array are used.
- Columns `A3` to `A8` provide the numbers of generalized words of lengths 3 to 8. (There are no words of shorter lengths, of course.)

It is possible to use the data frames `oacat` or `oacat3` directly for inspecting which arrays of a certain nature are available, for example for finding strength 2 32-run arrays which are regular but not strictly regular:

```
R> oacat$name[oacat$nruns == 32 & oacat$regular & !oacat$regular.strict]
```

```
[1] "L32.2.28.4.1"      "L32.2.25.4.2"      "L32.2.24.8.1"
[4] "L32.2.22.4.3"      "L32.2.21.4.1.8.1"  "L32.2.19.4.4"
[7] "L32.2.18.4.2.8.1"  "L32.2.16.4.5"      "L32.2.16.16.1"
[10] "L32.2.15.4.3.8.1"  "L32.2.13.4.6"      "L32.2.12.4.4.8.1"
[13] "L32.2.10.4.7"      "L32.2.9.4.5.8.1"  "L32.2.7.4.8"
[16] "L32.2.6.4.6.8.1"  "L32.2.4.4.9"      "L32.2.3.4.7.8.1"
[19] "L32.4.8.8.1"
```

The author prefers non-regular arrays for many situations, at least for the creation of screening designs. Looking at regular arrays in package **DoE.base** may nevertheless be of interest, if function `regular.design` of package **planor** runs for a long time without indicating failure for a run size that is in the scope of package **DoE.base**: If there is a regular array of the desired size in **DoE.base**, this array can be inspected and perhaps used after column optimization (see Section 5.2). Also, in principle, function `regular.design` can be expected to eventually succeed, unless estimability requirements make the existing array unsuitable.

The function `show.oas` allows inspection of the available arrays in a more convenient way; it shows arrays from `oacat`, unless option `Rgt3 = TRUE` requests arrays from `oacat3`. Suppose, for example, that a design for three 2-level factors, two 3-level factors and one 6-level factor is to be created, and between 20 and 54 runs are affordable (a full factorial would have 432). The following statement allows to inspect the candidate arrays, displaying also the quality metrics:

```
R> show.oas(nruns = c(20, 54), nlevels = c(2, 3, 3, 2, 2, 6),
+           showmetrics = TRUE)
```

5 arrays found

		name	nruns	lineage	GR	GRind	regular	SCones	A3	A4
78	L36.2.13.3.2.6.1		36		3.00	3.00	FALSE	6	45.3	158.4
81	L36.2.10.3.8.6.1		36		3.18	3.18	FALSE	0	130.3	737.2
83	L36.2.9.3.4.6.2		36		3.00	3.00	FALSE	17	82.3	338.2
87	L36.2.3.3.9.6.1		36		3.18	3.00	FALSE	12	73.8	300.2
88	L36.2.3.3.2.6.3		36		3.00	3.00	FALSE	37	35.6	73.1

	A5	A6	A7	A8
78	426	1010	1753.2	2306.8
81	3063	11096	31380.8	68828.1
83	1025	2828	5507.5	7780.0
87	912	2404	4354.2	5793.4
88	120	125	63.5	13.9

Only arrays in 36 runs have been found; these are quite different in the available patterns of numbers of levels, and therefore the quality metrics (especially the A_k) are not directly comparable. Nevertheless, the values for GR and GR_{ind} suggest that one might try the array L36.2.10.3.8.6.1, which is the only one without any completely confounded degree of freedom.

5.2. Optimization methods for function `oa.design`

Function `oa.design` allows to select an array, and to specify columns from that array either manually or by an optimization approach. As was mentioned earlier, if no array is specified, the function picks the first (and thus smallest) array in the catalog that is able to accommodate the requested factors. If no column selection approach is specified, the function simply takes the first available columns (from left to right).

The code below compares three designs:

- an unoptimized default design (i.e., the left-most suitable columns of the first array encountered, which is the L36.2.13.3.2.6.1);
- an optimized default design with optimization option `columns = "min34"` (i.e., an optimized choice of columns from the array L36.2.13.3.2.6.1, optimization being with respect to the number of generalized words of length 3, and subsequently length 4);
- and an optimized design obtained by selecting columns from the array L36.2.10.3.8.6.1 selected in Section 5.1 because of its GR_{ind} value.

```
R> oa.default <- oa.design(nlevels = c(2, 3, 3, 2, 2, 6))
R> GWLP(oa.default, digits = 2)
```

```
0   1   2   3   4   5   6
1.00 0.00 0.00 5.78 2.11 2.44 0.67
```

```
R> oa.optimized <- oa.design(nlevels = c(2, 3, 3, 2, 2, 6),
+     columns = "min34")
R> GWLP(oa.optimized, digits = 2)
```

```
0   1   2   3   4   5   6
1.00 0.00 0.00 4.11 3.61 2.78 0.50
```

```
R> oa.manualoptimized <- oa.design(L36.2.10.3.8.6.1,
+     nlevels = c(2, 3, 3, 2, 2, 6), columns = "min34")
R> GWLP(oa.manualoptimized, digits = 2)
```

0	1	2	3	4	5	6
1.00	0.00	0.00	2.44	6.44	1.78	0.33

Clearly, optimization improves the design from the default array, and the optimized pre-selected array is even better.

There are various methods for optimizing column allocation (see the documentation of function `oa.design` in the manual); `columns = "min34"` is the most important one among these. Depending on the number of columns on offer and the number of columns to be selected, optimization can take a very long time; in the above examples, $\binom{13}{3} \cdot \binom{2}{2} = 286$ or $\binom{10}{3} \cdot \binom{8}{2} = 3360$ column choices had to be checked, which is doable in reasonably short time. For larger designs, the resource implications of the numbers of available columns of the required lengths may also be considered in selecting an array to use.

5.3. Blocking general orthogonal arrays

Suppose a design with the above factors is to be run, and the 6-level factor is a blocking factor. In that case, the design should be randomized such that runs are randomized within blocks. Function `oa.design` does not directly allow to block randomization. However, the function `rerandomize.design` allows a post-hoc randomization within a single design factor declared as the block factor.

```
R> blockedoa1 <- rerandomize.design(oa.manualoptimized, seed = 24652,
+   block = "F")
R> blockedoa1

  run.no run.no.std.rp F A B C D E
  1      1      22.5.4 5 2 1 1 1 2
  2      2      5.5.2 5 1 2 3 2 2
  3      3      15.5.3 5 1 3 3 1 1
  4      4      1.5.1 5 1 1 2 1 1
  5      5      35.5.6 5 2 2 1 2 1
  6      6      33.5.5 5 2 3 2 2 2
  run.no run.no.std.rp F A B C D E
  7      7      32.3.5 3 2 2 1 2 2
  8      8      14.3.3 3 1 2 2 1 1
  9      9      34.3.6 3 2 1 3 2 1
  10     10     4.3.2 3 1 1 2 2 2
  11     11     3.3.1 3 1 3 1 1 1
  12     12     24.3.4 3 2 3 3 1 2
  run.no run.no.std.rp F A B C D E
  13     13     18.4.3 4 1 3 3 2 2
  14     14     26.4.5 4 2 2 2 2 1
  15     15     30.4.6 4 2 3 1 1 1
  16     16     11.4.2 4 1 2 1 1 2
  17     17     7.4.1 4 1 1 3 2 1
  18     18     19.4.4 4 2 1 2 1 2
  run.no run.no.std.rp F A B C D E
```

```

19      19      10.2.2 2 1 1 3 1 2
20      20      21.2.4 2 2 3 1 1 2
21      21      29.2.6 2 2 2 3 1 1
22      22      17.2.3 2 1 2 2 2 2
23      23      9.2.1 2 1 3 2 2 1
24      24      25.2.5 2 2 1 1 2 1
run.no run.no.std.rp F A B C D E
25      25      20.6.4 6 2 2 3 1 2
26      26      16.6.3 6 1 1 1 2 2
27      27      8.6.1 6 1 2 1 2 1
28      28      12.6.2 6 1 3 2 1 2
29      29      28.6.6 6 2 1 2 1 1
30      30      27.6.5 6 2 3 3 2 1
run.no run.no.std.rp F A B C D E
31      31      31.1.5 1 2 1 3 2 2
32      32      36.1.6 1 2 3 2 2 1
33      33      13.1.3 1 1 1 1 1 1
34      34      2.1.1 1 1 2 3 1 1
35      35      23.1.4 1 2 2 2 1 2
36      36      6.1.2 1 1 3 1 2 2
class=design, type= oa.blocked
NOTE: columns run.no and run.no.std.rp are annotation,
not part of the data frame

```

Function **GWLP** per default ignores confounding with the block factor, but can be requested to include it using the **with.block = TRUE** option:

```
R> GWLP(blockedoa1, digits = 2)
```

0	1	2	3	4	5
1.00	0.00	0.00	0.28	0.33	0.39

```
R> GWLP(blockedoa1, with.block = TRUE, digits = 2)
```

0	1	2	3	4	5
1.00	0.00	0.00	2.44	6.44	1.78

For designs that can also be created with function **FrF2** (package **FrF2**) and/or function **regular.design** (package **planor**), using one of the latter two may be a better choice, since they allow a more direct control over design quality via minimum aberration or estimable effects: The code below (results not shown) creates and inspects a 16-run design with eight 2-level factors in eight blocks of size 2 with all three methods; in all three cases, the design is resolution IV in terms of the experimental factors, which is systematically requested in both **FrF2** (by the minimum aberration approach of the function) and **regular.design** (by the **model** and **estimate** options). For function **oa.design**, the design quality is not as finely tunable, apart from the overall optimization that may precede usage of an array for which not all columns are used. For the designs below, however, the quality is the same for all three

designs: The GWLPs, starting with A_3 , are (28, 14, 56, 0, 28, 1) including the block factor and (0, 14, 0, 0, 0, 1) for the experimental factors alone.

```
R> planFrF2 <- FrF2(16, 8, blocks = 8, alias.block.2fis = TRUE)
R> planDoEbase <- oa.design(L16.2.8.8.1,
+   nlevels = c(2, 2, 2, 2, 2, 2, 2, 8),
+   factor.names = c(Letters[1:8], "Block"))
R> planDoEbase <- rerandomize.design(planDoEbase, seed = 31525,
+   block = "Block")
R> planplanor <- regular.design(factors = c(Letters[1:8], "Block"),
+   nlevels = c(rep(2, 8), 8),
+   model = ~ (A + B + C + D + E + F + G + H) ^ 2 + Block,
+   estimate = ~ A + B + C + D + E + F + G + H,
+   nunits = 16, randomize = ~ Block / UNITS)
R> GWLP(planFrF2, with.block = TRUE)
R> GWLP(planFrF2)
R> GWLP(planDoEbase, with.block = TRUE)
R> GWLP(planDoEbase)
R> GWLP(planplanor@design)
R> GWLP(planplanor@design[, 1:8])
```

5.4. Inspection methods for factorial designs

Any experimental plan should be carefully checked before using it for experimentation, since experimentation usually involves a lot of effort: Adverse consequences from mistakes in design creation may be severe, but can often be prevented without much trouble, if attended to at the design creation stage. For example, when implementing a self-chosen blocking, mistakes can easily happen:

```
R> planstupid <- FrF2(16, 8, blocks = c("AB", "AC", "BCD"),
+   alias.block.2fis = TRUE)
R> GWLP(planstupid, with.blocks = TRUE)
R> plot(data2design(undesign(planstupid)), select = "all2")
```

A simple check of the GWLP can serve as a first indication whether the design behaves as expected; here we find $A_2 = 4$ (not shown), which is of course undesirable. Further checks can involve plotting. The plot command in the code above reveals (not shown) that factors D , F , G and H are aliased with the block factor for the design `planstupid`, which explains the four words of length 2. Instead of looking at all two-factor sets with option `select = "all2"`, one can also look at a percentage of worst cases (see below). Note that the `data2design(undesign())` construction is used for incorporating the block factor, since the `plot` method for class ‘`design`’ does not have a `with.blocks` option and works on the design factors only.

We now visualize the worst case confounding of the two optimized designs discussed in Section 5.2:

```
R> plot(oa.optimized, select = "worst", selprop = 0.05)
```

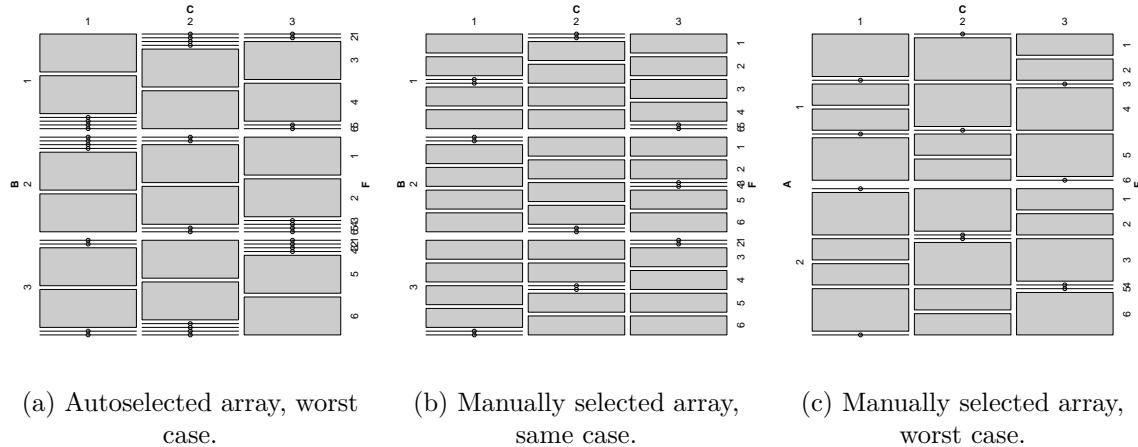


Figure 2: Mosaic plots of triples of factors.

```
R> plot(oa.manualoptimized, select = c(2, 3, 6))
R> plot(oa.manualoptimized, select = "worst", selprop = 0.05)
```

The code above generates the three mosaic plots of Figure 2. The optimized design based on the array that was automatically picked severely confounds the interaction of the two 3-level factors with the 6-level factor (factors 2, 3 and 6 of the design): A mosaic plot of that projection is created by the first line of the above code and is shown in Figure 2(a). It shows that the interaction of the two 3-level factors restricts the 6-level factor to a third of its possibilities only, which is the most severe form of aliasing possible for a triple of factors with this level combination; of the 54 possible level combinations in this triple of factors, the design contains 18 distinct variants only (instead of the possible 36). In comparison, a mosaic plot of the same projection for the optimized manually selected design shows a much better picture: Now, the design has 36 distinct level combinations, the maximum possible (see Figure 2(b)). The worst case confounding for that design is shown in Figure 2(c). It is distinctly less severe than that of Figure 2(a); however, apart from the one worst case shown in Figure 2(a), the optimized automatically-selected design is also quite reasonable.

Mosaic plots concentrate on individual low order projections. Package **DoE.base** offers another diagnostic plot for looking at the entire design at once: Function `corrPlot` inspects the correlations between model matrix columns. The creation of function `corrPlot` was inspired by function `colormap` of package **daewr** (Lawson 2016). Figure 4 in Section 6.1 illustrates its default version applied to a real life example. This function also offers a possibility to include the design's run order among the effects to be considered, since there are applications for which the run order might be of relevance. For such situations, a simple approach could be to re-randomize the design (with function `rerandomize.design`) until a randomization is encountered for which the correlation of the run order to relevant effect columns (as shown in the plot of absolute correlations) is reasonably small.

Numeric quality criteria can also be obtained; the function names for obtaining them were already mentioned in Section 3.2. These advanced criteria might deter practitioners but may be useful for experts: The function `GRind` calculates the metrics introduced by Grömping and Xu (2014) with the additional detail proposed in Grömping (2017a). For example, the code below shows that the optimized design based on the manually selected array has a clearly

better overall behavior regarding factor-specific worst case confounding ($GR_{tot,i}$ and $GR_{ind,i}$; see Section 3.2) than the optimized design based on the automatically selected first array.

```
R> GRind1 <- GRind(oa.optimized)
R> GRind2 <- GRind(oa.manualoptimized)
R> print(cbind(rep(c("oa.optimized", "oa.manualoptimized"), each = 2),
+   rbind(GRind1$GR.i, GRind2$GR.i)), quote = FALSE)

          A     B     C     D     E     F
GRtot.i oa.optimized 3.667 3    3    3.423 3.423 3.368
GRind.i oa.optimized 3.667 3    3    3.423 3.423 3
GRtot.i oa.manualoptimized 3.184 3.5 3.423 3.423 3.667 3.635
GRind.i oa.manualoptimized 3.184 3.5 3.423 3.423 3.667 3.423
```

6. An example from plant biotechnology

Vasilev *et al.* (2014) investigated cultivation factors for geraniol production by plant cells. There were four quantitative 2-level factors, two 3-level factors (one quantitative and one qualitative) and one qualitative 4-level factor. The data, including response values, have been published with the paper and are also included in package **DoE.base** as object **VSGFS**. The 2-level factors are coded in $-1/+1$ coding, the other factors in R's default dummy coding.

6.1. Creating and inspecting the design

For this experiment, a full factorial design would have had 576 runs. It would certainly have been necessary to conduct it in blocks, say in eight blocks of size 72 each. Such a design could have been created by function **fac.design** or by function **regular.design** of package **planor**, as shown in Section 4.

A full factorial design appeared neither feasible nor appropriate for the screening situation of the experiment. Instead, the design was conducted using a 72-run orthogonal array, which was generated with function **oa.design**, using automatic optimization (option **columns = "min34"**) of the manually pre-selected array **L72.2.43.3.8.4.1.6.1**. The optimization took quite a long time and resulted in the selection of columns 4, 22, 37, 41 for the 2-level factors, 46 and 48 for the 3-level factors, and 52 for the 4-level factor. The design is available as object **VSGFS**, and the code for reproducing it can be found in the manual.

When constructing the experimental plan for **VSGFS**, the array was pre-selected by intuition and trial and error from all 33 72-run designs that can accommodate the requested factors, using the versions of function **show.oas** and data frame **oacat** available at the time (no metrics information was contained in **oacat**, and thus, metric-based filtering was not possible with function **show.oas**). A key driver for the ad-hoc selection of designs within which to select columns was the number of choices a design offered for accommodating the factors; this criterion was used intuitively rather than systematically; post-hoc, note that the actually chosen design offers 3455480 choices, ranking fourth behind three designs offering 8959566, 5151510 and 3455760 possibilities. After the resource-intensive optimization of the **L72.2.43.3.8.4.1.6.1**, some further designs with distinctly fewer possibilities were additionally tried; all of these yielded worse A_3 .

With the present possibilities of function `show.oas`, the `GRgt3 = "ind"` option (i.e., requesting that $GR_{\text{ind}} > 3$ which means absence of any instance of complete confounding for individual degrees of freedom), one can pick arrays that are particularly promising for screening purposes. The following command shows the available arrays without any complete confounding for the example situation, together with the GR metrics (requested by the `showGRs` option):

```
R> show.oas(nlevels = c(2, 2, 2, 3, 2, 3, 4), GRgt3 = "ind", showGRs = TRUE)

4 arrays found
      name nruns          lineage   GR
366   L72.2.53.3.2.4.1    72           3.18
380   L72.2.44.3.12.4.1    72           3.18
382 L72.2.43.3.8.4.1.6.1    72           3.18
431   L72.2.13.3.25.4.1 72 3~24;24~1;:(24~1!2~13;3~1;4~1;) 3.18
      GRind
366  3.18
380  3.18
382  3.18
431  3.13
```

The array used in actual experimentation is the third array in this list; thus, the intuitive approach was lucky to pick a particularly promising array in this case. It can be expected that all four arrays listed above are reasonably suited for screening the experimental factors. As the designs are far from saturated, optimal column allocation can substantially improve the worst case confounding (and thus GR and GR_{ind}): For the eventual design, the output below shows $GR = GR_{\text{ind}} = 3.6$; ARFT and SCFT have been suppressed, because their background is outside the scope of this paper.

```
R> GRind(VSGFS, arft = FALSE, scft = FALSE)

$GRs
  GR GRind
3.667 3.667

$GR.i
  Light ShakFreq InocSize FilledVol     CM Sugar CDs
GRtot.i 3.667    3.667    3.667    3.864 3.667 3.864 3.808
GRind.i 3.667    3.667    3.667    3.808 3.667 3.808 3.667

attr("class")
[1] "GRind" "list"
```

For illustrating the worst case degree of confounding, Figure 3 shows the triple 3, 5, 7 in quite reasonable balance (this is one of the two worst case triples); the option `sub = "A"` in the code requests that the A_3 value for the triple is printed as a subtitle. Figure 4 shows for the entire design that the main effect model matrix columns have reasonably low correlations with

```
R> plot(VSGFS, select = list(c(3, 5, 7)), sub = "A")
```

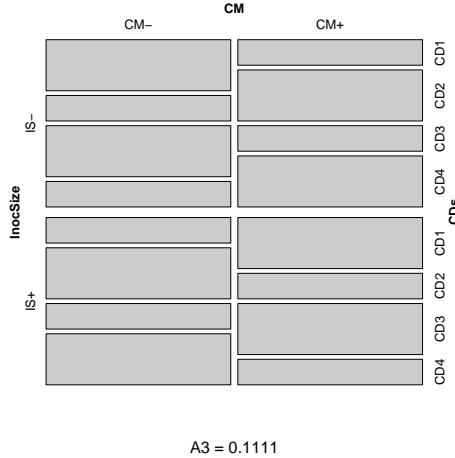


Figure 3: Mosaic plot for the worst case triple in the VSGFS example.

```
R> corrPlot(VSGFS)
```

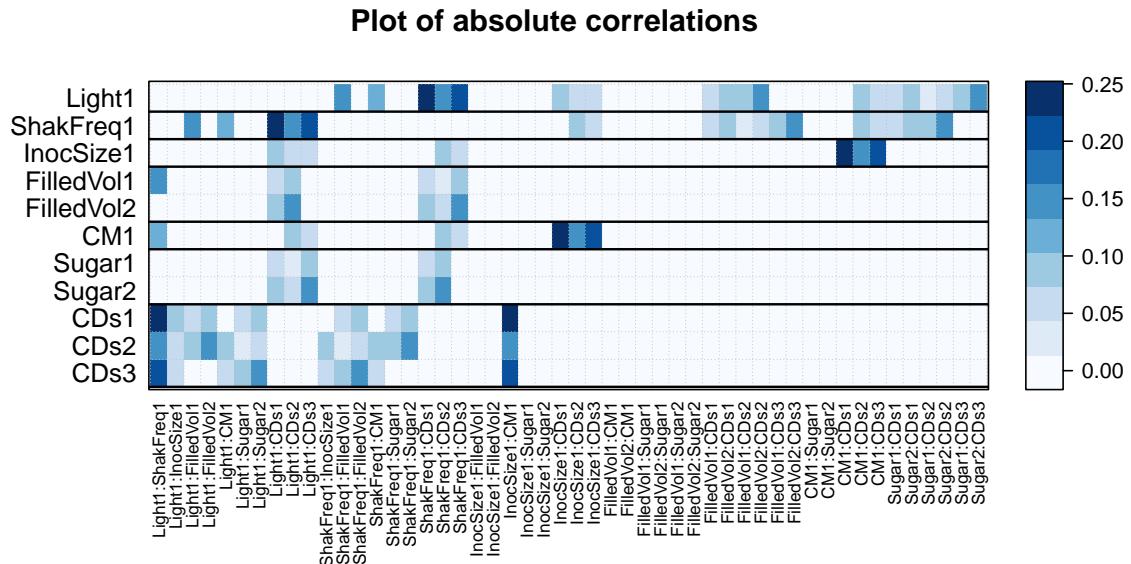


Figure 4: Plot of absolute correlations of main effects model matrix columns with two factor interaction model matrix columns for the VSGFS example (default normalized orthogonal coding is the Xu Wu coding).

the two-factor interaction model matrix columns; note that function `corrPlot` defaults to recoding all factors in normalized orthogonal coding (see also Section 7), in order to eliminate avoidable confounding.

The team was happy with the design and used it for collecting the data. Software-wise, data collection happened in **Excel**, exporting the randomized design with the `export.design`

function and re-importing it after data entry using the exported RDA file together with a CSV file with response values added to the data rows using function `add.response`.

6.2. Analyzing experimental data

The data frame `VSGFS` contains the experiment with response data. Package **DoE.base** offers a few functions for analysis purposes:

- The `plot` method for class ‘`design`’, in case of data with responses, creates simple main effects plots, by invoking the `plot` method from package **graphics**.
- The `lm` method for class ‘`design`’ runs a linear model with a modifiable default degree. For orthogonal arrays created with function `oa.design`, the default degree is one, i.e., a model with main effects only. Linear models can of course also be run by the `lm` function from the core package **stats**, and analysis of variance functionality can also be used (see below).
- The `halfnormal` method for classes ‘`lm`’ or ‘`design`’ creates a half-normal effects plot (see Section 7).

Of course, in general, other R functionality can also be used, for example interaction plots or functionality for mixed model analysis (not applicable for the example data). The functionality for handling repeated measurement data or replicated data (also not needed for this example) has been described for package **FrF2** in Sections 5.3 and 5.9 of Grömping (2014c) and is analogous here.

Main effects plots can be obtained by the simple command `plot(VSGFS)`, which creates these plots for all three response variables. Figure 5 shows the plots with default labeling, arranged with three plots on one page and reduced margin sizes. Of course, one would usually adapt the annotation for final reports or publications. The plot shows that the sugar sucrose is very beneficial for the biomass, not very good for the content, but nevertheless, because of the strong effect on biomass, beneficial for the yield. The content apparently can be increased by choosing the sugar mannitol, level one of factor CD and the + -level of light. Apart from the + -level of light, the other settings for high content are not helpful for overall yield.

An assessment of significance can be obtained from a linear model analysis. This can be obtained separately for each response, for example for the content. Here, the analysis confirms the findings from the main effects plot.

```
R> summary(lm(VSGFS, response = "Content"))

Number of observations used: 72
Formula:
Content ~ Light + ShakFreq + InocSize + FilledVol + CM + Sugar +
    CDs

Call:
lm.default(formula = fo, data = model.frame(fo, data = formula))

Residuals:
```

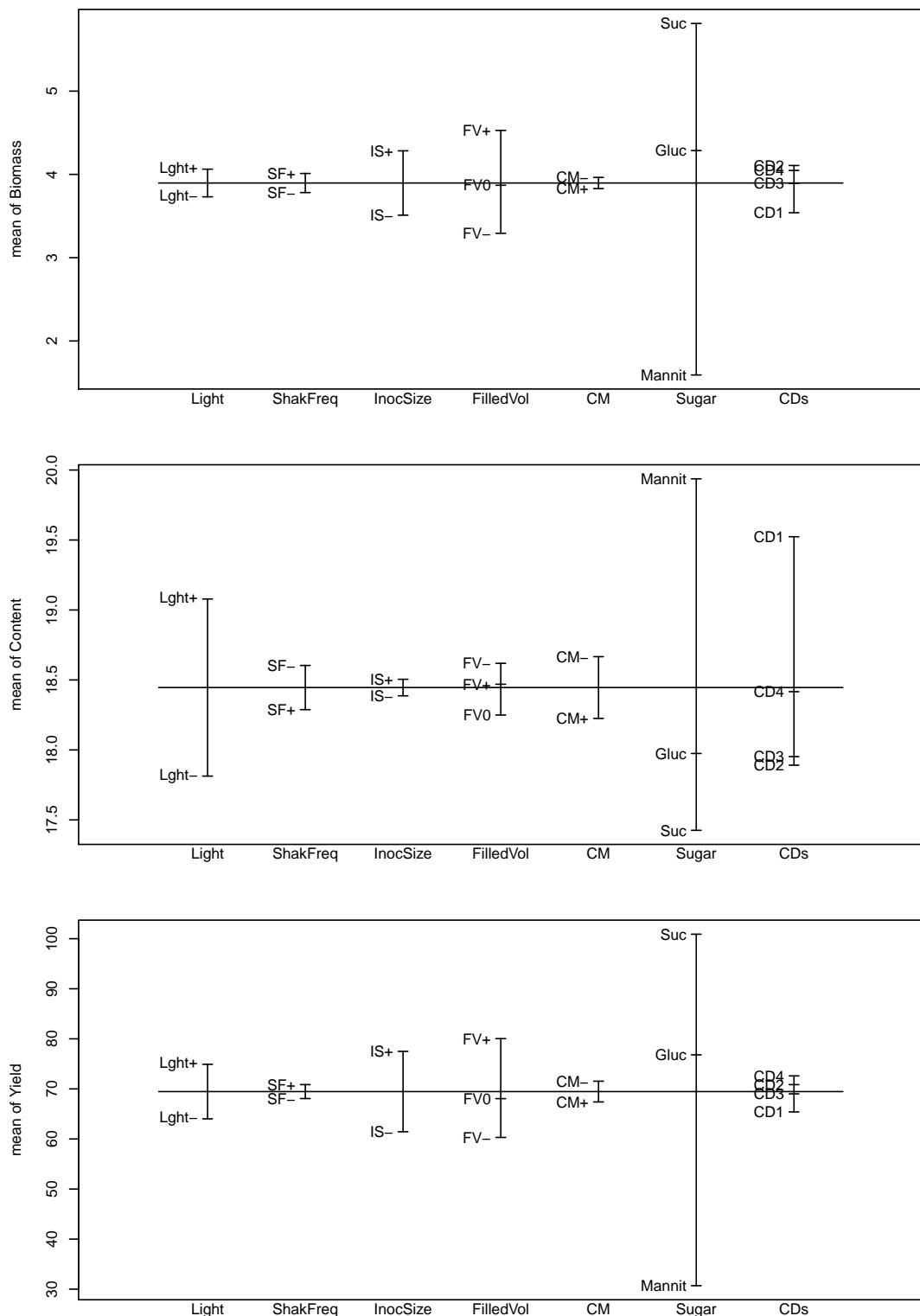


Figure 5: Main effects plots for all three responses.

	Min	1Q	Median	3Q	Max
	-2.5472	-1.1746	-0.1052	0.7008	4.6994

Coefficients:

	Estimate	Std. Error	t value	Pr(> t)
(Intercept)	18.67611	0.55106	33.891	< 2e-16 ***
Light1	0.63264	0.19483	3.247	0.00191 **
ShakFreq1	-0.15792	0.19483	-0.811	0.42083
InocSize1	0.05903	0.19483	0.303	0.76296
FilledVolFV0	-0.36958	0.47723	-0.774	0.44172
FilledVolFV+	-0.14958	0.47723	-0.313	0.75503
CM1	-0.22069	0.19483	-1.133	0.26182
SugarGluc	0.54917	0.47723	1.151	0.25441
SugarMannit	2.51167	0.47723	5.263	2e-06 ***
CDsCD2	-1.63222	0.55106	-2.962	0.00438 **
CDsCD3	-1.57111	0.55106	-2.851	0.00597 **
CDsCD4	-1.10722	0.55106	-2.009	0.04901 *

Signif. codes:	0 '***'	0.001 '**'	0.01 '*'	0.05 '.'
	0.1 ''	1		

Residual standard error: 1.653 on 60 degrees of freedom
 Multiple R-squared: 0.4787, Adjusted R-squared: 0.3831
 F-statistic: 5.008 on 11 and 60 DF, p-value: 1.75e-05

The model explains less than 50% of the response variability, which is far from perfect; the relatively large unexplained variation can be due both to random error and to the presence of interaction effects not captured in the main effects model. Nevertheless, the main effects results are trustworthy. The strong heredity principle – interactions are only active if both their component factors are active – suggests to look at a model with the three active main effect factors and their interactions, which leaves a somewhat confusing picture (not shown).

For the data at hand, there are enough degrees of freedom to run an Anova analysis with the full degree 2 model. Of course, while main effects are orthogonal to each other, two-factor interactions can be slightly confounded with main effects and severely confounded with other two-factor interactions. However, at least, the theory tells us that the estimable effects can be estimated without bias (unless effects of order higher than two bias them). As Anova analyses sums of squares, the analysis is invariant with respect to factor coding. In order to obtain an order-invariant assessment of significance, the function **Anova** from package **car** ([Fox and Weisberg 2011](#)) can be used; contrary to function **anova** from package **stats**, **Anova** avoids order-dependence by using type II sums of squares, which condition on all other effects except for the ones that contain the effect under investigation. The results point to the main effects that were also identified before, and a liberal look at *p* values additionally indicates some marginal two-factor interactions (sugar with each of **InocSize**, **FilledVol**, **CDs**, **Light**, and **ShakFreq:CM**, which is completely unrelated to the active main effects).

```
R> library("car")
R> Anova(lm(VSGFS, response = "Content", degree = 2))
```

Anova Table (Type II tests)

Response: Content	Sum Sq	Df	F value	Pr(>F)
Light	16.345	1	8.2606	0.0165515 *
ShakFreq	0.483	1	0.2441	0.6319195
InocSize	1.604	1	0.8105	0.3891387
FilledVol	0.889	2	0.2247	0.8026463
CM	0.805	1	0.4069	0.5378565
Sugar	67.692	2	17.1048	0.0005921 ***
CDs	29.468	3	4.9642	0.0230814 *
Light:ShakFreq	1.672	1	0.8450	0.3795988
Light:InocSize	0.320	1	0.1616	0.6961803
Light:FilledVol	3.296	2	0.8330	0.4628103
Light:CM	2.874	1	1.4524	0.2558948
Light:Sugar	11.387	2	2.8774	0.1030201
Light:CDs	6.073	3	1.0231	0.4231004
ShakFreq:InocSize	1.162	1	0.5871	0.4612301
ShakFreq:FilledVol	0.420	2	0.1061	0.9003686
ShakFreq:CM	6.303	1	3.1856	0.1046054
ShakFreq:Sugar	3.074	2	0.7767	0.4857813
ShakFreq:CDs	11.717	3	1.9738	0.1819413
InocSize:FilledVol	1.626	2	0.4109	0.6737704
InocSize:CM	0.752	1	0.3802	0.5512885
InocSize:Sugar	20.475	2	5.1739	0.0286697 *
InocSize:CDs	7.710	3	1.2989	0.3280819
FilledVol:CM	2.976	2	0.7521	0.4962887
FilledVol:Sugar	24.172	4	3.0539	0.0692876 .
FilledVol:CDs	6.333	6	0.5334	0.7719232
CM:Sugar	3.387	2	0.8559	0.4538120
CM:CDs	10.566	3	1.7799	0.2144014
Sugar:CDs	33.075	6	2.7858	0.0734918 .
Residuals	19.787	10		
<hr/>				
Signif. codes: 0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 ' ' 1				

Even though this design does not require analysis with half-normal effects plots, the next section will illustrate application of half-normal effects plots for this example.

7. Half-normal effects plots

In (almost) saturated designs, conventional analysis of variance methods are not very successful, because there are too few degrees of freedom for error. If one assumes a screening design, for which most effects are inactive, the inactive effects actually represent experimental error; however, it is not known a-priori, which are the active effects. Daniel (1959) proposed to use half-normal effects plots for diagnosing which effects are active, and Lenth (1989) pro-

posed a numerical activity check for these, known as “Lenth’s method”. The critical values proposed by Lenth were later found to be conservative, and it was proposed to use simulated ones instead. Half-normal effects plots with Lenth’s method and simulated critical values are implemented in function `halfnormal` of package **DoE.base**.

7.1. The principle

The standard use for such plots is with 2-level factors which are conventionally coded in $-1/+1$ coding (see, e.g., Grömping 2014c). Function `halfnormal` from package **DoE.base** covers not only these standard situations, but also offers half-normal plots for:

- 2-level designs with a few error degrees of freedom. For these, it automatically augments the estimated effects with error effects, distinguishing these into lack-of-fit and pure error. Significance assessment can be done with Lenth’s method on the augmented set of estimates, or with other methods proposed in the literature (Larntz and Whitcomb 1998; Edwards and Mee 2008). Note that error effects are not necessarily uniquely determined.
- 2-level designs with partially confounded effects. For these, it projects out all preceding effects from the remaining ones (thus, the plotting points depend on the model order for such situations).
- Mixed level designs, for which there is no unique coding and the plotting points are coding dependent. Mixed level designs can also have error degrees of freedom or partially confounded effects.

The strategy chosen in function `halfnormal` seems to be similar to that applied in the **JMP** software (SAS Institute Inc. 2018) screening platform (see Chapter 8 of SAS Institute Inc. 2012), both regarding the treatment of error points and the single degree of freedom representations for factors with more than two levels. On the contrary, **Design-Expert** (Stat-Ease Inc. 2017) does not plot individual degrees of freedom, but scaled Chi-squared values for effects with more than one degree of freedom. This avoids the coding dependence, but has the adverse effect that the number of plotting points is small so that effect sparsity is not easily achieved.

The steps for augmenting the estimated effects with error degrees of freedom are described, e.g., in Grömping (2015) and are very similar also to the suggestions by Langsrud (2001) in a different context. A coarse overview works as follows:

- Make sure the model matrix \mathbf{X} has orthogonal columns all of which have the same Euclidean length; if this is not the case, \mathbf{X} has to be pre-treated (see below).
- For N observations, a trivial saturated model matrix is the N -dimensional identity matrix \mathbf{I}_N . If a distinction between lack-of-fit and pure error is sought, one can replace this matrix by a matrix \mathbf{S} of dummy variables for *distinct runs*, and additionally include appropriately scaled orthogonal contrast matrices for replicated runs. In the following, for simplicity, \mathbf{I}_N is used.
- Residualize the matrix \mathbf{I}_N by projecting out the model matrix \mathbf{X} , i.e., calculate the residual matrix $\mathbf{R} = \mathbf{I}_N - \mathbf{X}(\mathbf{X}^\top \mathbf{X})^{-1} \mathbf{X}^\top$.

- Create the half-normal effects plot for the augmented model matrix ($\mathbf{X}|\mathbf{R}$), which has been created such that it has orthogonal columns where all have the same length.

The pre-treatment mentioned in the first bullet is as follows: If \mathbf{X} has orthogonal columns of varying Euclidean lengths, one simply has to normalize all columns to a common length. The case of non-orthogonal columns is more demanding and will be discussed using the model matrix \mathbf{X} for a full model in an unreplicated full factorial design for one 2-level and one 3-level factor, both in dummy coding, as given in Equation 1:

$$\mathbf{X} = \begin{pmatrix} \text{Int} & A_2 & B_2 & B_3 & A_2B_2 & A_2B_3 \\ 1 & 0 & 0 & 0 & 0 & 0 \\ 1 & 1 & 0 & 0 & 0 & 0 \\ 1 & 0 & 1 & 0 & 0 & 0 \\ 1 & 1 & 1 & 0 & 1 & 0 \\ 1 & 0 & 0 & 1 & 0 & 0 \\ 1 & 1 & 0 & 1 & 0 & 1 \end{pmatrix}. \quad (1)$$

The columns of this matrix are not orthogonal, which can be easily verified by inspecting $\mathbf{X}^\top \mathbf{X}$. As the intercept estimates the mean for the combination A_1B_1 and the coefficients for the main effects columns estimate deviations from that level, there is an obvious dependency. With the interaction coefficients also measuring deviations of particular cells from additivity, there is another clear dependence. It is well-known which effects are estimable in factorial models: the overall means and the contrasts. A model matrix formulated to directly estimate these has orthogonal columns.

Xu and Wu (2001) showed that it is particularly useful to code all main effects columns using an orthogonal coding normalized to squared length N , where N is the number of runs: Such a coding yields orthogonal columns of the model matrix for all effects up to degree s for any projection of s factors from a strength s design, if the interaction columns are created in the usual way as products of the normalized main effects contrast columns; the interaction columns will then also have squared Euclidean length N and will be orthogonal to each other and to the main effects columns. Such coding will be called Xu Wu coding in the sequel. It is available in two versions in package **DoE.base**: There are `contr.XuWu` and `contr.XuWuPoly` contrasts. The model matrix below is obtained by coding both factors A and B with `contr.XuWu` contrasts.

$$\mathbf{X} = \begin{pmatrix} \text{Int} & A_2 & B_2 & B_3 & A_2B_2 & A_2B_3 \\ 1 & -1 & -\sqrt{1.5} & -\sqrt{0.5} & \sqrt{1.5} & \sqrt{0.5} \\ 1 & 1 & -\sqrt{1.5} & -\sqrt{0.5} & -\sqrt{1.5} & -\sqrt{0.5} \\ 1 & -1 & \sqrt{1.5} & -\sqrt{0.5} & -\sqrt{1.5} & \sqrt{0.5} \\ 1 & 1 & \sqrt{1.5} & -\sqrt{0.5} & \sqrt{1.5} & -\sqrt{0.5} \\ 1 & -1 & 0 & \sqrt{2} & 0 & -\sqrt{2} \\ 1 & 1 & 0 & \sqrt{2} & 0 & \sqrt{2} \end{pmatrix} \quad (2)$$

If the design has strength s , a model matrix in a Xu Wu coding for main effects with up to s -factor interactions can also be obtained post-hoc from a model matrix \mathbf{X} based on non-orthogonal coding by sequential orthogonalization and normalization: Project out the first column (intercept column) from the second, which is just subtraction of the column mean. Project out the first two columns from the third and so on. In addition, one also has to

normalize all columns to squared length N (or any other common Euclidean length) in order to satisfy the requirements for a half-normal effects plot.

If the model has truly confounded effects that cannot be orthogonalized by simply applying Xu Wu coding, the same orthogonalization strategy can be applied, but the consequence is not only a recoding of in principle the same information, but an order-dependent removal of earlier confounded effects from later confounded effects.

7.2. Example application

The example design is an orthogonal array, i.e., main effect *contrasts* can be estimated independently of each other. However, depending on the coding, the actual *estimated coefficients* may be correlated, as was discussed above. For example, for 2-level factors, if $-1/+1$ coding is used, the estimates are uncorrelated, with 0/1 (dummy) coding, however, they are correlated. It is advisable to explicitly choose an orthogonal factor coding, ideally the Xu Wu variant discussed above. The code below creates an orthogonal main effects model matrix with squared Euclidean length $N = 72$ (output not shown).

```
R> VSGFS.XuWuPoly <- change.contr(VSGFS, "contr.XuWuPoly")
R> round(crossprod(model.matrix(lm(VSGFS.XuWuPoly))), 2)
```

Per default, function `halfnormal` refuses to work in case of correlated main effects (except in case of the perfect confounding of regular designs). The option `ME.partial = TRUE` can be used to change that; if the partial aliasing among main effects estimates is due to non-orthogonal coding in an orthogonal array, use of `ME.partial = TRUE` is acceptable in the light of the previous considerations (see the first command to create `hnAuto` below), although it seems preferable to decide on an orthogonal coding explicitly.

For the example design, a main effects analysis is quite well-protected against bias from two-factor interactions. However, two-factor interactions may be quite heavily confounded with each other. Nevertheless, we will now consider an almost saturated array in order to demonstrate the most general usage of the function `halfnormal`.

```
R> hnAuto <- halfnormal(lm(VSGFS, response = "Content", degree = 2),
+   ME.partial = TRUE, cex.text = 0.9, cex = 0.9, xlim = c(0, 1.1),
+   ylim = c(0, 2.8), main = "Half-normal effects plot for Content")
R> hnXuWuPoly <- halfnormal(lm(VSGFS.XuWuPoly, response = "Content",
+   degree = 2), cex.text = 0.9, cex = 0.9, xlim = c(0, 1.1),
+   ylim = c(0, 2.8), main = "Half-normal effects plot for Content")
R> hnXuWuPolyreordered <- halfnormal(lm(Content ~
+   (CDs + Sugar + CM + FilledVol + InocSize + ShakFreq + Light)^2,
+   VSGFS.XuWuPoly), cex.text = 0.9, cex = 0.9, xlim = c(0, 1.1),
+   ylim = c(0, 2.8), main = "Half-normal effects plot for Content")
```

Figure 6 shows half-normal effects plots from a model with all main effects and two-factor interactions for the response variable `Content`. In order to demonstrate the coding dependence of half-normal effects plots in case of factors with more than two levels, plot (a) shows the automatic coding obtained by orthogonalizing the model matrix that results from the dummy coding of the 3- and 4-level factors in `VSGFS`, plot (b) `contr.XuWuPoly` coding. Plot (c) shows

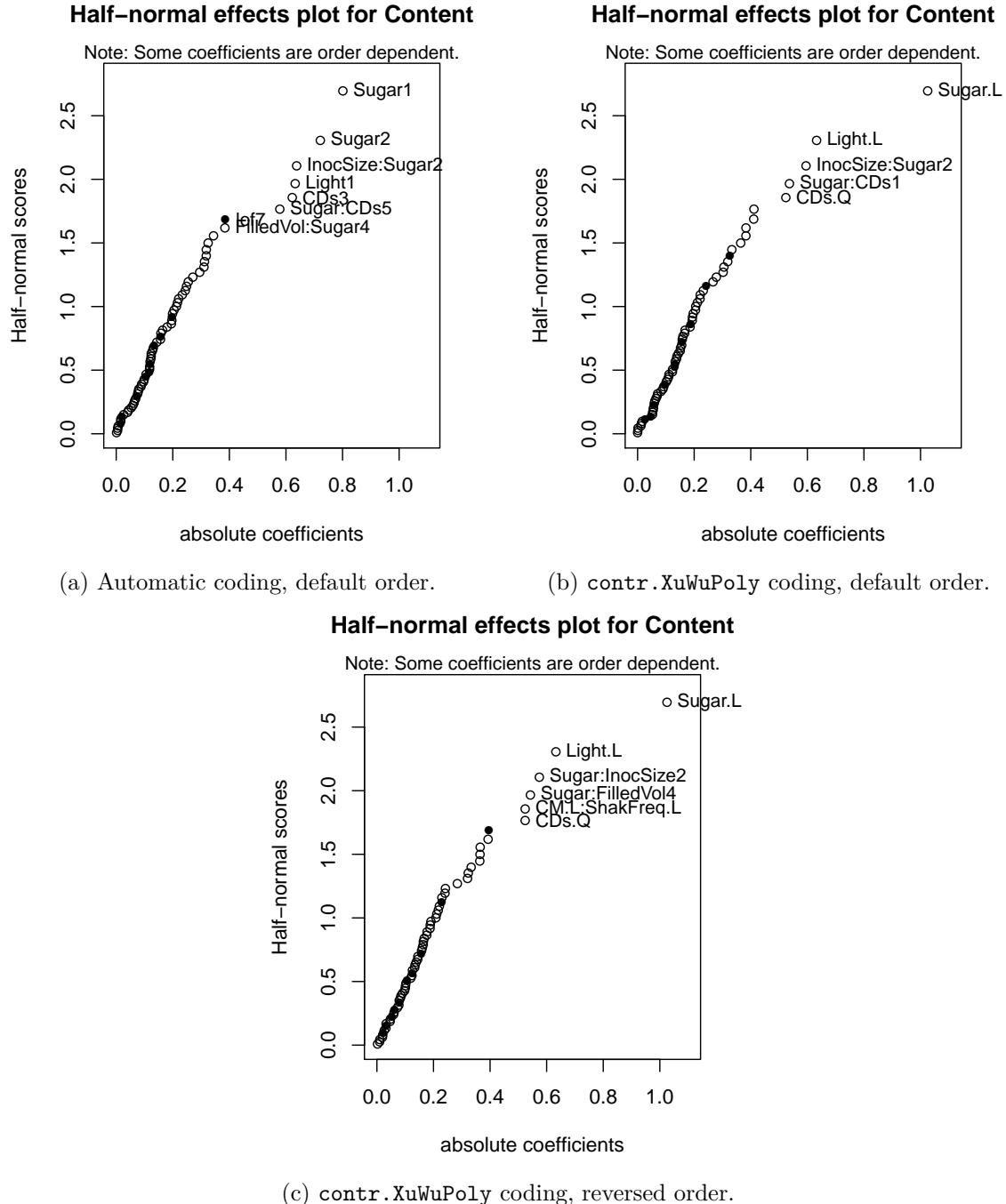


Figure 6: Half-normal effects plots for a model with all two-factor interactions, labeling from Lenth's method with $\alpha = 0.05$.

the analysis for `contr.XuWuPoly` coding with the effects in different order. The actual codings between (a) and (b) differ in the level ordering for the 3-level factors; for the 4-level factor, the coding differences are more complicated. Note that plot points from effects with more than one degree of freedom are not directly comparable between plots from different codings

(e.g., plots (a) and (b) of Figure 6); the comparable message is restricted to the presence or absence of individual effect components among the top-ranked absolute coefficients (and even that may be coding dependent). For the example data, while there are clear visual differences between the differently-coded plots, the message is more or less the same: As already seen in the linear main effects model, light, sugar and CDs are active factors. Furthermore, the sugar by CDs interaction and the inoculum size by sugar interaction seem to have active components. However, the interaction effects are order dependent. The plot with different interaction orders brings up three further interactions into the possibly active range and drops the sugar by CDs interaction; thus, clearly, one has to be cautious with statements on these effects from half-normal effects plots. However, all effects that show up in the half-normal effects plots have also been at least marginal in the Anova analysis of the previous section.

The calculation results have been stored in the following list objects: `hnAuto`, `hnXuWuPoly`, and `hnXuWuPolyreordered`. The list element `res` indicates which effects have been projected out from which other effects; furthermore, the results contain the model matrix after orthogonalization, and details about the orthogonalization itself. Printed output of orthogonalization steps has been suppressed here. The reader is encouraged to check the printed output of the simple command:

```
R> hnDemo <- halfnormal(VSGFS, ME.partial = TRUE, plot = FALSE)
```

which explains how columns of the main effects model matrix stored in

```
R> mm <- hnDemo$mm[, 2:11]
```

are obtained from the original model matrix stored (without intercept column) in

```
R> desnum(VSGFS) [, 1:10]
```

8. Further developments

Package **DoE.base** tries a balancing act of offering tools both for practitioners with a relatively weak statistical background and for statistical experts. In order to save the former from avoidable grave mistakes, the package takes a cautious strategy issuing many warnings, where designs might be improvable or analyses might be inadequate. In various cases, it would be desirable to avoid unnecessary warnings; for example, there are arrays for which it is known from theoretical work (Butler 2005) that all choices for certain columns are GMA. For these arrays, warnings for array optimization should be eliminated, which requires some slightly tedious work.

The current design catalog covers many situations, but is also limited especially with respect to availability of non-regular designs. Augmenting the catalog with further useful non-regular designs is a difficult task and should not be addressed before studying in more detail the relation between recently developed design quality criteria like *GR*, *GR_{ind}*, ARFT and SCFT and a design's usefulness for experimentation. Ideally, it should eventually be possible to automate the entire process of obtaining a good experimental design, combining array selection with the selection of columns within the array. It would also be good to have an option for

requesting a design whose main effects model matrix columns are (near) orthogonal to run order. The latter wishes are, however, quite resource intensive.

So far, the design catalog is used for selecting designs and optimizing column choices from them. [Kuhfeld \(2010\)](#) makes much more extensive use of the catalog within a SAS software macro suite for creating experimental designs from the arrays in the catalog, specifically created with marketing applications in mind ([Kuhfeld 2010](#)). Such use can certainly also be combined with the catalog in package **DoE.base**. The dependent package **support.CEs** ([Aizaki 2012](#)) implements choice experiments in R. However, implementing a functionality as flexible as the one offered in the SAS macros by [Kuhfeld \(2010\)](#) would require substantial additional effort.

Package **DoE.base** also serves as a tool for supporting further theoretical investigations into aspects like design regularity and new design quality criteria. For example, the work for [Grömping and Bailey \(2016\)](#) was supported by the package, and recent efforts at developing further criteria for checking arrays for combinatorial equivalence have been implemented in the package ([Grömping 2018a](#)).

Last but not least, the class ‘`design`’ and infrastructure for it are also available for use by other packages. Authors of other S3-based packages are invited to use it; the methods within package **DoE.base** for its class ‘`design`’ can be adjusted to accommodate further types of designs; please inform the package maintainer, if you would like a new type of design included into the methods offered.

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A. Class ‘design’ and functionality for it

Generally, all design generating functions from packages **DoE.base** and **FrF2** (exceptions highlighted in the documentation) create an output design that is of S3 class ‘**design**’ and thus the output follows a certain structure (cf. Section A.1) and allows application of certain inspection, modification and analysis methods and functions. The first subsection describes the class ‘**design**’ itself, the second subsection functionality applicable to class ‘**design**’ objects.

Experimental design matrices or data frames created with tools that do not output class ‘**design**’ objects can be brought into that class by function **data2design**, in order to make some of the package functionality work on them.

A.1. Class ‘design’

An object of S3 class ‘**design**’ is a data frame with the three attributes “**desnum**”, “**run.order**” and “**design.info**”.

- Attribute “**desnum**” can be used for a numeric version of the data frame, which may be useful for users who want to do manual matrix calculations. The package functionality itself makes little use of that attribute.
- Attribute “**run.order**” has the main purpose of always being able to switch back and forth between a standard order and the randomized run order.
- Attribute “**design.info**” is a list that contains all the important information on the design and is heavily used by the methods and functions discussed in the following subsection. This attribute will be described in some detail in the rest of this subsection.

The “**design.info**” attribute has some mandatory elements that have to be present for all class ‘**design**’ objects and many elements that are needed for some types of designs only. The author’s website contains a large table that details which types of designs need which elements of the “**design.info**” attribute. Tables 2, 3 and 4 list the elements and their meaning/context.

Apart from the elements listed in the table, there are further elements for class ‘**design**’ objects created by design combination functions. These are not discussed here. For such designs, the entry types given in the table can also be replaced by lists of several such entries.

A.2. Functionality for class ‘design’ objects

Package **DoE.base** offers inspection functionality (**print**, **summary**, **plot** as methods for their generics, **corrPlot** as a standalone function) and analysis methods (**plot**, **lm**, **halfnormal**) for this class, as well as a sub-setting method using **[**, which is useful for reordering experimental runs (to switch between standard order and randomized order, to re-randomize).

Besides these major functions, there are various further functions tailored to the need of class ‘**design**’ objects, either as methods for generic functions or just as functions:

Element	Data type	Role
<i>Mandatory elements</i>		
type	character string	identifies the type of design
nruns	number	number of runs (replications not counted)
nfactors	number	number of factors
factor.names	named list	factor names and factor levels (or scale ends for quantitative factors)
replications	number	number of replications or repeated measurements per run
repeat.only	logical	if TRUE, the number given in the replications element refers to repeated measurements only
randomize	logical	if TRUE, run order has been randomized
seed	number	the seed used for randomization
creator	call or list of menu settings in GUI	the creation history of the object
<i>Optional general elements</i>		
response.names	vector of character strings	names of the response columns (column names from the data frame)
<i>Elements for blocked designs (from functions <code>fac.design</code>, <code>rerandomize.design</code>, or <code>FrF2</code>)</i>		
block.name	character string	name of block variable
nblocks	number	number of blocks
block.gen	number	Yates matrix column number(s) of factor(s) used for blocking (<code>FrF2</code>) or block generator matrix (<code>fac.design</code>)
blocksize	number	run size of each block (without replications)
bbreps	number	number of between block replications (identical to replications)
wbreps	number	number of within block replications (these can be proper replications or repeated measurements only)
<i>Further element for full factorial designs and orthogonal array designs (functions <code>fac.design</code> or <code>oa.design</code>)</i>		
nlevels	numeric vector with nfactors elements	number of levels for each factor
<i>Further elements for general orthogonal array based designs (function <code>oa.design</code>)</i>		
generating.oa	character string	name of the OA used (from catalog of orthogonal arrays or user-provided)
selected.columns	numeric vector with nfactors elements	column number from <code>generating.oa</code> for each factor
origin	character string	origin specification from <code>generating.oa</code> (empty, if there is none)
comment	character string	comment from <code>generating.oa</code> (empty, if there is none)
residual.df	number	residual degrees of freedom for main effects only analysis

Table 2: Elements of the "design.info" attribute of class 'design' – Part I.

Element	Data type	Role
<i>Further element for designs created by function pb</i>		
ndummies	number	number of columns not assigned to experimental factors
<i>Further elements for designs created by function FrF2</i>		
aliased	list with character elements	information on the alias structure of the design up to degree 2
FrF2.version	character string	version number of package FrF2 , when design was created
generators	character vector or design key matrix	design generators in the format “D = ABC” etc. for FrF2 designs or design key matrix for designs created with package planor
catlg.name	character string	name of the catalog used for design creation
catlg.entry	list of length 1 of class ‘ catlg ’	the catalog entry used for the design
ntreat	number	identical to nfactors , present for blocked designs for historical reasons
aliased.with.blocks	character vector	lists two-factor interactions that are aliased with the block main effect
base.design	character string	element of design catalog used for creating a blocked or split plot design from
nfac.WP	number	number of whole plot factors
nfac.SP	number	number of split plot factors
nWPs	number	number of whole plots
plotsize	number	run size of each plot (without replications)
res.WP	number	resolution of the whole plot portion of the design
map	numeric vector with <i>k</i> elements	mapping of base factors so that estimability or randomization restriction requirement is fulfilled
orig.fac.order	numeric vector with nfactors elements	order of original factors from function call for split plot designs
clear	logical	if TRUE, the design is clear (for estimability requirement)
res3	logical	if TRUE, resolution III has been permitted for estimability request
quantitative	logical vector with nfactors elements	TRUE elements indicate quantitative factors
ncube	number	number of cube points in a design with center points
ncenter	number	number of center points
coding	list of formulae	coding of quantitative factors (for use with package rsm ; Lenth 2009)

Table 3: Elements of the "design.info" attribute of class ‘**design**’ – Part II.

Element	Data type	Role
<i>Further elements for designs created by function Dopt.design</i>		
		The function also outputs the split plot related elements and element quantitative described under specific elements for FrF2 above.
plot.name	character string	name of whole plot factor in split plot request
digits	number	number of digits to which quantitative factors are to be rounded
formula	formula	model formula for which <i>D</i> -optimality is to be achieved
constraint	logical expression	constraint applied to the experimental factors
optimality.criteria	named numeric vector	performance of design on several optimality criteria
<i>Further elements for designs created by function lhs.design</i>		
subtype	character string	type of latin hypercube design
The function also outputs digits and optimality.criteria elements described under specific elements for Dopt.design above.		

Table 4: Elements of the "design.info" attribute of class 'design' – Part III.

- Functions **design.info**, **run.order**, and **desnum** extract or set the respective attributes of a class 'design' object.
- Functions **factor.names** and **response.names** get or set the respective element from the "design.info" attribute.
- Function **generators** extracts generating information for designs of type **FrF2** or **planor**.
- Function **getblock** extracts block information for replicated designs (see also Section 2).
- Function **rerandomize.design** allows to re-randomize a class 'design' object without response data; this also allows explicit blocking of designs created with functions **oa.design** and **pb** (see also Section 5.3).
- Functions for designs that can be both in wide or in long format, i.e., parameter designs and designs with repeated measurements, can change between long and wide format or aggregate wide format designs: Function **rep2wide** brings repeated measurement designs into wide format, **reptolong** does the opposite; function **paramtowide** brings a parameter design to wide format (irreversible, they are created in long format); the **aggregate** method for 'design' objects aggregates designs in wide format into designs with a single response.
- Functions **add.response** and **col.remove** add responses or remove columns; function **response.names** can also be used to remove response columns.
- Function **qua.design** influences, which design columns are quantitative, and function **change.contr** changes the contrasts of design columns.

- Functions `undesign` and `redesign` remove the class ‘`design`’ properties from an object or reinstate them. Using function `data2design` instead of `redesign` on the result of `undesign` allows to treat a block factor as a normal experimental factor without a special role (see Section 5.3).
- Functions `cross.design` and `param.design` combine class ‘`design`’ objects into crossed designs or parameter designs, respectively.
- The `fix` method for ‘`design`’ objects has been adapted from package `utils`; it allows to edit ‘`design`’ objects; however, its use is not recommended.
- Function `export.design` can export a design in HTML or CSV format, together with an R workspace for the design. After entering response data in a spreadsheet program, responses can be added to the design itself using the `add.response` function.

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