Another Look at Supersaturated Designs

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ABSTRACT

A designed experiment in which the number of factors are at least as large as the number of runs is referred to as a supersaturated (SS) design. Recently these designs have received increased attention. Construction of such designs and analysis of data from these designs have been discussed by several authors. Our objective in this paper is to examine these designs and inform practitioners that the correlation structure inherent in SS designs can obscure real effects or promote non-real effects. Hence one should be cautious with the use of SS designs.

Key Words: Hadamard matrices, Misspecification, Plackett and Burman Designs, Rubber data

^{*}Work was done while on leave at the University of Waterloo.

1 Introduction

In industrial experimentation it is natural to look at many factors simultaneously. Cost considerations will sometimes make it impractical to design experiments so that effects of all these factors are simultaneously estimated. Assuming that only a few factors are active (having large effects), there have been attempts to identify them with high probability. The supersaturated (SS) designs used for this purpose have fewer observations than the number of factors considered (see, for example, Booth and Cox (1962), Lin (1993)). The objective usually is to identify all or most of the active factors. The intention of this paper is to investigate these designs and their ability to identify active factors.

It is assumed that there is no interaction between the factors and that the effect is linear in the level of each factor. Hence the basic model for observations y_i is given by

$$y_i = \beta_0 + \sum_{j=1}^m \beta_j X_{ij} + \epsilon_i , \ i = 1, 2, \dots, n \qquad n < m$$
 (1.1)

where m is the number of factors and n is the number of observations (runs). Here X_{ij} is the level of factor X_j in run i and is restricted to two values (+1 or -1). The parameter β_j (j = 1, ..., m) represents the effect of factor j and it is assumed that only a few of these parameters are different from 0. As the number of parameters is more than the number of observations, usual least square methods are not applicable for analyzing the data and some sequential procedure is needed. Lin (1993) suggests forward stepwise regression, while a second possibility is an exhaustive search over all models with a given number of predictors.

There are several methods of constructing SS designs (choosing "X" matrix). These include Wu (1993), Lin (1993), Nguyen (1996), Tang and Wu (1997), and Yamada and Lin (1997). Lin (1993) suggested using half fractions of Hadamard matrices. We will refer to this as the half replicate (HR) method. In Section 2 we will consider this procedure and discuss a set of data (the Rubber Data) which was previously considered by Williams (1968). Section 3 discusses the issue of misspecification and gives a simulation study to investigate the merits of the HR supersaturated design. Section 4 gives some concluding remarks.

2 Rubber Data

In order to illustrate the performance of the HR supersaturated design, Lin (1993) considered a set of data originally reported in Williams (1968) who, in the context of a rubber making process, used a Plackett-Burman (PB) design with 24 factors and 28 runs. The design and the data are given in Table 6 in the Appendix. As was pointed out by Box and Draper (1987) there was a typographical error in one of the entries of the original data (8th element from top in column 20) and this is corrected

in our table. Also columns 13 and 16 are identical and hence we eliminate column 16; however we keep the labeling of the other factors the same.

The HR method can be summarised as follows: In a standard Hadamard matrix whose first column entries are all +1, take any other column as a branching column. Consider the matrix formed by the rows which have the same sign in the branching columns. Assign factors to the columns other than the first and the branching columns. The resulting matrix is recommended as an SS design.

The HR construction procedure implies that the design obtained is not unique. In the current example since 23 factors are being investigated we consider a Hadamard matrix of order 28 with +1 in the first column. The 23 factors given in Table 6 can be assigned to any 23 of the other 27 columns of the Hadamard matrix. The remaining 4 columns of the Hadamard matrix can be considered as branching columns and can be used to produce 8 designs (see Vijayan (1976)). The rows corresponding to these 8 designs are given in Table 7 in the Appendix. For the analysis, we assume that the data for each run in the different designs are the same as those obtained by Williams. Lin (1993) proposed design number 5 for this situation and used a stepwise regression procedure to identify the active factors. We adopted the same procedure for each of these 8 designs (the stepwise procedure in SAS - Statistical Analysis System). For inclusion and exclusion the default option ("significance" level .15) was used first. Then we considered the level .075 to get results for design 5 to be close to those of Lin (1993). Table 1 shows the factors which were picked as active by each design. For instance, in design 1 with significance level (SL) = .075, factor 17 was selected as active in the first step and factors 15, 4, 22 and 10 were selected in subsequent steps. As can be seen from the table, different factors are chosen in different designs.

In the original study Williams (1968) selected factors 15, 20, and 17 as having major effects and 4, 22, 14 and 8 as having moderate effects. Combining process knowledge with the experimental results Williams chose factors 15, 10, 20 and 4 for subsequent use. Design 5 leads to factors 15, 12, 20, 4 and 10 as active. This set is somewhat similar to the one Williams (1968) used. However, factors 17 and 22 were not chosen as active in design 5; instead 12 and 10 were selected.

We adopted the 'best subset selection' procedure also to see how the results would compare. Procedures available in the usual packages such as Splus and SAS need to be modified for this situation since there are more columns (factors) than rows (runs) in the "X" matrix. Hence we restricted our search for the best subsets of size 5 or less factors; note that there are $\binom{23}{i}$ subsets of size i (i=1,2,3,4,5). The results are shown in Table 2. For example in design 1, factors 1, 15 and 20 provide the "best subset" of size 3; 1, 14, 15, and 20 provide the "best subset" of size 4; 1, 3, 14, 15, and 20 lead in subsets of size 5. Here best is in the sense of R^2 . From Table 2 we note the following.

(i) Different designs lead to different choices with subsets of size i = 3, 4, 5. No single design identifies the same five factors as an analysis of the full 28 runs.

Table 1: Rubber Data: Factors included after Five Steps in a Stepwise Procedure

	Steps	1	2	3	4	5	R^2
Design							
1	SL = .15	17	15	4	22	10	.90
	SL=.075	17	15	4	22	10	.90
2	SL=.15	15	24	18	13	8	.97
	SL = .075	15	24	18	13	8	.97
3	SL=.15	15	20	3	4	22	.92
	SL=.075	15	20	_	-	-	.73
4	SL=.15	2	13	8	3	20	.84
	SL=.075	2	13	8	-	-	.69
5*	SL=.15	15	12	20	4	10	.97
	SL=.075	15	12	20	4	10	.97
6	SL=.15	4	22	23	18	24	.88
	SL=.075	4	22	23	_	-	.74
7	SL=.15	14	12	11	23	-	.89
	SL=.075	14	12	11	23	-	.89
8	SL=.15	15	22	8	17	1	.94
	SL=.075	15	22	8	17	1	.94

Notes:

* Design 5 is the same as that used in Lin (1993)

"-" indicates that the procedure stopped before

- (ii) Stepwise regression often finds poorer models than best subsets regression. In designs 2 and 5 the two methods identify the same factors (up to step 5). In design 7 the two procedures lead to the same factors up to step 4 and in design 8 only up to step 3. In other designs they lead to different factors.
- (iii) If R^2 is used as a criterion and subsets of size 5 are considered, then design 3 is the "winner" with factors 1, 5, 8, 15 and 21; design 5 is a close contender with factors 4, 10, 12, 15, and 20.
- (iv) When the best subset selection procedure was used with the original PB design, factors 4, 15, 17, 20, and 22 were selected as the best five variable model. A half normal plot of effects for the original PB design identifies the same five factors as having the largest effects. The plot indicates that factor 15 is important; it is difficult to make concrete statements about other factors.

The important point is that different designs lead to the identification of different factors as active, which in turn would lead to different conclusions.

Table 2: Rubber Data: factors from 'best subset selection' procedure. The row marked "all" refers to results using all 28 runs. Each row corresponds to a model identified by best subsets, with a • indicating an active factor.

Design	1	2	3	4	5	6	7	8	9	10	11	12	13	14	15	17	18	19	20	21	22	23	24	R^2
all				•											•	•			•		•			0.6881
	•										<u></u>				•				•					0.7696
1	•													•	•				•					0.8736
	•			•										•	•				•					0.9360
															•		•						•	0.8869
2													•		•		•						•	0.9348
								•					•		•		•						•	0.9666
					•			•							•									0.8544
3	•				•			•							•									0.9568
	•				•			•			_				•					•				0.9741
		•				•													•					0.6938
4		•				•										•			•					0.8505
		•				•								•		•			•					0.9733
												•			•				•					0.8705
5				•								•			•				•					0.9548
				•						•	_	•			•				•					0.9730
	•	•																				•		0.7922
6	•	•													•							•		0.8954
	•	•									_				•			•				•		0.9549
											•	•		•										0.7747
7											•	•		•								•		0.8929
								•			•	•		•								•		0.9170
								•							•						•			0.7889
8															•				•	•	•			0.8734
	•							•							•	•					•			0.9399

3 Misspecification

The discussion in the previous section points to two issues: (i) Choice of a design. (ii) Misspecification. We now discuss the second issue more fully.

When the number of factors is very large it is inevitable that some of the inactive factors would be selected with high probability. This is a serious problem but may be manageable as long as active factors are also selected with high probability. However, sometimes the procedures fail to select active factors and select inactive factors instead (see, for instance, Chipman et. al. (1997), Hamada and Wu (1992) and Wu (1993)). The nonorthogonality of the columns of the design matrix X is the root of the problem.

To highlight the misspecification issue we conducted a simulation study. To verify some of the features of the supersaturated designs, we consider the two designs (14 run, 12 run) given in Lin (1993, 1995) and the model given in (1.1).

3.1 14 run design with 23 factors used in Lin (1993)

Step 1: Data generation:

We generate n = 14 N(0,1) random variates, $\epsilon_i, i = 1, ..., 14$ and then using the model (1.1) generate y_1, \dots, y_{14} for specific sets of values of the β 's.

Simulated Models:

(1) Pure noise: $\beta_j = 0$ for all j.

(2)
$$\beta_j = \begin{cases} .5, 1, 20 & j = 1 \\ 0 & j \neq 1 \end{cases}$$

(3)
$$\beta_2 = \beta_7 = 1$$
, other β 's zero $\beta_2 = \beta_7 = 20$, other β 's zero

(4)
$$\beta_{1} = \beta_{2} = \beta_{3} = 1,$$
 other β 's zero $\beta_{2} = \beta_{7} = \beta_{13} = .5,$ other β 's zero $\beta_{2} = \beta_{7} = \beta_{13} = 1,$ other β 's zero $\beta_{2} = 5, \beta_{7} = 10, \beta_{13} = 20,$ other β 's zero $\beta_{2} = 14, \beta_{7} = 20, \beta_{13} = 20,$ other β 's zero $\beta_{2} = \beta_{7} = \beta_{13} = 20,$ other β 's zero other β 's zero

Step 2: Analysis for factor selection.

We used forward selection with a specific SL for including a factor, stop after step 5 and record the selected factors.

Step 3: Repeat Steps 1 and 2, N = 200 times and record the number of times each factor was selected.

For this particular design, correlations between columns are either ± 0.14 or ± 0.43 .

 X_1 has a correlation of 0.43 with each of X_2, X_7, X_{13} , making case (4) a challenging problem. In models with single predictors, an effect will have t-ratio $\beta_j/\text{se}(\beta_j) = \beta_j/(\sigma/\sqrt{n})$. For n = 14 and $\sigma = 1$, even a coefficient of $\beta_1 = 1$ is large, since the expected t-ratio would be $\sqrt{14} = 3.74$. Coefficient values of 14 and 20 are very large indeed.

In the pure noise case with SL = .05, 79% of the times at least one factor was selected as active in 5 steps. This was not unexpected. With 23 contrasts tested to zero individually and assuming them to be independent (this is not quite true in our case) there is a chance $1 - (1 - \alpha)^{23}$ of selecting at least one significant contrast where α is the SL for inclusion. When $\alpha = .05$ this probability = .70 which is in the vicinity of the simulated value. We cannot avoid this situation even if we have a large number of observations. On the other hand if the present experiment is considered as a device to reduce the number of factors which need to be carefully examined in a later experiment, then it is important to know whether the procedure leads to the active factors with high probability. For this purpose we consider cases 2-4 where one, two, or three β 's are different from zero. The results from a stepwise procedure are presented in Table 3. We note the following from the table:

- (i) When only one β is different from zero or it is much larger than the rest, the corresponding factor gets selected in the first 5 steps. For instance, with a $SL=.1,~X_1$ gets selected 100% of the cases if $\beta_1=20,~87\%$ of the cases when $\beta_1=1$ and only 36.5% when $\beta_1=.5$. When SL=.05 and $\beta_1=.5,~X_1$ gets selected only in 36% of cases.
- (ii) When two β 's are important and very large ($\beta_2 = \beta_7 = 20$), the corresponding factors get selected about 78% of the time. This drops to 67% when $\beta_2 = \beta_7 = 1$. In both cases the inactive factor X_1 is selected in the first forward step about 22% of the time.
- (iii) When more than two β 's are important, conclusions are different depending on which β 's are non zero. For example when $\beta_1 = \beta_2 = \beta_3 = 1$ and other β 's are zero (situation 4), with SL = .1 the active factors X_1, X_2, X_3 get selected about 55% of the time in five steps. However, if $\beta_2 = \beta_7 = \beta_{13} = 1$ and other β 's are zero (active factors are now assigned to columns 3, 8, and 14 but the design is the same), then with SL = .1 the combination (X_2, X_7, X_{13}) get selected only 4% of the time in five steps and .5% of the time in three steps; X_1 gets selected 91.5% of the time in the first step. The same value (=1) was used for both cases but the active factors were in different columns of the matrix and the non orthogonality of the columns creates the difficulty.

Even when effects are very large, forward selection can be mislead. When $\beta_2 = 14$, $\beta_7 = \beta_{13} = 20$ or $\beta_2 = \beta_7 = \beta_{13} = 20$, the combination (X_2, X_7, X_{13}) is never selected while X_1 is selected 100% of the time. On the other hand, if $\beta_2 = 5$, $\beta_7 = 10$, $\beta_{13} = 20$, the combination is selected in 100% of the cases.

Table 3: Simulation Results: Selection Pattern of Factors, Lin's 14 Run Design

	Factors Corresp- Percent of times real effect factors selected Percent of time X_1													
Factors	Corresp-	Percent of	times real eff	ect factors	selected	Percent of	time X_1							
with Real	onding β	\mid in k steps		in 5 steps		selected in	the first							
Effects	values	k = # fac	tors in col. 1			step								
		SL=0.10	SL = 0.05	SL=0.10	SL = 0.05	SL = 0.10	SL=0.05							
	0.5	36.5	27.5	56.5	38									
X_1	1	87	85.5	93	90.5									
	20	100	100	100	100									
				1										
X_2, X_7	1,1	55	56	67	62.5	24.5	22							
	20, 20	78	74	78	74	22	26							
$X_1, X_2,$	1, 1, 1	40.5	38	55	45.5									
X_3														
	0.5,0.5,0.5	0	0	1	1	43.5	46.5							
$X_2, X_7,$	1, 1, 1,	.5	.5	4	2	91.5	91							
	5,10,20	100	100	100	100	0	0							
X_{13}	14, 20, 20	0	0	0	0	100	100							
	20,20,20	0	0	0	0	100	100							

The major point is that factors selected as active depend on which columns of the design matrix are assigned to the real active factors and that is difficult to do in advance. This is a result of the correlation structure of the columns of the design matrix.

We also performed simulations for cases (2) and (4) and used the best subset selection procedure to see how the selection pattern would be affected. The results are given in Table 4. We note the following.

- (i) Best subsets performs notably better than stepwise in the more difficult problem involving X_2, X_7 and X_{13} . It finds the correct model more than a third of the time when all three coefficients are 1. In the cases of $(\beta_2, \beta_7, \beta_{13}) = (14, 20, 20)$ and (20, 20, 20) where stepwise fails to find the right model, best subsets identifies the model 100% of the time.
- (ii) Unlike stepwise, the best subsets algorithm can identify different models of a given size with good fit. For this reason, Table 4 also gives the percentage of times that the correct model is among the five best of a given size. So for example, when $\beta_2 = \beta_7 = \beta_{13} = 1$, this model was ranked first among 3 term models by best subsets 33% of the time. It appeared in the top five models of size three 57% of the time.

Table	4:	Simulation	Results	with	Best	\mathbf{Subset}	Proceed	lure,	Lin's	14	Run	Design
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	Percent times correct factors													
Factors with	β -values	selected in	in subsets of size k *											
Real Effects		k = 3	k=4	k=5										
X_1, X_2, X_3	1,1,1	52 (77)	50 (77)	51 (70)										
X_2, X_7, X_{13}	.5,.5,.5	1 (8)	4 (9)	5 (15)										
	$1,\!1,\!1$	33 (57)	44 (65)	43 (67)										
	$5,\!10,\!20$	100 (100)	100 (100)	100 (100)										
	$14,\!20,\!20$	100 (100)	100 (100)	100 (100)										
	$20,\!20,\!20$	100 (100)	100 (100)	100 (100)										

^{*} Note: The first number gives frequency that the highest ranked model contained all active factors. The number in parentheses gives the frequency with which one of the five highest ranked models contained all the active factors.

Table 5: A Hadamard Matrix of Order 12

1	2	3	4	5	6	7	8	9	10	11	12
1	1	$-\mathbf{u_3}$	1	$-\mathbf{u_3}$	$-\mathbf{u_3}$	$-\mathbf{u_2}$	$-\mathbf{u_2}$	$-\mathbf{u_2}$	$\mathbf{u_1}$	\mathbf{u}_1	\mathbf{u}_1
1	1	$\mathbf{u_3}$	-1	$\mathbf{u_1}$	$\mathbf{u_2}$	$\mathbf{u_2}$	\mathbf{u}_1	$\mathbf{u_3}$	$-\mathbf{u_3}$	$-\mathbf{u_2}$	$-\mathbf{u_1}$
1	- 1	1	$\mathbf{u_1}$	$\mathbf{u_3}$	$\mathbf{u_2}$	$-\mathbf{u_2}$	$-\mathbf{u}_1$	$\mathbf{u_2}$	$-\mathbf{u_3}$	\mathbf{u}_1	$\mathbf{u_3}$
1	-1	-1	$-\mathbf{u}_1$	$-\mathbf{u_2}$	$-\mathbf{u_3}$	$\mathbf{u_1}$	$\mathbf{u_2}$	$-\mathbf{u_2}$	$\mathbf{u_3}$	$-\mathbf{u_3}$	$-\mathbf{u}_1$

Since SS designs are intended to screen variables, it is quite likely that practitioners will examine more than one model. The findings of this simulation study indicates that even when SS designs are used in this fashion, it is possible that the correct factors will not be identified.

3.2 12 run design

Lin (1995) presented an algorithm to generate certain SS designs and designs for 12 runs. General construction methods are available to produce these designs with number of factors as high as 66. Basically the design is constructed from the columns of a standard Hadamard matrix. Using the notation $\mathbf{1}' = (1, 1, 1), \mathbf{u}'_1 = (1, -1, -1), \mathbf{u}'_2 = (-1, 1, -1),$ and $\mathbf{u}'_3 = (-1, -1, 1)$ we can write down one such Hadamard matrix of order 12 as given in Table 5.

The different columns of the design matrix would be obtained by taking componentwise products of two columns of this basic matrix leading to a design matrix with

66 columns. We denote the componentwise product of the columns i and j by " $i \times j$ ". Lin (1995) demonstrated the usefulness of such SS designs by carrying out a simulation study using the above design. In the simulation study factors corresponding to columns $1 \times 4, 1 \times 8$, and 8×10 were given large effects (β values 17, 24 and 15 respectively), those corresponding to 1×11 , and 8×11 were given moderate effects (β values 3), and others were taken as zero. The results of the simulation indicated that the 5 active factors were always selected in the first five steps of a forward selection procedure.

We wish to demonstrate that the results depend on which columns of the X matrix correspond to the large and moderate effect factors. In our simulation study we used the same design as in Lin (1995) and we assigned the β values 17, 24, 15 (large effects in Lin's simulation) to columns 8×10 , 6×12 , and 9×11 and the β value 3 to columns 3×5 and 4×7 , and zero to the remaining columns. 12 observations were generated from model (1.1) using the β -values indicated before and a forward selection procedure as in Lin (1995) was used to select the active factors. In 200 repetitions, none of the active factors corresponding to columns 6×12 , 9×11 , 3×5 and 4×7 were picked up in the first five steps while the factor corresponding to column 8×10 was selected only 29% of the time. The factor corresponding to column 1×2 (inactive) was always the choice in the first step.

4 Additional Discussion and Concluding Remarks

Data analysis in Section 2, and the simulations and discussion in Section 3 indicate that one should be very cautious with the use of SS designs. The simulations, in particular show that there is a high chance of missing the real active factors and selecting the inactive ones instead. The assignment of factors to columns of these designs is crucial because of the correlation structure among the columns of the design. If SS designs are to be used, the results of Section 3 indicate that best subsets variable selection should be used rather than stepwise regression.

In Section 2 we commented that some additional criterion is necessary to select a design. One suggestion made in this regard is to consider a design in which the average value of S_{ij}^2 , where $S_{ij} = X_i X_j'$ (say, $E(S^2)$), is made as small as possible (for example see Booth and Cox (1962), Wu (1993)). In the case of a 12-run design with 13 factors it can be shown that the minimum value of $E(S^2) = 48$ and is attained by a design given in Wu (1993). In the eight different 14-run designs considered in Section 2, $E(S^2)$ is either 7.79 (for subsets 1 and 3) or 7.92 (for other subsets), leaving little theoretical grounds for distinguishing between the eight designs. The dramatically different conclusions found in the eight different subsets appears to be mainly due to random variation in the data.

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Appendix

Run											Fa	cto	r N	um	ber										Response
Number	1	2	3	4	5	6	7	8	9	10	11	12	13	14	15	16	17	18	19	20	21	22	23	24	y
1*	+	+	+	_	_	_	+	+	+	+	+	_	+	_	_	+	+	_	_	+	_	_	_	+	133
2	_	+	_	_	_	_	+	+	+	+	_	+	_	_	+	_	+	+	+	+	_	+	+	_	49
3*	+	_	_	_	_	_	+	+	+	_	_	_	+	+	+	+	_	+		_	+	+		_	62
4*	+	+	_	+	+	_	_	_	_	+	_	+	+	+	+	+	+	_	_	_	_	+	+	_	45
5	+	+	_	_	+	+	_	_	-	-	_	_	_	_		_	+	+	_	+	+	+	_	+	88
6*	+	+	_	+	_	+	_	_	_	+	+	_	+	_	+	+	_	+	+	+	_	_	_	_	52
7	_	_	+	+	+	+	+	+	_	+	_	+	+	_	_	+	_	+	_	_	_	_	_	_	300
8*	_	_	+	+	+	+	_	+	+	_	_	_	+	_	+	+	+	_		+	_	+	+	+	56
9*	_	_	+	+	+	+	+	_	+	+	+	_	_	+	+	_	+	+	+	+	+	+	_	_	47
10*	_	_	_	_	+	_	_	+	_	+	_	+	+	+	_	+	+	+	+	+	+	_	_	+	88
11	+	_	+	_	_	+	_	_	+	_	_	_	+	+	_	+	+	+	+	_	_		+	_	116
12	_	+	+	+	_	_	+	_	_	+	+		+	+	_	+	+	+	_	_	+	+	+	+	83
13*	_	+	+	_	-	+	_	+	_	+	_	_	_		_	_	_		+	_	+	+	+	_	193
14	_	_	_	+	_	_	_	_	+	+	_		_		+		_	_	_	_	+	_	_	+	230
15	+	_	+	_	+	_	+	_	_	+	_	_	_	+	+	_	_	_	+	+	_	_	+	+	51
16	_	+	+	_	+	_	_	_	+	_	+	+	+	_	+	+	+	_	+	_	+	_		_	82
17*	_	_	_	_	-	+	+	_	_		+	+	_	_	+	_	+	+	_	_	_		+	+	32
18	+	_	+	+	-	_	_	+	_	-	+	+	+	_	+	+	-	+	+	+	+	+	+	+	58
19	+	_	_	+	+	_	+	+	_	_	+		_	_	_	_	+	_	+	_	_	+	_	_	201
20	+	+	+	_	+	+	_	+	+	+	+	+	_	+	+	_	_	+	_	_	_	+	_	+	56
21	_	+	_	+	_	+	+	_	+		_	+	+	+		+	_	_	+	+	_	+	_	+	97
22^*	+	+	+	+	_	+	+	+		_	_	+		+	+	_	+	_	+		+		_	+	53
23^*	_	+	_	+	+	_	_	+	+	_	+	_	_	+	_	_	_	+	+	_	_		+	+	276
24^*	+	_	_	_	+	+	+	_	+	+	+	+	+			+		_	+	_	+	+	+	+	145
25^*	+	+	+	+	+	_	+		+	_	_	+	_	_	_	_	_	+	_	+	+	_	+	_	130
26	_	+	_	_	+	+	+	+		_	+	_	+	+	+	+	_	_	_	+	+		+	_	55
27	+	_	_	+	_	+	_	+	+	+	+	+	_	+	_	_	+	_	_	+	+	_	+	_	160
28*	-	_	+	_		_	_	_	_	-	+	+	_	+	_	_	_	-		+	_	+	-	_	127

Table 6: Rubber data, from Williams (1968). Runs used in Lin's (1993) design are marked with a *. The level of run 8, factor 20 is here shown as +. In the source reference, the lower level appears: that seems to be a typographical error.

Design															
Number	Rows from the PB design														
1 (2)	1	2	4	5	7	9	11	14	18	22	23	24	26	28	
3 (4)	1	3	4	5	7	9	13	15	16	17	18	21	23	27	
5 (6)	1	3	4	6	8	9	10	13	17	22	23	24	25	28	
7 (8)	1	4	9	10	11	13	14	17	18	19	20	21	25	26	

Table 7: 14-run supersaturated designs from a Hadamard matrix of order 28. Even numbered designs (in parentheses) are the complement of the rows listed.