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# Techniques for sensitivity analysis of simulation models

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Publication date: 1990

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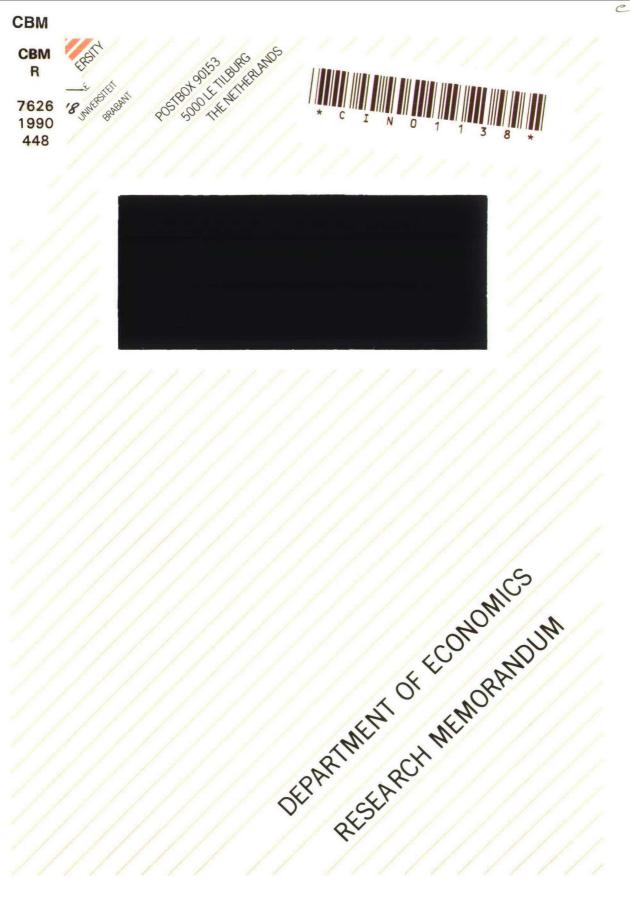
*Citation for published version (APA):* Kleijnen, J. P. C., van Ham, G., & Rotmans, J. (1990). *Techniques for sensitivity analysis of simulation models: A case study of the CO2 greenhouse effect.* (Research memorandum / Tilburg University, Department of Economics; Vol. 448). Unknown Publisher.

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TECHNIQUES FOR SENSITIVITY ANALYSIS OF SIMULATION MODELS: A CASE STUDY OF THE CO<sub>2</sub> GREENHOUSE EFFECT

Jack P.C. Kleijnen and Greet van Ham

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# ERRATUM

The cover page should be:

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FEW 448

Tilburg, July 27, 1990

Kenmerk: 320:90.225

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a case study of the  $\rm CO_2$  greenhouse effect

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Techniques for sensitivity analysis of simulation models:

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#### Abstract

Sensitivity analysis is needed for validation, what-if analysis, and optimization of complicated simulation models. One set of techniques for sensitivity analysis are least squares curve fitting, regression analysis, and statistical designs such as factorial designs. These techniques are applied to several modules of a large integrated assessment model for the greenhouse effect, developed in The Netherlands. The regression models turn out to be valid approximations to the simulation models. Some estimated effects are quite surprising for the simulation users.

Keywords: What-if analysis, validation, optimization, least squares, regression analysis, experimental design, factorial design, ecology, greenhouse effect, simulation, metamodel.

#### Introduction

Complicated simulation models have been constructed in many disciplines. All these models confront the analysts with the problem of sensitivity analysis; that is, what are the effects of changing the parameters and input variables of the simulation model? That question arisis in 'what if' analysis, validation, optimization, and so on. This article introduces and illustrates the application of *simple* techniques that originated in the discipline of mathematical statistics. These techniques are least squares curve-fitting, regression analysis, and statistical designs such as  $2^{k-p}$  designs. The techniques are applied to an integrated assessment model for the greenhouse effect. This model has been developed at a large Dutch institute called National Institute of Public Health and Environmental Protection (or RIVM in Dutch).

One of the major imminent ecological threats of the world is the 'enhanced greenhouse problem': the earth and the lower layers of its atmosphere have shown rising temperatures over the past hundred years. This heating phenomenon is probably caused by an increase of greenhouse gases (such as carbon dioxyde, methane, and ozon) that absorb the earth's heat radiation, so the global average temperature rises. Mankind is largely this responsible for increased 'greenhouse' gas concentration. Temperatures are expected to rise, but with different amounts in different regions of the earth (the tropics will be less affected probably). Higher temperatures will cause thermal expansion of the oceans and melting of artic ice, which raise the sea level. Many more processes, however, are involved; see [IPCC, 1990]. One consequence of a higher sea level is the need to raise the level of the dikes in the Netherlands. A survey of the effects for society is given in [Gezondheidsraad, 1987].

To gain quantitative insight into the greenhouse problem and develop long-term strategies for coping with climatic changes, RIVM developed the Integrated Model for the Assessment of the Greenhouse Effect or IMAGE. This model is a deterministic simulation (but most of the sensitivity techniques applied to this model can also be used in random simulation models). The state of the dynamic biospheric system is computed per half year, up to the year 2100, while starting at the year 1900. The model is composed of modules, which treat specific parts of the greenhouse problem. Modules get inputs from other modules. Also see figure 1 and the references [Rotmans, 1990; Rotmans, et al. 1990; den Elzen & Rotmans, 1988].

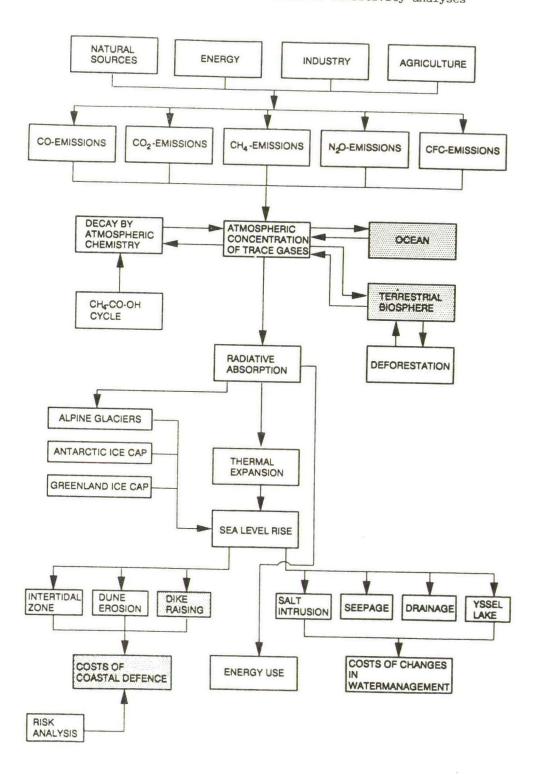
The sensitivity analysis techniques are applied to several modules. This paper concentrates on the carbon-cycle module; the dike raising modules are briefly discussed; see the shaded modules in figure 1.

Note that there are alternative techniques for sensitivity analysis. Latin Hypercube sampling is a Monte Carlo method, which is discussed at length in [Iman and Helton, 1988] and criticized in [Easterling, 1986]; also see [Kleijnen, 1987, p. 143-145]. This technique was applied to several IMAGE modules in [Lammerts, 1989]; it gave results similar to the results of this paper. More sophisticated techniques do not treat the simulation model as a black box; they use analytical differential analysis; see [McRae, 1989].

This article is organized as follows. First the need for sensitivity analysis is discussed, and the greenhouse case study is introduced. Then metamodels, which explain the input/output behavior of the underlying simulation model, are explained. The coefficients of the metamodel are estimated by least squares regression analysis. The resulting metamodel can be validated. Closely related to the metamodel specification is the selection of an efficient experimental design. All techniques are demonstrated by their application to several modules of the greenhouse simulation model.

#### Metamodeling through regression analysis

A simulation model maps its inputs into one or more outputs; hence a simulation model is a mathematical function (say) s( ). The inputs are parameters, input variables, and behavioral relationships (or submodules); Figure 1: The modules of *IMAGE*; the shaded modules are submitted to sensitivity analyses



see [Kleijnen, 1987, p. 136]. These inputs are called *factors* in the statistical design of experiments. They may be represented by  $z_j$  with  $j = s1, \ldots, k$  and  $k \ge 1$ . In the greenhouse model all factors are quantitative, but the techniques also apply to qualitative factors. The case study concentrates on a single output variable (say) y, namely the global average atmospheric CO<sub>2</sub> concentration in the year 2100. If there are several outputs, the technique can be applied per output. This yields

$$y = s(z_1, ..., z_j, ..., z_k).$$
 (1)

A mathematical function may be approximated by a Taylor series, under certain mathematical conditions. Suppose the initial approximation is

$$\hat{y} = \gamma_0 + \Sigma_{j=1}^k \gamma_j z_j + \Sigma_{j=1}^{k-1} \Sigma_{h=j+1}^k \gamma_{jh} z_j z_h,$$
(2)

where  $\gamma_0$  is called the overall or grand mean;  $\gamma_j$  is the first-order or main effect of factor j; and  $\gamma_{jh}$  is the interaction between the factors j and h, that is, the effect of factor j depends on the level of factor h.

Note that the variables in approximation (2) may be functions of the variables in the simulation model (1); for example,  $\log(z_j)$  or  $1/z_j$ . The approximation remains linear in the parameters  $\gamma$ ; so linear regression analysis still applies; see [Kleijnen, 1987, pp. 160-161].

The approximation in (2) is called a *metamodel*: it is a model of the input/output behavior of the underlying simulation model. The Taylor series argument may be one inspiration for the specification of a metamodel. Because the mathematical conditions of the Taylor series do not hold in complicated simulation models, the validity of the metamodel must be checked. In other words, the metamodel is only an approximation. Before that model can be validated, it must be calibrated, that is, its coefficients or parameters  $\gamma$  must be estimated. Moreover there is a scaling problem. These issues are discussed now. For simplicity's sake the interactions in the metamodel (2) are ignored temporarily. If the input variable  $z_j$  increases by one unit, then the output changes by  $\gamma_j$  units. We assume, however, that sensitivity analysis is meant to determine the effects of changes of the inputs over the *whole* experimental area, in order to detect the importance of those inputs. (Next those important factors are further investigated to validate and optimize the simulation model. If only optimization were the goal, then local marginal effects would suffice.) So the importance of factor j is measured by the difference between the outputs at the lowest and the highest value of that factor. Denoting those two extreme values by  $L_j$  and  $H_j$  respectively (so the 'experimental area' is a k-dimensional rectangle), the original variables  $z_j$  yield the *standardized variables*  $x_j$ , which range between -1 and +1:

$$z_{j} = a_{j}x_{j} + b_{j}$$
 with  $a_{j} = \frac{H_{j}-L_{j}}{2}$  and  $b_{j} = \frac{H_{j}+L_{j}}{2}$ . (3)

The simple transformation (3) together with the original metamodel (2) yields the standardized metamodel:

$$\hat{y} = \beta_0 + \Sigma_{j=1}^k \beta_j x_j + \Sigma_{j=1}^{k-1} \Sigma_{h=j+1}^k \beta_{jh} x_j x_h .$$
(4)

It is simple to prove that  $\beta_j$  reflects the *importance* of factor j:  $\beta_j = \gamma_j (H_j - L_j)/2$ , ignoring interactions. See [Bettonvil and Kleijnen, 1990].

Note that to search for the optimum combination of the input factors, Response Surface Methodology (RSM) combines a first-order metamodel with the steepest ascent technique. That search should not use the original or standardized model but a centered model:

$$\hat{y} = \delta_0 + \Sigma_{j=1}^k \delta_j (z_j - \bar{z}_j) + \Sigma_{j=1}^{k-1} \Sigma_{h=j+1}^k \delta_{jh} (z_j - \bar{z}_j) (z_h - \bar{z}_h), \quad (5)$$

where  $\bar{z}_j = \sum_{i=1}^n z_{ij}/n$  assuming that n combinations of input factors are simulated. See [Bettonvil and Kleijnen, 1990].

Calibration means that the parameters of the model are quantified. So the metamodel's parameters  $\beta$  in (4) are estimated. Therefore the metamodel is fitted to the simulation data. Let q denote the number of parameters in the metamodel; for example, in (4) q equals 1 + k + k(k-1)/2. To get estimated parameter values  $\hat{\beta}$ , n combinations of the factor values are simulated. That set of simulated combinations yields the n × q matrix of independent variables X corresponding to the metamodel of (4):

$$X = \begin{bmatrix} 1, x_{11}, \dots, x_{1k}, x_{11}, x_{12}, \dots, x_{1,k-1}, x_{1k} \\ 1, x_{11}, \dots, x_{1k}, x_{11}, x_{12}, \dots, x_{1,k-1}, x_{1k} \\ 1, x_{n1}, \dots, x_{nk}, x_{n1}, x_{n2}, \dots, x_{n,k-1}, x_{nk} \end{bmatrix}.$$
 (6)

*Example*: Suppose there are three factors (k=3), which in combination i have the values +1, -1 and -1 respectively. (Remember that standardization means that factor 1 is at its highest level  $H_1$ , factor 2 is at its lowest level  $L_2$  in this combination, and so on; see equation 3.) Then the interaction variable  $x_1x_2$  has the value (+1) (-1) = -1 in this combination, and so on. Obviously  $\beta_0$  corresponds to the 'variable' that is +1 in all combinations. So row i of X equals  $x_1' = (+1, +1, -1, -1, -1, -1, +1)$ .

The output of combination i is  $y_i$ ; see (1). Fitting the metamodel to the simulation data, using the *least squares* citerion, yields the estimated parameters  $\hat{\beta}$ :

$$\hat{\underline{\beta}} = (\mathbf{X}'\mathbf{X})^{-1} \mathbf{X}' \mathbf{y} . \tag{7}$$

The least squares criterion yields unique estimates only if X is non-singular so that the inverse of (X'X) exists. A necessary condition is  $n \ge q$  (the number of simulated factor combinations is not smaller than the number of parameters in the metamodel). This condition, however, is not sufficient. For example, if the factors 1 and 2 are changed simultaneously in the n combinations, then their two columns are identical ( $x_{i1} = x_{i2}$  for i

= 1,...,n) and X is singular. Obviously X is not singular if all its columns are *orthogonal* ( $\Sigma_{i=1}^{n} \times_{ij} \times_{ih} = 0$ , etc.). Under certain statistical assumptions, an orthogonal matrix X is optimal; see the next section.

The calibrated metamodel can now be *validated*. One aspect is how well this metamodel fits the simulation data. One overall criterion is  $R^2$ :

$$R^{2} = 1 - \frac{\Sigma_{i=1}^{n} (\hat{y}_{i} - y_{i})^{2}}{\Sigma_{i=1}^{n} (y_{i} - \bar{y})^{2}}, \qquad (8)$$

where  $\bar{y} = \sum_{i=1}^{n} y_i/n$ . A 'perfect' fit means that  $y_i = \hat{y}_i$  for all i, so the upper limit for R<sup>2</sup> is 1. Unfortunately, a lower threshold for R<sup>2</sup> is hard to give. Therefore we propose to compute the relative errors  $(y_i - \hat{y}_i)/y_i$ , which can be 'eyeballed' by the user.

Validating a model, however, usually means that the model is used to forecast the output; next that forecast is confronted with the true output. Therefore *cross-validation* should be used. Delete one combination  $(x'_i, y_i)$  from the old data set (X,y); denote the remaining set by  $(X_{-i}, y_{-i})$ . Reestimate the metamodel's parameters  $\beta$ , analogous to (7):

$$\hat{\beta}_{-i} = (X'_{-i} X_{-i})^{-1} X'_{-i} y_{-i}$$
(9)

Predict the output of combination i, not using the data of combination i:

$$\hat{\mathbf{y}}_{\mathbf{i}} = \mathbf{x}_{\mathbf{i}}' \, \hat{\underline{\beta}}_{-\mathbf{i}} \, . \tag{10}$$

Compute the forecast errors

$$e_{-i} = \hat{y}_{-i} - y_{i}$$
 (11)

The user may again evaluate the relative errors  $e_{-i}/y_i$ . This procedure is repeated for all i (i=1,...,n).

The computation of the errors e<sub>i</sub> is possible without applying the least squares criterion n times (to n-1 combinations). First no data are eliminated; see (7). Next the so-called 'hat' matrix H is computed:

$$H = X(X'X)^{-1} X'$$
(12)

with diagonal elements h .... Then

$$e_{-i} = e_i / (1 - h_{ii})$$
 (13)

Many modern regression analysis packages give those 'leave one out residuals'. See [Kleijnen, 1987, p. 178] and [Atkinson, 1985, p. 13].

The mathematical analysis can be refined if a *statistical* (sub)model is added for the *fitting errors* e. Kleijnen (1987, p. 164) assumes that these errors are normally and independently distributed with common variance (say)  $\sigma^2$ . Then the least squares algorithm yields the Best Linear Unbiased Estimators (BLUE); that is, the estimators have minimum variances and correct expected values. Those variances equal the main-diagonal elements of the variance-covariance matrix of  $\hat{\beta}$ :

$$\operatorname{cov} (\hat{\boldsymbol{\beta}}) = (\mathbf{X}' \mathbf{X})^{-1} \sigma^2 . \tag{14}$$

The parameter  $\sigma^2$  can be estimated through the Mean Squared Error:

$$\hat{\sigma}^{2} = \sum_{i=1}^{n} (\hat{y}_{i} - y_{i})^{2} / (n - q) .$$
(15)

The estimated variances (or standard errors) yield the Student t statistic with n - q degrees of freedom:

$$t_{n-q} = (\hat{\beta}_j - \beta_j) / \hat{\sigma}_j \quad (j=1,...,q) ,$$
 (16)

where  $\hat{\sigma}_{j}$  denotes  $\{var(\hat{\beta}_{j})\}^{\frac{1}{2}}$  and  $\beta_{j}$  is the j th element of  $\hat{\beta}$ ; so  $\beta_{1}$  in (16) is identical to  $\beta_{0}$  in (4),  $\beta_{2}$  in (16) is  $\beta_{1}$  in (4),...,  $\beta_{q}$  in (16) is  $\beta_{k-1,k}$  in (4). The *significance* of  $\hat{\beta}_{j}$  can be tested statistically, using the t table for a given significance level or type-I error (say)  $\alpha$ ; for

example, for  $\alpha = 0.05$  and v = 12 the critical t value is 2.18 in a twosided test, which considers the absolute value of  $t_{n-\alpha}$ .

Note that a more sophisticated model for the fitting errors is used by Sachs et al. (1989). They assume that the errors are not independent but form a stationary process with a specific correlation function. Also see [Kleijnen, 1990].

#### Statistical design of experiments

The metamodel determines the experimental design;, for example, a model with interactions such as (4) cannot be calibrated through a design that changes one factor at a time; see [Kleijnen, 1987, pp. 266-267]. If purely quadratic terms  $\beta_{jj} x_j^2$  are added to (4), then the variable  $x_j$  cannot be observed at only two levels (-1 and +1).

Given the metamodel, there is more than one design to calibrate that model. A necessary condition for the design is that the resulting matrix of independent variables X is non-singular; see (7). Consider, for example, a first-order model: in (4) the double summation term vanishes. That model has q = 1 + k effects; so a necessary condition is that the number of combinations satisfies:  $n \ge k + 1$ . For k = 3 table 1 gives two designs that give a non-singular X; X was defined in (6).

Table 1: Two designs for a first order model with k = 3

Combination	One factor at a time design			2 <sup>3-1</sup> design		
	×1	*2	<b>*</b> 3	<b>x</b> <sub>1</sub>	×2	<b>x</b> 3
1	-1	-1	-1	-1	-1	+1
2	+1	-1	-1	+1	-1	-1
3	-1	+1	-1	-1	+1	-1
4	-1	-1	+1	+1	+1	+1

The  $2^{3-1}$  design is *balanced*: each column of X has an equal number of plus and minus signs, and each pair of columns has an equal number of the four combinations (-, -), (-, +), (+, -), (+, +). If the classical statistical model for the errors is assumed, then the covariance matrix of  $\hat{\beta}$  is given by (14). An orthogonal X minimizes the variances of  $\hat{\beta}_{j}$ ; see [Kleijnen, 1987, p.335].

If the metamodel includes interactions, then the number of effects increases considerably. To keep the number of combinations relatively small, the user may specify which interactions may be important; the remaining interactions are assumed to be negligible. Examples will be presented later.

The metamodel may be expanded with *purely quadratic* effects:  $\Sigma_{j=1}^{k} \beta_{jj} x_{j}^{2}$  is added to (4). These quadratic effects quantify the curvature of the response surface. Then more than two values per factor must be simulated (otherwise all columns for  $x_{j}^{2}$  are identical to the column for  $x_{0}$ ). A classical design is the *central composite* design: each factor is observed not only at -1 and +1 but also at the 'center point' 0 and at two other values, for example, -2 and +2; together five values. The  $2^{k-p}$  designs that is used to estimate main effects and interactions, is augmented with an observation at the center plus the following 2k combinations:

( 0,	0,	•••,	0)
(-2,	0,	••••	0)
(+2,	0,	,	0)
( 0,	-2,	••••	0)
( 0	+2,	••••	0)
	÷		
( 0,		••••	-2)
		,	

Next applications of metamodeling and experimental design will be presented. First the results for a relatively simple module of IMAGE are discussed; then results and technical details for a more complicated module are presented.

#### Dike raising in IMAGE

One module of IMAGE estimates the magnitude of the necessary dike raise and the resulting costs; see the lower part of figure 1. Eleven factors are examined (k=11); for example, the "unit dike raising cost", which is the cost of increasing the dike level by one meter. Finding a valid metamodel takes several iterations; altogether nine different models (and their concomitant designs) are tried. In an early iteration the metamodel helped to detect a serious error in the underlying simulation model: the original module needed to be split into two modules; the first submodule yields the dike raise necessary to keep the flooding probability under a fixed safety level; the second submodule takes that raise as input and yields the costs as output. So metamodeling may serve verification of the simulation model. Moreover, metamodeling may show in which area the simulation model is valid; for factor combinations outside that experimental area the simulation is not a correct model.

To obtain a valid metamodel for the dike raising costs module, the ranges of the original input variables must be decreased. This makes sense mathematically: a Taylor approximation is better in a smaller area. The final metamodel yields relative forecast errors smaller than 10%, which is acceptable for the IMAGE analysts. Nine of the eleven factors are significant, and so is one interaction. The most important factor is the "unit dike raising cost", as the analysts expected. The order of importance of the other factors was surprising, and gives more insight into the simulation model; for details see [Van Ham et al., 1990].

#### The carbon-dioxide cycle in IMAGE: ocean module

There are two modules for the  $\text{CO}_2$  cycle in IMAGE: one for the oceans and one for the terrestrial biosphere; also see the upper part of figure 1. This section covers the first module; the next section discusses the second module.

The oceans show three  $CO_2$  processes, described in [Goudriaan & Ketner, 1984]. For these processes ten factors are investigated; for example, thickness of ocean layers is factor 5. For each factor a range is specified by the analysts; for example, factor 5 varies between 3,000 and 4,000 meters; factor 2 (diffusion coefficient) ranges between 3,716 and 5,984 cm<sup>2</sup>/second. These variables are standardized, as described by (3). The analysts list eleven specific interactions that might be important; the remaining 34 interactions are neglected. To verify the design the reader should know that the following interactions may be important: 1 3, 1 5, 2 3, 2 5, 3 6, 3 7, 5 6, 5 7, 5 8, 6 8 and 7 8, where 1 3 stands for  $\beta_1$  3, and so on. So the metamodel is given by (4) with k = 10 and only eleven specific interactions  $\beta_{ib}$ .

A 'full factorial' design requires  $2^{10}$  combinations, which takes too much computer time. The number of effects in the metamodel is: q = 1 + 10 + 11 = 22. Hence a classical  $2^{k-p}$  design with enough combinations requires:  $n = 2^{10-p} \ge 22$  or  $p \le 5$  (least squares applied to the whole data set requires  $n \ge q$ , whereas cross validation requires n > q). There are many  $2^{10-5}$  designs. Accounting for the eleven specific interactions, the following design is selected; details are given in [Kleijnen, 1987, pp. 295-300]. Write down all  $2^{k-p} = 2^5$  combinations of the factors 1,2,4,9, and 10. Write down element i of the column for factor 3 as the product of the elements i in the columns for the factors 9 and 10; that is,  $x_{13} = x_{19}$  $x_{110}$  for  $i = 1, \ldots, n$  and  $n = 2^{k-p} = 2^5$  or in short-hand: 3 = 9 10. This is called a 'generator' of the design. The  $2^{10-5}$  design is fully specified by its p = 5 generators:

$$3 - 9 = 10 \quad 5 - 4 = 10 \quad 6 - 1 \quad 9 \quad / - 2 \quad 9 \quad 8 - 1 \quad 2 \quad 4,$$
 (18)

where  $\mathbf{8} = \mathbf{1} \ \mathbf{2} \ \mathbf{4}$  means  $\mathbf{x}_{\mathbf{18}} = \mathbf{x}_{\mathbf{11}} \ \mathbf{x}_{\mathbf{12}} \ \mathbf{x}_{\mathbf{14}}$ . The generator  $\mathbf{3} = \mathbf{9} \ \mathbf{10}$  means that the main effect of factor 3 is confounded or aliased with the interaction between the factors 9 and 10; that is,  $\hat{\beta}_3 = \hat{\beta}_{9,10}$  and  $\mathbf{E}(\hat{\beta}_3) = \beta_3 + \beta_{9,10}$ . If indeed the interaction  $\beta_{9,10}$  is negligible, then this confounding is acceptable. Analogously, the generator  $\mathbf{8} = \mathbf{1} \ \mathbf{2} \ \mathbf{4}$  means:  $\mathbf{E}(\hat{\beta}_8) = \beta_8 + \beta_{1,2,4}$  where  $\beta_{1,2,4}$  is a 'three factor' interaction, which was not yet

defined in this paper; such high-order interactions, however, are assumed negligible in metamodeling.

The combinations of the  $2^{10-5}$  design are simulated, and the outputs are compared with the predictions of the calibrated metamodel. This results in relative errors exceeding 10% in eight combinations, which is considered unacceptable. Shrinking the ranges of the original variables does not help. Next the metamodel is expanded with purely quadratic effects. The central composite design of (17) requires 1 + 2k extra combinations. To save computer time, only five of the ten factors are investigated, namely those five factors that are significant in the previous metamodel. (Because that metamodel is not valid, it is dangerous to use it for the selection of factors; the resulting new metamodel, however, will be validated again.)

The quadratic model is used for the factors 3,4,5,7 and 10. Only four (not ten) interactions between these five factors are conjectured to be important. So the number of effects excluding purely quadratic effects, is: 1 + 5 + 4 = 10. So the  $2^{k-p}$  design, which is part of the central composite design, must satisfy:  $n = 2^{5-p} \ge 10$  or  $p \le 1$ . So one generator is selected; namely 4 = 3 7 10. These sixteen combinations are augmented with eleven more combinations, following (17). This experiment yields a calibrated metamodel, which is cross-validated. Six more combinations are selected randomly, simulated, and compared with the predicted outcomes. Finally, a 'base' combination is examined; this combination is not the center combination  $(0,0,\ldots,0)$  of (17), but is close to it; it is a combination intuitively specified by the analysts. All validation results are acceptable: the errors are smaller than 10%. The individual effects of this accepted metamodel are discussed next.

Because a statistical model for the fitting errors e is assumed, (16) gives the Student t statistic for effects, where the degrees of freedom is n-q = (16+11) - (10+5) = 12. For  $\alpha$  = 0.05 the critical t value is  $t_{12}^{\alpha/2}$  = 2.18. Table 2 shows significant effects only. Note that if no such statistical model were assumed, then the last column should be ignored. If the design were orthogonal, then 'significance' and importance would coincide: (14) through (16) show that  $t_{n-q} = \hat{\beta}_j / (\hat{\sigma} / \sqrt{n})$ . So if effects are sorted in order of magnitude  $(|\beta_j|)$ , they are sorted in order of significance  $(|t_{n-q}|)$ . The central composite design, however, is not orthogonal: quadratic effects and the overall mean are not orthogonal.

Summarizing, originally ten factors are investigated for the ocean module. Because the metamodel without quadratic effects cannot be accepted, a model including such effects is specified. That model, however, is restricted to five factors. The latter model can be accepted, and yields only four important factors. These factors have significant main effects, one significant quadratic effect, and three significant interactions.

Effect	Estimate	t Statistic	
β <sub>0</sub>	1074.66	154.08	
B <sub>5</sub>	-244.95	-139.35	
B4	158.3/	98.10	
B <sub>3</sub>	51.77	32.07	
β <sub>55</sub>	21.79	14.05	
β <sub>1</sub>	18.78	11.63	
B <sub>45</sub>	- 17.65	- 8.93	
B <sub>47</sub>	- 12.26	- 6.20	
β <sub>35</sub>	- 8.35	- 4.22	

# Table 2: Significant effects of ocean module

# The carbon-dioxide cycle in IMAGE: terrestrial biosphere module

The terrestrial biosphere module is described in [Goudriaan & Ketner, 1984]. The analysis of this module is presented, because the mo-

dule contains many input variables: k = 62. These 62 variables are described in [Van Ham et al., 1990]. There are designs that yield estimators of main effects without being biased by possible interactions; moreover, these designs yield estimators of confounded interactions: so-called resolution-IV designs (see Kleijnen, 1987, p. 301). Such designs require at least 2k combinations (so  $k + 1 \ll n \ll 1 + k + k(k-1)/2$ ). For k = 62 a resolution IV  $2^{k-p}$  design satisfies:  $2^{62-p} \ge 124$  or p = 55. So 55 generators must be selected. Each selection yields a specific confounding pattern of estimated effects. The analysts give 26 interactions that might be important. Based on that list, 55 generators are selected; see [Van Ham et al., 1990]. First one estimated effect turns out to have the wrong sign: the effect is significantly positive whereas the analysts expect a negative effect. Next the ranges of the input variables are decreased, and the results become acceptable: the relative forecast errors are small; all (26 confounded) interactions are non-significant; all significant (unbiased) main effects have the correct sign. There are only 13 significant main effects (significance is measured by the t statistic with 128 - (1+62+26) = 39 degrees of freedom). Note that for validation purposes an experiment with twelve randomly selected extra combinations is executed; its relative forecast errors vary between -5.5% and -0.10%.

#### Conclusion

Any simulation model requires sensitivity analysis. That model can be treated as a black box, if the techniques of regression analysis and experimental design are applied. The regression model is a metamodel of the simulation model, and guides the experimental design. The design leads to efficient and effective experimentation.

The case study demonstrates that application of these statistical techniques requires knowledge of the underlying simulation model and real world system. For example, potentially important factors and their ranges must be given by the analysts. Some statistical expertise is needed to select the generators for the design. The case study was succesful. The metamodels give acceptable forecast errors. The significance of certain effects surprises the analysts. For example, quadratic effects in the ocean module were not expected; and the major importance of the 'biotic growth' factor in the terrestrial module is also surprising. Another surprise is the 'bug' in the dike raising module; the metamodel shows that this module must be split into two submodules.

The sensitivity analysis of IMAGE took quite some time and effort. Fortunately this investment in metamodeling is judged to be profitable. The conclusions of this analysis will guide the development of an interactive version of IMAGE.

Summarizing, regression metamodels and experimental designs are useful in the sensitivity analysis of simulation models, as the case study demonstrates. Details on the techniques can be found in [Kleijnen, 1987]; for sensitivity analysis of simulation models with hundreds of factors, special screening designs are presented in [Bettonvil, 1990].

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