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1989

document version Publisher's PDF, also known as Version of record

Link to publication in VU Research Portal

### citation for published version (APA)

Rietveld, P., & Janssen, R. (1989). Sénsitivity analysis in discrete multiple criteria decision problems: on the siting of nuclear power plants. (Research Memorandum; No. 1989-66). Faculty of Economics and Business Administration, Vrije Universiteit Amsterdam.

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# SERIE RESEARCH MEMORANDA

SENSITIVITY ANALYSIS IN DISCRETE MULTIPLE CRITERIA DECISION PROBLEMS: ON THE SITING OF **NUCLEAR POWER** PLANTS

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Researchmemorandum 1989-66

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october 1989

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VRIJE UNIVERSITEIT FACULTEIT DER ECONOMISCHE WETENSCHAPPEN EN ECONOMETRIE A M S T E R D A M

### ABSTRACT

Inclusion of evaluation methods in decision support systems gives way to extensive sensitivity analysis. In this article new methods for sensitivity analysis are developed and applied to the siting of **nuclear** power plants in the Netherlands.

### 1. INTRODUCTION

The use of decision support systems in public planning is slowly increasing following developments in private enterprise. This development results in a renewed interest in formal evaluation methods such as multicriteria analysis and cost benefit analysis. A major advantage of the integration of evaluation methods in a decision support system are the increased opportunities for sensitivity analysis. The availability of a wide range of procedures for sensitivity analysis allows the decision maker to investigate the limits of a decision problem (see figure 1).



Figure 1. Feed back loops in a decision procedure.

The **main** focus of this article is the use of procedures for sensitivity analysis on results from discrete evaluation problems. This type of evaluation **can** be performed by a wide range of evaluation methods **such** as **cost benefit** analysis and multicriteria methods.

In the first three steps of an evaluation procedure scores are assigned to **all** alternatives, weights to **all** criteria and a ranking of the **al**ternatives is **produced**. Especially in decisions that involve **negotiations** or public **debate** it is useful to know within which limits the derived rankings hold. This results in the following types of **questions**:

- to what extent **can** these scores of weights increase or decrease without changing this ranking (calculation of robustness **inter**-vals).
- how similar is the set of weights that produces the first rank re-versal.

We **will** describe procedures to deal with these questions followed by an application to the highly controversial decision on the location of two nuclear power plants in the Netherlands.

The procedures described in this article are included in our decision support system DEFINITE. This system is developed to support DEcisions based on a FINITE set of alternatives. This system contains a wide range of procedures to **assist all** steps in the evaluation procedure. This **makes** it possible to feed back results of sensivitity analysis directly to problem definition and evaluation. (See Herwijnen and Janssen, 1988, for a description of DEFINITE, and Rietveld, 1988, for a complete description of the procedures for sensitivity analysis in DEFINITE.)

Sections 2 and 3 of this article are devoted to a formal introduction of discrete multicriteria methods and procedures for sensitivity **analy**sis respectively. In **section** 4 the multicriteria methods are applied to rank locations for nuclear plants in the Netherlands. The derived **rank**ings are analyzed using various procedures for sensitivity analysis. Finally, the usefulness of this type of approach is **discussed** in **section** 5.

### 2. DISCRETE MULTIPLE CRITERIA METHODS

An important aim of discrete multiple criteria analysis is to **provide** a rational basis for ranking a number of alternatives on the basis of multiple criteria. There are **many** different discrete multiple criteria methods currently in use (see, e.g. Nijkamp, 1979; Rietveld, 1980, Voogd, 1983).

A major step in these methods is the construction of an impact (or evaluation) matrix representing the effect of a certain alternative on a decision criterion. In order to aggregate the information of the evaluation matrix usually a weighting scheme is necessary which expresses the relative importance of the various criteria. The impact matrix will be denoted by the symbol P. This matrix has elements pij which represent the impact of alternative i (i = 1, . . . , I) on the value of criterion j (j = 1, . . . J). The vector of weights is denoted as  $\underline{\lambda} = (\lambda_1, \ldots, \lambda_J)$ . It is often assumed that the criteria have been defined in such a way that all weights are positive. In addition, one may impose the restriction that the weights add up to unity. Thus, the set S of feasible weights can be defined as:

$$S = \{ \underline{\lambda} \mid 0 \le \lambda_{i} \le 1 \text{ for all } j = 1, \dots, J, \text{ and } \underline{\Sigma} \lambda_{i} = 1 \}$$
(1)

In many applications, part of the information on P and  $\underline{\lambda}$  is soft. For example, for some criteria, no **precise** quantitative values of impacts may be available. At best one may have a ranking of alternatives in such a case. Similarly, one may only have a ranking of criteria to indicate their relative importance. Therefore the development of multicriteria methods which can deal with these types of problems is important. A survey of such methods is contained in Nijkamp, Rietveld and Voogd (1989). In this paper we will only shortly discuss those approaches which will be used in the empirical application.

A relatively easy way of dealing with ordinal data is by interpreting them as unknown quantitative data which satisfy certain inequalities. For example, if J <u>criteria</u> are ranked in increasing order, one arrives at J unknown cardinal weights satisfying:

$$\begin{bmatrix} 0 \le \lambda_1 \le \lambda_2 \le \dots \le \lambda_J \\ \sum_{j} \lambda_j = 1 \end{bmatrix}$$
(2)

Every cardinal  $\underline{\lambda}$  satisfying (2) is consistent with the original ranking. When one assumes that all vectors  $\underline{\lambda}$  satisfying (2) are equally probable, i.e.  $\underline{\lambda}$  is uniformly distributed on (2), one can derive expected values of the weights in a relatively straightforward way. As shown in Rietveld (1984), the expected values of weights are:

$$E(\lambda_{1}) = 1/J^{2}$$

$$E(\lambda_{2}) = 1/J^{2} + 1/[J.(J-1)]$$

$$E(\lambda_{3}) = 1/J^{2} + 1/[J.(J-1)] + 1/[J.(J-2)]$$
etc.
(3)

Along similar lines one **may** deal with ordinal data on criterion scores **p**<sub>ij</sub>. A ranking of alternatives in increasing order of **attractive**ness according to a certain criterion j, **combined** with the assumption of a uniform distribution leads to the following expected values:

$$E(p_{ij}) = i/I$$
 for  $i = 1, \dots, 1$  and a certain j (4)

where the highest outcome for pij has been set equal to 1 by way of standardization<sup>1</sup>).

This stochastic approach **provides** a basis for translating ordinal data into cardinal **ones.** Of course its **relevance** depends on the **appropriate-**

ness of the assumption of uniformly distributed variables. By using this cardinalization step, one **can** employ standard multicriteria methods for cardinal data even if (part of) P **or**  $\underline{\lambda}$  are cardinal. An example of a standard multicriteria method is <u>weighted summation</u> which is based on an additive utility **structure**.

Another **class** of multicriteria methods has been specifically designed to deal with evaluation problems **where** qualitative data are used. Examples of **such** methods are **EVAMIX** (Voogd, 19831, QUALIFLEX (Paelinck, 1976), and the regime method (Hinloopen, Nijkamp and Rietveld, 1983).

A multicriteria method frequently used is ELECTRE (also called concordance analysis, cf. Roy, 1974, Crama and Hansen, 1983). This method is based on a pairwise comparison of alternatives, thus using only the metric interval characteristics of the various scores in the evaluation of the impact matrix. The **basic** idea is to measure the degree to which the scores and their associated weights conform or contradict the dominant pairwise relationships among alternatives. The differences in weights and the differences in evaluation scores are usually analyzed separately. The central concept in ELECTRE is the so-called concordance index c;;. This index represents the extent to which alternative i is better than alternative i'. This index may be defined as the sum of weights attached to the criteria included in the so-called concordance set  $C_{ii'}$ ; this is the set of all evaluation criteria for which alternative i in the impact matrtix P is at least equally attractive as alternative i'. Clearly, 'this set can be determined irrespective of the degree of information on the impact matrix. Hence, the concordance index can be defined as follows:

 $c_{ii'} = j \xi_{c_{ii}}, \lambda_j$  (5)

A dominating alternative **can** now be found by employing threshold values, relative **dominance** indicators, **or** other **concepts** from graph theory.

In an analogous way, one **may define** a discordance index. This index reflects the extent to which alternative i is worse than i'. Instead of using weights in this index, the corresponding relative pairwise **dif**ferences from the impact matrix are then taken into consideration. By combining the results from the **concordance** and discordance approach, **final** inferences on the ranking of alternatives **may** be made (see e.g. Nijkamp, 1979). Most of the methods presented in this section will be applied in section 4.

### 3. SENSITIVITY ANALYSIS

### 3.1. Introduction

Results of multicriteria analysis **depend** on various **factors such** as the choice for a particular multicriteria method, the choice of weights, methods for standardizing criteria or methods for dealing with **uncer**-tainty on **effects** of alternatives. In this **section**, we **will pay atten**-tion to methods for investigating the sensitivity of evaluation results for the choice of criterion weights.

One way of dealing with the problem of uncertainty on weights is the Monte **Carlo** approach. In this case a **random** generator is used to **pro**duce a large number of weight **vectors centered** around a given weight vector. For **each** vector a multicriteria evaluation is carried **out**, **after** which the **result** is compared with the **result** for the given weight vector (see e.g. Nijkamp, 1979).

Another way is to formulate a limited number of policy views, **each** of which is represented by a certain weight vector. Then, in a **second** step, multicriteria analysis is carried **out** to find **out** whether the views lead to different options (see e.g. Voogd, 1983).

In the present paper another approach is proposed which gives a more detailed account of the sensitivity of, results of multicriteria **analy**-sis for the coice of weights. Our point of departure is a given weight vector  $\underline{\lambda}^{0}$ . For a certain multicriteria method, this **leads** to a ranking of alternatives, for example  $A_1 \ge A_2 \ge A_3 \ge \ldots$  where Ai  $\ge$  Ak means that alternative Ai performs equal to or better than Ak. Depending on the multicriteria method **chosen**, such a ranking is not necessarily complete, i.e., certain pairs of alternatives may be incomparable.

The method we will discuss is on the sensitivity of the result for an arbitrary pair of alternatives (e.g.,  $A_1 \ge A_2$ ) for changes in  $\underline{\lambda}^{\circ}$ . The question addressed is: "how far must  $\underline{\lambda}$  be from  $\underline{\lambda}^{\circ}$  before  $A_1 \ge A_2$  does no longer hold true"

This question **can** be approached in various ways. One way is to focus on the weight for one particular criterion and assume that the ratios **be**tween other weights remain unaltered. Another approach would be that the weights of **all** criteria are allowed to change freely (the only **condition** being that they add up to 1). Both methods **will** be **discussed** below.

### 3.2. The nearest tuming point: focus on one criterion

For the ease of presentation we start with te assumption that the **mul-**ticriteria method yields complete rankings. Let Sik be the set **of** 

weights for which Ai  $\geq$  Ak. Similarly, let Tik be the set of weights according to which the alternatives Ai and Ak perform equally well (T<sub>ik</sub> = S<sub>ik</sub>  $\bigcap$  S<sub>ki</sub>).

Suppose that  $\underline{\lambda}^{0}$  is an element of  $S_{ik}$ . We want to know how much a particular criterion weight (for example:  $\lambda_{1}$ ) has to change in order to make the weights vector an element of  $T_{ik}$ . Since we impose the restriction that the sum of the weights equals one, a change in  $\lambda_{1}$  implies that other weights will also change. We will assume that their ratios remain unchanged:

$$\begin{bmatrix} \lambda_j / \lambda_2 = \lambda_j^{\circ} / \lambda_2^{\circ} & \text{for } j = 3, \dots, J \end{bmatrix}$$
(6)

In Figure 2 an example is given for the case that J = 3. The turning point  $\underline{\lambda}^{\mathbf{A}}$  is found by extrapolating  $\underline{\lambda}^{\mathbf{0}}$ , using (1, 0, 0) as a reference point. In this case, one **finds** one turning point. It is not difficult to see that other examples could be given **where** there is no turning point at all. **Also**, the occurence of multiple turning points cannot be excluded. In that case the nearest turning point is the most relevant one.



Figure 2. Turning point in the rank order of two alternatives.

How can  $\underline{\lambda}^{\underline{A}}$  be found (if it exists)? If a utility based multicriteria method is used, the set Tik is defined by  $U(\underline{\lambda}, \underline{p_i}) = U(\underline{\lambda}, \underline{p_k})$ . In that case it is not difficult to determine the element of Tik which satisfies condition (6). However, if a method for multicriteria analysis is used which is not based on a utility concept, as is the case for example with concordance analysis, no such definition can be given of the set  $T_{\underline{ik}}$ . In that case, the turning point  $\underline{\lambda}^{\underline{A}}$  has to be found by a systematic inspection of the set of weights satisfying:

$$\begin{bmatrix} 0 \leq \lambda_{1} < 1 \\ \lambda_{j} = \lambda_{j}^{\circ} (1 - \lambda_{1}) / (j \neq 2 \lambda_{j}^{\circ}) & j = 2, \dots J \end{bmatrix}$$
(7)

It is not difficult to check that weights satisfying (7) are non-negative and add up to 1.

In order to find the turning point  $\underline{\lambda}^{\mathbf{A}}$ , we **propose** the method of **halv**ing. In terms of Figure 2, we first investigate whether a point exists on the line between D and E for which the original ranking Ai  $\geq$  Ak does not hold. If **such** a point appears to exist, an additional point is investigated which is in the middle of two points which are on different **sides** of the unknown line Tik. **After** a **sufficient** number of halvings one obtains a point which is **very** near to the turning point  $\underline{\lambda}^{\mathbf{A}}$ .

The first three steps of the following.algorithm are **carried out** to investigate whether a turning point exists. In addition, these steps aim at **solving** the problem of multiple turning points. If more than one turning point exists, it is the one nearest to  $\underline{\lambda}^{\circ}$  which has to be found. In septs 4 to 7 halving iterations are carried out.

In the algorithm, t stands for iteration. Further,  $a(\underline{\lambda})$  is used to **indicate** whether  $\underline{\lambda}$  is an element of the set  $S_{ik}$ :

 $\begin{bmatrix} \mathbf{a}(\underline{\lambda}) &= 1 & \text{if Ai} \geq Ak \\ a(\underline{\lambda}) &= 0 & \text{in all other cases.} \end{bmatrix}$ 

The algorithm consists of the following steps:

1. t = 0 A,(0) =  $\lambda_1^{\circ}$ Compute  $a(\underline{\lambda}(0))$  by means of multicriteria method, where  $\underline{\lambda}(0) = \underline{\lambda}^{\circ}$ .

2. t = t + 1If t = 12, stop: no turning point found.  $\lambda_1(t) = c_1(t)$ , where  $c_1(t)$  is defined below. Compute  $\underline{\lambda}(t)$ , by means of (7). Compute  $a(\underline{\lambda}(t))$  by means of multicriteria method. (For t = 1, . . . 13 c,(t) assumes the following values<sup>2</sup>: TEL λî .6λ<sup>9</sup><sub>1</sub>.4λ<sup>9</sup><sub>1</sub>.2λ<sup>9</sup><sub>1</sub> .8λº .0) 7.0 ু**∓**ে -~ 1 7:10 3. If  $a(\underline{\lambda}(t)) - a(\underline{\lambda}(t-1)) = 0$ , return to 2. If  $a(\underline{\lambda}(t)) - a(\underline{\lambda}(t-1)) \neq 0$ , go to 4. 4.  $y = \underline{\lambda}(t-1)$ )  $a(y) = a(\underline{\lambda}(t-1))$  $\underline{z} = \underline{\lambda}(t)$  $a(\underline{z}) = a(\lambda(t))$ 5.  $\underline{\mathbf{v}} = (\underline{\mathbf{y}} + \underline{\mathbf{z}})/2$ Compute  $\mathbf{a}(\underline{\mathbf{v}})$  by **means** of multicriteria method. 6. If  $\mathbf{a}(\mathbf{y}) - \mathbf{a}(\mathbf{v}) = 0$ , then  $\mathbf{y} = \mathbf{v}$  and  $\mathbf{a}(\mathbf{y}) = \mathbf{a}(\mathbf{v})$ , go to 7. If  $\mathbf{a}(\underline{z}) - \mathbf{a}(\underline{v}) = 0$ , then  $\underline{z} = \underline{v}$  and  $\mathbf{a}(\underline{z}) = \mathbf{a}(\underline{v})$ , go to 7.

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If |y<sub>j</sub> - z<sub>j</sub>| ≤ ε z<sub>j</sub> for all j where ε is a certain small value, stop: turning point found; otherwise: return to 5.

The algorithm **needs** some straightforward extentions when multicriteria methods are used which may yield incomplete rankings. As illustrated in Figure 4, the indifference line  $T_{ik}$  is replaced by a band of weights leading to the result that two alternatives are incomparable. In this case two turning points are found ( $\underline{\lambda}^B$  and  $\underline{\lambda}^C$ ). In a strict sense, the algorithm does not guarantee that a turning point is found if there exists one, even if  $T_{ik}$  is continuous. As shown in Figure 3, a turning point may be overlooked when the points investigated in Step 2 are too far removed from each other. The probability that this occurs can be made very small by increasing the number of iterations in Step 2.



Figure 3. Possibility of overlooking a turning point.

The method of halving provides a decision making unit (DMU) with an indication of the degree of sensitivity of a certain outcome Ai > Ak for changes in the value of a certain weight. For a DMU, the degree of uncertainty about the original weight vector  $\underline{\lambda}^{\circ}$  is usually considerable. By comparing  $\underline{\lambda}^{\circ}$  with  $\underline{\lambda}^{A}$ , the DMU is informed on how sensitive the results of multicriteria analysis are for changes in the weights vector. The method of halving can be used for any multicriteria method.



Figure 4. Turning points with incomparable alternatives.

## 3.3. The nearest turning point: all weights are allowed to vary

A limitation in the method of **halving** is that **all** weights **except** one are assumed to be proportional to the original weights. Thus, one **ar**rives at a point  $\underline{\lambda}^{\mathbf{A}}$ , which **may** be far removed from the point  $\underline{\lambda}^{\mathbf{Z}}$ which is the point on  $\mathbf{T}_{\mathbf{ik}}$  nearest to  $\underline{\lambda}^{\mathbf{0}}$  when **all** weights are allowed to move freely (see Figure 5).

If a linear utility function would be used,  $\underline{\lambda}^{z}$  can be found as the solution of a quadratic programming problem. Since it is our aim to develop a method which is applicable to a **much** more general class of multicriteria methods, a different approach has to be followed.



Figure 5. Search procedure for turning point **when all** weights are **al**-lowed to vary.

The **structure** of the algorithm reads as follows. Let  $\underline{\lambda}^{0}$  be an element of  $s_{ik}$ . Then the following steps have to be made in order to find the turning point  $\underline{\lambda}^{\mathbf{Z}}$  nearest to  $\underline{\lambda}^{0}$ .

- 1. Find a weight vector  $\underline{v}$  which is an element of  $\mathbf{S_{ki}}$  and which consists of nonnegative elements **adding** up to 1, If no **such** vector **can** be found, a turning point does not exist.
- 2. Use the method of halving based on the vectors  $\underline{\lambda}^{0}$  and  $\underline{v}$  to determine the vector  $\underline{w}$  which is an element of  $T_{ki}$ . Let  $d(\underline{\lambda}^{0}, \underline{w})$  be the distance between the vectors  $\underline{\lambda}^{0}$  and  $\underline{w}$ .
- 3. Find in a neighbourhood of  $\underline{w}$  a new point  $\underline{v}$  with  $d(\underline{\lambda}^{\circ}, \underline{v}) \leq d(\underline{\lambda}^{\circ}, \underline{w})$ and where  $\underline{v}$  is an element of  $s_{ki'}$

4. Return to step 2, but stop when subsequent results of  $\underline{v}$  come very near to each other.

Figure 3 gives an illustration of the algorithm for the first two iterations. We will now discuss the four steps in more detail.

Step 1. For each of the extreme points in the weights set (1, 0,...,0),(0,1,0,...0),... (0,...,0, 1) an investigation is made whether it is an element of  $S_{ki}$ . For those extreme points  $\underline{v}$  which are indeed an element of  $S_{ki}$ , the distance  $d(\underline{v}, \underline{\lambda}^{\circ})$  to  $\underline{\lambda}^{\circ}$  is measured, where  $d(\underline{v}, \underline{\lambda}^{\circ})$  is defined as:

 $d = \left[\sum_{j} (v_{j} - \lambda_{j}^{\circ})^{2}\right]^{5}$ 

The extreme point  $\underline{v}$  with minimum distance is  ${\tt selected}$  to be used in the  ${\tt second}$  step.

A difficulty is that there may be cases where feasible weight vectors in  $S_{ki}$  exist, but where there are no extreme points in  $S_{ki}$ (see Figure 6). Several approaches can be followed to counter this problem. First, one may examine in a systematic way points on the faces of the polyhedral set S. For example, when J=4, one can examine the points: (0, 1/3, 1/3, 1/3), (1/3, 0, 1/3, 1/3), (1/3, 1/3, 0, 1/3), (1/3, 1/3, 1/3, 0), (0, 0, 1/2, 1/2), (0, 1/2, 0, 1/2), (0, 1/2, 1/2, 0), (1/2, 0, 0, 1/2), (1/2, 0, 1/2, 0), (1/2, 1/2, 0, 0).

Another approach to generate in a random way a set of points in S after which for each point it is examined whether it is in  $S_{ki}$ . Appendix 1 contains a method to generate random weights which are uniform in S.

step 2 The method of halving as presented in section 3.2 can be used directly here.

step 3. A neighbourhood of  $\underline{w}$  is defined here as a polyhedral set around  $\underline{w}$  which is **contained** in the set S. Element  $\underline{x}$  of **such** a set **can** be generated by using the formula:

 $x = b\underline{w} + (1-b)\underline{\delta}$ ,

where  $\underline{\delta}$  is an arbitrary element of the set S. An example of such.a neighbourhood is given in Figure 7 (for J=3 and b=.5). The method presented in Appendix 1 can be used again for generating weights which are uniformly distributed in S. The generation of point  $\underline{x}$  continues until a point is found which is nearer to  $\underline{\lambda}^{0}$  than is  $\underline{w}$ , and which is an element of  $S_{ki}$ . If no such point  $\underline{x}$  is found,  $\underline{w}$  is the optimal solution.



Figure 6. Existence of turning point: a special case.

Step 4. The algorithm stops when for some subsequent iterations t:

## $|w_{j}(t) - w_{j}(t-1)| < \varepsilon w_{j}(t)$ ,

where  $\epsilon$  is a suitably choosen small number. In order to deal with the problem of local optima, one may return the algorithm with another starting point in Step 1.



Figure 7. A neighbourhood for  $\underline{w}$ .

The parameters to be fixed in the algorithm are b and  $\epsilon$ . When fixing b a compromise must be found between the probability of finding a nearer point, and the expected size of decrease of distance. Further, a maximum level must be set for the number of random vectors  $\underline{x}$  in Step 3 (and Step 1).

In a **strict** sense the algorithm does not guarantee that the turning point with minimum distance is found. One problem is related to fixing the number of points generated in Step 3 (or Step 1). If the maximum number of interations in Step 3 is small, the algorithm **may** stop at a point which is far removed from the nearest turning point. The **probabi**lity that this occurs **can** be made arbitrary small by increasing the maximum number of iterations. Another problem is that the optimum **solution** found is not the global one. This probability **can** be made **arbitra**ry **small** by restarting the algorithm for a sufficiently large number of starting points. Finally, the algorithm **needs** some straightforward **ex**tensions **when** multicriteria methods are used which **may** yield incomplete rankings (cf. **section** 3.2).

# 4. AN APPLICATION: THE SELECTION OF THE OPTIMAL LOCATION FOR NUCLEAR PLANTS IN THE NETHERLANDS

### 4.1. Introduction

The share of nuclear power in total power production is **very** small at present in the Netherlands. In 1985 the Dutch government expressed the intention to build two new nuclear power plants with a capacity of 1000 **MWe each.** Given this decision an important decision is **where** to **locate** these two plants. **After** some initial **scoping**, nine potential locations for the plants were **selected** (Tweede Kamer 188303, 43-44). These locations are shown in figure 8.

In this section these nine locations **will** be ranked using 15 appraisal criteria. Our support system for decisions on a finite set of **alter**-natives (**DEFINITE**) is used to **produce** this ranking. This **will** be done by using the following steps:

- 1. Problem definition.
- 2. Problem presentation.
- 3. Problem evaluation.
- 4. Sensitivity analysis.

Steps 1 to 3 are described only briefly. This section **concentrates** on the methods for sensitivity analysis **discussed** in the previous **sec-tion.** 



Figure 8. Potential sites for **nuclear** plants.

0 potential **site** 

1 Bath/Hoedekenskerke 6 Maasvlakte

2 Borssele

4 Flevo Noord 5 Ketelmeer

- 3 Eems
- 7 Moerdijk
- 8 West. NOP-dijk 9 Wieringermeer

0 20 km range around the potential **site** 

### 4.2. Problem definition

The impact matrix of this evaluation problem is shown in Table 1, where nine potential locations are **scored** according to 15 criteria. Only the score for population around a **site**, is measured on a cardinal scale. **All** other scores are measures on an ordinal scale: a score 1 is **as**signed to the best alternative, 2 to the **second** best, etc. (see Appendix 11 for a definition of the criteria).

Table 1 Impact matrix (Source: Tweede Kamer 18830, 43-44; **advice** to the government)

	Bath	Bors- <b>sele</b>	Eems	Flevo	Ketel	Maas vlak	Moer- dijk	NOPolder	<b>Wie-</b> ring
Population	51	49	16	27	30	43	100	19	21
Evacuation	1	2	1	1	1	2	1	1	1
Agricult at risk	2	2	2	2	2	1	3	2	2
Industry at risk	1	4	3	2	1	5	3	1	1
Fr water at risk	1	1	1	2	2	1	2	2	2
Cool-water quant	2	1	1	1	1	1	3	1	1
Cool-water <b>qual</b>	2	1	2	3	3	1	2	3	3
Air pollution	2	2	2	2	2	1	2	2	2
Thermal poll.	3	2	2	2	3	1	2	3	3
Indirect landuse	2	3	3	2	1	2	4	1	1
Landscape	3	1	1	1	3	1	2	3	3
Nat environment	3	1	3	1	2	1	1	1	1
National grid	2	2	3	1	1	2	2	2	3
Infrastructure	2	1	1	2	2	1	1	2	2
Coal-location	3	6	4	3	2	7	5	1	1

#### 4.3. Problem presentation

A graphical presentation of the impact matrix is shown in Figure 9. This figure is derived by standardising **all** criterion scores between 0 and 1 (see **section** 2). The highest bar in **each** row represents a score of 1 corresponding to the best alternative for that row. As a next step the criteria are ordered from most important (top> to least important (bottom). These priorities have been expressed in an ordinal way by experts of the government advisory board on physical planning (Tweede Kamer 18830, **43-44**). Using a combination of the expected value method for weights and the weighted summation method (see **section** 2) the **in**formation on priorities and scores **can** be used to sort the alternatives from best (**left**) to worst (right). It is **clear** from this figure that the impact matrix contains **many** tied scores and that differences **be**tween alternative locations are fairly small.





## 4.4. Problem evaluation

A variety of evaluation methods **can** be used to rank these alternatives (see for a short description **section 2**). In this application the **ex**pected value method is used to transform the priority ranking of the criteria to quantitative weights (Table 2). Using these weights both the weighted summation and the Electre method were applied to **generate** a ranking of the alternatives (Table 2). Both the weighted summation method and the Electre 2 method **result** in an **almost** complete ranking of the alternatives. In the results of the weighted summation method, **al**ternatives Maasvlakte and Borsele share the 7th and 8th position and the Electre method results in a tie for alternatives Flevo and **Ketel**meer. Note that the methods **generate** different rankings.

Weights Ranking Sco		Score	Weighted summation Ranking Score			<b>ELECTRE</b> 2 Ranking		
1: 2: 3: 5:	Population Industry at risk Agricult at risk Fr. water at risk Cool-water quant. Cool-water qual.	0.221 0.155 0.110 0.110 0.064 0.064	1: 2: 3: 4: 5: 6:	Eems NOPolder Flevo <b>Wiering</b> Ketel Bath	0.86 0.83 0.82 0.81 0.79 0.76	1: 2: 3: 5: 6:	NOPolder Eems Flevo Ketel Maasvlak Wiering	
9:	Thermal poll. Coal-location Air pollution Landscape Nat. environment	0.064 0.064 0.027 0.027 0.027	7: 9:	<b>Maasvlak</b> Borssele Moerdijk	0.75 0.75 0.52	7: 8: 9:	Borssele Bath Moerdijk	
14:	National grid Infrastructure Evacuation Indirect landuse	0.027 0.027 7.0E-03 7.0E-03						

Table 2. Ranking of the alternatives according to the weighted **summa-tion** and the ELECTRE-2 method.

### 4.5. Sensitivity analysis

The location of **nuclear** plants is a politically sensitive decision. It is therefore interesting to analyze the relationship between assigned priorities and the ranking of the locations. The general robustness of the derived ranking **can** be analyzed as a first step. In this case a Monte Carlo analysis shows that if the weights were allowed to vary by  $\pm$  5% and assuming the weights are normally distributed, the overall ranking proves uncertain (see also Rietveld 1988). However, the selection of the two best alternatives, proves sufficiently certain. Since the government aims to select two locations this is a useful result. More interesting than the overall stability of the derived ranking is, in cases like this, the sensitivity of the ranking to specific weights and the stability of the ranking of specific alternatives to changes in weights.

Since the governement **wishes** to select two sites we **will** try to analyze **how** firmly the alternatives Eems and Noord-Oost Polder hold the first two positions. Firstly we **will** use the methods as described in **section** 3.2 to **calculate** robustness intervals for these two alternatives and secondly we **will** use the methods as described in **section** 3.3 to find the nearest weight combination that brings one of the other **alter**-natives to the first or **second** position.

Robustness intervals are calculated for the weight of the most important criterion: population living **around** the site.

As shown in Table 2 weighted summation **selects** Eems as the best alternative followed by NO Polder. Alternative Flevo ranks on the third **place.** In Table 3 robustness intervals are calculated for the pairs Eems - Flevo and NO Polder - Flevo. Table 3 shows that Eems ranks **higher** than Flevo for **any** weight assigned to the population criterion. Table 3 **also** shows that if the weight assigned to population is reduced below 0.16 the ranking of Flevo and NO Polder is reversed.

Table 3 Robustness intervals; weighted summation method.

Ranking	Weight of criterion population
Eems ≥ Flevo	0 - > 1
Flevo ≥ Eems	empty
NO Polder ≥ Flevo	0.16 - > 1
Flevo ≥ NO Polder	0 - > 0.16

The ELECTRE-2 method **also** ranks Eems and NO Polder as the best two **al**ternatives, but in the reversed order. Table 4 shows the required **changes** in weights to move Eems **or** NO Polder from their first position. It is shown that with **any** weight assigned to population, alternative Flevo **will** not replace alternative Eems **or** NO Polder from their first two positions. The same **can** be shown for alternative Ketelmeer. It is interesting to note that alternative Maasvlakte, which is ranked at a fifth position in the initial ranking, ranks **higher** than NO Polder if the weight assigned to population is **lower** than 0.13, and **higher** than Eems if this weight is less than 0.12. A similar procedure **can** be applied to establish robustness intervals for criterion scores.

Table 4. robustness intervals; ELECTRE-2 method.

Ranking	Weight of criterion population
Eems ≥ Flevo	0 -> 1
Flevo ≥ Eems	empty
NO Polder ≥ Flevo	0 -> 1
Flevo ≥ NO Polder	empty
Eems > Maasvlak Maasvlak > Eems	$\begin{array}{cccccccccccccccccccccccccccccccccccc$
NO Polder ≥ Maasvlak	0.13 - > 1
Maasvlak ≥ NO Polder	0 - > 0.13

Only the relative weight of the population criterion to **all** other criteria was **changed** in calculating the weight intervals. If we allow **all** weights to change it becomes clear **how** sensitive the derived ranking is to overall changes in weights. Using the method described in **section** 3.3 for **each** pair of alternatives, the set of weights with the **smallest** Euclidean distance from the original weights that **reverses** the ranking of the alternatives **can** be calculated. Table 5 shows the results for **six** pairs of alternatives. The small values for distances **indicate how** sensitive the ranking is to changes in weights.

Table	5.	Weight	combinations	with	rank	reversal;	ELECTIE-2	method.
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Criterion	Original	Flevo > Eems	Flevo > NOPolder	Ketel > Eems	Ketel > NOPolder	Maas > Eems
Population	0.21	0.150	0.158	0.157	0.150	0.160
Evacuation	0.007	0.047	0.020	0.020	0.047	0.019
Agricult at risk	0.110	0.075	0.103	0.081	0.075	0.092
Industry at risk	0.155	0.105	0.106	0.230	0.105	0.119
Fr water at risk	0.110	0.075	0.078	0.088	0.075	0.088
Cool-water quant	0.064	0.043	0.070	0.057	0.043	0.061
Cool-water <b>qual</b>	0.064	0.043	0.058	0.051	0.043	0.060
Air pollution	0.027	0.018	0.024	0.040	0.018	0.039
Thermal poll.	0.064	0.043	0.129	0.061	0.043	0.136
Indirect landuse	0.007	0.047	0.032	0.037	0.047	0.020
Landscape	0.027	0.018	0.043	0.025	0.018	0.030
Nat environment	0.027	0.018	0.043	0.024	0.018	0.041
National grid	0.027	0.341	0.052	0.022	0.341	0.041
Infrastructure	0.027	0.018	0.040	0.031	0.018	0.043
Coal-location	0.064	0.043	0.047	0.076	0.043	0.054
Distance to original	0	0.043	0.119	0.113	0.332	0.111

It is **already** clear from Table 2 that different evaluation methods might **result** in different rankings of the alternatives. The results of both multicriteria methods are shown in Figure 10. It is clear that both methods rank Eems and NO Polder above **all** other alternatives. The ranking of the alternatives is similar for both methods. From Figure 10 it **can** be safely concluded that Eems and NO Polder should be **selected** as the best two locations.

As a last step DEFINITE offers conclusions on the usefulness of the derived rankings based on the results of sensitivity analysis. The **con**-clusions **read** as follows:

- The overall ranking is insufficiently certain.
- Eems and NO Polder are the best two alternatives.

This **result** should prevent the decision maker from publishing **the** complete ranking. Although our aim to derive a complete ranking has failed the **result** obtained is useful. Two locations had to be **selected:** the **difference** in ranking between the first two alternatives is in this special case irrelevant.



Figure 10. Ranking of five alternatives according to two multicriteria methods.

### 5. CONCLUDING REMARKS

The availability of various evaluation methods and methods for sensitivity analysis in a decision support system **such** as DEFINITE has clear advantages. It increases availability of these methods to various types of users; although the exact calculation procedures **will** not be clear to **all** users, the results are easy to interpret and unambiguous. In addition, it allows users to study the sensitivity of outcomes for the choice of a particular method of multicriteria analysis. In the case study presented in this paper, it appears that the ranking **produced** with the weighted summation method is **indeed** different from the ranking **produced** with ELECTRE. For the choice of the two highest **rank**ing alternatives, the methods appear to yield identical outcomes, **how**ever. The methods for sensitivity analysis available in DEFINITE allows one to carry **out** detailed studies for the sensitivity of evaluation results. For the choice of **nuclear** power plants, it appears that the choice of location is **rather** sensitive to the value of the weights. Sensitivity analysis of this type are important because the information on weights is **rather** soft in **many real** world applications. This is **also** the reason in this paper special attention is given to methods for dealing with ordinal information.

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### NOTES

- 1) The **difference** between (3) and (4) is **caused** by the different ways of standardizing weights and criterion scores.
- 2) By decreasing the step size in Step 2 one can increase the probability that in the case of multiple turning points it is the nearest one which is found.
- 3) Documents presented to the lower house of parliament for discussion.



### APPENDIX I

### Generating random weights

Consider a weight vector  $\underline{\lambda}$  which is uniformly distributed in S as defined in section 2:

 $s = \{\underline{\lambda} \mid 0 \le \lambda_j \le 1 \text{ for all } j=1,..., J, and \sum_{j} \lambda_j = 1\}$ 

One might be tempted to generate weight vectors  $\underline{\lambda}$  by drawing J random numbers  $x_1, \dots x_J$  from a uniform distribution on the interval  $0 \le x \le 1$ , and setting  $\lambda = x / \sum_{j=1}^{J} x_j$ .

However, this does not lead to weight vectors which are uniformly distributed on S. Therefore, another approach has to be followed.

The joint density of  $\underline{\lambda}$  is:

 $g(\lambda_1, \dots, \lambda_{J-1}) = (J-1)! \text{ for } 0 \le \lambda_1 \le 1$  $0 \le \lambda_2 \le 1 - \lambda_1$  $= 0 \qquad \qquad 0 \le \lambda_{J-1} \le 1 - \lambda_1 - \dots + \lambda_{J-2}$ elsewhere

On the basis of this joint density function one can derive for the density of  $\lambda_1 \colon$ 

$$g(\lambda_1) = (J-1)(1-\lambda_1)^{J-2} \text{ for } 0 \le \lambda_1 \le 1$$
  
= 0 elsewhere

=0

Further, the conditional density functions  ${\tt can}$  be shown to  ${\tt read}$  as follows for  $j{=}2,\ldots,J{-}1.$ 

$$g(\lambda_{j}|\lambda_{1},...,\lambda_{j-1})$$

$$= (J-j)(1-\lambda_{1}-...-\lambda_{j})^{J-j-1}(1-\lambda_{1}-...\lambda_{j-1})^{j-J}$$
for  $0 \le \lambda_{j} \le 1-\lambda_{1}-...-\lambda_{j-1}$ 

elsewhere

Then, a **random** weight vector **can** be generated by drawing a value for  $\lambda_1$  on the basis of  $g(\lambda_1)$ , followed by drawing a value for  $\lambda_2$  on the basis of  $g(\lambda_2 | \lambda_1)$ , etc. Finally,  $\lambda_J can$  be computed as  $1 - \lambda_1 - \ldots - \lambda_{J-1}$ .

The conditional distributions mentioned above are not included in standard statistical packages. Therefore, random weight vectors cannot be directly created by means of random generators. A solution for this problem is given by the theorema which says that if  $G(\mathbf{x})$  is the distribution function of x, then  $\mathbf{u}=G(\mathbf{x})$  is uniformly distributed on the interval  $0 < \mathbf{u} < 1$ . (Hogg and Craig, 1970, p. 349). For the latter uniform distribution, standard random generators are available. then, if  $\mathbf{u}_1$  is uniformly distributed on the interval (0,1),  $\lambda_1 = G^{-1}(\mathbf{u}_1)$  can be shown to be distributed according to the density function  $g(\lambda_1)$  corresponding with the distribution function  $G(\lambda_1)$ . Thus, random values for  $\lambda_1$  can be found by using the following transformation.

 $\lambda_1 = 1 - (1 - u_1)^{1/(J-1)}$ 

For  $\lambda_2,\ldots,\,\lambda_{J-1}$  the following transformation has to be used:

$$\begin{split} \lambda_{j} &= (1 - \lambda_{1} - \ldots - \lambda_{j-1}) (1 - (1 - u_{j})^{1/(J-j)}) & j = 2, \ldots, J-1 \\ \text{Finally, } \lambda_{j} \text{ can be computed as } 1 - \lambda_{1} - \ldots - \lambda_{J-1} \end{split}$$

APPENDIX11

### Definition of evaluation criteria

Population.	A weighted sum of population around a location
	was calculated to quantify this score. The weight
	assigned decreases with distance. The <b>result</b> is
	standardized by dividing by the maximum score. A
	minus sign is added to <b>indicate</b> that the <b>crite-</b>
	rion is a <b>cost</b> criterion.

- Evacuation. The score reflects the availability of sufficient transport infrastructure.
- Agriculture at risk. This score reflects the location of agricultural land in the vicinity.
- Industry at risk. This score reflects the **size** and **importance** of industry near the location.
- Fresh water at risk. This score reflects the quantity of fresh water that **may** be **affected** by a **nuclear** plant at **each** location.
- Cool-water quantity. This score represents the quantity of available water for cooling the **nuclear** plant.
- Cool-water quality. This score represents the capacity of the **coolant** to flush **out** pollution originating from a **nuclear** plant at **each** location.
- Air pollution. It is **assumed** that the **nuclear** plant is an **alter**native to a conventional **coal** power plant, and so has the most beneficial effect at the most **pol**luted location.
- Thermal pollution. The amount of pollution is lower if users of the waste heat are available. The score reflects the availability of **such** users.
- Indirect land use. This score reflects limitations on potential land uses around a **nuclear** plant.
- Landscape. This score reflects the visual **effects** of the landscape and the extent to which a **nuclear** plant fits in with existing activities.
- Natural environment. This score reflects expected damage to the **natu**ral environment.
- National grid. This score reflects the proximity of high voltage lines and connector stations.
- Infrastructure. This score reflects the availability of transport and other infrastructure near the site.
- Coal-location. It is assumed that the **nuclear** plant is an **alter**native to a conventional **coal** power plant. The score reflects the **cost** of the lost opportunity to build a **coal** plant at the **site** if a **nuclear** plant is constructed.