WARPLAM DSS: using Cluster Analysis as an approach to delineate IWRM regions

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Abstract. The lack of uniform and integrated water resources regions is a critical issue, especially in transboundary water regions and federative countries. Overlaying levels of planning and management, as a result of uncoordinated water resources regions, hamper Integrated Water Resources Management. In order to harmonize multiple objectives and better represent the interaction between environmental, socio-economic, political and historical aspects, it becomes imperative to define appropriate territorial limits for water resources planning and management. The present study introduces an approach to support the process of delineating water resources regions. It is based both on recognition of more comprehensive aspects and incorporation of those aspects into a decision support system. This paper describes how cluster analysis is applied in the model design. Dynamic Programming is selected as the suitable method to be combined with Cluster Analysis to improve the algorithm efficiency.

Key Terms: water resources planning and management, decision support systems, cluster analysis, dynamic programming

1. Introduction

The lack of uniform and integrated water resources regions is a critical issue, especially in transboundary water regions and federative countries (Matthews and Germain 2007; Ganoulis et. al. 1996; EC 2002). Overlaying levels of planning and management, as a result of uncoordinated water resources regions, hampers Integrated Water Resources Management (IWRM). In addition, the process of delineating these regions has often been executed without sufficient scientific support, usually resulting from political and historical circumstances. In spite of this, it is possible to improve results by using knowledge from prior experiences, modern techniques and decision support systems (DSS). In order to harmonize multiple objectives and better represent the interaction between environmental, socio-economic, political and historical aspects, it becomes imperative to define appropriate territorial limits for water resources planning and management.

The present study introduces an approach to support the process of delineating water resources regions based both on recognition of more comprehensive aspects and incorporation of those aspects into a DSS. The proposed Water Resources Planning and Management Regions (WARPLAM) DSS is designed to be used by federal and state governments, international commissions and water councils. Considering that river basins are the most suitable boundaries to attain IWRM goals (Dourojeanni et. al 2002; Wegerich 2008;

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Falkenmark 2004), the DSS simulation model offers the option for decision makers to include socio-economic, political and environmental aspects into the analysis, as suggested by Porto & Porto (2008). It intends to promote a better understanding about the reasoning related to this process, and to reinforce the principles of IWRM. WARPLAM DSS is also a very flexible solution to support the delineation of regions in multiple levels of subsidiarity and to be adaptable to regional circumstances.

This paper describes how cluster analysis is applied in the model design, combined with Dynamic Programming. Among the available techniques, Dynamic Programming has proven to be a valuable tool to support cluster analysis (Belman, 1973). It increases substantially the algorithm efficiency, considering the number of combinations in an exhaustive enumeration search can be too extensive.

This paper is organized in two main topics. The first topic describes how the DSS is developed, including a general overview of its structure and procedures. The second topic is the description of the algorithm that constitutes the model of the DSS. It includes the logic associated with combining Cluster Analysis and Dynamic Programming. Finally, conclusions and general recommendations are presented.

2. WARPLAM DSS: The Proposed Approach

Water Resources Planning and Management Regions Decision Support System is the proposed approach to address the issue of lack of uniform and integrated water resources regions. It constitutes a structured and instructive tool to help decision makers to delineate water resources regions, which is usually an ill structured task. Another important characteristic of the proposed approach is its ability to help harmonizing multiple interests from different stakeholders.

To describe the process of developing this approach, it is helpful to understand the main steps of the decision making process related to the exercise of delineating water resources planning and management regions. In this study, it is organized into five basic steps. The first step is the definition of a consistent basis over which to develop an aggregation process. This is a very important step because it represents the main aspect to be considered for the water resources regions. From the grouping of those smaller territorial units, for example natural drainage areas or municipalities, water resources planning and management regions will be created. The second step is the selection of criteria, beyond river basin boundaries, that reflect the main aspects related to IWRM principles. Those criteria represent the recognition of more comprehensive objectives and multiple interests into the analysis. A specific comparative analysis was performed, based on selected examples from European and American countries, in order to enhance the selection of criteria. This comparative study constitutes an Expert System, as proposed by Turban (1998), used to support the decision making process. This step also includes weights assignment for each criterion. The third step is the combination of selected criteria with the basis in order to define the ‘measure of closeness’ for each adjacent pair of territorial units contained in the basis. Each of those pairs constitutes one grouping alternative. The ‘measure of closeness’ for each alternative is defined taking into account overlaying area values of all the criteria. The fourth step is the application of compromise programming to sum up all weighted criteria values for each alternative, considering the different scale range or space dimensions of the criteria’ values. The fifth and last proposed step is the application of Cluster Analysis to define different grouping alternatives that represent ‘ideal’ IWRM regions.
WARPLAM DSS: using Cluster Analysis as an approach to delineate IWRM regions

After considering the main steps of the decision making process, the DSS procedures and structure are presented, followed by a description of its components, as well as the model outline.

2.1. DSS Procedures and Structure

WARPLAM DSS is structured using ESRI ArcGIS, Microsoft Excel and Visual Basic functionalities. The main procedures performed within the DSS are described below.

The first two steps of the decision making process are supported mainly by ArcGIS functionalities. The criteria and basis selection is facilitated through the use of GIS techniques. The Expert System is integrated to the GIS interface to provide the necessary understanding about the criteria selection process, based on heuristic rules derived from the comparative study. In such cases, the decision makers are able to learn from past experiences and to decide, based on their own preferences, which of those aspects are important in the specific context of the case in analysis.

As described before, the basis contains the territorial units to be grouped. For example, the adoption of a consistent basis considering natural drainage area limits represents the consideration of watershed boundaries as the basis for the analysis. Instead, the adoption of municipalities represents the consideration of political-administrative boundaries as the basis for the analysis. The selected criteria should reflect the main aspects related to IWRM. These criteria, as well as the basis, must be available in the format of spatial data, as the necessary input for the model. As soon as the criteria are selected, data can be easily imported to the DSS. The ESRI Geodatabase format is recommended, but data may also be imported using shapefile or coverage formats.

After data is imported, the Database Management System handles all pre-processing analysis, as part of Step 3 of the decision making process, in order to prepare the input data to the model system. Knowledge rules, imported from the Expert System are directly integrated into the database. The intersection among chosen criteria and the consistent basis is performed. In order to support the creation of a more functional and user-friendly interface, the Model Builder ArcGIS functionality is used. This tool allows all the repeated tasks, to be performed at one click, according to the selected functionalities. In such case, the calculation of all overlaying areas is performed by one click and the results are being incorporated into the model system through the use of a single workspace. Microsoft Visual Basic functionalities are also used in this stage of the process to perform some necessary data management tasks, in integration with ArcGIS. As a result of this pre-processing stage, all overlaying areas of selected criteria are calculated and combined with the knowledge rules from the Expert System. In addition, all adjacent pairs are listed as possible alternatives to be grouped. Therefore, the necessary input for the model system is ready and the algorithm can be started.

Steps 4 and 5 of the decision making process are basically performed inside the model system, which will be described in the following section. The algorithm is developed using Microsoft Excel Macros and Visual Basic Codes, which guarantee the necessary integration among the data management system and the model system. In addition, optimization techniques are applied to support the clustering process and to increase the algorithm’s efficiency. As soon as the data is read in the model system, the user needs to define the weights for each criterion and some parameters for the cluster analysis. This is also facili-
tated through a user-friendly interface. Finally, the results from the simulation are displayed into the GIS interface automatically. Figure 1 illustrates a summary of those procedures.

3. Model Outline and Algorithm Structure

The model structure is comprised of the algorithm developed to address the delineation of water resources regions. It is divided in two main modules, correspondent to the Steps 4 and 5 of the decision making process, as described above. The main input for the algorithm comes from the intersection between selected criteria and the basis, performed in Step 3. Each pair of adjacent units contained in the basis constitutes one alternative to be considered for the cluster analysis.

Cluster analysis is a set of procedures used to create classification and reorganize data into homogeneous groups (MOPU 1984; Kaufman & Rousseeuw 1990). The first stage of the Cluster Analysis is to define a numerical measure of homogeneity (Bellman 1973; Aldenderfer & Blashfield 1984). For that measure, this approach employs the concept of ‘measure of closeness’. It is defined, for each adjacent pair, taking into account the criteria overlaying area values over the basis. Considering that the calculations are performed based on area values, it is not necessary to standardize the data. As soon as the initial units are defined in the database management system, uniform outputs are provided. In addition, the compromise programming step handles different data dimensions. According to Coelho et. al. (2005) the ‘measure of closeness’ can be calculated through the size and proportion of the common criteria area overlaying one adjacent pair of the basis’ territorial units. Besides showing how relevant a common criterion is to the pair (size), the measure also needs to express how equal the two units are in reference to the criteria (proportion). In addition the common perimeter is also considered. By grouping these three aspects, the following vector-based equation was applied for each alternative and each criterion:

\[
C_{i,2} = \frac{2 \cdot CP_{i,2}}{P_{WS1} + P_{WS2}} \cdot \frac{A_{WS1} \cdot A_{WS2}}{A_{WS1} + A_{WS2}}
\] (1)
WARPLAM DSS: using Cluster Analysis as an approach to delineate IWRM regions

Considering:

\[ A_{Ci, WS1} = \text{Overlaying Area of Criteria } i \text{ over territorial unit } WS1 \ (i = 1, 2, \ldots, N) \]

\[ A_{Ci, WS2} = \text{Overlaying Area of Criteria } i \text{ over territorial unit } WS2 \ (i = 1, 2, \ldots, N) \]

N = number of criteria defined by the user

\[ A_{WS1} = \text{Area of territorial unit } WS1 \]

\[ P_{WS1} = \text{Perimeter of territorial unit } WS1 \]

\[ CP_{1,2} = \text{Common Perimeter between territorial units } 1 \text{ and } 2 \]

\[ C_{1,2} = \text{Measure of closeness between units } 1 \text{ and } 2, \text{ considering Criteria } i \]

\[ C_{1,2} \text{ ranges from } 0 \text{ to } 1 \]

As soon as the list of alternatives (adjacent pairs) and respective measures of closeness is ready, the algorithm is started. The first module of the algorithm is the application of compromise programming to sum up the measures of closeness of each criterion value for each alternative, resulting in the total measure of closeness \( C_{a,b} \) for each alternative. This method was considered the most adequate considering the different scale range of criteria values (different space dimensions) and its ability to rank alternatives according to their ‘closeness’ to certain ‘ideal’ criteria levels (Hajkowicz & Collins 2007; Labadie 2007). The scaling function was applied by the selection of the best and worst values of the alternatives for each criterion, according to Equation 02. The Total Measure of Closeness is assigned as a ‘link’ between adjacent basis’ territorial units, and represents the proximity between those units into each pair. Compromise solutions are the result of combining different L1, L2 and L\( \infty \) norms and different sets of weights.

\[
L_{\sigma}(j) = \sum_{i=1}^{P} \alpha_{i}^{\sigma} \left[ \frac{z_{i}^{*} - z_{i}(j)}{z_{i}^{*} - z_{i}^{**}} \right]^{\sigma} 
\]  
(2)

The second module of the algorithm is the application of Cluster Analysis over alternatives to define different grouping alternatives or clusters. The total measure of closeness between each pair is used as the input to the similarity matrix of elements to be clustered. It is considered the most adequate method to directly represent the relative distances between the elements to be clustered. Alternatives of groups with higher similarity will be formed in order to delineate the “ideal” regions for water resources planning and management. The partitioning method is applied according to the calculated overall proximity of each cluster created. The partitioning clustering method, in contrast to the hierarchical method, generally results in better patterns of similarities between elements of the groups, because the overall distance of the group is being considered (Kaufman & Rousseeuw 1990; Aldenderfer & Blashfield 1984). This overall distance is calculated using the average proximity between all elements of the group. The objective function is to maximize the overall proximity of all clusters (minimize intra-cluster variance). The constraints associated with the problem are derived from the knowledge rules existent at the Expert System, according to the decision maker preferences.

A significant drawback of this method is the infinite number of alternatives to be analyzed (Kaufman & Rousseeuw 1990). Depending on the number of elements to be grouped, the analysis may become too extensive. In such case, Dynamic Programming (DP) can be applied to support the evaluation of multiple alternatives. It speeds up the analysis consistently and is ideally suited to be applied with cluster analysis (Bellman 1973; Esogbue 1986).
Currently, the DP method is being tested to support the analysis, using the generalized dynamic programming software developed by Labadie (1990). A 9-element data set is adopted as the trial exercise of the method in study, assuming that it is a result of the Steps 1, 2 and 3 of the decision making process. The following distance matrix contains the ‘total measure of closeness’ for each of the ten pairs of alternatives in the analysis (Table 1):

**Table 1. Distance Matrix adopted as an example for the Cluster Analysis**

<table>
<thead>
<tr>
<th></th>
<th>a</th>
<th>b</th>
<th>c</th>
<th>d</th>
<th>e</th>
<th>f</th>
<th>g</th>
<th>h</th>
<th>i</th>
</tr>
</thead>
<tbody>
<tr>
<td>a</td>
<td>NA</td>
<td>0.2</td>
<td>0.4</td>
<td></td>
<td></td>
<td></td>
<td></td>
<td></td>
<td></td>
</tr>
<tr>
<td>b</td>
<td>0.2</td>
<td>NA</td>
<td>0.8</td>
<td></td>
<td></td>
<td></td>
<td></td>
<td></td>
<td></td>
</tr>
<tr>
<td>c</td>
<td>0.4</td>
<td>NA</td>
<td>0.7</td>
<td>0.5</td>
<td></td>
<td></td>
<td></td>
<td></td>
<td></td>
</tr>
<tr>
<td>d</td>
<td>0.8</td>
<td>0.7</td>
<td>NA</td>
<td></td>
<td></td>
<td></td>
<td></td>
<td></td>
<td></td>
</tr>
<tr>
<td>e</td>
<td>0.5</td>
<td>NA</td>
<td>0.6</td>
<td></td>
<td></td>
<td></td>
<td></td>
<td></td>
<td></td>
</tr>
<tr>
<td>f</td>
<td>0.6</td>
<td>NA</td>
<td>0.1</td>
<td>0.5</td>
<td></td>
<td></td>
<td></td>
<td></td>
<td></td>
</tr>
<tr>
<td>g</td>
<td></td>
<td>0.1</td>
<td>NA</td>
<td>0.7</td>
<td></td>
<td></td>
<td></td>
<td></td>
<td></td>
</tr>
<tr>
<td>h</td>
<td>0.5</td>
<td>NA</td>
<td>0.5</td>
<td></td>
<td></td>
<td></td>
<td></td>
<td></td>
<td></td>
</tr>
<tr>
<td>i</td>
<td></td>
<td></td>
<td>0.7</td>
<td>0.5</td>
<td>NA</td>
<td></td>
<td></td>
<td></td>
<td></td>
</tr>
</tbody>
</table>

The intra-cluster measure of homogeneity is calculated considering the overall average of the ‘measures of closeness’. For example, for the 4-element cluster ‘a-b-c-d’ it is equal to 0.525, taking into account the list of pairs and respective ‘measure of closeness’ contained in Table 2. The inter-cluster measure of homogeneity is then calculated by taking the average of the intra-cluster measure of homogeneity. For example, the nine available elements can be clustered in three groups of 2, 3 and 4 elements, respectively. The inter-cluster measure of homogeneity is then the average of the three intra-cluster measures of homogeneity, as shown in Table 3.

**Table 2. Intra-Cluster Measure of Homogeneity**

<table>
<thead>
<tr>
<th>ab</th>
<th>0.2</th>
</tr>
</thead>
<tbody>
<tr>
<td>ac</td>
<td>0.4</td>
</tr>
<tr>
<td>bd</td>
<td>0.8</td>
</tr>
<tr>
<td>cd</td>
<td>0.7</td>
</tr>
</tbody>
</table>

<table>
<thead>
<tr>
<th>ab</th>
<th>0.2</th>
</tr>
</thead>
<tbody>
<tr>
<td>ac</td>
<td>0.4</td>
</tr>
<tr>
<td>bd</td>
<td>0.8</td>
</tr>
<tr>
<td>cd</td>
<td>0.7</td>
</tr>
</tbody>
</table>

**Table 3. Inter-Cluster Measure of Homogeneity**

<table>
<thead>
<tr>
<th>ef</th>
<th>0.6</th>
</tr>
</thead>
<tbody>
<tr>
<td>fg</td>
<td>0.1</td>
</tr>
<tr>
<td>hi</td>
<td>0.3</td>
</tr>
</tbody>
</table>

It is assumed that if the cluster has one element, the intra-cluster measure of homogeneity is equal to zero. The objective is to reduce the inter-cluster measure of homogeneity if there are clusters containing just one element. This way, the best grouping alternatives —
or the ones containing the highest inter-cluster measure of homogeneity – are more homogeneous. For example: having clusters ‘e-f’ and ‘c-d’ (Inter-Cluster = 0.65, as the average of 0.6 and 0.7) is better than having clusters ‘c-d-e’ and ‘f’ (Inter-Cluster = 0.30, as the average of 0.60 and 0.00).

Therefore, the objective is to maximize the inter-cluster measure of homogeneity. For that, the DP analysis is divided in two parts, according to the method suggested by Bellman & Zadeh (1970), Bellman (1973), Esogbue (1986). It consists in dividing the set of alternatives into I groups, according to the intra-cluster measure of homogeneity (measure of closeness), and determining the optimal value of I according to the inter-cluster measure of homogeneity, and then the optimal subdivision. The additive objective function is to maximize the total benefits of allocating m_i objects to I clusters. The DP recursion relation and other related equations are defined as following:

\[ F_i^{'}(m_{i+1}) = \max \{ u_i^{'}(m_i) + F_{i-1}(m_i) \} \]  

(03)

\[ \text{S.T.} \]
\[ 0 \leq x_i \leq M \]
\[ 0 < m_i = x_i - x_{i+1} \leq M \text{ (no cluster with 0 elements)} \]
\[ m_i = 1, 2, \ldots, M \]

For all discrete \( x_{i+1} \): \( 0 \leq x_{i+1} \leq M-i \)

- Over stages \( i = 1, 2, \ldots, M \)
- \( x_1 = M; x_M = 0 \) (all element should be clustered at the end)

Optimal solution can be found in any stage when \( x_{i+1} = 0 \)

Starting with: \( F_0(m_1) = 0 \)

Max \( F_i(M) \)

Considering:
- \( M \) = total number of elements to be grouped
- \( m_i \) = number of elements in the cluster
- \( u_i^{'}(m_i) \) = best benefit of having \( m_i \) elements in the cluster
- \( i \) = number of clusters = number of stages in DP
- \( x_i \) = state variables
- \( m_i \) = decision variables

The decision variable is the number of elements to be included in the cluster in each stage. The state variables are the number of elements remaining to be allocated, using the concept of the resources allocation problem. They are both integer values, according to the nature of the problem. The benefit is equal to the intra-cluster measure of homogeneity (average of the measures of closeness). It is calculated in a pre-optimization step that returns the best possible benefit for a cluster having \( m_i \) elements. The forward DP recursion relation and the inverted form of the state dynamic equation are adopted.

The important concept that is added to Belman’s first proposed recursion relation is the ability to store the information calculated in the stage before and use it as an input for the sequence of the solution. The proposed method stores the best results in each stage to be used in the next stage in order to exclude the elements already clustered.
To do that, the binary string concept is also applied. This is a really efficient and unique way to organize the data. Considering all possible combinations among the elements, the position in a string determines if the element is included in the cluster or not. Therefore, in each stage and state variable, the algorithm returns a unique number that is associated with a string that represents the clustered elements. There is a unique number associated with each possible combination (Table 4). This unique number is used to guarantee that the elements previously clustered are not included in the current stage.

<table>
<thead>
<tr>
<th>b</th>
<th>128</th>
<th>0</th>
<th>1</th>
<th>0</th>
<th>0</th>
<th>0</th>
<th>0</th>
<th>0</th>
<th>0</th>
</tr>
</thead>
<tbody>
<tr>
<td>c-d</td>
<td>96</td>
<td>0</td>
<td>0</td>
<td>1</td>
<td>1</td>
<td>0</td>
<td>0</td>
<td>0</td>
<td>0</td>
</tr>
<tr>
<td>c-d-e</td>
<td>224</td>
<td>0</td>
<td>1</td>
<td>1</td>
<td>1</td>
<td>0</td>
<td>0</td>
<td>0</td>
<td>0</td>
</tr>
<tr>
<td>c-e-f</td>
<td>88</td>
<td>0</td>
<td>0</td>
<td>1</td>
<td>0</td>
<td>1</td>
<td>1</td>
<td>0</td>
<td>0</td>
</tr>
<tr>
<td>c-e-f-h</td>
<td>90</td>
<td>0</td>
<td>0</td>
<td>1</td>
<td>0</td>
<td>1</td>
<td>1</td>
<td>0</td>
<td>0</td>
</tr>
</tbody>
</table>

Table 4. Unique Number and Binary String for Different Combinations

In addition, the running average concept, as defined by Lee & Labadie (2007) is applied to calculate the objective function in each stage. The additive objective function, in principle, does not allow the calculation of the inter-cluster measure of homogeneity in each stage, because the average is required. To overcome this issue, a discount factor (DF) is added to both parts of the recursion equation. In the first part of the recursion equation – \( f^*(m_i - m_{i-1}) \) – the discount factor is equal to ‘1/i’. In the second part of the recursion equation - \( F_{i-1}(m_i) \) – the discount factor is equal to ‘(i-1)/i’. As a result, the objective function is adapted to the DP format and the ‘running average’ is calculated in each stage. CSUDP allows the user to define a Discount Factor for the objective function. The discount factor cannot be the same for all stages because the elements would have a different weight depending on the stage it is selected to be part of the cluster. The DP recursion relation, including the DFs, becomes:

\[
F_i(m_{i+1}) = \max \left\{ \frac{1}{i}f^*(m_i - m_{i-1}) + (\frac{i-1}{i})F_{i-1}(m_i) \right\}
\]

4. Conclusions and Recommendations

The presented approach constitutes a prototype of a decision support system to address the necessity of defining water resources regions. As demonstrated in this paper, it includes human intuition and judgment, e.g. subjective criteria selection and weighting processes. Through a user-end focus, it also provides easy access to information; interaction, which is supported by visualization of criteria; and flexibility, since it is open to aggregate other criteria in order to consider new aspects. In addition, it constitutes a learning process because decision makers can better understand the aspects related to water units delineation, using Expert Systems.
Regarding the potential results, it is possible to observe that the model can easily incorporate a set of criteria and weights. Multiple simulations can be simply performed in order to support the decision making process. Dynamic Programming has proven to increase the efficiency of the algorithm, especially when compared to the alternative exhaustive enumeration method. According to the results of the simulations being tested, for a data set containing five elements, 24 intra-cluster and 48 inter-cluster measure of homogeneity can be found in exhaustive enumeration. Dynamic Programming one-dimensional algorithm analyzes only 16. For the given 9-elements dataset presented in this paper, 90 intra-cluster and 1300+ inter-cluster valid measures of homogeneity can be found in exhaustive enumeration Dynamic Programming one-dimensional algorithm analyzes around 240.

![Figure 2. Possible Solutions and Respective Return Values in each stage.](image)

Table 5 presents the maximum benefit considering the best solution for the given 9-element data set adopted as the trial exercise of the method in study.

<table>
<thead>
<tr>
<th>Clusters</th>
<th>f(x)</th>
<th>m_i</th>
<th>X_i</th>
<th>X_{i+1}</th>
</tr>
</thead>
<tbody>
<tr>
<td>b-d</td>
<td>0.8</td>
<td>2</td>
<td>9</td>
<td>7</td>
</tr>
<tr>
<td>g-i</td>
<td>0.7</td>
<td>5</td>
<td>7</td>
<td>5</td>
</tr>
<tr>
<td>a-c-e-f-h</td>
<td>0.5</td>
<td>2</td>
<td>5</td>
<td>0</td>
</tr>
<tr>
<td>Max</td>
<td>0.67</td>
<td></td>
<td></td>
<td>3 clusters</td>
</tr>
</tbody>
</table>

The optimization of the best number of groups can be incorporated into the analysis, according to the user preferences. However, it is important to highlight the objective of this DSS is not to guarantee the optimum number of groups. Instead, different simulations of grouping alternatives seem to be more important for the decision maker in order to evaluate the problem.

Regarding the cluster analysis method, its key is to define ‘real groups’ instead of ‘imposed groups’ (Aldenderfer & Blashfield 1984). In this case, the combination of cluster analysis with dynamic programming and the adopted ‘closeness measure’ guarantees ‘ideal’ solutions. The combination of these techniques used to address the presented issue is the innovation proposed in this study. However, it is important to affirm that this paper refers to an ongoing study that may lead to future reviews and adjustments.
It is expected that the final algorithm to be developed will incorporate fuzzy analysis in order to indicate the membership function of each element to its final cluster, as suggested by Kaufman & Rousseeuw (1990). It will represent the uncertainty associated with defining element X as part of cluster Y. According to Hajkowicz & Collins (2007) the fuzzy set theory is excellent to handle uncertainty inherent in ill-structured problems. In addition, another advantage of the selected partitioning method for the cluster analysis is that it allows the representation of the results considering fuzzy logic.

In addition, it is possible that a multi-dynamic program analysis will be necessary to prove that the best solution is reached in DP. As an ongoing research project, some tests are being generated in order to evaluate it. Generic algorithm will also be assessed as a way to increase the algorithm’s efficiency. Another important aspect that needs special attention is the occurrence of ties when returning the benefit value, in the pre-optimization step.

Finally, the 2nd United Nations World Water Development Report: “Water, a shared responsibility” pointed out the need for an integrated and holistic approach to water resources management, highlighting the benefits from IWRM: 1) multiple uses and cooperation between different sectors; 2) coordinated management and development of land, water and other resources; and 3) balanced social, environmental and economic benefits (UNESCO, 2006). Therefore, to integrate political divisions within river basin units is one of the biggest challenges. Still according to this report, the difficulties of IWRM are directly related to the fact that political boundaries are not coincident with natural river basins units.

In this sense, the presented study reinforces the importance of defining IWRM units and demonstrates the DSS approach as a method to support multiple interest decision process, reflecting human judgment through easy access to information, education, interaction and flexibility. As demonstrated, it addresses the solution of such a complex and ill-structured problem. Future decisions related to water resources regions delineation may have increased quality by using knowledge from prior experience and modern techniques, instead of letting the process to be a result of political and historical circumstances only.

References


