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ROBUST AUTOMATIC METHODS FOR OUTLIER AND ERROR DETECTION

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within these subsets to identify errors. The second uses a robust regression tree modelling

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Key words: Survey data editing; representative outliers; robust regression; regression tree

model; gross errors; M-estimates

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1. Introduction

1.1 Overview

Outliers are a common phenomenon in business and economic surveys. These are data values that are so unlike values associated with other sample units that ignoring them can lead to wildly inaccurate survey estimates. Outlier identification and correction is therefore an important objective of survey processing, particularly for surveys carried out by national statistical agencies. In most cases these processing systems operate by applying a series of edits that identify data values outside bounds determined by the expectations of subject matter specialists. These outlier values are then investigated further, in many cases by re-contacting the survey respondents, to establish whether they are due to errors in the data capture process or whether they are in fact valid. Chambers (1986) refers to the latter valid values as representative outliers, insofar as there is typically no reason to believe that they are unique within the survey population. Outlier values that are identified as errors, on the other hand, are not representative, and it is assumed that they are corrected as part of survey processing. A common class of such errors within the business survey context is where the survey questionnaire asks for answers to be provided in one type of unit (e.g. thousands of pounds) while the respondent mistakenly provides the required data in another unit (e.g. single pounds). Sample cases containing this type of error therefore have true data values inflated by a factor of 1000. Left uncorrected, such values can seriously destabilise the survey estimates.

The standard approach to the type of situation described above is to use a large number of edits to identify as many outliers as possible during survey processing. These outliers are then followed up to establish their correct values. If the correct value is identical to the value that triggered the edit failure, then this value is not an error but corresponds to a representative outlier. In this case the usual strategy is to replace it by an imputed value,

typically one that is subjectively determined as "more typical". In continuing surveys this can be the previous value of the same variable, provided that value is acceptable.

There are two major problems with this approach. The first is that it can be extremely labour intensive. This is because the edit bounds are often such that a large proportion of the sample data values lie outside them. This leads to many unnecessary re-contacts of surveyed individuals or businesses, resulting in an increase in response burden. Secondly, the subjective corrections applied to representative outliers leads to biases in the survey estimates, particularly for estimates of change. Since there are often large numbers of such representative outliers identified by this type of strategy, the resulting biases from their "correction" can be substantial.

This paper describes research aimed at identifying an editing strategy for surveys that are subject to both outliers and errors that overcomes some of the problems identified above. In particular, the aim is to develop an automated outlier identification strategy that finds as many significant errors in the data as possible, while minimising the number of representative outliers also identified (and whose values are therefore incorrectly changed). In particular, the methods described below do not rely on specification of edit bounds and use modern robust methods to identify potential errors, including outliers, from the sample data alone.

1.2 The ABI Data

This research has been carried out as part of the *Euredit* project (Charlton *et al*, 2001). In particular, we use two data sets that were created within this project for the specific purpose of evaluating automatic methods for edit and imputation. Both contain data for 6099 businesses that responded to the UK Annual Business Inquiry (*ABI*) in the late 1990s. The values in the first data set have been thoroughly checked and there are no missing values. We refer to it as the *clean data* below. Note that these data are not necessarily the "truth". What

they represent are data of sufficient quality for use in official statistics. The second data set contains values for the same variables and businesses as in the first. However, these values now include both introduced errors and missing values, and can be considered as representing the type of "raw" data that are typically seen prior to editing. We refer to this data set as the *perturbed data* below. Note that the *clean data* and the *perturbed data* contain a significant number of common extreme values (i.e. representative outliers), reflecting the fact that the *clean data* may still contain errors.

Table 1. The ABI variables

Name	Definition
turnover	Total turnover
emptotc	Total employment costs
purtot	Total purchases of goods and services
taxtot	Total taxes paid
assacq	Total cost of all capital assets acquired
assdisp	Total proceeds from capital asset disposal
employ	Total number of employees

Table 1 lists the names and definitions of the variables collected in the *ABI* that we consider in this paper. The names are those used in the *Euredit* project, and these seven variables represent the major outcome variables for the *ABI*. In addition we assume that we have access to "complete" (no missing values and no errors) auxiliary information for the sampled businesses on the Inter-Departmental Business Register or *IDBR* (the sample frame for the *ABI*). The most important auxiliary variable is the estimated turnover of a business (*turnreg*, defined in terms of the *IDBR* value of turnover for a business, in thousands of pounds). Other auxiliary information on the *IDBR* relates to the estimated number of employees of a business and its industrial classification. Together, these define the sampling strata for the *ABI*. Below we assume that we have access to the strata affiliations of the sampled businesses.

Figures 1 and 2 show the relationship between the ABI variables and the auxiliary variable turnreg for the values contained in the $clean\ data$ and the $perturbed\ data$. Because these variables are extremely heteroskedastic, scatterplots of their raw values reveal little. Consequently all plots in Figures 1 and 2 are on the log scale. Also, for reasons of confidentiality, no scale is displayed on these plots. It is clear from Figures 1 and 2 that although the general relationship between turnreg and the ABI variables is linear in the log scale, comparison of the $clean\ data$ and the $perturbed\ data$ show there are a very large number of significant errors leading to outliers (these appear as Δ 's in Figure 2, but not in Figure 1) as well as large representative outliers (these appear in both Figure 1 and Figure 2).

Figures 1 and 2 about here

2. Outlier Identification Via Forward Search

The first automatic method we investigate was suggested by Hadi and Simonoff (1993). See also Atkinson and Riani (2000). The basic idea is simple. In order to avoid the well-known masking and swamping problems that can occur when there are multiple outliers in a data set (see Barnett and Lewis 1994, Rousseeuw and Leroy 1987), the algorithm starts from an initial subset of observations of size m < n that is chosen to be outlier free. Here n denotes the number of observations in the complete data set. In the regression context, a model for the variable of interest is estimated from this initial "clean" subset. Fitted values generated by this model are then used to generate n distances to the actual sample data values. The next step in the algorithm redefines the clean subset to contain those observations corresponding to the m+1 smallest of these distances and the procedure repeated. The algorithm stops when distances to all sample observations outside the clean subset are all too large or when this subset contains all n sample units.

In order to more accurately specify this forward search procedure, we assume values of a p-dimensional multivariate survey variable \mathbf{Y} and a q-dimensional multivariate auxiliary variable \mathbf{X} are available for the sample of size n. We denote an individual's value of \mathbf{Y} and \mathbf{X} by \mathbf{y}_i and \mathbf{x}_i respectively. The matrix of sample values of \mathbf{Y} and \mathbf{X} is denoted by $\mathbf{y} = (\mathbf{y}_1 \cdots \mathbf{y}_n)'$ and $\mathbf{x} = (\mathbf{x}_1 \cdots \mathbf{x}_n)'$ respectively. We seek to identify possible outliers in \mathbf{y} .

Generally identification of such outliers is relative to some assumed model for the conditional expectation of **Y** given **X**. Given the linear structure evident in Figures 1 and 2, we assume that a linear model $E(\mathbf{Y}) = \beta \mathbf{X}$ can be used to characterise this conditional expectation, where $\boldsymbol{\beta} = (\beta_1', \dots, \beta_p')'$ is a $p \times q$ matrix of unknown parameters. A large residual for one or more components of **Y** is typically taken as evidence that these values are outliers.

For p = 1 we let $\hat{\beta}_{(m)}$ and $\hat{\sigma}^2_{(m)}$ denote the regression model parameter estimates based on a clean subset of size m. For an arbitrary sample unit i, Hadi and Simonoff (1993) suggest the distance from the observed value y_i to the fitted value generated by these estimates be calculated as

$$d_{i(m)} = \frac{\left| y_i - \mathbf{x}_i' \hat{\boldsymbol{\beta}}_{(m)} \right|}{\hat{\sigma}_{(m)} \sqrt{1 - \lambda_i \mathbf{x}_i' (\mathbf{X}_{(m)}' \mathbf{X}_{(m)})^{-1} \mathbf{x}_i}}$$

where $\mathbf{X}_{(m)}$ denotes the matrix of values of \mathbf{X} associated with the sample observations making up the clean subset, and λ_i takes the value 1 if observation i is in this subset and -1 otherwise. The clean subset of size m+1 is then defined by those sample units with the m+1 smallest values of $d_{i(m)}$. For p>1, Hadi (1994) uses the squared Mahalanobis distance

$$D_{i(m)}^{2} = (\mathbf{y}_{i} - \hat{\mathbf{y}}_{i(m)})' \hat{\mathbf{S}}_{(m)}^{-1} (\mathbf{y}_{i} - \hat{\mathbf{y}}_{i(m)})$$

where $\hat{\mathbf{y}}_{i(m)}$ denotes the fitted value for \mathbf{y}_i generated by the estimated regression models for the components of this vector, and $\hat{\mathbf{S}}_{(m)} = (m-q)^{-1} \sum_{i} (\mathbf{y}_i - \hat{\mathbf{y}}_{i(m)}) (\mathbf{y}_i - \hat{\mathbf{y}}_{i(m)})'$ denotes the

estimated covariance matrix of the errors associated with these models. The summation here is over the observations making up the clean subset of size m.

For p=1 Hadi and Simonoff (1993) suggest stopping the forward search when the $(m+1)^{th}$ order statistic for the distances $d_{i(m)}$ is greater than the $1-\alpha/2(m+1)$ quantile of the t-distribution on m-q degrees of freedom. When this occurs the remaining n-m sample observations are declared as outliers. Similarly, when p>1, Hadi (1994) suggests the forward search be stopped when the (m+1)th order statistic for the squared Mahalanobis distances $D_{i(m)}^2$ exceeds the $1-\alpha/n$ quantile of the chi-squared distribution on p degrees of freedom.

Definition of the initial clean subset is important for implementing the forward search procedure. Since the residuals from the estimated fit based on the initial clean subset define subsequent clean subsets, it is important that the parameter estimates defining this estimated fit are unaffected by possible outliers in the initial subset. This can be achieved by selecting observations to enter this subset only after they have been thoroughly checked. Alternatively, we use an outlier robust estimation procedure applied to the entire sample to define a set of robust residuals, with the initial subset then corresponding to the *m* observations with smallest absolute residuals relative to this initial fit. Our experience is that this choice is typically sufficient to allow use of more efficient, but non-robust, least squares estimation methods in subsequent steps of the forward search algorithm.

Before ending this section, we should point out that the forward search method described above is based on a linear model for the non-outlier data, with stopping rules that implicitly assume that the error term in this model is normally distributed. Although Figures 1 and 2 indicate that these assumptions are not unreasonable for logarithmic transforms of the *ABI* data, it is also clear that they are unlikely to hold exactly. This may not be of practical consequence if these stopping rules are activated when the error distribution starts to deviate

from the normal. However, an alternative outlier detection method that does not assume either linearity or normality would appear to be worth considering.

3. Outlier Identification Via Robust Tree Modelling

Regression tree models (Breiman *et al*, 1984) are now widely used in statistical data analysis, especially in data mining applications. Here we use a tree modelling approach that is robust to the presence of outliers in data to identify gross errors and extreme outliers. The WAID software for regression and classification tree modelling that was used for this purpose was developed for missing data imputation under the *Autimp* project (Chambers *et al*, 2001). Under the *Euredit* project a toolkit of programs has been created that emulates and extends the capabilities of WAID. These programs work under R (Ihaka and Gentleman, 1996). Similar software products are CART (Steinberg and Colla, 1995), the S-PLUS (MathSoft, 1999) *tree* function (also available in R), the R function *rpart* and the CHAID program in SPSS AnswerTree (SPSS, 1998). The code for the WAID toolkit is available from the authors.

The basic idea behind WAID is to sequentially divide the original data set into subgroups or nodes that are increasingly more homogeneous with respect to the values of a response variable. The splits themselves are defined in terms of the values of a set of categorical covariates. The categories of these covariates do not need to be ordered. By definition, WAID is a nonparametric regression procedure. It also has the capacity to implement outlier robust splitting based on M-estimation methodology (Huber, 1981). In this case outliers are "locally" down-weighted when calculating the measure of within node heterogeneity (weighted residual sum of squares) used to decide whether a node should be split or not. The weights used for this purpose are themselves based on outlier robust influence functions.

3.1 The WAID Regression Tree Algorithm For Univariate Y

This assumes a rectangular data set containing n observations, values $\{y_i\}$ of a univariate response variable Y and values $\{x_{1i}, \dots x_{pi}\}$ of p categorical covariates X_1, \dots, X_p . No missing X-values are allowed in the current version of WAID. For scalar Y WAID builds a regression tree. If Y is categorical, WAID builds a classification tree. The only difference between these two types of trees is the heterogeneity measure used to determine tree-splitting behaviour. Since our focus is outlier identification, we are concerned with scalar response variables only and so we restrict consideration to WAID's regression tree algorithm.

The basic idea used in WAID (as well as other tree-based methods) is to split the original data set into smaller subsets or "nodes" in which Y-values are more homogeneous. In WAID this is accomplished by sequential binary splitting. At each step in the splitting process, all nodes created up to that point are examined in order to identify the one with maximum heterogeneity. An optimal binary split of this "parent" node is then carried out. This is based on identifying a set of values of one of the covariates $X_1, ..., X_p$ such that a split of the parent node into one "child" node containing only cases possessing these values and another child node containing the remaining cases minimises the heterogeneity of these child nodes. In searching for this optimal split, covariates are classified as monotone or nonmonotone. Candidate splits for a monotone X are determined by splits with respect to the ordered values of X in that node. Candidate splits for a non-monotone X are defined with respect to values of X sorted by their corresponding average Y value in the node. The splitting process continues until a suitable stopping criterion is met. At present this is when either (i) all candidate parent nodes are effectively homogeneous; (ii) all candidate parent nodes are too small to split further; or (iii) a user-specified maximum number of nodes is reached. Unlike some other tree modelling software packages (e.g. CART), there is no attempt to find an "optimal" tree. The set of nodes defining the final tree are typically referred to as the terminal nodes of the tree.

As the above description implies, the crucial step in the splitting process is the calculation of the heterogeneity for a particular node. For the k^{th} node created in the splitting process this is the weighted sum of squared residuals

$$WSSR_k = \sum_{i \in k} w_{ik} (y_i - \overline{y}_{wk})^2$$

where $i \in k$ denotes the cases making up the node, w_{ik} is the weight attached to i^{th} case in the node and \bar{y}_{wk} is the weighted mean of Y in the node,

$$\overline{y}_{wk} = \sum\nolimits_{i \in k} w_{ik} y_i / \sum\nolimits_{i \in k} w_{ik} \,.$$

The weight w_{ik} is calculated as the ratio

$$w_{ik} = \frac{\psi(y_i - \overline{y}_{wk})}{y_i - \overline{y}_{wk}}$$

where $\psi(x)$ denotes an appropriately chosen influence function. The S-PLUS/R toolkit version of WAID uses weights returned by the robust regression function rlm in the MASS robust statistics library (Venables and Ripley, 1994). These weights are rescaled within WAID to sum to the number n_k of cases within node k.

3.2 The WAID Regression Tree Algorithm For Multivariate Y

Unlike other regression tree algorithms, the WAID toolkit can also build a regression tree for a p-dimensional response variable Y, and so can be used for multivariate outlier detection. The only difference between the univariate and multivariate tree fitting procedures is the method used to calculate the heterogeneity of a candidate node. Three options are available in this regard. In what follows tree "stages" are indexed by k (k = 1 corresponds to the original data set and k = K denotes the final stage of the tree), and the candidate nodes for splitting at stage k are indexed by k.

Option 1: The program first grows p univariate trees, one for each component of the response vector. Each such tree is characterised by an $n \times K$ matrix of weights, where column k of this matrix contains the weights used to determine node heterogeneity for all nodes defined at stage k of the tree growing procedure.

Let $w_{ij}^{(hk)}$ denote the weight associated with case i in node h at stage k of the univariate tree defined by response variable j. WAID then builds a tree using the heterogeneity measure for candidate node h at stage k of the multivariate tree:

$$WRSS_{hk} = \sum_{i \in h} \sum_{j=1}^{p} w_{ij}^{(hk)} (y_{ij} - \overline{y}_{whj}^{(k)})^2$$

where

$$\overline{y}_{whj}^{(k)} = \sum_{i \in h} w_{ij}^{(hk)} y_{ij} / \sum_{i \in h} w_{ij}^{(hk)}.$$

We can think of this as an "average heterogeneity" approach. Note that it is not scale invariant – a component response variable that is much larger in scale than the other component response variables will dominate this heterogeneity measure and hence dominate the tree growing process. Consequently component variables that differ wildly in terms of scale should be first rescaled before this option is used to build a multivariate tree.

Option 2: Here again WAID grows p univariate trees to obtain the weights $w_{ij}^{(hk)}$. However, in this case the measure of heterogeneity for candidate node h at stage k in the multivariate tree is

$$WRSS_k = \sum_{i \in h} w_i^{(hk)} \sum_{i=1}^p (y_{ij} - \bar{y}_{whj}^{(k)})^2$$

where

$$\overline{y}_{whj}^{(k)} = \sum\nolimits_{i \in h} {w_i^{(hk)} y_{ij} / \sum\nolimits_{i \in h} {w_i^{(hk)}}}$$

and

$$w_i^{(hk)} = p^{-1} \sum_{j=1}^p w_{ij}^{(hk)}.$$

We can think of this approach as an "average weight" approach. It also is not scale invariant.

Option 3: This is the only truly multivariate tree growing option in WAID. The weight associated with observation i in candidate node h at stage k is calculated iteratively as

$$w_{i}^{(hk)} = \frac{\psi(\left\|\mathbf{y}_{i} - \overline{\mathbf{y}}_{wh}^{(k)}\right\|_{wh})}{\left\|\mathbf{y}_{i} - \overline{\mathbf{y}}_{wh}^{(k)}\right\|_{wh}}$$

where \mathbf{y}_i denotes the *p*-vector of response values for this case, $\overline{\mathbf{y}}_{wh}^{(k)} = \sum_{i \in h} w_i^{(hk)} \mathbf{y}_i / \sum_{i \in h} w_i^{(hk)}$, the function ψ corresponds to an influence function and $\|\mathbf{y}_i - \overline{\mathbf{y}}_{wh}^{(k)}\|_{wh} = \sqrt{\sum_{j=1}^p s_{whj}^{-2} (y_{ij} - \overline{y}_{whj}^{(k)})^2}$,

where
$$s_{whj}^2 = \frac{\sum_{i \in h} w_i^{(hk)} (y_{ij} - \overline{y}_{whj}^{(k)})^2}{\sum_{i \in h} w_i^{(hk)}}$$
.

3.3 Outlier Identification Using WAID

Each time WAID splits the data set to create two new nodes it creates a new set of weights for the cases making up those nodes. When these weights are based on a robust influence function, outliers within the node have weights close to zero and non-outliers have weights around one. These weights reflect distance from a robust estimate of location for the values in the node. Consequently a value that is not immediately identifiable as an outlier within larger nodes created earlier in the tree building process is more likely to become identified as such as it is classified into smaller and smaller nodes. In effect, the weights associated with such cases tend to move towards zero. Conversely, extreme points in the covariate space with corresponding extreme *Y* values are initially given small weights. However, such points are quickly isolated into terminal nodes in the tree splitting process, at which point the weights associated with these points increase back to values near one.

The WAID outlier identification algorithm defines an outlier as a case with an average weight over all node splits less than a specified threshold. An optimal threshold value w^* is

one that successfully identifies outliers due to errors while minimising identification of "true" (i.e. representative) outliers. Let N_{errors} equal to the total number of true errors in the data, and, for a given threshold w, put $N_{outliers}(w)$ equal to the total number of outliers identified by WAID on the basis of the specified threshold w, $n_{errors}(w)$ equal to the corresponding number of errors identified as outliers, and $n_{non-errors}(w)$ equal to the total number of non-errors identified as outliers. The proportion of error-generated outliers identified by WAID is

$$R_{\rm I}(w) = \frac{n_{errors}(w)}{N_{errors}}$$

while the proportion of non-errors identified as outliers using the threshold w is

$$R_2(w) = \frac{n_{non-errors}(w)}{N_{outliers}(w)} = 1 - \frac{n_{errors}(w)}{N_{outliers}(w)}.$$

The optimal threshold value is then

$$w^* = \arg \max_{w} [R_1(w)(1 - R_2(w))].$$

4. Identifying Errors And Outliers In The Perturbed Data

4.1 Error Detection Using Forward Search

Table 2 shows the incidence of errors and missing values for each of the *ABI* variables in the *perturbed data*. It also shows the incidence of "n/a" codes for these variables (indicating no response was required for that variable for a sampled business) and the incidence of zero values. Note the large number of zero values for the *assdisp* and *assacq* variables.

Table 2. Incidence of incorrect and non-standard data types in the perturbed data.

Data type	Variables										
	turnover	emptotc	purtot	taxtot	assacq	assdisp	employ				
n/a	0	0	0	0	908	1389	0				
missing	42	41	28	45	57	63	35				
errors	241	332	629	482	248	223	49				
zero	3	658	5	390	2106	3208	721				

We first applied the forward search algorithm described in Section 2 to these data treating each variable separately (i.e. a univariate forward search). In each case we fitted a linear model in the logarithm of the variable concerned, using the logarithm of turnreg as the covariate. Two types of model were investigated. The first (across stratum) was fitted using all cases in the data set. The second (stratum level) fitted a separate linear model within each sampling stratum in the data set. Cases with zero, n/a or missing values were excluded from the outlier search procedure. As described in Section 2, the initial subset for the forward search procedure was defined by the smallest absolute residuals from a robust regression fit to the entire data set. This regression fit was based on the bisquare influence function $\psi(t) = t(1 - \min(1, t^2/c^2))^2$, with c = 4.685. The size of this initial data was set at 70 per cent of the size of the overall data set, since smaller initial data sets greatly increased the execution time of the algorithm and led to no change in the set of identified outliers. The stopping rule suggested by Hadi and Simonoff (1993) was used, with $\alpha = 0.01$. Table 3 shows the results from this univariate outlier search. On the basis of these results the across stratum specification of the regression model used in the forward search seems preferable for the ABI data.

Table 3. Numbers of outliers detected (with numbers of errors detected in parentheses) using univariate forward search applied to the *perturbed data*.

Model type				Variables			
	turnover	emptotc	purtot	taxtot	assacq	assdisp	employ
Across stratum	349	226	361	224	8	14	11
	(225)	(219)	(281)	(219)	(8)	(14)	(5)
Stratum level	467	279	441	245	102	56	77
	(227)	(237)	(294)	(235)	(90)	(52)	(24)

A multivariate forward search was also carried out. In this case we restricted attention to the 5118 cases where *turnover*, *emptotc*, *purtot*, *taxtot* and *employ* were all strictly positive. We excluded the variables *assacq* and *assdisp* from consideration because the large number of zero values for these variables meant that only a small number of cases had strictly positive

values for all seven variables. In the columns labelled (a) in Table 4 we show the number of cases with errors detected by this method distributed according to the number of errors in each case. We see that out of a total of 824 cases with one or more errors, the multivariate forward search detected 385. It also identified 96 cases with no errors as outliers. If we restrict attention to those cases with "significant" errors in one or more of *turnover*, *emptotc*, *purtot*, *taxtot* and *employ*, i.e. those cases where the perturbed values of these variables differ by more than 100 per cent from their values in the *clean data*, then we obtain the results shown in the columns labelled (b) in Table 4. In this case the multivariate forward search procedure finds 336 out of the 541 cases with at least one "significant" error.

Table 4. Error detection performance of multivariate forward search procedure (across stratum model) applied to *perturbed data*. Numbers in columns labelled (a) refer to all cases, while numbers in columns labelled (b) refer to cases with "significant" errors.

Number of errors per record	Total number of records		Records declared as outliers		
	(a)	(b)	(a)	(b)	
0	4294	4577	96	145	
1	489	374	194	172	
2	165	8	31	5	
3	13	2	3	2	
4	154	155	154	155	
5	3	2	3	2	
Total	5118	5118	481	481	

In Table 5 we contrast the performance of the multivariate forward search with that of the individual univariate forward search procedures. In both cases the across stratum version of the model was used in the forward search procedure. Here we see that 36 records were identified as outliers by the multivariate search and were not identified as such by any of the univariate searches. Furthermore only 14 records were identified by one of the univariate searches as containing an outlier and were not identified as such by the multivariate search. All records containing two or more outliers (as identified by the univariate searches) were also identified as outliers by the multivariate search.

Table 5. Comparing the outlier detection performances of the univariate and multivariate forward searches (across stratum model) applied to the *perturbed data*.

Number of outliers	Total number of	Number of records	Number of records
detected by	records	not identified by	identified by
univariate search		multivariate search	multivariate search
0	4659	4623	36
1	254	14	240
2	39	0	39
3	16	0	16
4	146	0	146
5	4	0	4
Total	5118	4637	481

Given that the multivariate forward search is restricted to cases where all components are non-zero, and given the lack of a strong differentiation in the performance of these two methods with the *perturbed data*, we restrict attention to comparisons with the univariate forward search method in what follows.

4.2 Error Detection Using WAID

Initially we focus on a univariate approach, building robust trees for the individual variables in Table 1. To save space we provide results below only for turnover, emptotc and assacq, since these are representative of the results obtained for the other variables. In all cases we built trees for the logarithm of the variable value. As in the previous section, cases with zero, n/a or missing values were excluded from the tree building process. All trees used the register variable turnreg, categorised into its percentile classes, as the covariate. This covariate was treated as monotone in the tree building process and each tree was built using the default option of robust splitting based on the bisquare influence function (c = 4.685). All trees were grown to 50 terminal nodes, with no node containing less than 5 cases.

Comparison of the trees grown using the *clean data* and the *perturbed data* showed that they were virtually identical, indicating that the errors in the *perturbed data* has virtually

no impact on the tree growing process. Furthermore, these trees were substantially different from corresponding non-robust trees ($\psi(t) = t$) grown from these data sets.

Figure 3 shows the plots of $R_1(w)$ and $R_1(w)(1-R_2(w))$ generated by these trees for the variables *turnover*, *emptotc* and *assacq*. We see that for all three variables, $R_1(w)(1-R_2(w))$ attains a maximum early, then falls away steadily. This behaviour is reflected in the plots for $R_1(w)$, which shows that most of the errors in these variables are detected using small values of w, with few non-errors detected at the same time. As w increases the remaining errors are then gradually detected, at the cost of identifying increasing numbers of non-errors as outliers, evidenced by the increasing separation of $R_1(w)$ and $R_1(w)(1-R_2(w))$.

Figure 3 about here

In most cases, errors detected at larger values of w are "non-significant", reflecting small differences from corresponding true values. As in the previous section, we define a "significant" error as one where the relative difference between the *perturbed data* value and the *clean data* value is greater than one. Table 6 shows the values of $R_1(w)$ (denoted R_{sig}) when only significant errors are taken into account. Observe that for all three variables over 80 per cent of such errors are identified at the optimal value w^* .

Table 6 also contrasts the performance of the WAID procedure with the corresponding forward search procedures. Here we see that for *turnover* this approach does identify all the significant errors in the data, but at the cost of identifying many more outliers than the WAID-based procedure. Both approaches seem comparable for *emptotc*. However, for *assacq* the WAID-based procedure is clearly superior in terms of identifying significant errors, while at the same time keeping the number of non-errors identified as outliers to an acceptable level.

Table 6. Univariate WAID error detection performance (perturbed data) for turnover, emptotc and assacq compared with the corresponding forward search (FS) performance. Note that WAID results are for $w = w^*$.

Method	Outliers	Errors	Significant	R_1	R_{sig}	R_2	$R_1(1-R_2)$
1.1011100	detected	detected	errors		_		
	(N_{out})	(N_{error})	detected				
			(N_{sig})				
			turnover				
WAID	194	191	190	0.793	0.913	0.015	0.780
FS/Across	349	225	206	0.941	1.000	0.355	0.607
FS/Stratum	467	227	206	0.950	1.000	0.514	0.462
			emptotc				
WAID	249	236	234	0.711	0.876	0.052	0.674
FS/Across	226	219	218	0.664	0.823	0.031	0.643
FS/Stratum	279	237	229	0.718	0.864	0.151	0.577
			assacq				
WAID	210	195	195	0.799	0.837	0.071	0.742
FS/Across	8	8	8	0.033	0.035	0	0.033
FS/Stratum	102	90	90	0.372	0.390	0.118	0.328

The trees used in the preceding analysis all had 50 terminal nodes. A natural question to ask at this stage is therefore about the impact of tree size (as measured by numbers of terminal nodes) on outlier identification. Rather surprisingly, at least as far as the *perturbed data* is concerned, tree size turns out to have little impact. Table 7 shows the performance characteristics of trees of varying size for *turnover* and *assacq*. Provided a tree has 10 or more terminal nodes, there is little to be gained by increasing the size of the tree.

Table 7. Impact of tree size (numbers of terminal nodes) on outlier and error detection performance with the *perturbed data*. In all cases the optimal value *w** was used.

Size	W_{opt}	N_{out}	N_{error}	N_{sig}	R_1	R_{sig}	R_2	$R_1(1-R_2)$
				turnover				
5	0.00001	206	189	188	0.784	0.904	0.083	0.720
10	0.00161	195	191	190	0.793	0.913	0.021	0.776
25	0.00075	194	191	190	0.793	0.913	0.015	0.780
50	0.00053	194	191	190	0.793	0.913	0.015	0.780
100	0.00053	194	191	190	0.793	0.913	0.015	0.780
				assacq				
5	0.03600	213	187	187	0.766	0.803	0.122	0.673
10	0.05009	226	195	195	0.799	0.837	0.137	0.690
25	0.04525	211	195	195	0.799	0.837	0.076	0.739
50	0.04464	210	195	195	0.799	0.837	0.071	0.742
100	0.03705	209	194	194	0.795	0.833	0.072	0.738

So far we have investigated the performance of univariate trees for outlier and error detection. We now consider the use of a multivariate tree for the same purpose. As with the multivariate forward search procedure, we restrict attention to the five *ABI* variables *turnover*, *emptotc*, *purtot*, *taxtot* and *employ* where excessive zero values are not a problem. Again, all data values were transformed to the logarithmic scale.

Recollect that there are three options available for growing a multivariate tree, corresponding to the way the "multivariate heterogeneity" associated with a particular tree split is defined (see Section 3.2). In Figure 4 we show how the error detection performance for the *turnover* variable for the three different trees generated by these options varies with w. This clearly shows that option 1 (average heterogeneity) is the preferable approach for this variable.

Figure 4 about here

Table 8. Error detection performance for a multivariate tree based on the *perturbed data*. Results for the three multivariate options, with $w = w^*$, are set out below each other for each variable. Note that for option 1 (first line for each variable), the optimal value w^* is the optimal univariate value.

w*	N_{out}	N_{error}	N_{sig}	R_1	R_{sig}	R_2	$R_1(1-R_2)$					
	turnover											
0.0199	165	160	159	0.812	0.914	0.030	0.788					
0.2593	177	119	119	0.604	0.684	0.328	0.406					
0.0076	163	157	157	0.797	0.902	0.037	0.768					
			purto	ot								
0.0326	214	204	197	0.394	0.925	0.047	0.375					
0.4456	339	179	151	0.346	0.709	0.472	0.183					
0.1062	254	198	193	0.382	0.906	0.221	0.298					
			taxto	ot .								
0.1085	334	275	273	0.679	0.7358	0.177	0.559					
0.2593	177	121	121	0.299	0.3261	0.316	0.204					
0.0672	196	182	182	0.449	0.4906	0.071	0.417					
			empto	otc								
0.0571	209	200	199	0.702	0.901	0.043	0.672					
0.2593	177	119	119	0.418	0.539	0.328	0.281					
0.0076	163	158	158	0.554	0.715	0.030	0.537					
	employ											
0.1075	87	25	24	0.595	0.857	0.713	0.171					
0.0178	8	2	2	0.048	0.071	0.750	0.012					
0.4455	656	30	24	0.714	0.857	0.954	0.033					

In Table 8 we show the error detection performance diagnostics for each of the five component variables and for each option at the optimal value of w thus defined. This confirms our observation above that option 1 (average heterogeneity) is preferable to the two other multivariate options. However, comparing the results in Table 8 with those in Table 6, we see that none of the multivariate trees perform significantly better than the corresponding univariate trees. This is consistent with our earlier observation about the lack of a similar improvement using the forward search approach, and indicates that, for the *perturbed data* at least, virtually all the errors are in low dimensions and so are easily detected via a univariate outlier search procedure.

5. Discussion

In this paper we compare two approaches to identification of gross errors and outliers in survey data. The first uses a forward search procedure based on a parametric model for the data. The second is based on a robust nonparametric modelling approach, based on regression tree methodology. Software (WAID) for implementing this second approach is described, and both approaches are evaluated using a realistic business survey data set created to reflect the errors that are typically found in such data. Overall, the regression tree method performs rather well and, for the application we consider, seems preferable to the forward search procedure. Multivariate versions of both approaches were also evaluated. Unfortunately, the data set we used for the application did not appear to have a significant number of truly multivariate outliers, and so we are unable at this stage to recommend the multivariate version of the regression tree procedure.

The regression tree approach is dependent on choice of a tuning constant, corresponding to the optimal weight cutoff w^* . In the *Euredit* project the value of this constant is determined using a "completely edited" data set obtained from a previous survey.

This assumes that the value of w^* remains the same over time. This is unlikely to be the case. Further work therefore needs to be carried out to determine an optimum updating strategy for this parameter.

The problem of dealing with "special" values (e.g. zero) when carrying out error detection remains an open problem. In this paper we have implicitly assumed these values are recorded without error. However, this will typically not be true. Alternative models need to be constructed to predict when an observed special value is an error. Some experience in using these models in the *Euredit* project seems to show that if such errors are randomly distributed over the survey data base, then there is little that can be done to automatically identify them.

Finally, we note that this paper does not address the issue of what should be done about errors and outliers once they are identified. Automatic methods for the imputation of these values are a focus of the *Euredit* project. Here we just note that both the forward search and robust regression tree methods for outlier and error detection imply methods for correcting these detected values. In the forward search case this is through the use of fitted values generated from the final "clean" subset of the data. In the regression tree context this is through the use of within node imputation, using either the robust estimate of the node mean or by making a random draw from the non-outlier cases in the node (i.e. those cases with average weights greater than the weight cut-off value).

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Figure 1. Plot of the *clean data*, log scale. Starting from top left, variables are *turnover*, *emptotc*, *purtot*, *taxtot*, *assacq*, *assdisp* and *employ*. In all cases *x*-axis is *turnreg*.

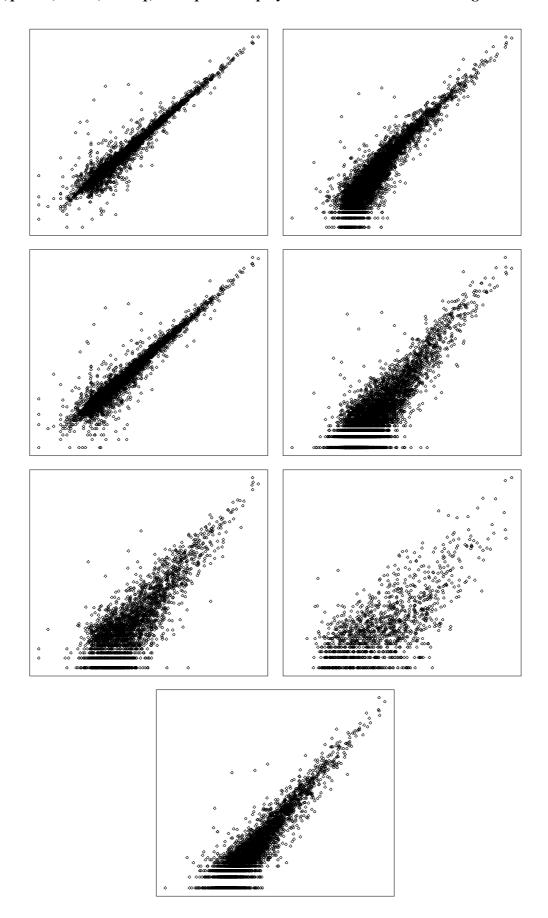


Figure 2. Plot of the *perturbed data*, log scale. " Δ " indicates an introduced error value. Starting from top left, variables are *turnover*, *emptotc*, *purtot*, *taxtot*, *assacq*, *assdisp* and *employ*. In all cases *x*-axis is *turnreg*.

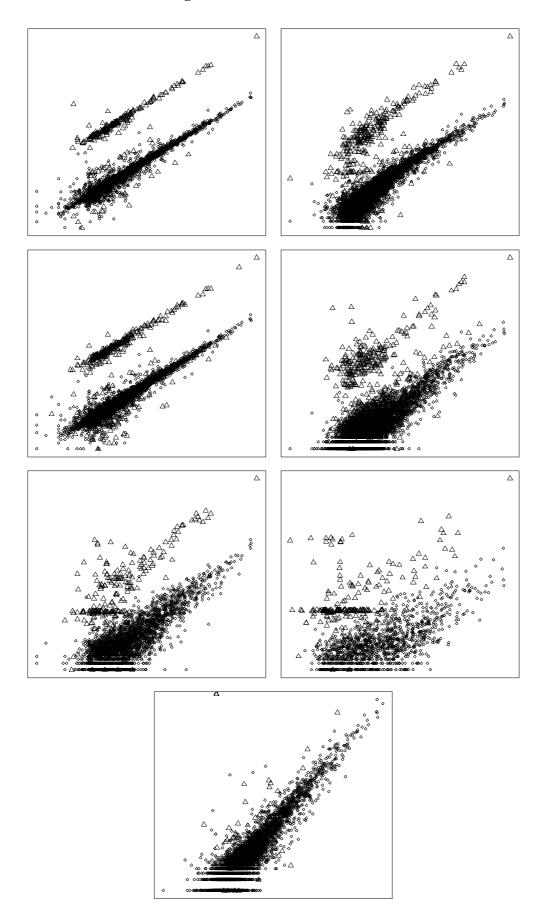


Figure 3. Plot of $R_1(w)$ and $R_1(w)(1-R_2(w))$ for univariate WAID trees. The x-axis is the value of w. The dashed line is $R_1(w)$. The solid line is $R_1(w)(1-R_2(w))$.

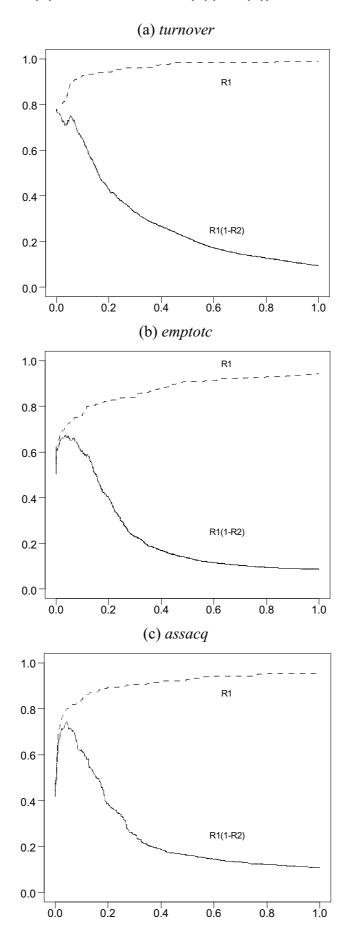


Figure 4. Plot of $R_1(w)$ and $R_1(w)(1-R_2(w))$ for multivariate WAID trees for *turnover*. The x-axis is the value of w. The dashed line is $R_1(w)$. The solid line is $R_1(w)(1-R_2(w))$.

