

NAG Library Function Document

nag_regsn_quant_linear (g02qgc)

Note: this function uses **optional parameters** to define choices in the problem specification and in the details of the algorithm. If you wish to use default settings for all of the optional parameters, you need only read Sections 1 to 10 of this document. If, however, you wish to reset some or all of the settings please refer to Section 11 for a detailed description of the algorithm, to Section 12 for a detailed description of the specification of the optional parameters and to Section 13 for a detailed description of the monitoring information produced by the function.

1 Purpose

nag_regsn_quant_linear (g02qgc) performs a multiple linear quantile regression. Parameter estimates and, if required, confidence limits, covariance matrices and residuals are calculated. nag_regsn_quant_linear (g02qgc) may be used to perform a weighted quantile regression. A simplified interface for nag_regsn_quant_linear (g02qgc) is provided by nag_regsn_quant_linear_iid (g02qfc).

2 Specification

```
#include <nag.h>
#include <nagg02.h>

void nag_regsn_quant_linear (Nag_OrderType order,
    Nag_IncludeIntercept intcpt, Integer n, Integer m, const double dat[],
    Integer pddat, const Integer isx[], Integer ip, const double y[],
    const double wt[], Integer ntau, const double tau[], double *df,
    double b[], double bl[], double bu[], double ch[], double res[],
    const Integer iopts[], const double opts[], Integer state[],
    Integer info[], NagError *fail)
```

3 Description

Given a vector of n observed values, $y = \{y_i : i = 1, 2, \dots, n\}$, an $n \times p$ design matrix X , a column vector, x , of length p holding the i th row of X and a quantile $\tau \in (0, 1)$, nag_regsn_quant_linear (g02qgc) estimates the p -element vector β as the solution to

$$\underset{\beta \in \mathbb{R}^p}{\text{minimize}} \sum_{i=1}^n \rho_{\tau}(y_i - x_i^T \beta) \quad (1)$$

where ρ_{τ} is the piecewise linear loss function $\rho_{\tau}(z) = z(\tau - I(z < 0))$, and $I(z < 0)$ is an indicator function taking the value 1 if $z < 0$ and 0 otherwise. Weights can be incorporated by replacing X and y with WX and Wy respectively, where W is an $n \times n$ diagonal matrix. Observations with zero weights can either be included or excluded from the analysis; this is in contrast to least squares regression where such observations do not contribute to the objective function and are therefore always dropped.

nag_regsn_quant_linear (g02qgc) uses the interior point algorithm of Portnoy and Koenker (1997), described briefly in Section 11, to obtain the parameter estimates $\hat{\beta}$, for a given value of τ .

Under the assumption of Normally distributed errors, Koenker (2005) shows that the limiting covariance matrix of $\hat{\beta} - \beta$ has the form

$$\Sigma = \frac{\tau(1-\tau)}{n} H_n^{-1} J_n H_n^{-1}$$

where $J_n = n^{-1} \sum_{i=1}^n x_i x_i^T$ and H_n is a function of τ , as described below. Given an estimate of the covariance matrix, $\hat{\Sigma}$, lower ($\hat{\beta}_L$) and upper ($\hat{\beta}_U$) limits for an $(100 \times \alpha)\%$ confidence interval can be calculated for each of the p parameters, via

$$\hat{\beta}_{Li} = \hat{\beta}_i - t_{n-p,(1+\alpha)/2} \sqrt{\hat{\Sigma}_{ii}}, \hat{\beta}_{Ui} = \hat{\beta}_i + t_{n-p,(1+\alpha)/2} \sqrt{\hat{\Sigma}_{ii}}$$

where $t_{n-p,0.975}$ is the 97.5 percentile of the Student's t distribution with $n - k$ degrees of freedom, where k is the rank of the cross-product matrix $X^T X$.

Four methods for estimating the covariance matrix, Σ , are available:

(i) Independent, identically distributed (IID) errors

Under an assumption of IID errors the asymptotic relationship for Σ simplifies to

$$\Sigma = \frac{\tau(1-\tau)}{n} (s(\tau))^2 (X^T X)^{-1}$$

where s is the sparsity function. `nag_regsn_quant_linear` (g02qgc) estimates $s(\tau)$ from the residuals, $r_i = y_i - x_i^T \hat{\beta}$ and a bandwidth h_n .

(ii) Powell Sandwich

Powell (1991) suggested estimating the matrix H_n by a kernel estimator of the form

$$\hat{H}_n = (nc_n)^{-1} \sum_{i=1}^n K\left(\frac{r_i}{c_n}\right) x_i x_i^T$$

where K is a kernel function and c_n satisfies $\lim_{n \rightarrow \infty} c_n \rightarrow 0$ and $\lim_{n \rightarrow \infty} \sqrt{nc_n} \rightarrow \infty$. When the Powell method is chosen, `nag_regsn_quant_linear` (g02qgc) uses a Gaussian kernel (i.e., $K = \phi$) and sets

$$c_n = \min(\sigma_r, (q_{r3} - q_{r1})/1.34) \times (\Phi^{-1}(\tau + h_n) - \Phi^{-1}(\tau - h_n))$$

where h_n is a bandwidth, σ_r , q_{r1} and q_{r3} are, respectively, the standard deviation and the 25% and 75% quantiles for the residuals, r_i .

(iii) Hendricks–Koenker Sandwich

Koenker (2005) suggested estimating the matrix H_n using

$$\hat{H}_n = n^{-1} \sum_{i=1}^n \left[\frac{2h_n}{x_i^T (\hat{\beta}(\tau + h_n) - \hat{\beta}(\tau - h_n))} \right] x_i x_i^T$$

where h_n is a bandwidth and $\hat{\beta}(\tau + h_n)$ denotes the parameter estimates obtained from a quantile regression using the $(\tau + h_n)$ th quantile. Similarly with $\hat{\beta}(\tau - h_n)$.

(iv) Bootstrap

The last method uses bootstrapping to either estimate a covariance matrix or obtain confidence intervals for the parameter estimates directly. This method therefore does not assume Normally distributed errors. Samples of size n are taken from the paired data $\{y_i, x_i\}$ (i.e., the independent and dependent variables are sampled together). A quantile regression is then fitted to each sample resulting in a series of bootstrap estimates for the model parameters, β . A covariance matrix can then be calculated directly from this series of values. Alternatively, confidence limits, $\hat{\beta}_L$ and $\hat{\beta}_U$, can be obtained directly from the $(1 - \alpha)/2$ and $(1 + \alpha)/2$ sample quantiles of the bootstrap estimates.

Further details of the algorithms used to calculate the covariance matrices can be found in Section 11.

All three asymptotic estimates of the covariance matrix require a bandwidth, h_n . Two alternative methods for determining this are provided:

(i) Sheather–Hall

$$h_n = \left(\frac{1.5(\Phi^{-1}(\alpha_b)\phi(\Phi^{-1}(\tau)))^2}{n(2\Phi^{-1}(\tau) + 1)} \right)^{\frac{1}{5}}$$

for a user-supplied value α_b ,

(ii) Bofinger

$$h_n = \left(\frac{4.5(\phi(\Phi^{-1}(\tau)))^4}{n(2\Phi^{-1}(\tau) + 1)^2} \right)^{\frac{1}{5}}$$

nag_regsn_quant_linear (g02qgc) allows optional arguments to be supplied via the **iopts** and **opts** arrays (see Section 12 for details of the available options). If the default values for these optional arguments are sufficient then **iopts** and **opts** can be set to **NULL**, otherwise prior to calling nag_regsn_quant_linear (g02qgc) the optional parameter arrays, must be initialized by calling nag_g02_opt_set (g02zkc) with **optstr** set to **Initialize = g02qgc**. If bootstrap confidence limits are required (**Interval Method = BOOTSTRAP XY**) then one of the random number initialization functions nag_rand_init_repeatable (g05kfc) (for a repeatable analysis) or nag_rand_init_nonrepeatable (g05kge) (for an unrepeatable analysis) must also have been previously called.

4 References

Koenker R (2005) *Quantile Regression* Econometric Society Monographs, Cambridge University Press, New York

Mehrotra S (1992) On the implementation of a primal-dual interior point method *SIAM J. Optim.* **2** 575–601

Nocedal J and Wright S J (1999) *Numerical Optimization* Springer Series in Operations Research, Springer, New York

Portnoy S and Koenker R (1997) The Gaussian hare and the Laplacian tortoise: computability of squared-error versus absolute error estimators *Statistical Science* **4** 279–300

Powell J L (1991) Estimation of monotonic regression models under quantile restrictions *Nonparametric and Semiparametric Methods in Econometrics* Cambridge University Press, Cambridge

5 Arguments

1: **order** – Nag_OrderType *Input*

On entry: the **order** argument specifies the two-dimensional storage scheme being used, i.e., row-major ordering or column-major ordering. C language defined storage is specified by **order = Nag_RowMajor**. See Section 2.3.1.3 in How to Use the NAG Library and its Documentation for a more detailed explanation of the use of this argument.

Constraint: **order = Nag_RowMajor** or **Nag_ColMajor**.

2: **intcpt** – Nag_IncludeIntercept *Input*

On entry: indicates whether an intercept will be included in the model. The intercept is included by adding a column of ones as the first column in the design matrix, X .

intcpt = Nag_Intercept

An intercept will be included in the model.

intcpt = Nag_NoIntercept

An intercept will not be included in the model.

Constraint: **intcpt = Nag_NoIntercept** or **Nag_Intercept**.

- 3: **n** – Integer *Input*
On entry: the total number of observations in the dataset. If no weights are supplied, or no zero weights are supplied or observations with zero weights are included in the model then **n** = *n*. Otherwise **n** = *n* + the number of observations with zero weights.
Constraint: **n** ≥ 2.
- 4: **m** – Integer *Input*
On entry: *m*, the total number of variates in the dataset.
Constraint: **m** ≥ 0.
- 5: **dat**[*dim*] – const double *Input*
Note: the dimension, *dim*, of the array **dat** must be at least
pddat × **m** when **order** = Nag_ColMajor;
pddat × **n** when **order** = Nag_RowMajor.
Where **DAT**(*i*, *j*) appears in this document, it refers to the array element
dat[(*j* – 1) × **pddat** + *i* – 1] when **order** = Nag_ColMajor;
dat[(*i* – 1) × **pddat** + *j* – 1] when **order** = Nag_RowMajor.
On entry: the *i*th value for the *j*th variate, for *i* = 1, 2, ..., **n** and *j* = 1, 2, ..., **m**, must be supplied in **DAT**(*i*, *j*)
The design matrix *X* is constructed from **dat**, **isx** and **intcpt**.
- 6: **pddat** – Integer *Input*
On entry: the stride separating row or column elements (depending on the value of **order**) in the array **dat**.
Constraints:
if **order** = Nag_ColMajor, **pddat** ≥ **n**;
otherwise **pddat** ≥ **m**.
- 7: **isx**[**m**] – const Integer *Input*
On entry: indicates which independent variables are to be included in the model.
isx[*j* – 1] = 0
The *j*th variate, supplied in **dat**, is not included in the regression model.
isx[*j* – 1] = 1
The *j*th variate, supplied in **dat**, is included in the regression model.
Constraints:
isx[*j* – 1] = 0 or 1, for *j* = 1, 2, ..., **m**;
if **intcpt** = Nag_Intercept, exactly **ip** – 1 values of **isx** must be set to 1;
if **intcpt** = Nag_NoIntercept, exactly **ip** values of **isx** must be set to 1.
- 8: **ip** – Integer *Input*
On entry: *p*, the number of independent variables in the model, including the intercept, see **intcpt**, if present.
Constraints:
1 ≤ **ip** < **n**;
if **intcpt** = Nag_Intercept, 1 ≤ **ip** ≤ **m** + 1;
if **intcpt** = Nag_NoIntercept, 1 ≤ **ip** ≤ **m**.

- 9: **y[n]** – const double *Input*
On entry: y , the observations on the dependent variable.
- 10: **wt[n]** – const double *Input*
On entry: optionally, the diagonal elements of the weight matrix W .
 If weights are not provided then **wt** must be set to **NULL**.
 When
Drop Zero Weights = YES
 If **wt**[$i - 1$] = 0.0, the i th observation is not included in the model, in which case the effective number of observations, n , is the number of observations with nonzero weights.
 If **Return Residuals = YES**, the values of **res** will be set to zero for observations with zero weights.
Drop Zero Weights = NO
 All observations are included in the model and the effective number of observations is **n**, i. e., $n = \mathbf{n}$.
Constraints:
 the effective number of observations ≥ 2 ;
wt[i] = 0.0, for all i .
- 11: **ntau** – Integer *Input*
On entry: the number of quantiles of interest.
Constraint: **ntau** ≥ 1 .
- 12: **tau[ntau]** – const double *Input*
On entry: the vector of quantiles of interest. A separate model is fitted to each quantile.
Constraint: $\sqrt{\epsilon} < \mathbf{tau}[j - 1] < 1 - \sqrt{\epsilon}$ where ϵ is the *machine precision* returned by `nag_machine_precision` (X02AJC), for $j = 1, 2, \dots, \mathbf{ntau}$.
- 13: **df** – double * *Output*
On exit: the degrees of freedom given by $n - k$, where n is the effective number of observations and k is the rank of the cross-product matrix $X^T X$.
- 14: **b[ip × ntau]** – double *Input/Output*
Note: where **B**(i, l) appears in this document, it refers to the array element **b**[($l - 1$) × **ip** + $i - 1$].
On entry: if **Calculate Initial Values = NO**, **B**(i, l) must hold an initial estimates for $\hat{\beta}_i$, for $i = 1, 2, \dots, \mathbf{ip}$ and $l = 1, 2, \dots, \mathbf{ntau}$. If **Calculate Initial Values = YES**, **b** need not be set.
On exit: **B**(i, l), for $i = 1, 2, \dots, \mathbf{ip}$, contains the estimates of the parameters of the regression model, $\hat{\beta}$, estimated for $\tau = \mathbf{tau}[l - 1]$.
 If **intcpt = Nag_Intercept**, **B**($1, l$) will contain the estimate corresponding to the intercept and **B**($i + 1, l$) will contain the coefficient of the j th variate contained in **dat**, where **isx**[$j - 1$] is the i th nonzero value in the array **isx**.
 If **intcpt = Nag_NoIntercept**, **B**(i, l) will contain the coefficient of the j th variate contained in **dat**, where **isx**[$j - 1$] is the i th nonzero value in the array **isx**.
- 15: **bl[dim]** – double *Output*
Note: the dimension, dim , of the array **bl** must be at least **ntau** when **Interval Method** \neq NONE.
 Where **BL**(i, l) appears in this document, it refers to the array element **bl**[($l - 1$) × **ip** + $i - 1$].

On exit: if **Interval Method** \neq NONE, **BL**(i, l) contains the lower limit of an $(100 \times \alpha)\%$ confidence interval for **B**(i, l), for $i = 1, 2, \dots, \mathbf{ip}$ and $l = 1, 2, \dots, \mathbf{ntau}$.

If **Interval Method** = NONE, **bl** is not referenced and can be set to **NULL**.

The method used for calculating the interval is controlled by the optional parameters **Interval Method** and **Bootstrap Interval Method**. The size of the interval, α , is controlled by the optional parameter **Significance Level**.

16: **bu**[dim] – double *Output*

Note: the dimension, dim , of the array **bu** must be at least **ntau** when **Interval Method** \neq NONE.

Where **BU**(i, l) appears in this document, it refers to the array element **bu**[($l - 1$) \times **ip** + $i - 1$].

On exit: if **Interval Method** \neq NONE, **BU**(i, l) contains the upper limit of an $(100 \times \alpha)\%$ confidence interval for **B**(i, l), for $i = 1, 2, \dots, \mathbf{ip}$ and $l = 1, 2, \dots, \mathbf{ntau}$.

If **Interval Method** = NONE, **bu** is not referenced and can be set to **NULL**.

The method used for calculating the interval is controlled by the optional parameters **Interval Method** and **Bootstrap Interval Method**. The size of the interval, α is controlled by the optional parameter **Significance Level**.

17: **ch**[dim] – double *Output*

Note: the dimension, dim , of the array **ch** must be at least

if **Interval Method** \neq NONE and **Matrix Returned** = COVARIANCE, **ip** \times **ip** \times **ntau**;
if **Interval Method** \neq NONE, IID or BOOTSTRAP XY and
Matrix Returned = H INVERSE, **ip** \times **ip** \times (**ntau** + 1).

Where **CH**(i, j, l) appears in this document, it refers to the array element **ch**[($l - 1$) \times **ip** \times **ip** + ($j - 1$) \times **ip** + $i - 1$].

On exit: depending on the supplied optional parameters, **ch** will either not be referenced, hold an estimate of the upper triangular part of the covariance matrix, Σ , or an estimate of the upper triangular parts of nJ_n and $n^{-1}H_n^{-1}$.

If **Interval Method** = NONE or **Matrix Returned** = NONE, **ch** is not referenced.

If **Interval Method** = BOOTSTRAP XY or IID and **Matrix Returned** = H INVERSE, **ch** is not referenced.

Otherwise, for $i, j = 1, 2, \dots, \mathbf{ip}$, $j \geq i$ and $l = 1, 2, \dots, \mathbf{ntau}$:

If **Matrix Returned** = COVARIANCE, **CH**(i, j, l) holds an estimate of the covariance between **B**(i, l) and **B**(j, l).

If **Matrix Returned** = H INVERSE, **CH**($i, j, 1$) holds an estimate of the (i, j)th element of nJ_n and **CH**($i, j, l + 1$) holds an estimate of the (i, j)th element of $n^{-1}H_n^{-1}$, for $\tau = \mathbf{tau}[l - 1]$.

The method used for calculating Σ and H_n^{-1} is controlled by the optional parameter **Interval Method**.

In cases where **ch** is not going to be referenced it can be set to **NULL**.

18: **res**[$n \times \mathbf{ntau}$] – double *Output*

Note: the (i, j)th element of the matrix is stored in **res**[($j - 1$) \times **n** + $i - 1$].

On exit: if **Return Residuals** = YES, **res**[($l - 1$) \times **n** + $i - 1$] holds the (weighted) residuals, r_i , for $\tau = \mathbf{tau}[l - 1]$, for $i = 1, 2, \dots, \mathbf{n}$ and $l = 1, 2, \dots, \mathbf{ntau}$.

If **wt** is not **NULL** and **Drop Zero Weights** = YES, the value of **res** will be set to zero for observations with zero weights.

If **Return Residuals** = NO, **res** is not referenced and can be set to **NULL**.

19: **iopts** $[dim]$ – const Integer *Communication Array*

Note: the dimension, dim , of this array is dictated by the requirements of associated functions that must have been previously called. This array **MUST** be the same array passed as argument **iopts** in the previous call to nag_g02_opt_set (g02zkc).

On entry: if the default values of the optional arguments are sufficient, then **iopts** can be set to **NULL**, otherwise the optional parameter array, as initialized by a call to nag_g02_opt_set (g02zkc) must be supplied.

20: **opts** $[dim]$ – const double *Communication Array*

Note: the dimension, dim , of this array is dictated by the requirements of associated functions that must have been previously called. This array **MUST** be the same array passed as argument **opts** in the previous call to nag_g02_opt_set (g02zkc).

On entry: if the default values of the optional arguments are sufficient, then **opts** can be set to **NULL**, otherwise the optional parameter array, as initialized by a call to nag_g02_opt_set (g02zkc) must be supplied.

21: **state** $[dim]$ – Integer *Communication Array*

Note: the dimension, dim , of this array is dictated by the requirements of associated functions that must have been previously called. This array **MUST** be the same array passed as argument **state** in the previous call to nag_rand_init_repeatable (g05kfc) or nag_rand_init_nonrepeatable (g05kge).

If **Interval Method** = BOOTSTRAP XY, **state** contains information about the selected random number generator. Otherwise **state** is not referenced and can be set to **NULL**.

22: **info** $[ntau]$ – Integer *Output*

On exit: **info** $[i]$ holds additional information concerning the model fitting and confidence limit calculations when $\tau = \mathbf{tau}[i]$.

Code Warning

0	Model fitted and confidence limits (if requested) calculated successfully
1	The function did not converge. The returned values are based on the estimate at the last iteration. Try increasing Iteration Limit whilst calculating the parameter estimates or relaxing the definition of convergence by increasing Tolerance .
2	A singular matrix was encountered during the optimization. The model was not fitted for this value of τ .
4	Some truncation occurred whilst calculating the confidence limits for this value of τ . See Section 11 for details. The returned upper and lower limits may be narrower than specified.
8	The function did not converge whilst calculating the confidence limits. The returned limits are based on the estimate at the last iteration. Try increasing Iteration Limit .
16	Confidence limits for this value of τ could not be calculated. The returned upper and lower limits are set to a large positive and large negative value respectively as defined by the optional parameter Big .

It is possible for multiple warnings to be applicable to a single model. In these cases the value returned in **info** is the sum of the corresponding individual nonzero warning codes.

23: **fail** – NagError * *Input/Output*

The NAG error argument (see Section 2.7 in How to Use the NAG Library and its Documentation).

6 Error Indicators and Warnings

NE_ALLOC_FAIL

Dynamic memory allocation failed.

See Section 2.3.1.2 in How to Use the NAG Library and its Documentation for further information.

NE_ARRAY_SIZE

On entry, **pddat** = $\langle value \rangle$ and **m** = $\langle value \rangle$.

Constraint: **pddat** \geq **m**.

On entry, **pddat** = $\langle value \rangle$ and **n** = $\langle value \rangle$.

Constraint: **pddat** \geq **n**.

NE_BAD_PARAM

On entry, argument $\langle value \rangle$ had an illegal value.

NE_INITIALIZATION

On entry, either the option arrays have not been initialized or they have been corrupted.

NE_INT

On entry, **m** = $\langle value \rangle$.

Constraint: **m** \geq 0.

On entry, **n** = $\langle value \rangle$.

Constraint: **n** \geq 2.

On entry, **ntau** = $\langle value \rangle$.

Constraint: **ntau** \geq 1.

NE_INT_2

On entry, **ip** = $\langle value \rangle$ and **n** = $\langle value \rangle$.

Constraint: $1 \leq \mathbf{ip} < \mathbf{n}$.

NE_INT_ARRAY

On entry, **isx**[$\langle value \rangle$] = $\langle value \rangle$.

Constraint: **isx**[i] = 0 or 1 for all i .

NE_INTERNAL_ERROR

An internal error has occurred in this function. Check the function call and any array sizes. If the call is correct then please contact NAG for assistance.

An unexpected error has been triggered by this function. Please contact NAG.

See Section 2.7.6 in How to Use the NAG Library and its Documentation for further information.

NE_INVALID_STATE

On entry, **state** vector has been corrupted or not initialized.

NE_IP_INCOMP_SX

On entry, **ip** is not consistent with **isx** or **intcpt**: **ip** = $\langle value \rangle$, expected value = $\langle value \rangle$.

NE_NEG_WEIGHT

On entry, **wt**[$\langle value \rangle$] = $\langle value \rangle$.

Constraint: **wt**[i] \geq 0.0 for all i .

NE_NO_LICENCE

Your licence key may have expired or may not have been installed correctly.
See Section 2.7.5 in How to Use the NAG Library and its Documentation for further information.

NE_OBSERVATIONS

On entry, effective number of observations = $\langle value \rangle$.
Constraint: effective number of observations $\geq \langle value \rangle$.

NE_REAL_ARRAY

On entry, $\mathbf{tau}[\langle value \rangle] = \langle value \rangle$ is invalid.

NW_POTENTIAL_PROBLEM

A potential problem occurred whilst fitting the model(s).
Additional information has been returned in **info**.

7 Accuracy

Not applicable.

8 Parallelism and Performance

`nag_regsn_quant_linear` (g02qgc) is threaded by NAG for parallel execution in multithreaded implementations of the NAG Library.

`nag_regsn_quant_linear` (g02qgc) makes calls to BLAS and/or LAPACK routines, which may be threaded within the vendor library used by this implementation. Consult the documentation for the vendor library for further information.

Please consult the x06 Chapter Introduction for information on how to control and interrogate the OpenMP environment used within this function. Please also consult the Users' Note for your implementation for any additional implementation-specific information.

9 Further Comments

`nag_regsn_quant_linear` (g02qgc) allocates internally approximately the following elements of double storage: $13n + np + 3p^2 + 6p + 3(p + 1) \times \mathbf{ntau}$. If **Interval Method** = BOOTSTRAP XY then a further np elements are required, and this increases by $p \times \mathbf{ntau} \times \mathbf{Bootstrap Iterations}$ if **Bootstrap Interval Method** = QUANTILE. Where possible, any user-supplied output arrays are used as workspace and so the amount actually allocated may be less. If **order** = Nag_RowMajor, **wt** is NULL, **intcpt** = Nag_NoIntercept and **ip** = **m** an internal copy of the input data is avoided and the amount of locally allocated memory is reduced by np .

10 Example

A quantile regression model is fitted to Engels 1857 study of household expenditure on food. The model regresses the dependent variable, household food expenditure, against two explanatory variables, a column of ones and household income. The model is fit for five different values of τ and the covariance matrix is estimated assuming Normal IID errors. Both the covariance matrix and the residuals are returned.

10.1 Program Text

```

/* nag_regsn_quant_linear (g02qgc) Example Program.
 *
 * NAGPRODCODE Version.
 *
 * Copyright 2016 Numerical Algorithms Group.
 *
 * Mark 26, 2016.
 */
/* Pre-processor includes */
#include <stdio.h>
#include <string.h>
#include <nag.h>
#include <nag_stdlib.h>
#include <nagg02.h>
#include <nagg05.h>
#include <nagx04.h>

#define DAT(i,j)    dat[(order==Nag_RowMajor) ? (i*pddat+j) : (j*pddat+i)]

#define LOPTSTR 80

int main(void)
{
    /* Integer scalar and array declarations */
    Integer lseed = 1, liopts = 100, lopts = 100, lcvalue = LOPTSTR;
    Integer exit_status = 0;
    Integer i, ip, ivalue, j, l, lc, lstate, loptstr,
           m, n, ntau, subid, tdch, pddat;
    Integer *info = 0, *iopts = 0, *isx = 0, *state = 0;
    Integer seed[1];
    Nag_BaseRNG genid;

    /* NAG structures */
    NagError fail;
    Nag_OrderType order;
    Nag_IncludeIntercept intcpt;
    Nag_Boolean weighted;
    Nag_VariableType optype;

    /* Double scalar and array declarations */
    double df, rvalue;
    double *b = 0, *bl = 0, *bu = 0, *ch = 0, *dat = 0,
           *opts = 0, *res = 0, *tau = 0, *wt = 0, *y = 0;

    /* Character scalar and array declarations */
    char semeth[30], *poptstr, *cvalue = 0;
    char optstr[LOPTSTR], corder[40], cintcpt[40], cweighted[40], cgenid[40];
    char *clabs = 0, **clabsc = 0;

    /* Initialize the error structure to print out any error messages */
    INIT_FAIL(fail);

    printf("nag_regsn_quant_linear (g02qgc) Example Program Results\n\n");
    fflush(stdout);

    /* Skip heading in data file */
#ifdef _WIN32
    scanf_s("%*[\n] ");
#else
    scanf("%*[\n] ");
#endif

    /* Read in the problem size */
#ifdef _WIN32
    scanf_s("%39s%*[\n]", corder, (unsigned)_countof(corder));
#else
    scanf("%39s%*[\n]", corder);
#endif
#endif

```

```

scanf_s("%39s%39s%*[\n]", cintcpt, (unsigned)_countof(cintcpt),
        cweighted, (unsigned)_countof(cweighted));
#else
scanf("%39s%39s%*[\n]", cintcpt, cweighted);
#endif
#ifdef _WIN32
scanf_s("%" NAG_IFMT "%" NAG_IFMT "%" NAG_IFMT "%*[\n]", &n, &m, &ntau);
#else
scanf("%" NAG_IFMT "%" NAG_IFMT "%" NAG_IFMT "%*[\n]", &n, &m, &ntau);
#endif
order = (Nag_OrderType) nag_enum_name_to_value(corder);
intcpt = (Nag_IncludeIntercept) nag_enum_name_to_value(cintcpt);
/* weighted is a Nag_Boolean flag used in this example program to indicate
 * whether weights are being supplied (weighted=Nag_TRUE)
 * or not (weighted=Nag_FALSE)
 */
weighted = (Nag_Boolean) nag_enum_name_to_value(cweighted);

pddat = (order == Nag_RowMajor) ? m : n;

/* Allocate memory for input arrays */
if (!(y = NAG_ALLOC(n, double)) ||
    !(tau = NAG_ALLOC(ntau, double)) ||
    !(isx = NAG_ALLOC(m, Integer)) ||
    !(dat = NAG_ALLOC(m * n, double)) ||
    !(cvalue = NAG_ALLOC(lcvalue, char)) ||
    !(clabs = NAG_ALLOC(10 * 10, char)) || !(clabsc = NAG_ALLOC(10, char *)))
{
printf("Allocation failure\n");
exit_status = -1;
goto END;
}

if (weighted) {
/* Data includes a weight */
if (!(wt = NAG_ALLOC(n, double)))
{
printf("Allocation failure\n");
exit_status = -1;
goto END;
}
for (i = 0; i < n; i++) {
#ifdef _WIN32
for (j = 0; j < m; j++)
scanf_s("%lf", &DAT(i, j));
#else
for (j = 0; j < m; j++)
scanf("%lf", &DAT(i, j));
#endif
scanf_s("%lf%lf", &y[i], &wt[i]);
#else
scanf("%lf%lf", &y[i], &wt[i]);
#endif
}
#ifdef _WIN32
scanf_s("%*[\n] ");
#else
scanf("%*[\n] ");
#endif
}
else {
/* No weights supplied */
for (i = 0; i < n; i++) {
#ifdef _WIN32
for (j = 0; j < m; j++)
scanf_s("%lf", &DAT(i, j));
#else
for (j = 0; j < m; j++)
scanf("%lf", &DAT(i, j));

```

```

#endif
#ifdef _WIN32
    scanf_s("%lf", &y[i]);
#else
    scanf("%lf", &y[i]);
#endif
    }
#ifdef _WIN32
    scanf_s("%*[\n] ");
#else
    scanf("%*[\n] ");
#endif
    }

    /* Read in variable inclusion flags and calculate IP */
    ip = (intcpt == Nag_Intercept) ? 1 : 0;
    for (j = 0; j < m; j++) {
#ifdef _WIN32
        scanf_s("%" NAG_IFMT, &isx[j]);
#else
        scanf("%" NAG_IFMT, &isx[j]);
#endif
        if (isx[j] == 1)
            ip++;
    }
#ifdef _WIN32
    scanf_s("%*[\n] ");
#else
    scanf("%*[\n] ");
#endif

    /* Read in the quantiles required */
#ifdef _WIN32
    for (l = 0; l < ntau; l++)
        scanf_s("%lf", &tau[l]);
#else
    for (l = 0; l < ntau; l++)
        scanf("%lf", &tau[l]);
#endif
#ifdef _WIN32
    scanf_s("%*[\n] ");
#else
    scanf("%*[\n] ");
#endif

    /* Allocate memory for option arrays */
    if (!(opts = NAG_ALLOC(lopts, double)) ||
        !(iopts = NAG_ALLOC(liopts, Integer)))
    {
        printf("Allocation failure\n");
        exit_status = -1;
        goto END;
    }

    /* Initialize the optional argument array with nag_g02_opt_set (g02zkc) */
    nag_g02_opt_set("INITIALIZE = G02QG", iopts, liopts, opts, lopts, &fail);
    if (fail.code != NE_NOERROR) {
        printf("Error from nag_g02_opt_set (g02zkc).\n%s\n", fail.message);
        exit_status = 1;
        goto END;
    }

    /* Read in any optional arguments. Reads in to the end of
       the input data, or until a blank line is reached */
    for (;;) {
        if (!fgets(optstr, LOPTSTR, stdin))
            break;

        /* Left justify the option */
        popostr = (optstr + strspn(optstr, " \n\t"));
        /* Get the string length */

```

```

    loptstr = strlen(poptstr);
    if (poptstr[loptstr - 1] == '\n') {
        /* Remove any trailing line breaks */
        poptrstr[loptstr - 1] = '\setminus 0';
    }
    else {
        /* Clear the rest of the line */
#ifdef _WIN32
        scanf_s("%*[\n] ");
#else
        scanf("%*[\n] ");
#endif
    }

    /* Break if read in a blank line */
    if (!*(poptstr))
        break;

    /* Set the supplied option (g02zkc) */
    nag_g02_opt_set(optstr, iopts, liopts, opts, lopts, &fail);
    if (fail.code != NE_NOERROR) {
        printf("Error from nag_g02_opt_set (g02zkc).\n%s\n", fail.message);
        exit_status = 1;
        goto END;
    }
}

/* Allocate memory for the output arrays */
if (!(b = NAG_ALLOC(ip * ntau, double)) ||
    !(info = NAG_ALLOC(ntau, Integer)))
{
    printf("Allocation failure\n");
    exit_status = -1;
    goto END;
}

/* Query optional arguments via nag_g02_opt_get (g02zlc) and calculate which
 * of the optional arrays are required and their sizes
 * ...
 */
nag_g02_opt_get("INTERVAL METHOD", &ivalue, &rvalue, cvalue, lvalue,
                &octype, iopts, opts, &fail);
if (fail.code != NE_NOERROR) {
    printf("Error from nag_g02_opt_get (g02zlc).\n%s\n", fail.message);
    exit_status = 1;
    goto END;
}
#ifdef _WIN32
    strcpy_s(semeth, (unsigned)_countof(semeth), cvalue);
#else
    strcpy(semeth, cvalue);
#endif
if (strcmp(semeth, "NONE") != 0) {
    /* Require the intervals to be output */
    if (!(b1 = NAG_ALLOC(ip * ntau, double)) ||
        !(bu = NAG_ALLOC(ip * ntau, double)))
    {
        printf("Allocation failure\n");
        exit_status = -1;
        goto END;
    }
}

/* Decide whether the state array is required, and initialize if it is */
if (strcmp(semeth, "BOOTSTRAP XY") == 0) {
    /* Read in the generator ID and a seed */
#ifdef _WIN32
    scanf_s("%39s% NAG_IFMT \"% NAG_IFMT \"%*[\n] ", cgenid,
            (unsigned)_countof(cgenid), &subid, &seed[0]);
#else
    scanf("%39s% NAG_IFMT \"% NAG_IFMT \"%*[\n] ", cgenid, &subid,
            &seed[0]);
#endif
}

```

```

#endif
    genid = (Nag_BaseRNG) nag_enum_name_to_value(cgenid);

    /* Query the length of the state array (g05kfc) */
    lstate = 0;
    nag_rand_init_repeatable(genid, subid, seed, lseed, state, &lstate,
                             &fail);
    if (fail.code != NE_NOERROR) {
        printf("Error from nag_rand_init_repeatable (g05kfc).\n%s\n",
              fail.message);
        exit_status = 1;
        goto END;
    }

    /* Allocate memory to state */
    if (!(state = NAG_ALLOC(lstate, Integer)))
    {
        printf("Allocation failure\n");
        exit_status = -1;
        goto END;
    }

    /* Initialize the RNG (g05kfc) */
    nag_rand_init_repeatable(genid, subid, seed, lseed, state, &lstate,
                             &fail);
    if (fail.code != NE_NOERROR) {
        printf("Error from nag_rand_init_repeatable (g05kfc).\n%s\n",
              fail.message);
        exit_status = 1;
        goto END;
    }
}

/* Calculate the size of the covariance matrix, ch. */
tdch = 0;
nag_g02_opt_get("MATRIX RETURNED", &ivalue, &rvalue, cvalue, lcvalue,
               &optype, iopts, opts, &fail);
if (fail.code != NE_NOERROR) {
    printf("Error from nag_g02_opt_get (g02zlc).\n%s\n", fail.message);
    exit_status = 1;
    goto END;
}

if (strcmp(cvalue, "COVARIANCE") == 0) {
    tdch = ntau;
}
else if (strcmp(cvalue, "H INVERSE") == 0) {
    /* NB: If we are using bootstrap or IID errors then any request for
       H INVERSE is ignored */
    if (strcmp(semeth, "BOOTSTRAP XY") != 0 && strcmp(semeth, "IID") != 0)
        tdch = ntau + 1;
}
if (tdch > 0) {
    /* Need to allocate ch */
    if (!(ch = NAG_ALLOC(ip * ip * tdch, double)))
    {
        printf("Allocation failure\n");
        exit_status = -1;
        goto END;
    }
}

/* Calculate the size of the residual array, res */
nag_g02_opt_get("RETURN RESIDUALS", &ivalue, &rvalue, cvalue, lcvalue,
               &optype, iopts, opts, &fail);
if (fail.code != NE_NOERROR) {
    printf("Error from nag_g02_opt_get (g02zlc).\n%s\n", fail.message);
    exit_status = 1;
    goto END;
}
}

```

```

if (strcmp(cvalue, "YES") == 0) {
    /* Need to allocate res */
    if (!(res = NAG_ALLOC(n * ntau, double)))
    {
        printf("Allocation failure\n");
        exit_status = -1;
        goto END;
    }
}
}
/* ...
* end of handling the optional arguments, and allocating optional arrays
*/

/* Call the model fitting routine (nag_regsn_quant_linear (g02qgc)) */
nag_regsn_quant_linear(order, intcpt, n, m, dat, pddat, isx, ip, y, wt,
    ntau, tau, &df, b, bl, bu, ch, res, iopts, opts,
    state, info, &fail);
if (fail.code != NE_NOERROR) {
    printf("Error from nag_regsn_quant_linear (g02qgc).\n%s\n", fail.message);
    if (fail.code == NW_POTENTIAL_PROBLEM) {
        printf("Additional error information: ");
        for (i = 0; i < ntau; i++)
            printf("%" NAG_IFMT " ", info[i]);
        printf("\n");
    }
    else {
        printf("Error from nag_regsn_quant_linear (g02qgc).\n%s\n",
            fail.message);
        exit_status = -1;
        goto END;
    }
}

/* Display the parameter estimates */
for (l = 0; l < ntau; l++) {
    printf(" Quantile: %6.3f\n\n", tau[l]);
    if (bl && bu) {
        printf("          Lower   Parameter   Upper\n");
        printf("          Limit   Estimate   Limit\n");
        for (j = 0; j < ip; j++)
            printf(" %3" NAG_IFMT "%10.3f%10.3f%10.3f\n", j + 1, bl[l * ip + j],
                b[l * ip + j], bu[l * ip + j]);
    }
    else {
        printf("          Parameter\n");
        printf("          Estimate\n");
        for (j = 0; j < ip; j++)
            printf(" %3" NAG_IFMT "%10.3f\n", j + 1, b[l * ip + j]);
    }
    printf("\n\n");
    fflush(stdout);
    if (ch) {
        lc = l * ip * ip;
        if (tdch == ntau) {
            /* nag_gen_real_mat_print_comp (x04cbc).
            * Print real general matrix (comprehensive).
            */
            nag_gen_real_mat_print_comp(Nag_ColMajor, Nag_UpperMatrix,
                Nag_NonUnitDiag, ip, ip, &sch[lc], ip,
                "%9.2e", "Covariance matrix",
                Nag_NoLabels, 0, Nag_NoLabels, 0, 80,
                0, 0, &fail);
        }
        else {
            if (l == 0) {
                nag_gen_real_mat_print_comp(Nag_ColMajor, Nag_UpperMatrix,
                    Nag_NonUnitDiag, ip, ip, ch, ip,
                    "%9.2e", "J", Nag_NoLabels, 0,
                    Nag_NoLabels, 0, 80, 0, 0, &fail);
            }
            printf("\n");
        }
    }
}

```

```

    }
    lc = lc + ip * ip;
    nag_gen_real_mat_print_comp(Nag_ColMajor, Nag_UpperMatrix,
                               Nag_NonUnitDiag, ip, ip, &sch[lc], ip,
                               "%9.2e", "H inverse",
                               Nag_NoLabels, 0, Nag_NoLabels, 0, 80,
                               0, 0, &fail);
}
if (fail.code != NE_NOERROR) {
    printf("Error from nag_gen_real_mat_print_comp (x04cbc).\n%s\n",
           fail.message);
    exit_status = 1;
    goto END;
}
printf("\n");
}
}

if (res) {
    printf(" First 10 Residuals\n");
    fflush(stdout);
    /* set up column labels for matrix printer */
#ifdef _WIN32
    for (l = 0; l < ntau; l++)
        sprintf_s(&clabs[10 * l], 10, "%6.3f", tau[l]);
#else
    for (l = 0; l < ntau; l++)
        sprintf(&clabs[10 * l], "%6.3f", tau[l]);
#endif
    for (l = 0; l < ntau; l++)
        clabsc[l] = &clabs[l * 10];
    /* nag_gen_real_mat_print_comp (x04cbc).
     * Print real general matrix (comprehensive).
     */
    nag_gen_real_mat_print_comp(Nag_ColMajor, Nag_GeneralMatrix,
                               Nag_NonUnitDiag, MIN(10, n), ntau, res, n,
                               "%10.5f", "
                               Quantile",
                               Nag_IntegerLabels, NULL, Nag_CharacterLabels,
                               (const char **) clabsc, 80, 2, NULL, &fail);
    if (fail.code != NE_NOERROR) {
        printf("Error from nag_gen_real_mat_print_comp (x04cbc).\n%s\n",
               fail.message);
        exit_status = 1;
        goto END;
    }
}
else {
    printf(" Residuals not returned\n");
}

END:

NAG_FREE(info);
NAG_FREE(iopts);
NAG_FREE(isx);
NAG_FREE(state);
NAG_FREE(b);
NAG_FREE(bl);
NAG_FREE(bu);
NAG_FREE(ch);
NAG_FREE(dat);
NAG_FREE(opts);
NAG_FREE(res);
NAG_FREE(tau);
NAG_FREE(wt);
NAG_FREE(y);
NAG_FREE(cvalue);

```



```

NAG_FREE(clabs);
NAG_FREE(clabsc);

return (exit_status);
}

```

10.2 Program Data

```

nag_regsn_quant_linear (g02qgc) Example Program Data
Nag_ColMajor                :: sorder
Nag_Intercept   Nag_FALSE  :: intcpt, weighted
235                1                5                :: n, m, ntau
420.1577  255.8394      800.7990  572.0807      643.3571  459.8177
541.4117  310.9587      1245.6964  907.3969      2551.6615  863.9199
901.1575  485.6800      1201.0002  811.5776      1795.3226  831.4407
639.0802  402.9974      634.4002  427.7975      1165.7734  534.7610
750.8756  495.5608      956.2315  649.9985      815.6212  392.0502
945.7989  633.7978      1148.6010  860.6002      1264.2066  934.9752
829.3979  630.7566      1768.8236  1143.4211     1095.4056  813.3081
979.1648  700.4409      2822.5330  2032.6792     447.4479  263.7100
1309.8789  830.9586      922.3548  590.6183      1178.9742  769.0838
1492.3987  815.3602     2293.1920  1570.3911     975.8023  630.5863
502.8390  338.0014      627.4726  483.4800      1017.8522  645.9874
616.7168  412.3613      889.9809  600.4804      423.8798  319.5584
790.9225  520.0006     1162.2000  696.2021     558.7767  348.4518
555.8786  452.4015     1197.0794  774.7962     943.2487  614.5068
713.4412  512.7201     530.7972  390.5984     1348.3002  662.0096
838.7561  658.8395     1142.1526  612.5619     2340.6174 1504.3708
535.0766  392.5995     1088.0039  708.7622     587.1792  406.2180
596.4408  443.5586     484.6612  296.9192     1540.9741  692.1689
924.5619  640.1164     1536.0201 1071.4627     1115.8481  588.1371
487.7583  333.8394     678.8974  496.5976     1044.6843  511.2609
692.6397  466.9583     671.8802  503.3974     1389.7929  700.5600
997.8770  543.3969     690.4683  357.6411     2497.7860 1301.1451
506.9995  317.7198     860.6948  430.3376     1585.3809  879.0660
654.1587  424.3209     873.3095  624.6990     1862.0438  912.8851
933.9193  518.9617     894.4598  582.5413     2008.8546 1509.7812
433.6813  338.0014     1148.6470  580.2215     697.3099  484.0605
587.5962  419.6412     926.8762  543.8807     571.2517  399.6703
896.4746  476.3200     839.0414  588.6372     598.3465  444.1001
454.4782  386.3602     829.4974  627.9999     461.0977  248.8101
584.9989  423.2783     1264.0043  712.1012     977.1107  527.8014
800.7990  503.3572     1937.9771  968.3949     883.9849  500.6313
502.4369  354.6389     698.8317  482.5816     718.3594  436.8107
713.5197  497.3182     920.4199  593.1694     543.8971  374.7990
906.0006  588.5195     1897.5711 1033.5658     1587.3480  726.3921
880.5969  654.5971     891.6824  693.6795     4957.8130 1827.2000
796.8289  550.7274     889.6784  693.6795     969.6838  523.4911
854.8791  528.3770     1221.4818  761.2791     419.9980  334.9998
1167.3716  640.4813     544.5991  361.3981     561.9990  473.2009
523.8000  401.3204     1031.4491  628.4522     689.5988  581.2029
670.7792  435.9990     1462.9497  771.4486     1398.5203  929.7540
377.0584  276.5606     830.4353  757.1187     820.8168  591.1974
851.5430  588.3488     975.0415  821.5970     875.1716  637.5483
1121.0937  664.1978     1337.9983 1022.3202     1392.4499  674.9509
625.5179  444.8602     867.6427  679.4407     1256.3174  776.7589
805.5377  462.8995     725.7459  538.7491     1362.8590  959.5170
558.5812  377.7792     989.0056  679.9981     1999.2552 1250.9643
884.4005  553.1504     1525.0005  977.0033     1209.4730  737.8201
1257.4989  810.8962     672.1960  561.2015     1125.0356  810.6772
2051.1789 1067.9541     923.3977  728.3997     1827.4010  983.0009
1466.3330 1049.8788     472.3215  372.3186     1014.1540  708.8968
730.0989  522.7012     590.7601  361.5210     880.3944  633.1200
2432.3910 1424.8047     831.7983  620.8006     873.7375  631.7982
940.9218  517.9196     1139.4945  819.9964     951.4432  608.6419
1177.8547  830.9586     507.5169  360.8780     473.0022  300.9999
1222.5939  925.5795     576.1972  395.7608     601.0030  377.9984
1519.5811 1162.0024     696.5991  442.0001     713.9979  397.0015
687.6638  383.4580     650.8180  404.0384     829.2984  588.5195
953.1192  621.1173     949.5802  670.7993     959.7953  681.7616

```

```

953.1192 621.1173 497.1193 297.5702 1212.9613 807.3603
953.1192 621.1173 570.1674 353.4882 958.8743 696.8011
939.0418 548.6002 724.7306 383.9376 1129.4431 811.1962
1283.4025 745.2353 408.3399 284.8008 1943.0419 1305.7201
1511.5789 837.8005 638.6713 431.1000 539.6388 442.0001
1342.5821 795.3402 1225.7890 801.3518 463.5990 353.6013
511.7980 418.5976 715.3701 448.4513 562.6400 468.0008
689.7988 508.7974 800.4708 577.9111 736.7584 526.7573
1532.3074 883.2780 975.5974 570.5210 1415.4461 890.2390
1056.0808 742.5276 1613.7565 865.3205 2208.7897 1318.8033
387.3195 242.3202 608.5019 444.5578 636.0009 331.0005
387.3195 242.3202 958.6634 680.4198 759.4010 416.4015
410.9987 266.0010 835.9426 576.2779 1078.8382 596.8406
499.7510 408.4992 1024.8177 708.4787 748.6413 429.0399
832.7554 614.7588 1006.4353 734.2356 987.6417 619.6408
614.9986 385.3184 726.0000 433.0010 788.0961 400.7990
887.4658 515.6200 494.4174 327.4188 1020.0225 775.0209
1595.1611 1138.1620 776.5958 485.5198 1230.9235 772.7611
1807.9520 993.9630 415.4407 305.4390 440.5174 306.5191
541.2006 299.1993 581.3599 468.0008 743.0772 522.6019
1057.6767 750.3202 :: (x[1..m],y)[i] for i = 0...n-1
1 :: isx[1..m]
0.10 0.25 0.50 0.75 0.90 :: tau[1..ntau]
Return Residuals = Yes
Matrix Returned = Covariance
Interval Method = IID

```

10.3 Program Results

nag_regsn_quant_linear (g02qgc) Example Program Results

Quantile: 0.100

	Lower Limit	Parameter Estimate	Upper Limit
1	74.946	110.142	145.337
2	0.370	0.402	0.433

Covariance matrix
3.19e+02 -2.54e-01
2.59e-04

Quantile: 0.250

	Lower Limit	Parameter Estimate	Upper Limit
1	64.232	95.483	126.735
2	0.446	0.474	0.502

Covariance matrix
2.52e+02 -2.00e-01
2.04e-04

Quantile: 0.500

	Lower Limit	Parameter Estimate	Upper Limit
1	55.399	81.482	107.566
2	0.537	0.560	0.584

Covariance matrix
1.75e+02 -1.40e-01
1.42e-04

Quantile: 0.750

	Lower	Parameter	Upper
--	-------	-----------	-------

	Limit	Estimate	Limit
1	41.372	62.396	83.421
2	0.625	0.644	0.663

Covariance matrix

1.14e+02	-9.07e-02
	9.23e-05

Quantile: 0.900

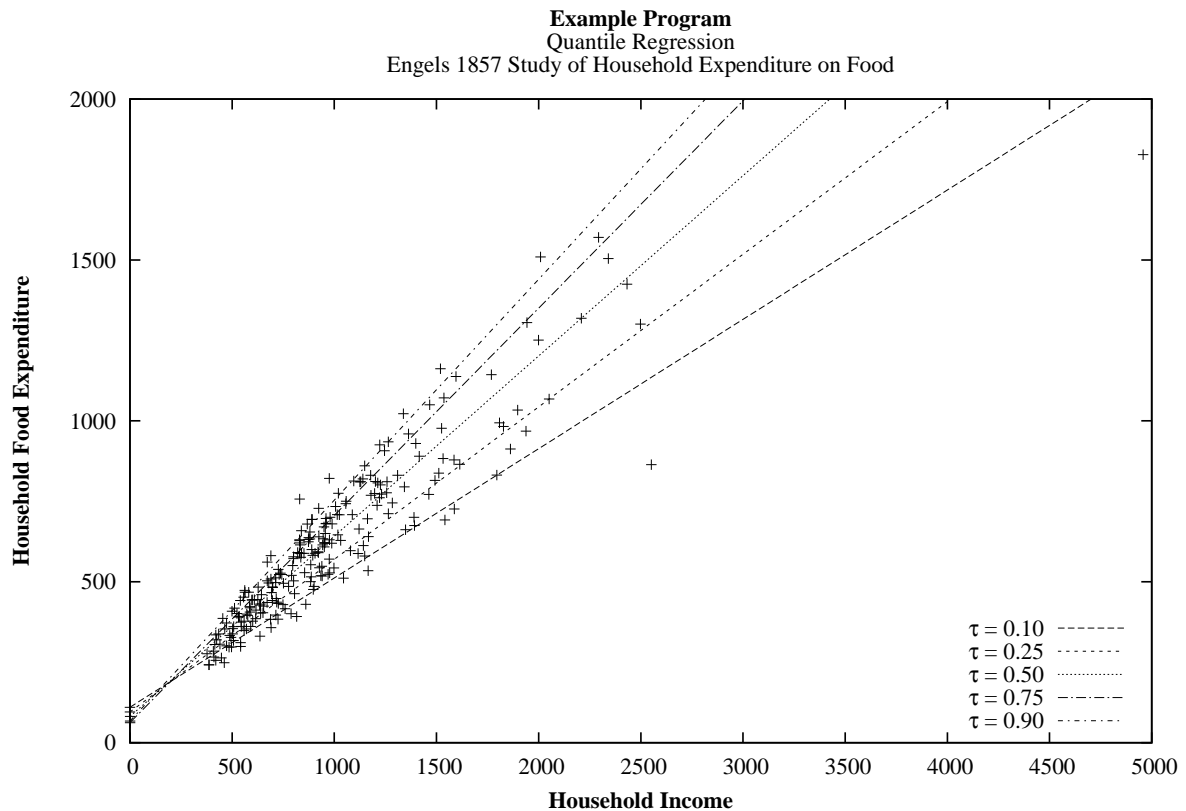
	Lower Limit	Parameter Estimate	Upper Limit
1	26.829	67.351	107.873
2	0.650	0.686	0.723

Covariance matrix

4.23e+02	-3.37e-01
	3.43e-04

First 10 Residuals

	Quantile				
	0.100	0.250	0.500	0.750	0.900
1	-23.10718	-38.84219	-61.00711	-77.14462	-99.86551
2	140.20549	96.93582	42.00636	-6.04177	-44.85812
3	91.19725	59.31654	17.93924	-16.90993	-49.06884
4	-16.70358	-41.20981	-73.81193	-100.11463	-127.96277
5	296.77717	221.32470	128.09970	42.75414	-14.87476
6	-271.39185	-441.31464	-646.95350	-841.78309	-954.63488
7	13.48419	-37.04518	-100.61322	-157.07478	-200.13481
8	218.91527	146.69601	57.31834	-24.28017	-80.01908
9	0.00000	-115.21109	-255.74639	-387.16920	-468.03911
10	36.09526	4.52393	-36.48522	-70.97584	-102.95390



11 Algorithmic Details

By the addition of slack variables the minimization (1) can be reformulated into the linear programming problem

$$\underset{(u,v,\beta) \in \mathbb{R}_+^n \times \mathbb{R}_+^n \times \mathbb{R}^p}{\text{minimize}} \quad \tau e^T u + (1 - \tau) e^T v \quad \text{subject to} \quad y = X\beta + u - v \quad (2)$$

and its associated dual

$$\underset{d}{\text{maximize}} \quad y^T d \quad \text{subject to} \quad X^T d = 0, d \in [\tau - 1, \tau]^n \quad (3)$$

where e is a vector of n 1s. Setting $a = d + (1 - \tau)e$ gives the equivalent formulation

$$\underset{a}{\text{maximize}} \quad y^T a \quad \text{subject to} \quad X^T a = (1 - \tau)X^T e, a \in [0, 1]^n. \quad (4)$$

The algorithm introduced by Portnoy and Koenker (1997) and used by `nag_regsn_quant_linear` (`g02qgc`), uses the primal-dual formulation expressed in equations (2) and (4) along with a logarithmic barrier function to obtain estimates for β . The algorithm is based on the predictor-corrector algorithm of Mehrotra (1992) and further details can be obtained from Portnoy and Koenker (1997) and Koenker (2005). A good description of linear programming, interior point algorithms, barrier functions and Mehrotra's predictor-corrector algorithm can be found in Nocedal and Wright (1999).

11.1 Interior Point Algorithm

In this section a brief description of the interior point algorithm used to estimate the model parameters is presented. It should be noted that there are some differences in the equations given here – particularly (7) and (9) – compared to those given in Koenker (2005) and Portnoy and Koenker (1997).

11.1.1 Central path

Rather than optimize (4) directly, an additional slack variable s is added and the constraint $a \in [0, 1]^n$ is replaced with $a + s = e, a_i \geq 0, s_i \geq 0$, for $i = 1, 2, \dots, n$.

The positivity constraint on a and s is handled using the logarithmic barrier function

$$B(a, s, \mu) = y^T a + \mu \sum_{i=1}^n (\log a_i + \log s_i).$$

The primal-dual form of the problem is used giving the Lagrangian

$$L(a, s, \beta, u, \mu) = B(a, s, \mu) - \beta^T (X^T a - (1 - \tau)X^T e) - u^T (a + s - e)$$

whose central path is described by the following first order conditions

$$\begin{aligned} X^T a &= (1 - \tau)X^T e \\ a + s &= e \\ X\beta + u - v &= y \\ S U e &= \mu e \\ A V e &= \mu e \end{aligned} \quad (5)$$

where A denotes the diagonal matrix with diagonal elements given by a , similarly with S, U and V . By enforcing the inequalities on s and a strictly, i.e., $a_i > 0$ and $s_i > 0$ for all i we ensure that A and S are positive definite diagonal matrices and hence A^{-1} and S^{-1} exist.

Rather than applying Newton's method to the system of equations given in (5) to obtain the step directions $\delta_\beta, \delta_a, \delta_s, \delta_u$ and δ_v , Mehrotra substituted the steps directly into (5) giving the augmented system of equations

$$\begin{aligned} X^T(a + \delta_a) &= (1 - \tau)X^T e \\ (a + \delta_a) + (s + \delta_s) &= e \\ X(\beta + \delta_\beta) + (u + \delta_u) - (v + \delta_v) &= y \\ (S + \Delta_s)(U + \Delta_u)e &= \mu e \\ (A + \Delta_a)(V + \Delta_v)e &= \mu e \end{aligned} \quad (6)$$

where $\Delta_a, \Delta_s, \Delta_u$ and Δ_v denote the diagonal matrices with diagonal elements given by $\delta_a, \delta_s, \delta_u$ and δ_v respectively.

11.1.2 Affine scaling step

The affine scaling step is constructed by setting $\mu = 0$ in (5) and applying Newton's method to obtain an intermediate set of step directions

$$\begin{aligned} (X^T W X) \delta_\beta &= X^T W (y - X\beta) + (\tau - 1) X^T e + X^T a \\ \delta_a &= W (y - X\beta - X\delta_\beta) \\ \delta_s &= -\delta_a \\ \delta_u &= S^{-1} U \delta_a - U e \\ \delta_v &= A^{-1} V \delta_s - V e \end{aligned} \quad (7)$$

where $W = (S^{-1} U + A^{-1} V)^{-1}$.

Initial step sizes for the primal ($\hat{\gamma}_P$) and dual ($\hat{\gamma}_D$) parameters are constructed as

$$\begin{aligned} \hat{\gamma}_P &= \sigma \min \left\{ \min_{i, \delta_{a_i} < 0} \{a_i / \delta_{a_i}\}, \min_{i, \delta_{s_i} < 0} \{s_i / \delta_{s_i}\} \right\} \\ \hat{\gamma}_D &= \sigma \min \left\{ \min_{i, \delta_{u_i} < 0} \{u_i / \delta_{u_i}\}, \min_{i, \delta_{v_i} < 0} \{v_i / \delta_{v_i}\} \right\} \end{aligned} \quad (8)$$

where σ is a user-supplied scaling factor. If $\hat{\gamma}_P \times \hat{\gamma}_D \geq 1$ then the nonlinearity adjustment, described in Section 11.1.3, is not made and the model parameters are updated using the current step size and directions.

11.1.3 Nonlinearity Adjustment

In the nonlinearity adjustment step a new estimate of μ is obtained by letting

$$\hat{g}(\hat{\gamma}_P, \hat{\gamma}_D) = (s + \hat{\gamma}_P \delta_s)^T (u + \hat{\gamma}_D \delta_u) + (a + \hat{\gamma}_P \delta_a)^T (v + \hat{\gamma}_D \delta_v)$$

and estimating μ as

$$\mu = \left(\frac{\hat{g}(\hat{\gamma}_P, \hat{\gamma}_D)}{\hat{g}(0, 0)} \right)^3 \frac{\hat{g}(0, 0)}{2n}.$$

This estimate, along with the nonlinear terms ($\Delta u, \Delta s, \Delta a$ and Δv) from (6) are calculated using the values of $\delta_a, \delta_s, \delta_u$ and δ_v obtained from the affine scaling step.

Given an updated estimate for μ and the nonlinear terms the system of equations

$$\begin{aligned} (X^T W X) \delta_\beta &= X^T W (y - X\beta + \mu(S^{-1} - A^{-1})e + S^{-1} \Delta_s \Delta_u e - A^{-1} \Delta_a \Delta_v e) + (\tau - 1) X^T e + X^T a \\ \delta_a &= W (y - X\beta - X\delta_\beta + \mu(S^{-1} - A^{-1})e) \\ \delta_s &= -\delta_a \\ \delta_u &= \mu S^{-1} e + S^{-1} U \delta_a - U e - S^{-1} \Delta_s \Delta_u e \\ \delta_v &= \mu A^{-1} e + A^{-1} V \delta_s - V e - A^{-1} \Delta_a \Delta_v e \end{aligned} \quad (9)$$

are solved and updated values for $\delta_\beta, \delta_a, \delta_s, \delta_u, \delta_v, \hat{\gamma}_P$ and $\hat{\gamma}_D$ calculated.

11.1.4 Update and convergence

At each iteration the model parameters (β, a, s, u, v) are updated using step directions, $(\delta_\beta, \delta_a, \delta_s, \delta_u, \delta_v)$ and step lengths $(\hat{\gamma}_P, \hat{\gamma}_D)$.

Convergence is assessed using the duality gap, that is, the differences between the objective function in the primal and dual formulations. For any feasible point (u, v, s, a) the duality gap can be calculated from equations (2) and (3) as

$$\begin{aligned} \tau e^T u + (1 - \tau) e^T v - d^T y &= \tau e^T u + (1 - \tau) e^T v - (a - (1 - \tau) e)^T y \\ &= s^T u + a^T v \\ &= e^T u - a^T y + (1 - \tau) e^T X \beta \end{aligned}$$

and the optimization terminates if the duality gap is smaller than the tolerance supplied in the optional parameter **Tolerance**.

11.1.5 Additional information

Initial values are required for the parameters a, s, u, v and β . If not supplied by the user, initial values for β are calculated from a least squares regression of y on X . This regression is carried out by first constructing the cross-product matrix $X^T X$ and then using a pivoted QR decomposition as performed by `nag_dgeqp3` (f08bfc). In addition, if the cross-product matrix is not of full rank, a rank reduction is carried out and, rather than using the full design matrix, X , a matrix formed from the first p -rank columns of XP is used instead, where P is the pivot matrix used during the QR decomposition. Parameter estimates, confidence intervals and the rows and columns of the matrices returned in the argument **ch** (if any) are set to zero for variables dropped during the rank-reduction. The rank reduction step is performed irrespective of whether initial values are supplied by the user.

Once initial values have been obtained for β , the initial values for u and v are calculated from the residuals. If $|r_i| < \epsilon_u$ then a value of $\pm\epsilon_u$ is used instead, where ϵ_u is supplied in the optional parameter **Epsilon**. The initial values for the a and s are always set to $1 - \tau$ and τ respectively.

The solution for δ_β in both (7) and (9) is obtained using a Bunch–Kaufman decomposition, as implemented in `nag_dsytrf` (f07mdc).

11.2 Calculation of Covariance Matrix

`nag_regsn_quant_linear` (g02qgc) supplies four methods to calculate the covariance matrices associated with the parameter estimates for β . This section gives some additional detail on three of the algorithms, the fourth, (which uses bootstrapping), is described in Section 3.

(i) Independent, identically distributed (IID) errors

When assuming IID errors, the covariance matrices depend on the sparsity, $s(\tau)$, which `nag_regsn_quant_linear` (g02qgc) estimates as follows:

- Let r_i denote the residuals from the original quantile regression, that is $r_i = y_i - x_i^T \hat{\beta}$.
- Drop any residual where $|r_i|$ is less than ϵ_u , supplied in the optional parameter **Epsilon**.
- Sort and relabel the remaining residuals in ascending order, by absolute value, so that $\epsilon_u < |r_1| < |r_2| < \dots$.
- Select the first l values where $l = h_n n$, for some bandwidth h_n .
- Sort and relabel these l residuals again, so that $r_1 < r_2 < \dots < r_l$ and regress them against a design matrix with two columns ($p = 2$) and rows given by $x_i = \{1, i/(n - p)\}$ using quantile regression with $\tau = 0.5$.
- Use the resulting estimate of the slope as an estimate of the sparsity.

(ii) Powell Sandwich

When using the Powell Sandwich to estimate the matrix H_n , the quantity

$$c_n = \min(\sigma_r, (q_{r3} - q_{r1})/1.34) \times (\Phi^{-1}(\tau + h_n) - \Phi^{-1}(\tau - h_n))$$

is calculated. Dependent on the value of τ and the method used to calculate the bandwidth (h_n), it is possible for the quantities $\tau \pm h_n$ to be too large or small, compared to **machine precision** (ϵ). More specifically, when $\tau - h_n \leq \sqrt{\epsilon}$, or $\tau + h_n \geq 1 - \sqrt{\epsilon}$, a warning flag is raised in **info**, the value is truncated to $\sqrt{\epsilon}$ or $1 - \sqrt{\epsilon}$ respectively and the covariance matrix calculated as usual.

(iii) Hendricks–Koenker Sandwich

The Hendricks–Koenker Sandwich requires the calculation of the quantity $d_i = x_i^T (\hat{\beta}(\tau + h_n) - \hat{\beta}(\tau - h_n))$. As with the Powell Sandwich, in cases where $\tau - h_n \leq \sqrt{\epsilon}$, or $\tau + h_n \geq 1 - \sqrt{\epsilon}$, a warning flag is raised in **info**, the value truncated to $\sqrt{\epsilon}$ or $1 - \sqrt{\epsilon}$ respectively and the covariance matrix calculated as usual.

In addition, it is required that $d_i > 0$, in this method. Hence, instead of using $2h_n/d_i$ in the calculation of H_n , $\max(2h_n/(d_i + \epsilon_u), 0)$ is used instead, where ϵ_u is supplied in the optional parameter **Epsilon**.

12 Optional Parameters

Several optional parameters in `nag_regsn_quant_linear` (g02qgc) control aspects of the optimization algorithm, methodology used, logic or output. Their values are contained in the arrays **iopts** and **opts**; these must be initialized before calling `nag_regsn_quant_linear` (g02qgc) by first calling `nag_g02_opt_set` (g02zkc) with **optstr** set to **Initialize** = g02qgc.

Each optional parameter has an associated default value; to set any of them to a non-default value, use `nag_g02_opt_set` (g02zkc). The current value of an optional parameter can be queried using `nag_g02_opt_get` (g02zlc).

The remainder of this section can be skipped if you wish to use the default values for all optional parameters.

The following is a list of the optional parameters available. A full description of each optional parameter is provided in Section 12.1.

Band Width Alpha

Band Width Method

Big

Bootstrap Interval Method

Bootstrap Iterations

Bootstrap Monitoring

Calculate Initial Values

Defaults

Drop Zero Weights

Epsilon

Interval Method

Iteration Limit

Matrix Returned

Monitoring

QR Tolerance

Return Residuals

Sigma

Significance Level

Tolerance

Unit Number

12.1 Description of the Optional Parameters

For each option, we give a summary line, a description of the optional parameter and details of constraints.

The summary line contains:

the keywords, where the minimum abbreviation of each keyword is underlined (if no characters of an optional qualifier are underlined, the qualifier may be omitted);

a parameter value, where the letters a , i and r denote options that take character, integer and real values respectively;

the default value, where the symbol ϵ is a generic notation for *machine precision* (see `nag_machine_precision` (X02AJC)).

Keywords and character values are case and white space insensitive.

Band Width Alpha *r* Default = 1.0

A multiplier used to construct the parameter α_b used when calculating the Sheather–Hall bandwidth (see Section 3), with $\alpha_b = (1 - \alpha) \times \mathbf{Band\ Width\ Alpha}$. Here, α is the **Significance Level**.

Constraint: **Band Width Alpha** > 0.0.

Band Width Method *a* Default = 'SHEATHER HALL'

The method used to calculate the bandwidth used in the calculation of the asymptotic covariance matrix Σ and H^{-1} if **Interval Method** = HKS, KERNEL or IID (see Section 3).

Constraint: **Band Width Method** = SHEATHER HALL or BOFINGER.

Big *r* Default = 10.0²⁰

This parameter should be set to something larger than the biggest value supplied in **dat** and **y**.

Constraint: **Big** > 0.0.

Bootstrap Interval Method *a* Default = QUANTILE

If **Interval Method** = BOOTSTRAP XY, **Bootstrap Interval Method** controls how the confidence intervals are calculated from the bootstrap estimates.

Bootstrap Interval Method = T

t intervals are calculated. That is, the covariance matrix, $\Sigma = \{\sigma_{ij} : i, j = 1, 2, \dots, p\}$ is calculated from the bootstrap estimates and the limits calculated as $\beta_i \pm t_{(n-p, (1+\alpha)/2)} \sigma_{ii}$ where $t_{(n-p, (1+\alpha)/2)}$ is the $(1 + \alpha)/2$ percentage point from a Student's t distribution on $n - p$ degrees of freedom, n is the effective number of observations and α is given by the optional parameter **Significance Level**.

Bootstrap Interval Method = QUANTILE

Quantile intervals are calculated. That is, the upper and lower limits are taken as the $(1 + \alpha)/2$ and $(1 - \alpha)/2$ quantiles of the bootstrap estimates, as calculated using `nag_double_quantiles` (g01amc).

Constraint: **Bootstrap Interval Method** = T or QUANTILE.

Bootstrap Iterations *i* Default = 100

The number of bootstrap samples used to calculate the confidence limits and covariance matrix (if requested) when **Interval Method** = BOOTSTRAP XY.

Constraint: **Bootstrap Iterations** > 1.

Bootstrap Monitoring *a* Default = NO

If **Bootstrap Monitoring** = YES and **Interval Method** = BOOTSTRAP XY, then the parameter estimates for each of the bootstrap samples are displayed. This information is sent to the unit number specified by **Unit Number**.

Constraint: **Bootstrap Monitoring** = YES or NO.

Calculate Initial Values *a* Default = YES

If **Calculate Initial Values** = YES then the initial values for the regression parameters, β , are calculated from the data. Otherwise they must be supplied in **b**.

Constraint: **Calculate Initial Values** = YES or NO.

Defaults

This special keyword is used to reset all optional parameters to their default values.

Drop Zero Weights *a* Default = YES

If a weighted regression is being performed and **Drop Zero Weights** = YES then observations with zero weight are dropped from the analysis. Otherwise such observations are included.

Constraint: **Drop Zero Weights** = YES or NO.

Epsilon *r* Default = $\sqrt{\epsilon}$

ϵ_u , the tolerance used when calculating the covariance matrix and the initial values for u and v . For additional details see Section 11.2 and Section 11.1.5 respectively.

Constraint: **Epsilon** \geq 0.0.

Interval Method *a* Default = IID

The value of **Interval Method** controls whether confidence limits are returned in **bl** and **bu** and how these limits are calculated. This parameter also controls how the matrices returned in **ch** are calculated.

Interval Method = NONE

No limits are calculated and **bl**, **bu** and **ch** are not referenced.

Interval Method = KERNEL

The Powell Sandwich method with a Gaussian kernel is used.

Interval Method = HKS

The Hendricks–Koenker Sandwich is used.

Interval Method = IID

The errors are assumed to be identical, and independently distributed.

Interval Method = BOOTSTRAP XY

A bootstrap method is used, where sampling is done on the pair (y_i, x_i) . The number of bootstrap samples is controlled by the parameter **Bootstrap Iterations** and the type of interval constructed from the bootstrap samples is controlled by **Bootstrap Interval Method**.

Constraint: **Interval Method** = NONE, KERNEL, HKS, IID or BOOTSTRAP XY.

Iteration Limit *i* Default = 100

The maximum number of iterations to be performed by the interior point optimization algorithm.

Constraint: **Iteration Limit** $>$ 0.

Matrix Returned *a* Default = NONE

The value of **Matrix Returned** controls the type of matrices returned in **ch**. If **Interval Method** = NONE, this parameter is ignored and **ch** is not referenced. Otherwise:

Matrix Returned = NONE

No matrices are returned and **ch** is not referenced.

Matrix Returned = COVARIANCE

The covariance matrices are returned.

Matrix Returned = H INVERSE

If **Interval Method** = KERNEL or HKS, the matrices J and H^{-1} are returned. Otherwise no matrices are returned and **ch** is not referenced.

The matrices returned are calculated as described in Section 3, with the algorithm used specified by **Interval Method**. In the case of **Interval Method** = BOOTSTRAP XY the covariance matrix is calculated directly from the bootstrap estimates.

Constraint: **Matrix Returned** = NONE, COVARIANCE or H INVERSE.

Monitoring *a* Default = NO

If **Monitoring** = YES then the duality gap is displayed at each iteration of the interior point optimization algorithm. In addition, the final estimates for β are also displayed.

The monitoring information is sent to the unit number specified by **Unit Number**.

Constraint: **Monitoring** = YES or NO.

QR Tolerance *r* Default = $\epsilon^{0.9}$

The tolerance used to calculate the rank, k , of the $p \times p$ cross-product matrix, $X^T X$. Letting Q be the orthogonal matrix obtained from a QR decomposition of $X^T X$, then the rank is calculated by comparing Q_{ii} with $Q_{11} \times$ **QR Tolerance**.

If the cross-product matrix is rank deficient, then the parameter estimates for the $p - k$ columns with the smallest values of Q_{ii} are set to zero, along with the corresponding entries in **bl**, **bu** and **ch**, if returned. This is equivalent to dropping these variables from the model. Details on the QR decomposition used can be found in nag_dgeqp3 (f08bfc).

Constraint: **QR Tolerance** > 0.0.

Return Residuals *a* Default = NO

If **Return Residuals** = YES, the residuals are returned in **res**. Otherwise **res** is not referenced.

Constraint: **Return Residuals** = YES or NO.

Sigma *r* Default = 0.99995

The scaling factor used when calculating the affine scaling step size (see equation (8)).

Constraint: $0.0 < \mathbf{Sigma} < 1.0$.

Significance Level *r* Default = 0.95

α , the size of the confidence interval whose limits are returned in **bl** and **bu**.

Constraint: $0.0 < \mathbf{Significance Level} < 1.0$.

Tolerance *r* Default = $\sqrt{\epsilon}$

Convergence tolerance. The optimization is deemed to have converged if the duality gap is less than **Tolerance** (see Section 11.1.4).

Constraint: **Tolerance** > 0.0.

Unit Number *i* Output sent to stdout

The unit number to which any monitoring information is sent. See nag_open_file (x04acc) for details on how to assign a file to a unit number. If no unit number is specified then any monitoring information will be sent to standard output (stdout).

Constraint: **Unit Number** > 1.

13 Description of Monitoring Information

See the description of the optional argument **Monitoring**.
