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# An Interior Point Algorithm for Nonlinear Quantile Regression 

Roger Koenker*<br>and<br>Beum J. Park ${ }^{*}$

April 1992


#### Abstract

A new algorithm for computing quantile regression estimates for problems in which the response function is nonlinear in parameters is described. The nonlinear $l_{1}$ estimation problem is a special (median) case. The algorithm is closely related to recent developments on interior point methods for solving linear programs. Performance of the algorithm on a variety of test problems including the censored linear quantile regression problem of Powell (1986) is reported.


Keywords: Quantile Regression, Nonlinear Regression, Linear Programming, Interior Point Algorithms, Nonlinear Programming

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## 1. Introduction

About a century ago Edgeworth observed that methods of estimation based upon minimizing sums of absolute residuals could be far superior to least-squares methods under nonGaussian error conditions. Laplace had drawn similar conclusions a century earlier. See Stigler (1986) on this early history. But computation of $l_{1}$-estimators, even for linear regression, remained a major impediment to applications until the emergence of the simplex algorithm for linear programming in the 1940's. Papers by Charnes, Cooper and Ferguson (1955) and Wagner (1959) provided a foundation for modern algorithms for linear $l_{1}$-regression by Barrodale and Roberts (1973), Bartels and Conn (1980) and others. These algorithms are readily extended to linear quantile regression, Koenker and Bassett (1978), of which $l_{1}$-regression is an important (median) special case.

The current state of algorithms for nonlinear quantile regression is far less satisfactory. Certainly nothing comparable to the venerable Gauss-Newton algorithm for nonlinear least squares problems has emerged. Despite a flurry of interest by prominent numerical analysts in the 1970's and early 1980's, see, e.g., Osborne and Watson (1971), Murray and Overton (1981) and Bartels and Conn (1982), occasional applications of nonlinear quantile regression have relied on the Nelder and Mead (1965) algorithm and other generic optimization methods. An excellent statement of the current state-of-the-art is provided in the thesis of Busovaca (1985).

In contrast, the statistical theory of nonlinear quantile regression has developed rapidly in recent years. Powell (1986) has emphasized its value in the analysis of censored and truncated responses. Asymptotic theory for independent errors has been developed by Oberhofer(1982), Dupacova(1987), and Powell(1991). Theoretical developments by Weiss (1991) and White (1991) have stressed applications to time-series analysis. Applications of Horowitz and Neu-
mann (1987), Chamberlain (1990), and others have demonstrated its value in applied econometrics.

In this paper we will describe a new approach to the computation of nonlinear quantile regression estimators based on recent interior point methods for solving linear programs. In the next section we review interior point methods for strictly linear problems. Section 3 describes our approach to nonlinear problems, and Section 4 describes our computational experience.

## 2. Interior Point Methods for Linear Programs

In this section we provide a brief discussion of interior point methods for solving strictly linear programs including the linear quantile regression problem. Our exposition will follow closely that of Vanderbei, Meketon, and Freedman (1986) and Meketon (1986). We should emphasize that in our experience interior point algorithms for linear quantile regression do not appear to be competitive in efficiency terms with existing simplex method algorithms. See Koenker and d'Orey(1987) for a description of a simplex based algorithm for linear quantile regression. However, unlike simplex based methods they do appear to offer a natural extension to nonlinear problems. Thus a clear understanding of the linear case is an essential first step in our exposition.

### 2.1. A Canonical LP

Consider the equality constrained linear program

$$
\begin{equation*}
\min \left\{c^{\prime} \omega \mid \omega \in \Omega \equiv\left\{\omega \in \mathbf{R}_{+}^{n}, A \omega=b\right\}\right\} \tag{2.1}
\end{equation*}
$$

where $\mathbf{R}_{+}^{n}$ denotes the positive orthant of $\mathbf{R}^{n}$. Given a feasible point in the interior of the constraint set, $\omega \in \operatorname{int}(\Omega)$, interior point methods proceed in two steps. First we transform coordinates to reposition $\omega$ so it is centered relative to the set $\Omega$. Then a (projected) gradient step is
taken toward the boundary of $\Omega$. Repeating this process brings us arbitrarily close to a solution, and a stopping criterion is eventually invoked.

To flesh out this brief description, let $D=\operatorname{diag}(\omega)$ and consider the transformation

$$
\omega \rightarrow D^{-1} \omega
$$

We have $D^{-1} \omega=1_{n}$, an $n$-vector of ones, so the transformation $D$ has the effect of centering $\omega$ relative to the orthant boundaries of $\Omega$. Correspondingly, we may define $\tilde{A}=A D$ and $\tilde{c}=D c$. In the transformed coordinates we wish to move in the gradient direction $-\tilde{c}$, but to preserve feasibility we should instead project $\tilde{c}$ onto the null space of $\tilde{A}$ to insure that the equality constraints are satisfied.

Let $\hat{c}$ denote this projection, i.e.,

$$
\hat{c}=\left(I-\tilde{A}^{\prime}\left(\tilde{A^{\prime}} \tilde{A}\right)^{-1} \tilde{A}\right) \tilde{c}
$$

Clearly, $\hat{c}$ is a direction of descent; and we now move toward the boundary of $\Omega$ in this direction.
Let

$$
\hat{\alpha}=\max _{i=1, \ldots, n}\left\{e_{i}^{\prime} \hat{c}\right\}
$$

where $e_{i}$ is the $i^{\text {th }}$ unit basis vector for $\mathbf{R}^{n}$. For some fixed $\eta \in(0,1)$, consider

$$
\omega \leftarrow \omega-(\eta / \alpha) D \hat{c}
$$

which defines a sequence of iterations $\omega_{k+1}=T\left(\omega_{k}\right)$. Since at each iteration

$$
c^{\prime} \omega_{k+1}=c^{\prime} \omega_{k}-(\eta / \alpha) c^{\prime} D \hat{c}=c^{\prime} \omega_{k}-(\eta / \alpha)\|\hat{c}\|
$$

we expect to see an improvement in the objective function at each iteration.

Proposition. If $\hat{c} \leq 0$ the problem (2.1) is unbounded, unless $\hat{c}=0$ in which case every $\omega \in \Omega$ is
optimal. Otherwise, the problem is bounded and the sequence $\left\{c^{\prime} \omega_{k}\right\}$ is strictly decreasing. Proof. Since the proof of this proposition, found in Vanderbei, Meketon and Freedman (1986), is both elementary and revealing we repeat it here for the sake of completeness. If $\hat{c}=0$, there exists a vector $z$ such that $\tilde{c}=\tilde{A}^{\prime} z$, hence $D \tilde{c}=D \tilde{A} z$ and since $\omega \in \operatorname{int}(\Omega)$ it follows that $c=A z$. But then for any $\omega \in \Omega$,

$$
c^{\prime} \omega=z^{\prime} A \omega=z^{\prime} b
$$

which is independent of $\omega$, establishing that $c^{\prime} \omega$ is constant on all of $\Omega$. Next consider $\hat{c} \leq 0$. Note that

$$
\begin{equation*}
c^{\prime} \omega_{1} \equiv c^{\prime} \omega-\gamma(\omega) \tilde{c}^{\prime} \hat{c}=c^{\prime} \omega-\gamma(\omega)\|\hat{c}\|_{2}^{2} \tag{2.4}
\end{equation*}
$$

where $\gamma(\omega)=\eta / \alpha$. Since $\hat{c} \leq 0$,

$$
\omega_{\rho}=\omega-\rho D \hat{c}
$$

is feasible for any $\rho>0$ and

$$
c^{\prime} \omega_{\rho}=c^{\prime} \omega-\rho\|\hat{c}\|_{2}^{2}
$$

implies that $c^{\prime} \omega_{\rho} \rightarrow-\infty$ as $\rho \rightarrow \infty$. Finally, if $\hat{c} \geq 0$, then since $\gamma(\omega)>0$ and $c^{\prime} \omega_{1}<c^{\prime} \omega$ follows from (2.4), establishing that the step is a direction of descent.
2.2. Linear $l_{1}$-regression

In the linear model

$$
y_{i}=x_{i}^{\prime} \beta+u_{i} \quad i=1, \cdots, n
$$

as noted in the introduction, the $l_{1}$-estimator of $\beta$ which minimizes

$$
R(b)=\sum_{i=1}^{n}\left|y_{i}-x_{i}^{\prime} b\right|
$$

may be formulated as a linear program. The dual problem may be written as

$$
\max \left\{y^{\prime} d \mid d \in \Omega=\left\{d \in[-1,1]^{n}, X^{\prime} d=0\right\}\right\}
$$

where $y$ is the $n$-vector of responses, and $X$ is the $n \times p$ design matrix. To solve the dual problem we proceed as before, except that the centering is slightly altered to accommodate the altered form of $\Omega$. For any initial feasible point $d$, e.g., $d=0$, following Meketon (1986), set

$$
D=\operatorname{diag}\left(\min \left\{1+d_{i}, 1-d_{i}\right\}\right)
$$

In the transformed coordinates $D^{-1} d$ the projected gradient is

$$
D \hat{u}=\left(I-D X^{\prime}\left(X^{\prime} D^{2} X\right)^{-1} X^{\prime} D\right) D y=D(y-X b)
$$

where $b=\left(X^{\prime} D^{2} X\right)^{-1} D^{2} y$. Note that as in the former case the transformation has the effect of centering the point $d$ in the feasible set $\Omega$. Now let

$$
\alpha=\max _{i}\left\{\max \left\{\frac{e_{i}^{\prime} D^{2} \hat{u}}{1+d_{i}}, \frac{-e_{i}^{\prime} D^{2} \hat{u}}{1-d_{i}}\right\}\right\}
$$

and again for $\eta \in(0,1)$ we take the step

$$
d \leftarrow d+(\eta / \alpha) D^{2} \hat{u}
$$

Note the change in sign since we are now maximizing. The iteration sequence $d_{k+1}=T\left(d_{k}\right)$ in the dual vector implicitly defines a corresponding primal sequence with

$$
b_{k}=\left(X^{\prime} D_{k}^{2} X\right)^{-1} X^{\prime} D_{k}^{2} y
$$

As Meketon notes, the duality theory yields a natural stopping criterion. Since

$$
y^{\prime} d_{k} \leq \sum\left|y_{i}-x_{i}^{\prime} b_{k}\right|
$$

with optimality if and only if equality holds, it is reasonable to stop iterating when the difference between the dual and primal values is less than a specified tolerance.

### 2.3 Linear Quantile Regression

If we replace the (symmetric) $l_{1}$-criterion with an asymmetric linear criterion so we minimize

$$
R_{\theta}(b)=\sum_{i=1}^{n} \rho_{\theta}\left(y_{i}-x_{i} b\right)
$$

$\rho_{\theta}(u)=(\theta-I(u<0)) u$, we obtain the regression quantiles of Koenker and Bassett (1978). The dual problem is now,

$$
\max \left\{y^{\prime} d \mid d \in \Omega=\left\{d \in[\theta-1, \theta]^{n}, X^{\prime} d=0\right\}\right\}
$$

This leads to an algorithm identical to the $l_{1}$ special case except that now

$$
D=\operatorname{diag}\left(\min \left(\theta-d_{i}, 1-\theta+d_{i}\right)\right)
$$

and

$$
\alpha=\max _{i}\left(\max \left[\frac{e_{i}^{\prime} D \hat{c}}{\theta-d_{i}}, \frac{-e_{i}^{\prime} D \hat{c}}{1-\theta+d_{i}}\right]\right)
$$

## 3. Nonlinear Quantile Regression

To extend these ideas to the case of nonlinear response functions we begin by considering the nonlinear $l_{1}$ problem

$$
\begin{equation*}
\min _{t \in \mathbf{R}^{p}} \sum\left|f_{i}(t)\right| \tag{3.1}
\end{equation*}
$$

where, for example,

$$
f_{i}(t)=y_{i}-f_{0}\left(x_{i}, t\right)
$$

As noted by El Attar, et al (1979) a necessary condition for $t^{*}$ to solve (3.1) is that there exists a vector $d \in[-1,1]^{n}$ such that

$$
\begin{gather*}
J\left(t^{*}\right)^{\prime} d=0  \tag{3.2}\\
f\left(t^{*}\right)^{\prime} d=\sum\left|f_{i}\left(t^{*}\right)\right| \tag{3.3}
\end{gather*}
$$

where $f(t)=\left(f_{i}(t)\right)$ and $J(t)=\left(\partial f_{i}(t) / \partial t_{j}\right)$
Thus, as proposed by Osborne and Watson (1971), one approach to solving (3.1) is to solve a succession of linearized $l_{1}$ problems minimizing

$$
\sum\left|f_{i}(t)-J_{i}(t)^{\prime} \delta\right|=\|f-J \delta\|_{1},
$$

choosing a step length, $\lambda$, at each iteration, by line search in the resulting directions $\delta$. The difficulty, as we see it, with this approach is not only that we must expend the effort to solve an $l_{1}$ linear program at each iteration, but, perhaps more significantly, the resulting directions may actually be inferior to directions determined by incomplete solutions to the sequence of linearized problems.

Let $t$ be the value of the parameter at the current iteration, and consider the dual problem

$$
\begin{equation*}
\max \left\{f^{\prime} d \in[-1,1]^{n}, J^{\prime} d=0\right\} . \tag{3.4}
\end{equation*}
$$

If the model were linear so

$$
f(s)=f(t)-K(s-t)
$$

for some fixed matrix $K$, then a solution can be found by applying Meketon's algorithm to find $d^{*}$ to solve (3.4), computing

$$
\delta^{*}=\left(K^{\prime} D^{2} K\right)^{-1} K^{\prime} D^{2} f
$$

where $D=\operatorname{diag}\left(\min \left(1-d_{i}^{*}, 1+d_{i}^{*}\right\}\right)$ and setting $t^{*}=t+\delta^{*}$. When $f$ is nonlinear there is no longer a compelling argument for fully solving (3.4) at each iteration, indeed, in our experience only a few iterations to refine the dual vector is preferable. In the version of the algorithm we have implemented to conduct the tests reported in the next section we take two dual steps between successive updates of $f$ and $J$. A detailed description of the algorithm is now provided.

### 3.1. Dual Step

For any feasible $d$ in the interior of the constraint set of (3.4) we refine $d$, following Meketon, as follows. Let

$$
\begin{aligned}
& D=\operatorname{diag}\left(\min \left\{1-d_{i}, 1+d_{i}\right\}\right) \\
& s=D^{2}\left(I-J\left(J^{\prime} D^{2} J\right)^{-1} J D^{2}\right) f . \\
& d \leftarrow d+(\eta / \alpha) s
\end{aligned}
$$

where

$$
\alpha=\max \left\{\max \left\{s_{i} /\left(1-d_{i}\right),-s_{i} /\left(1+d_{i}\right)\right\}\right\}
$$

and $\eta \in(0,1)$ is the constant chosen to insure feasibility. Following Meketon, we use $\eta=.97$. Updating $D, s$, and the new $d$ continues the iteration. This process is embedded in a sequence of primal iterations in which we update $f$ and $J$ as follows.

### 3.2. Primal Step

The dual step yields the primal direction

$$
\delta=\left(J^{\prime} D^{2} J\right)^{-1} J^{\prime} D^{2} f^{\prime}
$$

which we explore by line search. Our current implementation uses Brent's (1973) algorithm from the PORT3 library Fox (1984). Updating we have

$$
t \leftarrow t+\lambda^{*} \delta
$$

where $\lambda^{*}=\underset{\lambda}{\operatorname{argmin}}\|f(t+\lambda \delta)\|_{1}$, and we then update $f$ and $J$. However before returning to the dual step we must adjust the current $d$ so that it is feasible for the new value of $J$. This is accomplished, somewhat naively, by projecting the current $d$ onto the null space of the new $J$, i.e. $\hat{d}=\left(I-J\left(J^{\prime} J\right)^{-1} J^{\prime}\right) d$ and then shrinking it to insure that it lies in $[-1,1]^{n}$, so

$$
d \leftarrow \hat{d} /\left(\max _{i}\left\{\left|\hat{d}_{i}\right|\right\}+\varepsilon\right)
$$

for some tolerance parameter $\varepsilon>0$. Obviously, when the problem is linear, so $J$ is fixed, this "adjustment" is nugatory since $d$ is already in the null space of $J$, and the algorithm is essentially like Meketon's.

### 3.3. Stopping

The algorithm currently terminates when the new iterate fails to improve the objective function by a specified tolerance. Exploration of alternative stopping rules is a topic for future research.

### 3.4 Related Literature

Gill, Murray, Saunders, Tomlin, and Wright (1986) and Bayer and Lagarias (1991) have recently pointed out the close relationship of "projected Newton barrier" methods (see Fiacco and McCormick (1965)) and interior point methods. Algorithms closely related to the one described above could presumably be formulated employing logarithmic barrier functions in the
dual vector $d$.

### 3.5. Quantile Regression

As in the case of the linear problem the generalization of the $l_{1}$ problem to other quantiles is straightforward involving only a modification of the constraint set $[-1,1]^{n}$ to $[\theta-1, \theta]^{n}$ for some $\theta \in(0,1)$.

## 4. Numerical Experience

In this section we describe our computational experience with a variety of test problems. To facilitate comparison with existing results in the literature we have chosen problems from Busovaca(1985) and Wormersley(1986). We focus exclusively on the $l_{1}$ case since there are no comparable results in the literature for other quantiles. The problems used are described in detail in Appendix A. We have attempted to investigate all of the problems reported on by Busovaca, however in several cases we were unable find a complete description of the problem. The problem taken from Wormersley is included to explore the important special case of piecewise linear reponse functions which arise in Powell's(1986) formulation of the quantile regression problem for censored data.

In Appendix B we provide explicit versions of our algorithm in both S, Becker, Chambers and Wilks(1988), and Gauss. All of the reported tests were carried out in $S$ on a Sun $3 / 50$. Note that the line search algorithm in the $S$ and Gauss versions are different. To implement a simple version of the Osborne and Watson(1971) algorithm in S we employ the $S$ function l1fit which does $l_{1}$ regression using the Barrodale and Roberts(1973) algorithm. The $S$ function lsfit carries out the corresponding weighted least squares computations for the interior point algorithm.

A summary of our experience on the test problems appears in Tables 5.1 and 5.2.. For Wormersley's(1986) censored regression problem (Problem 1) our version of the interior point algorithm converges to Wormersley's reported solution. However, it should be noted that the solution to this problem is notoriously nonunique. Busovaca's algorithm cannot be employed on Problem 1 due to the fact that the Hessian of the response function is identically zero almost everywhere. The remaining problems are all taken from Busovaca, and generally our interior point solutions correspond closely to his. In Problems 7 and 13 there are small discrepancies favoring Busovaca; in Problem 9 there is a larger descrepancy favoring the interior point method. Results for our implementation of the Osborne and Watson algorithm are somewhat less satisfactory. It fails completely on Problems 11 and 12, performs poorly in Problems 1, 4b, and 13, but does slightly better than the interior point method on Problem 5. All three algorithms fail for Problem 4 a which is highly degenerate at the specified initial point. At an alternative starting point, the interior point algorithm performs well.

## 5. Some Concluding Remarks

We have described a simple approach to computing quantile regression estimates for problems with nonlinear response functions. The approach is based on recent developments on interior point methods for linear programming, but may be viewed as a variant of well-known iteratively reweighted least squares. While the algorithm seems to perform well on a variety of test problems, there is considerable room for improvement. Handling rank deficiency in the model Jacobian is critical. Alternative stopping criteria also should be explored.

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Figure 5.1
Algorithmic Performance on Several Test Problems

| Example | Starting <br> Point | Interior Point Algorithm |  | Osborne-Watson Algorithm |  | Busovaca Algorithm |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
|  |  | Optimal Objective | Number of Iterations | Optimal Objective | Number of Iterations | Optimal Objective | Number of Iterations |
| 1.Wormersley | $(0,0)$ | 3.032544 | 3 | 5.234825 | 3 | NA |  |
| 2.Bard | (1,1,1) | 0.1243555 | 6 | 0.1243383 | 5 | 0.1243406 | 13 |
| 3.Beale | (1,0.1) | $0.2928905 \mathrm{e}-07$ | 6 | $0.1547611 \mathrm{e}-06$ | 6 | 0.3695488e-05 | 8 |
| 4.a.Biggs | (1,1,1,1,1,1) | F |  | F |  | F |  |
| b.Biggs | (1,8,2,2,2,2) | 0.0 | 11 | 0.7289559 | 45 | NA |  |
| 5.Brown\&Dennis | (25,5,-5,-1) | 903.3648 | 29 | 903.2406 | 113 | 903.2343 | 2 |
| 6.El-Attar 5.1 | $(1,2)$ | 0.47042 | 10 | 0.4704267 | 6 | 0.4704247 | 8 |
| 7.El-Attar 5.2 | (1,1,1) | 7.902733 | 13 | 7.904731 | 22 | 7.894227 | 5 |
| 8.Madsen | $(3,1)$ | 1.0 | 16 | 1.000010 | 11 | 1.000002 | 13 |
| 9. Osborne 1 | (0.5, 1.5,-1,0.01, 0.02) | 0.0293912 | 14 | 0.0293914 | 10 | 0.8203727 | 55 |
| 10.Osborne 2 | $\begin{gathered} \text { (1.3,0.65,0.65,0.7,0.6 } \\ 3,5,7,2,4.5,5.5) \end{gathered}$ | F |  | F |  | - F |  |
| 11.Powell | (3,-1, 0, 1) | $0.1272 \mathrm{e}-07$ | 15 | F |  | 0.29039e-08 | 3 |
| 12.Rosenbrock | (-1.2,1) | 0.0 | 15 | F |  | $0.506642 \mathrm{e}-06$ | 51 |
| 13.Watson | (1,1,1,1) | 0.6487749 | 25 | 1.278432 | 5 | 0.6018584 | 24 |
| 14.Wood | (0,0,0,0) | 0.0 | 7 | 0.000003 | 10 | 0.0 | 25 |

See Appendix A for a detailed description of the test problems.

Findicates the algorithm failed to meet convergence criteria for the problem.

NA indicates results are not available for this entry.

Table 5.2
Optimal Points for Several Test Problems

| Example | Interior Point Algorithm | Osborne-Watson Algorithm | Busovaca Algorithm |
| :---: | :---: | :---: | :---: |
| 1.Wormersley | -6.74166, 4.59299 | -7.29804, 4.74178 | NA |
| 2.Bard | $0.10095,1.52545,1.97182$ | $0.10094,1.52516,1.97211$ | $0.10094,1.52513,1.97214$ |
| 3.Beale | 3, 0.5 | 3, 0.5 | 2.99999, 0.49999 |
| 4a.Biggs | F | F | F |
| b.Biggs | $1,10,1,5,4,3$ | $1.82143,81.94978,2.27882$ | NA |
| 5.Brown\&Dennis | $\begin{gathered} -9.70273,11.74096,-0.44204 \\ 0.55827 \end{gathered}$ | $\begin{gathered} -10.0227,11.91354,-0.44026 \\ 0.55823 \end{gathered}$ | $\begin{gathered} -10.2236,11.90843,-0.45804 \\ 0.58032 \end{gathered}$ |
| 6.El-Attar 5.1 | $2.84250,1.92018$ | 2.84250, 1.92018 | 2.84250, 1.92018 |
| 7.El-Attar 5.2 | 0.53558, -0.00139, 0.02871 | 0.53148, $-0.00004,0.02751$ | $0.53606,0.0,0.00319$ |
| 8.Madsen | $0.0,0.00016$ | 0.0,0.0022 | 0.0, -0.00205 |
| 9.Osborne 1 | $\begin{gathered} 0.37706,2.19244,-1.72549 \\ 0.01332,0.02129 \end{gathered}$ | $\begin{gathered} 0.37706,2.19246,-1.72552 \\ 0.01332,0.02129 \end{gathered}$ | $\begin{gathered} 1.06716,1.80257,-1.80731 \\ 0.00345,0.00109 \end{gathered}$ |
| 10.Osborne 2 | F | F | F |
| 11.Powell | $0.7268 \mathrm{e}-04,-0.7268 \mathrm{e}-05$ | F | $0.5588 \mathrm{e}-08,-0.3725 \mathrm{e}-09$ |
|  | $0.1162 e-04,0.1163 \mathrm{e}-04$ |  | 0.1250e-08, $0.1716 \mathrm{e}-08$ |
| 12.Rosenbrock | 1.0, 1.0 | F | 0.99999, 0.99999 |
| 13.Watson | $\begin{aligned} & -0.37526,1.14089 \\ & -0.42239,0.39683 \end{aligned}$ | $\begin{aligned} & -0.23584,1.03241 \\ & -0.22747,0.41384 \end{aligned}$ | $\begin{aligned} & -0.44271,1.19321 \\ & -0.47676,0.38449 \end{aligned}$ |
| 14.Wood | $1.0,1.0,1.0,1.0$ | $1.0,1.0,1.0,1.0$ | $1.0,1.0,1.0,1.0$ |

## Appendix A

Test problem 1 (Wormersley, 1986)
The results of temperature accelerated life tests on electrical insulation in 40 motorettes are recorded in Table A.1. This data is originally from Schmee and Hahn (1979). Ten motorettes were tested at each of four temperatures. Testing was terminated at different times at each temperature. The model used to fit the data is

$$
\log _{10} H=x_{1}+\frac{1000 x_{2}}{(T+273.2)}+\epsilon
$$

where $H$ is the failure time and $T$ is the temperature.
Table A. 1 : Data for motorettes example

|  |  | Test temperature $T^{\circ} \mathrm{C}$ |  |  |
| :---: | :---: | :---: | :---: | :---: |
|  | 150 | 170 | 190 | 200 |
| Failure times H in hours |  | 1764 | 408 | 408 |
|  |  | $2772^{*}$ | 408 | 408 |
|  |  | 3444 | 1344 | 504 |
|  |  | 3542 | 1344 | 504 |
|  |  | 3780 | 1440 | 504 |
|  |  | 4860 |  |  |
|  |  | 5196 |  |  |
| Termination time $\bar{H}$ | 8064 | 5448 | 1680 | 528 |
|  | 10 units | 3 units | 5 units | 5 units |
|  |  |  |  |  |

* Wormersley gives the second failure time at $170^{\circ}$ as 2722 , but his results are consistent with the values recorded here from Schmee and Hahn.

At each temperature there is an upper bound $\bar{H}$ (the time at which testing was stopped) on the observed failure times, so the algorithms of the observed failure times are given by

$$
\min \left(\log _{10} \bar{H}, x_{1}+\frac{1000 x_{2}}{(T+273.2)}+\epsilon\right) .
$$

Test problem 2 (Bard, 1970)

$$
f_{i}(x)=y_{i}-\left(x_{1}+\frac{u_{i}}{v_{i} x_{2}+w_{i} x_{3}}\right)
$$

where $i=1,2, \cdots, 15, u_{i}=i, v_{i}=16-i, w_{i}=\min \left(u_{i}, v_{i}\right)$, and

| i | $y_{i}$ | i | $y_{i}$ | i | $y_{i}$ |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 1 | 0.14 | 6 | 0.32 | 11 | 0.73 |
| 2 | 0.18 | 7 | 0.35 | 12 | 0.96 |
| 3 | 0.22 | 8 | 0.39 | 13 | 1.34 |
| 4 | 0.25 | 9 | 0.37 | 14 | 2.10 |
| 5 | 0.29 | 10 | 0.58 | 15 | 4.39 |

Test problem 3 (Beale, 1958)

$$
f_{i}(x)=y_{i}-x_{1}\left(1-x_{2}^{i}\right),
$$

where $i=1,2,3, y_{1}=1.5, y_{2}=2.25$ and $y_{3}=2.625$.

Test problem 4 (Biggs, 1971)

$$
f_{i}(x)=x_{3} \exp \left(-t_{i} x_{1}\right)-x_{4} \exp \left(-t_{i} x_{2}\right)+x_{6} \exp \left(-t_{i} x_{5}\right)-y_{i}
$$

where $i=1, \cdots, 13, t_{i}=(0.1) i$ and

$$
y_{i}=\exp \left(-t_{i}\right)-5 \exp \left(-10 t_{i}\right)+3 \exp \left(-4 t_{i}\right)
$$

Test problem 5 (Brown and Dennis, 1971)

$$
f_{i}(x)=\left(x_{1}+t_{i} x_{2}-\exp \left(t_{i}\right)\right)^{2}+\left(x_{3}+x_{4} \sin \left(t_{i}\right)-\cos \left(t_{i}\right)\right)^{2}
$$

Test problem 6 (El-Attar 5.1, 1979)

$$
\begin{aligned}
& f_{1}(x)=x_{1}^{2}+x_{2}-10 \\
& f_{2}(x)=x_{1}+x_{2}^{2}-7 \\
& f_{3}(x)=x_{1}^{2}-x_{2}^{3}-1
\end{aligned}
$$

Test problem 7 (El-Attar 5.2)

$$
\begin{aligned}
& f_{1}(x)=x_{1}^{2}+x_{2}^{2}+x_{3}^{2}-1 \\
& f_{2}(x)=x_{1}^{2}+x_{2}^{2}+\left(x_{3}-2\right)^{2} \\
& f_{3}(x)=x_{1}+x_{2}+x_{3}-1 \\
& f_{4}(x)=x_{1}+x_{2}-x_{3}+1 \\
& f_{5}(x)=2 x_{1}^{3}+6 x_{2}^{2}+2\left(5 x_{3}-x_{1}+1\right)^{2} \\
& f_{6}(x)=x_{1}^{2}-9 x_{3}
\end{aligned}
$$

Test problem 8 (Madsen, see Overton and Murray, 1981)

$$
\begin{aligned}
& f_{1}(x)=x_{1}^{2}+x_{2}^{2}+x_{1} x_{2} \\
& f_{2}(x)=\sin \left(x_{1}\right) \\
& f_{3}(x)=\cos \left(x_{2}\right)
\end{aligned}
$$

Test problem 9 (Osborne 1, 1972)

$$
f_{i}(x)=y_{i}-\left(x_{1}+x_{2} \exp \left(-t_{i} x_{4}\right)+x_{3} \exp \left(-t_{i} x_{5}\right)\right)
$$

where $i=1,2, \cdots, 33, t_{i}=10(i-1)$, and

| i | $y_{i}$ | i | $y_{i}$ | i | $y_{i}$ |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 1 | 0.844 | 12 | 0.718 | 23 | 0.478 |
| 2 | 0.908 | 13 | 0.685 | 24 | 0.467 |
| 3 | 0.932 | 14 | 0.658 | 25 | 0.457 |
| 4 | 0.936 | 15 | 0.628 | 26 | 0.448 |
| 5 | 0.925 | 16 | 0.603 | 27 | 0.438 |
| 6 | 0.908 | 17 | 0.580 | 28 | 0.431 |
| 7 | 0.881 | 18 | 0.558 | 29 | 0.424 |
| 9 | 0.850 | 19 | 0.538 | 30 | 0.420 |
| 10 | 0.818 | 20 | 0.522 | 31 | 0.414 |
| 11 | 0.784 | 21 | 0.506 | 32 | 0.411 |
|  | 0.751 | 22 | 0.490 | 33 | 0.406 |

Test problem 10 (Osborne 2)

$$
\begin{aligned}
f_{i}(x)= & y_{i}-\left(x_{1} \exp \left(-t_{i} x_{5}\right)+x_{2} \exp \left(-\left(t_{i}-x_{9}\right)^{2} x_{6}\right)\right. \\
& \left.+x_{3} \exp \left(-\left(t_{i}-x_{10}\right)^{2} x_{7}\right)+x_{4} \exp \left(-\left(t_{i}-x_{11}\right)^{2} x_{8}\right)\right)
\end{aligned}
$$

where $i=1,2, \cdots, 65, t_{i}=(i-1) / 10$, and

| i | $y_{i}$ | i | $y_{i}$ | i | $y_{i}$ |
| :---: | :---: | :---: | :---: | :---: | :--- |
| 1 | 1.366 | 23 | 0.694 | 45 | 0.672 |
| 2 | 1.191 | 24 | 0.644 | 46 | 0.708 |
| 3 | 1.112 | 25 | 0.624 | 47 | 0.633 |
| 4 | 1.013 | 26 | 0.661 | 48 | 0.668 |
| 5 | 0.991 | 27 | 0.612 | 49 | 0.645 |
| 6 | 0.885 | 28 | 0.558 | 50 | 0.632 |
| 7 | 0.831 | 29 | 0.533 | 51 | 0.591 |
| 8 | 0.847 | 30 | 0.495 | 52 | 0.559 |
| 9 | 0.786 | 31 | 0.500 | 53 | 0.597 |
| 10 | 0.725 | 32 | 0.423 | 54 | 0.625 |
| 11 | 0.746 | 33 | 0.395 | 55 | 0.739 |
| 12 | 0.679 | 34 | 0.375 | 56 | 0.710 |
| 13 | 0.608 | 35 | 0.372 | 57 | 0.729 |
| 14 | 0.655 | 36 | 0.391 | 58 | 0.720 |
| 15 | 0.616 | 37 | 0.396 | 59 | 0.636 |
| 16 | 0.606 | 38 | 0.405 | 60 | 0.581 |
| 17 | 0.602 | 39 | 0.428 | 61 | 0.428 |
| 18 | 0.626 | 40 | 0.429 | 62 | 0.292 |
| 19 | 0.651 | 41 | 0.523 | 63 | 0.162 |
| 20 | 0.724 | 42 | 0.562 | 64 | 0.098 |
| 21 | 0.649 | 43 | 0.607 | 65 | 0.054 |
| 22 | 0.649 | 44 | 0.653 |  |  |
|  |  |  |  |  |  |

Test problem 11 (Powell, 1962)

$$
\begin{aligned}
& f_{1}(x)=x_{1}+10 x_{2} \\
& f_{2}(x)=5^{1 / 2}\left(x_{3}-x_{4}\right) \\
& f_{3}(x)=\left(x_{2}-2 x_{3}\right)^{2} \\
& f_{4}(x)=10^{1 / 2}\left(x_{1}-x_{4}\right)^{2}
\end{aligned}
$$

Test problem 12 (Rosenbrock, 1960)

$$
\begin{aligned}
& f_{1}(x)=10\left(x_{2}-x_{1}^{2}\right) \\
& f_{2}(x)=1-x_{1}
\end{aligned}
$$

Test problem 13 (Watson, see Kowalik and Osborne, 1968)

$$
f_{i}(x)=\sum_{j=2}^{n}(j-1) x_{j} t_{i}^{j-2}-\left(\sum_{j=1}^{n} x_{j} t_{i}^{j-1}\right)^{2}-1
$$

where $i=1, \cdots, 29, t_{i}=i / 29, f_{30}(x)=x_{1}$ and $f_{31}(x)=x_{2}-x_{1}^{2}-1$.

Test problem 14 (Wood, see Colville, 1968)

$$
\begin{aligned}
& f_{1}(x)=10\left(x_{2}-x_{1}^{2}\right) \\
& f_{2}(x)=1-x_{1} \\
& f_{3}(x)=90^{1 / 2}\left(x_{4}-x_{3}^{2}\right) \\
& f_{4}(x)=1-x_{3} \\
& f_{5}(x)=10^{1 / 2}\left(x_{2}+x_{4}-2\right) \\
& f_{6}(x)=10^{-1 / 2}\left(x_{2}-x_{4}\right)
\end{aligned}
$$

## Appendix B

```
"nlrq" <-
function(x, y, model, t, theta, k = 2, eps = le-06, big = le+20, eta = 0.97)
(
This is a function to compute nonlinear l1 estimate.
    Input
            model - user-provided function which returns components
                        f=(f_i (x_i , t)
                        J=(grad f_i )
            t - vector of initial values of the unknown parameters
            theta - desired quantile
            k - number of Meketon's iterations used to calculate the
                        dual step
            eps - small positive number
            big - big positive number
            eta - 0.97
    Output
            t - vector of estimated parameters
            f - function value at minimum
            n <- length(y)
            zero <- rep(0, n)
            w <- zero
            d <- rep (l, n)
            m <- model (x, y, t, theta)
            snew <- sum(abs(m$f))
            sold <- big
            while(sold - snew > eps) (
                z <- mekrq(m$J, m$f, w, theta, k, int = F, eps, big, eta)
                step <- z$coef
                # Calculate an optimal step length lambda
                lambda <- step.length(t,step)
                t <- t + lambda * step
                m <- model(x, y, t, theta)
                sold <- snew
                snew <- sum(abs(m$f))
                w <- z$w
                w <- lsfit(m$J, w, int = F)$resid
                if(max(abs(w)) >= l)
                                    w <- w/(max(abs (w)) + eps)
            }
            f <- snew
            return(t, f)
)
"mekrq* <-
function(x,y, w, theta, kmax = 1000, int = T, eps, big, eta)
{
# Compute linear regression quantile estimate (Meketon, 1985).
# However, note that }W\mathrm{ is not initialized and the maximum number of
# iteration is given by kmax.
#
    if(int == T) x <- cbind(1, x)
    sr <- big
    k <- l
    while(k <= kmax & sr - crossprod(y, w) > eps) {
        d <- pmin(theta - w, l - theta + w)
        z<- lsfit(x, y, d^2, int = F)
```


## sr <- sum(abs(z\$resid))

$k<-k+1$
$s<-\quad$ \$resid * $d^{\wedge} 2$
alpha <- max $(\operatorname{pmax}(s /(t h e t a-w) .-s /(1-t h e t a+w)))$
$w<-w+(e t a / a l p h a) * s$
\}
coef $<-$ z\$coef
return(coef. W)

This procedure is a function to compute nonlinear regression quantile estimate.

## PROC NLRQ

FORMAT
$\{t, f\}=n l r q(x, y, \& m o d e l, s t a r t$, theta,k,eps,big,eta )
INPUT
model - user-provided function which returns components $\mathrm{f}=\left(\mathrm{f}\right.$ _i $\left.\left(\mathrm{x} \_i, \mathrm{t}\right)\right)$ $J=\left(g r a d f_{-}^{-i}\right)$
start - vector of initial values of the unknown parameters
theta - desired quantile
$k$ - number of Meketon's iterations used to calcuate the dual step
eps - small positive number (le-06)
big - big positive number (le+20)
eta - 0.97
OUTPUT
$t$ - vector of estimated parameters
f - function value at minimum

The following procedure is a function to compute linear regression quantile estimate (Meketon, 1985). However, note that $W$ is not initialized and the maximum number of iterations is given by kmax.

```
proc(2) = mekrq(x,y,w,theta,kmax,inter,eps,big,eta);
```

    local sr,k,wy,wx,d,t,r,s,alpha;
        if inter eq 1;
            \(\mathrm{x}=\) ones(rows(x),1)-x;
        endif;
        sr = big;
        \(\mathrm{k}=1\);
        do while \(k\) <= kmax and sr - \(y^{\prime *}\) w > eps;
            \(\mathrm{d}=\operatorname{minc}\left((\mathrm{theta}-\mathrm{w})^{\prime} \mid\left(1-\mathrm{theta}^{\prime} \mathrm{w}\right)^{\prime}\right)\);
            \(\mathrm{d}=\mathrm{vec}(\mathrm{d})\);
            \(\mathrm{wx}=\mathrm{x} . * \mathrm{~d}\);
            wy = \(\mathrm{y} . * \mathrm{~d}\);
            \{ t \} = olsqr(wy,wx);
            \(\mathrm{r}=\mathrm{y}-\mathrm{x}\) *;
            sr \(=\) sumc(abs(r));
            \(\mathrm{k}=\mathrm{k}+1\);
            s = r.*d^2;
            alpha \(=\operatorname{maxc}\left(\operatorname{maxc}\left(\left(s . /(\text { theta-w) })^{\prime} \mid\right.\right.\right.\)
                        (-s./(1-theta+w))');
            w = w + eta/alpha*s;
        endo;
        retp(t,w);
    endp;
    ```
proc(2) = nlrq(x,y,&model,t,theta,kmax,eps,big,eta);
    local n,d,w,f,J,snew,sold,step,lambda,t,t1,p,inter,
    model:proc;
    n = rows(y);
    w = zeros(n,1);
    d = ones(n,1);
    {f,J } = model(x,y,t);
    snew = sumc(abs(f));
    sold = big;
    inter = 0;
    do while sold - snew > eps;
        { step,w } = mekrq(J,f,w,theta,kmax,inter,eps,
                    big,eta);
        /* Calculate an optimal step length lambda */
        { lambda } = stepl(J,f,t,step,&model,eps);
        t = t + lambda*step;
        { f,J } = model(x,y,t);
        sold = snew;
        snew = sumc(abs(f));
        { t1,w,p } = olsqr2(w,J);
        if maxc(abs(w)) >= 1;
            w = w/(maxc(abs(w))+eps);
        endif;
    endo;
    retp(t,snew);
endp;
```


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