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AN IMPROVED DIRECT SEARCH NUMERICAL OPTIMIZATION PROCEDURE. (U)

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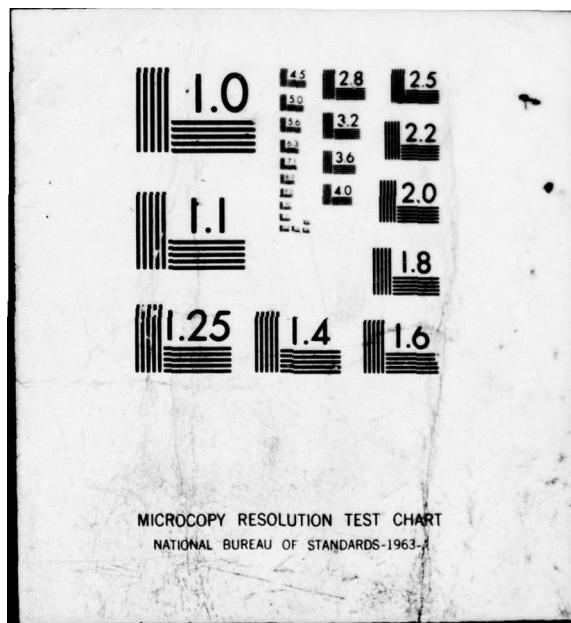


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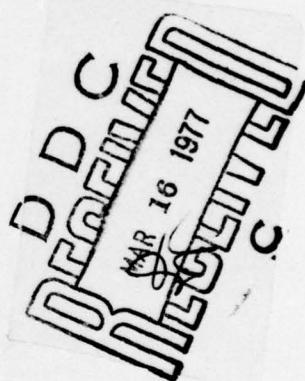
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AN IMPROVED DIRECT SEARCH  
NUMERICAL OPTIMIZATION PROCEDURE

by

Michael Pappas

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New Jersey Institute of Technology  
323 High Street  
Newark, New Jersey 07102

## ABSTRACT

An improved, nonlinear, constrained mathematical programming optimization algorithm is presented in this report. It couples a rotating coordinate pattern search with a feasible direction finding procedure used at points of pattern search termination. The procedure is compared with nineteen algorithms, including most of the popular methods, on ten test problems. These problems are such that the majority of codes failed to solve more than half of them. The new method proved superior to all others in the overall generality and efficiency rating, being the only one solving all problems. It was particularly effective on constrained problems where it was best in all rating categories.



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## NOMENCLATURE

$B_j$  = behavior of constraints  $g_j(x_i)$

$e_1, e_2, e_3$  = small arbitrary constants defining finite difference step size used to calculate the gradient of the objective function, nearness to constraints, and convergence criterion respectively.

$f_a, \bar{f}_a$  = efficiency ratings (eqs. 30 and 31)

$f(x_i)$  = objective function

$F(x_i)$  = composite objective function

$g_j(x_i)$  = constraint function

$I$  = number of variables

$J$  = number of constraints

$K_A^-, K_A^+$  = set of active lower and upper regional, constraints, respectively

$L_j$  = lower limit on the behavior of  $B_j$

$n_a, N_A$  = generality ratings (eq. 29)

$P(x_i)$  = penalty function used when a constraint is violated

$s_i$  = best feasible direction vector

$T_A$  = generality and efficiency rating (eq. 32)

$U_j$  = upper limit on the behavior of  $B_j$

$x_i$  = problem variables or variable vector

$\bar{x}_i$  = optimal variable values

$w_j$  = weighting factor for constraint  $g_j(x_i)$  used in the feasible direction finding problem

$\alpha$  = step size  
 $\alpha_{\min}$  = minimum step size  
 $\epsilon_j$  = constraint activity limits for the direction finding problem  
 $\sigma$  = objective function and variable in the feasible direction finding problem  
 $\nabla$  = gradient operator  
 $|\phi|$  = magnitude of  $\phi$

**Superscripts**

$l$  = lower limit  
 $r$  = at rth iteration  
 $t$  = local exploration base  
 $T$  = transpose of vector  
 $u$  = upper limit  
 $*$  = comparison quantity

## INTRODUCTION

A variety of methods involving ordinary and variational calculus, mathematical programming, and a number of special techniques, such as the fully stressed concept used in structural design, are available to treat optimal design problems. Among these methods, the mathematical programming procedures appear to have the broadest range of application [1]<sup>1</sup>. Such methods are flexible, easy to adapt, and can offer "automatic" optimal computer solutions [2].

Mathematical programming methods solve, or approach the solution to, the problem: Find those values of the variables,  $x_i$ , that minimize (or maximize) the objective function

$$f(x_i) \quad i = 1, 2, \dots I \quad (1)$$

such that all constraints

$$g_j(x_i) \leq 0 \quad j = 1, 2, \dots J \quad (2)$$

are satisfied. The objective function (often called the merit or payoff function) quantitatively defines the merit of the design as a function of the design variables,  $x_i$ . Inequality constraints are used to control or define the behavior of the design (behavior constraints).

Inequality constraints can also be used to limit the values of the variables within specified limits (regional constraints).

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<sup>1</sup>Numbers in brackets designate References at end of report.

It is usually convenient to treat such constraints somewhat differently than behavior constraints. Regional constraints can be given in the form

$$x_i^l \leq x_i \leq x_i^u \quad (3)$$

where  $x_i^l$  and  $x_i^u$  are the lower and upper limits respectively on the variable  $x_i$ .

A number of effective methods exist for special forms of this problem, such as the unconstrained problem, the linear programming problem, and many quadratic programming problems. Relatively few effective methods are available, however, to treat the form most often encountered in design, the nonlinear constrained optimization problem. Even the better nonlinear programming methods have limitations on the size and complexity of problems they can solve within a justifiable amount of computer time. Large order, highly nonlinear systems requiring lengthy design computations can be quite formidable. A typical nonlinear mathematical programming procedure requires several hundred to several thousand sets of objective and constraint function evaluations to approach the optimum on most multivariable problems. Furthermore, reliability is a major problem with these methods [3].

The new Direct Search-Feasible Direction (DSFD) algorithm appears to offer superior performance with respect to speed and reliability [4]. It is the only procedure studied that solved all of Eason and Fenton's test problems [3-4].

No attempt was made in the original work on this algorithm, as reported in [4], to maximize its effectiveness. Rather the performance of the basic procedure was studied and presented.

It later became apparent that it was possible to substantially improve the speed of the procedure by eliminating needless computations and by other detailed improvements.

This report presents and evaluates the latest version of the DSFD algorithm where such improvements have been incorporated.

#### THE OPTIMIZATION ALGORITHM

##### The basic search

The "rotating coordinate" (RC) pattern search used in [5] and described in detail in [6] is employed as the basic optimal search technique. The usual pattern search strategy [7] is utilized in this procedure except that after establishing a pattern move, the first local exploration step is taken in the direction of the move and all other local steps are taken normal to this direction rather than in the direction of the variable coordinates, as in the usual pattern search procedure. The movement of this search on a hypothetical composite objective surface using the penalty function of [5] is shown in Fig. 1. It may be seen from Path 1 that in the feasible region the search tends to move essentially in the direction of the gradient, while near the feasible-infeasible boundary it moves easily along the ridge. The local steps taken

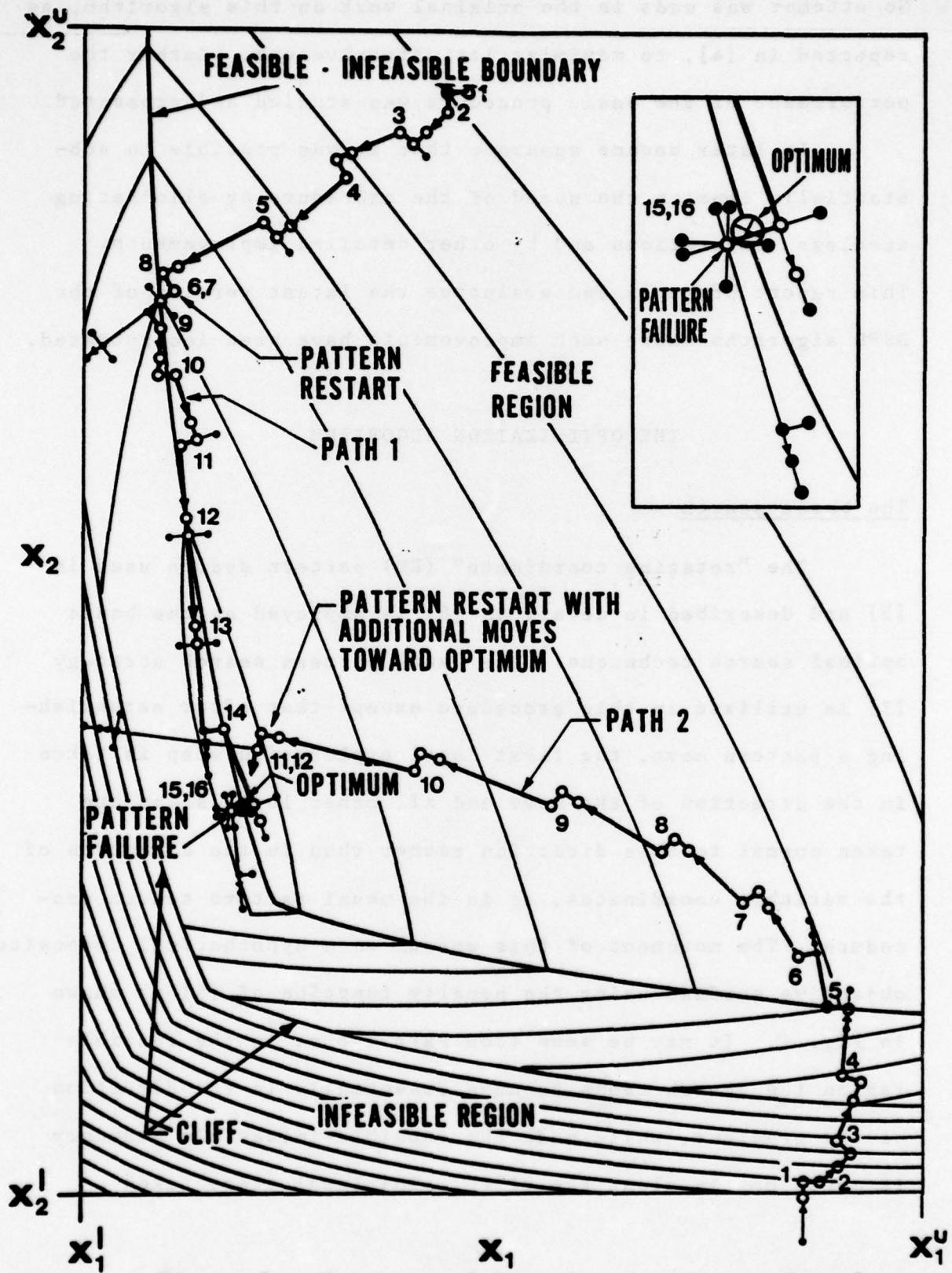


Figure 1. Movement of Rotating Coordinate Search

normal to the pattern move provide the turning component necessary to maintain efficient movement. It should be noted that the method does not require a feasible starting point as do most procedures (see Path 2). The closed dots in this figure indicate unsuccessful steps while the open dots are successful moves. The digits, associated with some open dots, identify the pattern base numbers.

This basic search procedure was found to be substantially superior to the original Hooke and Jeeves search in earlier comparison studies [5-7].

Since the basic search scheme is formulated to treat unconstrained optimization problems, the constrained problem of (1) - (3) is transformed to the form; Find

$$F(\bar{x}_i) = \min F(x_i) \quad (4)$$

where

$$F(x_i) = f(x_i) + P(x_i). \quad (5)$$

Here  $P(x_i)$  is the largest of the penalties  $p_j(x_i)$  which are of the form

$$p_j(x_i) = -\lambda_j[g_j(x_i)] \quad (6)$$

where

$$\begin{aligned} [\phi] &= \phi & \phi < 0 \\ [\phi] &= 0 & \phi > 0. \end{aligned} \quad (7)$$

When  $0 > g_j > -e_2$ , then

$$\lambda_j = 2[f(x_i) - f(x_i + \Delta x_i)]/[g_j(x_i) - g_j(x_i + \Delta x_i)] \quad (8)$$

where

$$\Delta x_1 = e_1 \nabla f(x_1^t) / |\nabla f(x_1^t)|. \quad (9)$$

The quantity  $e_1$  is an arbitrary small real number representing the size of the step,  $\Delta x_1$ , taken in the direction of the gradient of  $f(x_1)$  where this gradient is evaluated at the local exploration base point  $x_1^t$ . It may be seen that a penalty is used only if a constraint is violated. If the violation of any of these constraints is greater than an arbitrarily specified amount  $e_2$ , then  $\lambda_j$  is not calculated from eqn (8) but is made arbitrarily large. That is, when

$$\left. \begin{array}{l} g_j(x_1) < -e_2 \\ \lambda_j = K \end{array} \right\} \quad (10)$$

where  $K$  is an arbitrary large number.

The justification for the use of this penalty function form is discussed in detail in Ref. [5].

Violations of eqns (3) are handled by prohibiting moves that violate any of these equations.

All of the constraints are evaluated only at each temporary base (points reached after a pattern move). For all other points in the search only nearby constraints, this is those where

$$g_j(x_f) \geq e_2 \quad (11)$$

are evaluated since it is presumed that only these constraints need be considered.

### The Secondary Search

Unfortunately, the pattern search, even with the rotating coordinate improvement, is not particularly reliable [5,6]. Thus, some method is needed to determine if the point of pattern search termination is optimal and if it is not, to determine a direction in which to restart the search. Pappas and Amba-Rao [8] and Siddal [9] apply a secondary search at points of pattern search termination. Unfortunately, the earlier secondary search strategies although more powerful locally than the pattern search do not provide a rigorous procedure for confirming optimality.

The direction finding procedure of Zoutendijk [10] can, however, be used to confirm the optimality of a point of pattern search termination and determine a direction in which to restart the pattern search if the point is not optimal. The direction finding problem can be stated: Given the set  $x_i$ , find the set  $s_i$  that results in a

$$\max \sigma \quad (12)$$

for which

$$\sigma > 0 \quad (13)$$

$$(s_i)^T \nabla f(x_i) + \sigma < 0 \quad (14)$$

$$(s_i)^T \nabla g_j(x_i) + w_j \sigma \leq 0 \quad j \in J_a \quad (15)$$

$$-1 \leq s_k \leq 0 \quad k \in K_a^- \quad (16)$$

$$0 \leq s_k \leq 1 \quad k \in K_a^+. \quad (17)$$

Here  $(s_i)^T$  indicates the transpose of vector  $s_i$ ,  $w_j$  is a weighing parameter, the set  $J_a$  contains the active constraints

$$g_j(x_1) > -\epsilon_j \quad (18)$$

where  $\epsilon_j$ , an array of small arbitrary positive constants defining "activity," and  $K_a^- K_a^+$  constitute the active upper and lower side constraints, respectively.

Zoutendijk [10] shows that if the solution  $s_i$  is a null vector, then the point is a local optimum and if not then the direction of  $s_i$  is the best feasible direction.

Equations (12)-(17) constitute a linear programming problem with the variable  $s_i$  and  $\sigma$ . Equation (12) is the objective function and the remaining equations the constraints. The solution  $s_i$  can be obtained reliably and efficiently using any suitable linear programming method such as the simplex procedure.

#### The Optimization Procedure.

The general procedure here is similar to that used in [3] and [6] except the pattern search and optimality checks used in [5] are replaced by those described previously. An arbitrary initial starting point  $x_i^0$  is selected,  $F^*$  defined as equal to  $F(x_i^0)$ , and a quantity  $D^*$  set equal to unity. The optimal search is then started using the RC procedure with a step size  $\alpha^0$  and continued until it terminates at some point  $x_i^r$ , where the secondary search is invoked.

If the pattern search terminates at a point in the feasible region, that is, where equation (3) is satisfied for all  $j$ , the direction finding problem is formulated with  $w_j = 1$  and solved at the point  $x_i^r$ . The components of the gradients are

generated using simple forward differences and thus no analytical derivatives are employed.

If the pattern search terminates in the infeasible region at a point near the feasible-infeasible boundary, that is, if some

$$e_2 \geq g_j(x_i^r) \geq 0 \quad (19)$$

(see Base 15 of Path 1 in Fig. 1) a weighting parameter of  $w_j = 100$  is used to help drive the search to the feasible region.

In the case where the pattern search terminates in the infeasible region away from the boundary, that is, where any constraint equation violates the left side of (19), the search is abandoned since in this instance the pattern search has failed to generate a feasible point, even with the use of a large penalty parameter.

Once the direction vector  $s_i$  is computed, the quantity  $F(x_i^{r+1})$  is evaluated, where

$$x_i^{r+1} = x_i^r + \Delta x_i \quad (20)$$

and

$$\Delta x_i = \alpha^r s_i / \left| \left[ \sum_{m=1}^I s_m^2 \right]^{1/2} \right| \quad \text{if } s_i \neq 0 \quad (21)$$

Here  $\alpha^r$  is the step size used for the pattern search just terminated. If any constraint not considered in the formulation of the direction finding problem has been violated at  $x_i^{r+1}$ , then the direction finding problem is reformulated considering these constraints and a new  $s_i$  and  $F(x_i^{r+1})$  computed.

If  $s_i = 0$  the activity definition limits  $\epsilon_j$  and the step size  $\alpha^r$  are halved. If the set  $J_a$  changes as a result of the

change in  $\epsilon_j$  due to one of the constraints becoming deactivated as a result of the reduction in the activity specification the direction finding problem is reformulated and solved. Otherwise the activity limit and step size is further reduced. This process is repeated until a nonzero solution of  $s_i$  is obtained or until

$$\alpha^r \leq \alpha_{\min} \quad (22)$$

where  $\alpha_{\min}$  is a specified minimum basic step size.

Now if

$$F(x_n^{r+1}) < F(x_n^r) \quad (23)$$

the RC search is restarted using  $x_n^{r+1}$  as the new base point in the pattern search strategy. If equation (22) is not satisfied, then  $\alpha$  is refined as

$$\alpha^{r+1} = \alpha^r / 2 \quad (24)$$

Now if

$$\alpha^{r+1} < \alpha_{\min} \quad (25)$$

or if

$$\left. \begin{array}{l} D^r \leq e_3 \\ \text{and} \\ D^r < D^* \end{array} \right\} \quad (26)$$

where  $D^r = [F^* - F(x_1^r)]/F^*$

and where  $\alpha_{\min}$  and  $e_3$  are arbitrary small positive variable and objective function convergence constants, respectively, the search is abandoned and the point is assumed to be sufficiently near optimum. Otherwise a new  $x_i^{r+1}$  is defined by equations (20-21) with  $\alpha^r$  replaced by  $\alpha^{r+1}$  until equation (23), (25), or (26) is satisfied.

Once equation (23) is satisfied, then the RC search is restarted, with  $x_i^{r+1}$  as a new base point in the pattern search strategy, with step size  $\alpha^{r+1}$ ,  $F^* = F(x_i^r)$ , and  $D^* = D^r$ . If, however, two successive step size reductions fail to satisfy equation (23), then the entire optimization procedure is restarting from point  $x_i^r$ , with step size  $\alpha^{r+1}$ ,  $F^* = F(x_i^r)$ , and  $D^* = D^r$  and repeated until equation (25) or (26) is satisfied.

It may be seen that the step defined by equation (21), taken when  $s_i$  is nonzero, is a move in the best feasible direction. Thus, if equation (23) is not satisfied by this move, it indicates that the step size is too large. The step size is then reduced in an effort to locate a better point, first by taking a smaller step in the  $s_i$  direction, and in the event this fails, by attempting to restart the pattern using this smaller step. The process is repeated until convergence of the objective function is achieved or a minimum step size is reached.

A zero  $s$  vector indicates the presence of an optimum. Since, however, in the consideration of the active constraints,  $\epsilon_j \neq 0$ , the point  $x_i^r$  may be merely near, rather than at, the optimum. Thus, the activity limit is reduced when a null  $s$  vector solution is encountered and the Feasible Direction problem reexamined to determine if the point is indeed an optimum. Since in this procedure

$$\epsilon_j = C_1 \alpha^r / \alpha^0 \quad (27)$$

when equation (22) is satisfied it is assumed that the  $\epsilon_j$  are sufficiently small to allow the point  $x_i^r$  to be considered an optimum.

This procedure is the same as that of reference [4] except for the following differences. In reference [4];

1. the values of  $\nabla f(x_i)$  were evaluated at each point calling for the computation of a penalty parameter  $\lambda_j$ . Thus  $\nabla f(x_i)$  is often calculated several times during a local exploration. Since, however, only an estimate of  $\nabla f$  is required here,  $\nabla f(x_i^T)$  is computed here only once during local explorations. A substantial reduction in the number of required objective function evaluations required for penalty evaluations is thereby achieved.
2. all constraints were evaluated at all points. Since, however, only those constraints which are active or near active influence movement toward the optimum at a given point, the procedure used here only considers such constraints. All constraints are evaluated only periodically (after each pattern jump) to monitor constraint activity.
3. at points of pattern search termination in the unfeasible region near the feasible-infeasible boundary the pattern search was restarted using  $\lambda_j = K$ , the large penalty parameter, in an attempt to drive the search to the feasible region. In the strategy used here, movement in a direction finding problem with  $W_j = 100$ , is utilized.
4. slightly different version of the Rotating Coordinate Pattern Search is used. The version employed here contains several improvements.

## COMPARISON STUDY

The work of Eason and Fenton [3] forms the primary basis for the comparison between the algorithm presented here and other optimization procedures. A FORTRAN IV code called CADOP3 based on this algorithm was developed and used on the test problems given in [3]. The code is operational with IBM FORTRAN G and H and the UNIVAC TDOS and TSOS systems. This code required the user to supply only the objective and constraint functions, the initial values of the variables, and regional constraint values, if used. The convergence criteria and initial step size may also be specified.

If no step size is specified, the size is internally generated. In this program the initial step size is selected such that at the starting point a step of size  $\alpha^0$  in the  $V_f$  direction produces a one percent change in the value of the objective function. The internally generated step size is constrained so that  $\alpha^0 > 0.005$ . The minimum step size is defined as  $\alpha_{\min} = \alpha^0/1000$  unless specified otherwise by the user.

Nondimensional constraint equations of the form

$$\left. \begin{aligned} g_j(x_i) &= (B_j - U_j)/|U_j| \leq 0 \text{ when } U_j \neq 0 \text{ and } B_j \neq 0 \\ \text{or} \\ g_j(x_i) &= (L_j - B_j)/|L_j| \leq 0 \text{ when } L_j \neq 0 \text{ and } B_j \neq 0 \end{aligned} \right\} \quad (28)$$

are used. Otherwise a dimensional form is used.

Here  $B_j$  represents the constraint "behavior" and  $U_j$  and  $L_j$  the upper and lower limits on this behavior, respectively. The user supplies the expressions for  $f(x_i)$  and the expressions or parameters defining  $B_j$ ,  $U_j$ , and  $L_j$ .

A constraint given as  $0 \leq b(x_i) \leq A$  would be written as two constraint equations. For example, one could be of the form of the first of equations (28) where  $B_1 = b(x_n)$  and  $U_1 = A$ . The second would be of the form  $g_2 = -b(x_n)$  since  $L_1 = 0$ .

The program uses  $e_1 = 0.0001$ ,  $e_2 = 0.10$  and  $e_3 = 1 \times 10^{-7}$ . Multiple starting points can be tried in a single computer run to provide additional optimality confirmation. Such a technique is recommended even though the algorithm presented here seems reasonably reliable, since no nonlinear method can be considered absolutely reliable.

All ten problems of reference [3] were run with the CADOP3 code from the starting points specified in reference [3]. The internally generated step sizes were used for all problems and no algorithm control constants were changed during the study. Thus, there was no tuning of the code to individual problems. All problems were run on an IBM 370 model <sup>168</sup>91 using the "G" level compiler and double precision. Thus, the runs closely simulate those of the Eason and Fenton study which likewise used an IBM machine with the same level compiler and precision specification. Thus the comparisons can be considered accurate.

Table 1 briefly describes the codes studied in reference [3] with the addition of the DSDA code, the earlier CADOP2 code, and the CADOP3 code studied here. Additional details on these codes can be found in references [4] and [11]. Table 2 presents results of the performance of these codes on the ten test problems in the form of a success matrix. The numerical entries indicate the normalized time required for solution [3], the symbol

TABLE 1  
CODE DESCRIPTIONS

| Name   | Description   | Algor class* |
|--------|---|--------------|
| ADRANS | Random search followed by pattern moves   | DS           |
| CLIMB  | Rosenbrock search   | DS           |
| DAVID  | Davidon-Fletcher-Powell with numerical derivatives                              | GF           |
| DFMCG  | Fletcher-Reeves conjugate gradient method with secant approximation derivatives | G            |
| DFMFP  | Davidon-Fletcher-Powell with secant approximation derivatives                   | G            |
| FMIND  | Hook & Jeeves pattern search  | DS           |
| GRADA4 | Steepest descent method   | G            |
| GRID1  | Grid and star network search, with shrinkage                                    | AR           |
| MEMGRD | Davidon-Fletcher-Powell with retained step length information                   | GF           |
| NMSERS | Simplex search  | DS           |
| PATSH  | Modified pattern search   | DS           |
| PATRNO | Modified pattern search, dome strategy  | DS           |
| PATRN1 | Modified pattern search, ridge strategy   | DS           |
| RANDOM | Random search with shrinkage  | AR           |
| SEEK1  | Pattern search followed by random search  | DS           |
| SEEK3  | Modified pattern search   | DSF          |
| SIMPLX | Modified simplex search   | DSF          |
| DSDA   | Modified pattern search followed by Mugele's search                             | DS           |
| CADOP2 | Direct search-feasible direction algorithm                                      | DS           |
| CADOP3 | Refined direct search-feasible direction algorithm                              | DS           |

\*DS = direct search, DSF = direct search employing SUMT strategy and penalty function, G = gradient procedure, GF = gradient procedure using SUMT strategy and penalty function. AR = area reduction method.

TABLE 2  
PERFORMANCE OF OPTIMIZATION CODES\*

| Problem No.   | 1     | 2     | 3     | 4     | 5     | 6     | 7     | 8     | 9    | 10   |
|---------------|-------|-------|-------|-------|-------|-------|-------|-------|------|------|
| Variables     | 5     | 3     | 5     | 4     | 2     | 2     | 3     | 2     | 2    | 4    |
| Constraints** | 10    | 2     | 6     | 0     | 0     | 1     | 2     | 0     | 6    | 0    |
| Code names    |       |       |       |       |       |       |       |       |      |      |
| ADRANS        | 2.64  | 0.458 | 1.65  | P     | 0.100 | 0.654 | 0.159 | 0.069 | P    | P    |
| CLIMB         |       | 0.188 | 0.84  | 0.015 | 0.005 |       |       | 0.007 |      |      |
| DAVID         |       |       | P     | 1.132 | 0.075 | 0.046 |       |       |      |      |
| DFMCG         |       |       | P     | 0.015 | P     |       |       |       |      |      |
| DFMFP         |       |       | P     | 0.038 | 0.250 |       | P     | 0.004 |      |      |
| FMIND         | 0.004 | P     |       | 0.140 |       | 0.003 | P     | 0.387 | P    | 3.74 |
| GRAD4         |       | P     | 0.283 | 0.393 |       | P     | 0.004 |       |      | P    |
| GRID1         |       | P     | 0.033 | P     |       | P     | 0.004 |       |      |      |
| MEMGRD        | 0.143 | 0.91  | 0.059 | 0.066 | 0.067 | P     | 0.037 | P     |      | P    |
| NMSERS        | 0.019 | 0.045 | 0.060 | 0.002 | 0.012 | P     | 0.007 | 0.39  | P    | P    |
| PATSH         | 1.78  | 0.220 | 1.00  | 0.013 | 0.018 | 0.020 | 0.010 | 0.009 | P    | 3.54 |
| PATRNO        | P     | P     | P     | 0.021 | P     |       | P     | 0.002 | P    | 1.95 |
| PATRN1        | P     | P     | P     | 0.008 | 0.002 | P     | 0.001 | 1.02  | P    | 1.20 |
| RANDOM        | P     | P     | P     | P     | 0.024 | 1.20  | 0.013 |       |      |      |
| SEEK1         | P     | P     | P     | P     | 0.010 | 0.010 | 0.007 | P     |      | 1.53 |
| SEEK3         | 0.102 | 0.14  | 0.212 | 0.035 | 0.191 | 0.141 | 0.013 | 4.20  | P    |      |
| SIMPLX        | 2.97  | 0.297 | 1.31  | 0.196 | 0.035 | 0.191 | 0.262 | 0.050 | P    |      |
| DSDA          | P     | 0.021 | 0.047 | 0.104 | 0.003 | 0.008 | 0.145 | 0.004 | 1.94 | 1.35 |
| CADOP2        | 0.307 | 0.016 | 0.090 | 0.075 | 0.008 | 0.011 | 0.069 | 0.003 | 1.55 | 1.73 |
| CADOP3        | 0.119 | 0.006 | 0.030 | 0.080 | 0.004 | 0.011 | 0.025 | 0.005 | 1.00 | 1.60 |

\*Numerical entry indicates normalized time required for solution, and P indicates progress toward a solution [3, 12].

\*\*Excluding regional constraints.

"p" indicates progress toward a solution, and a blank indicates failure. Definitions of "solution" and "progress" vary for each problem and are given in reference [11] as well as a detailed description of each problem.

The data from Table 2 may be applied to a number of rating schemes for comparing the codes tested. Table 3, presents relative rankings of the codes using some of the rating criteria of reference [3]. The rating equations used are as follows:

The number of problems solved by code "a" is  $n_a$  and

$$N_a = n_a + n'_a / 2 \quad (29)$$

where  $n'_a$  is the number of problems in which a P rating was achieved. The efficiency ratings are given by

$$f_a = \left[ \sum_{p=1}^{10} b_{ap} t_{ap} / \min(t_{ap}) \right] / n_a \quad (30)$$

where  $b_{ap} = 1$  if code "a" solved problem "p" and zero otherwise,  $t_{ap}$  is the normalized time required for solution (cpu time divided by cpu time required to run a timing standardization program) [3], and  $\min(t_{ap})$  is the shortest time required by any of the codes to solve problem "p." The other efficiency criterion is

$$\bar{f}_a = \left[ \sum_{p=1}^{10} b_{ap} t_{ap} / \text{mean}(t_{ap}) \right] / n_a \quad (31)$$

where  $\text{mean}(t_{ap})$  is the average time required by the codes to solve problem "p." Here the overall rating number for generality (reliability) and efficiency (speed) is given by

$$T_a = \sum_{p=1}^{10} t_{ap} \quad (32)$$

TABLE 3  
RELATIVE RANKING OF OPTIMIZATION CODES

|    |         | Generality |                    | Efficiency |                   | Generality and Efficiency |                         |
|----|---------|------------|--------------------|------------|-------------------|---------------------------|-------------------------|
|    |         | $n_a$      | $N_a$              | $f_a$      | $\bar{f}_a$       | $T_a$                     | $\frac{T_a}{\bar{f}_a}$ |
| 10 | CADOP 3 | 10         | CADOP 3<br>CADOP 2 | 1.3<br>1.0 | PATRN1            | 0.17                      | NMSERS                  |
|    | CADOP 2 |            | CADOP 2            |            |                   | 2.5                       | CADOP 3                 |
| 9  | DSDA    | 9.5        | DSDA               | 3.4<br>3.5 | CADOP 3<br>NMSERS | 0.26                      | PATRN1                  |
|    | PATSH   |            | PATSH              | 5.6        | CADOP 2           |                           | DSDA                    |
| 8  | SEEK 3  | SEEK 3     |                    | 8.9        | DSDA              | 0.34                      | CADOP 2                 |
|    | SIMPLX  | 8.5        | SIMPLX<br>ADRANS   |            |                   | 0.41                      | DSDA                    |
| 7  | ADRANS  |            |                    |            |                   |                           | PATRN1                  |
|    | NMSERS  | 8.0        | NMSERS             | 15         | PATSCH            | 0.77                      | FMIND<br>PATSCH         |
| 5  | PATRN1  | 7.0        | PATRN1             | 16         | FMIND             | 0.90                      | NEMGRD                  |
|    | FMIND   | 6.0        | FMIND              | 20         | SEEK 3            |                           | SIMPLX                  |
| 4  | NEMGRD  | 6.0        | SEEK 1             | 22         | NEMGRD            |                           | ADRANS                  |
|    | DAVID   |            |                    |            |                   |                           | 24                      |
| 3  | SEEK 1  | 5.5        | NEMGRD<br>PATRNO   | 42         | SIMPLX            | 1.0                       | SEEK 3                  |
|    | CLIMB   |            |                    |            |                   |                           | 25                      |
| 2  | DFMFP   |            | DAVID              | 51         | DAVID             | 1.5                       | SIMPLX                  |
|    | GRAD 4  |            | GRAD 4             |            |                   | 2.4                       | DAVID                   |
| 1  | PATRNO  | 5.0        | GRID 1             | 60         | ADRANS            | 2.5                       | ADRANS                  |
|    | RANDOM  |            | RANDOM             |            |                   |                           |                         |
| 0  | DFMCG   | 4.0        | DFMFP              |            |                   |                           |                         |
|    | GRID 1  |            | CLIMB              |            |                   |                           |                         |
| -1 | DFMCG   | 3.0        | DFMCG              |            |                   |                           |                         |

where  $t_{ap}$  is set equal to twice the time used by the slowest code solving a problem "p" that code "a" could not solve. This penalty time is used to penalize code unreliability. Only codes that solved half or more of the problems are rated for efficiency.

The tables presented here are essentially similar to those given in reference [3], except that the DSDA, CADOP2 and CADOP3 codes are included. Thus, both the rating schemes and presentation of results are essentially those of [3].

Based on its performance on the ten problems of reference [3] the new algorithm appears to be a relatively fast and reliable optimization procedure. Codes based on this method were the only ones solving all problems attempted. CADOP3 is comparable in speed to the faster codes (NMSERS, and PATRN1). It was significantly faster than average in all problems and was the fastest code solving problems 1 and 2. Compared to the relatively reliable codes ( $n_a \geq 8$ ), CADOP3 was faster than DSDA on seven of nine problems solved by DSDA, faster than CADOP2 on seven of ten, faster than PATSH on seven of nine, and faster than SEEK3 and SIMPLX in all problems solved by these schemes. As may be seen that DSDA and CADOP2 are the only reasonably reliable codes comparable in speed to CADOP2. CADOP3 appears significantly faster than PATSCH, SEEK3, and SIMPLX. Only the straight (PATRN1) and simplex (NMSERS) codes were generally faster. Of this group, only NMSERS appears to be sufficiently reliable to merit serious consideration. The differences in speed between

NMSERS and CADOP3 is, however, overshadowed by the superior reliability of the latter. Thus, in the overall generality and efficiency rating, CADOP3 stands alone. Viewed on the basis of these comparisons, CADOP3 appears to be a superior nonlinear MP code.

It should be noted that problems 4, 5, 8 and 10 are unconstrained at the optimum points. Thus, even though problem 10 has a regional constraint which is active at the start of the search, these problems can be considered essentially unconstrained.

Those codes (PATRN1 and (NMSER) which are faster overall than CADOP3 obtain their superior efficiency ratings primarily from faster solutions of such problems. The DSFD algorithm is designed, however, for constrained problems. Thus, although its overall performance on constrained and unconstrained problems, or its performance on unconstrained problems, is of interest, a direct evaluation of its performance on constrained problems is also appropriate. Thus a comparison on those problems with behavior constraints (problems 1, 2, 3, 6, 7, and 9) is given in Table 4.

In Table 4 only those codes that solved three or more of the constrained problems are evaluated for efficiency. In this comparison CADOP3 is comparable to NMSERS in speed with these codes being substantially more efficient than any other code. CADOP3 is ranked much higher in the overall rating since NMSERS failed on two of the six problems. CADOP3 is faster than CADOP2 and DSDA on all but one problem. This problem has only one constraint, however, and thus the reduction associated with checking only nearby rather than all constraints is lost.

TABLE 4

RELATIVE RANKING OF OPTIMIZATION CODES ON PROBLEMS  
WITH BEHAVIOR CONSTRAINTS

| <u>Generality</u>    |   | <u>Efficiency</u>    |                                  |                  | <u>Generality and Efficiency</u> |                      |                         |   |
|----------------------|---|----------------------|----------------------------------|------------------|----------------------------------|----------------------|-------------------------|---|
| <u>n<sub>a</sub></u> |   | <u>N<sub>a</sub></u> | <u>f<sub>a</sub></u>             |                  | <u>-f<sub>a</sub></u>            | <u>T<sub>a</sub></u> |                         |   |
| 6                    | CADOP3<br>CADOP2                                    | 6                    | CADOP3<br>CADOP2                 | 2.2<br>2.6       | NMSERS<br>CADOP3                 | 0.14<br>0.16         | NMSERS<br>CADOP3        | 1.2 CADOP3<br>2.0 CADOP2                    |
| 5                    | DSDA<br>PATSH<br>ADRANS<br>SEEK3                    | 5                    | DSDA<br>PATSH<br>ADRANS<br>SEEKS | 6.3<br>12.<br>22 | CADOP2<br>DSDA<br>PATSH<br>SEEK3 | 0.30<br>0.44<br>0.93 | CADOP2<br>DSDA<br>PATSH | 8.1 DSDA<br>8.8 NMSERS<br>11 PATSH<br>SEEK3 |
| 4                    | NMSERS  | 4.5                  | NMSERS                           |                  |                                  | 1.1                  | NEMGRD<br>DAVID         | 13 SIMPLX<br>NEMGRD                         |
| 3                    | MEMGRD<br>SIMPLX<br>DAVID                           | 3.5                  | SIMPLX<br>MEMGRD<br>DAVID        | 25<br>27<br>51   | NEMGRD<br>DAVID<br>SIMPLX        | 1.2<br>1.9           | SEEK3<br>SIMPLX         | 14 ADRANS<br>18 DAVID                       |
| 2                    | FMIND<br>RANDOM                                     | 3.0                  | FMIND<br>SEEK1                   | 65               | ADRANS                           | 2.9                  | ADRANS                  |   |
| 1                    | PATRN1<br>SEEK1                                     | 2.0                  | PATRNO<br>GRID1                  |                  |                                  |                      |                         |   |
| 0                    | PATRNO<br>GRID1<br>GRAD4<br>DFMCG<br>DFMFD<br>CLIMB | 1.0<br>0.5<br>0.0    | DFMFD<br>DFMCG<br>CLIMB          |                  |                                  |                      |                         |   |

CADOP3 is faster than NMSERS on three of four problems although both codes are essentially similar in speed. It is faster on all problems than the remainder of the codes rated for efficiency. CADOP3 is first in the generality and efficiency rating since it is faster than CADOP2 and since all other codes were penalized in this rating for failure on one or more problems. It may be seen therefore that the DSFD improvements cited here did indeed increase performance.

Thus the improved DSFD procedure seems, on the basis of this comparison, clearly superior to the other procedures studied on constrained problems.

Table 5 lists the number of function evaluations required by CADOP3 to reach solution. The comparison data of reference [3] does not provide these values for the other codes. These values are provided here since it is felt that for most engineering problems the number of function evaluations is a better criterion of performance than nondimensional time.

The time required for solution can be viewed as the sum of the function evaluation and algorithm execution time. For very simple functions most of the required time may be associated with algorithm execution particularly where the algorithms are complex. For such problems, however, time is usually of little importance since any reasonably effective procedure can generate a relatively low cost solution. On the other hand, efficiency becomes important where function evaluation is computationally demanding. Here algorithm execution time may be quite small

TABLE 5  
NUMBER OF FUNCTION EVALUATIONS FOR SOLUTION

| <u>Problem No.</u> | <u>Objective Function Evaluations</u> | <u>Constraint Function Evaluations</u> |
|--------------------|---------------------------------------|--|
| 1                  | 2088                                  | 4860                                   |
| 2                  | 266                                   | 190                                    |
| 3                  | 622                                   | 993                                    |
| 4                  | 4910                                  | --                                     |
| 5                  | 347                                   | --                                     |
| 6                  | 378                                   | 322                                    |
| 7                  | 556                                   | 737                                    |
| 8                  | 217                                   | --                                     |
| 9                  | 488                                   | 447                                    |
| 10                 | 849                                   | --                                     |

compared to function evaluation time. In such cases the number of function evaluations is a reasonably accurate criterion for performance.

Preliminary experiments indicate that although nondimensional time is a valid comparison criterion where codes are run using the similar compilation software it is not valid otherwise. It is felt therefore that the number of function evaluations is a superior comparison criterion since it is a reliable indicator of performance on problems where efficiency is important. Thus this information is provided here so as to allow future comparisons.

The principal apparent disadvantage of CADOP3 is its relative complexity compared to some of the other reasonably reliable methods. It contains 713 FORTRAN statements, while, for example, DSDA contains 372 and PATSH only about 75. Thus for simple problems of relatively low dimensionality, one of the simpler codes may be preferable.

It may be seen that the gradient-based procedures, including those employing the SUMT strategy and penalty function [12], so widely used in engineering problems, performed rather poorly. The best gradient-based code DAVID solved only half the problems. The direct search procedures in general provided much better performance although only four, CADOP3, CADOP2, DSDA, and PATSH, solved or approached the solution to all ten problems with only CADOP3 and CADOP2 solving all problems.

It should be noted, however, that this comparison employed rather small test problems (two to five variables with zero to ten behavior constraints). It is not clear that the direct search procedures would also possess reliability and speed superior to the gradient methods on large problems. Still, the superiority in this study of the better direct search procedures is dramatic while the performance of the gradient schemes is rather dismal. Furthermore, although the better direct search procedures may not be as reliable on larger problems, one would certainly not expect the reliability of the gradient procedures to improve on such problems.

## CONCLUSION

Although it is recognized that such favorable findings on the CADOP3 code presented by its developer can be viewed with some skepticism, it should be noted that the performance cited is based primarily on rating schemes and problems selected by Eason and Fenton and not by the author. Furthermore, this performance was achieved with a conscious effort to avoid any tuning of the code to individual problems. Thus, these results indicate that the DSFD algorithm appears to be a superior nonlinear mathematical programming procedure at least on relatively small problems on which it was tested.

It should be noted, however, that this algorithm combines the strength of the direct search procedures (speed) and the Zoutendijk method (ability to confirm optimality and identify a feasible direction) in such a way that the weakness of the direct search in treating problems of large dimensionality is minimized. Thus, one would expect this new procedure to also be quite effective on larger problems. This potential has, however, yet to be demonstrated.

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## **APPENDIX A AND B**

### **USER INSTRUCTIONS**

**and**

### **PROGRAM LISTING**

## APPENDIX A

### USER INSTRUCTIONS FOR CADOP3

#### 1. Instruction

This code treats inequality constrained or unconstrained linear or nonlinear minimization problems by means of a direct search mathematical programming method. The method starts from a user specified starting or initial point and generates a sequence of better points until, hopefully, an optimal, or near optimal, point is reached. The code is intended primarily for constrained nonlinear problems. Linear problems are much more effectively handled by one of many linear programming methods. Unconstrained problems are better treated by UNCDP, a simplified version of CADOP3, or one of several other unconstrained nonlinear problem codes. CADOP3 will, however, treat both linear and unconstrained problems.

It should be noted that nonlinear mathematical programming methods generally do not guarantee an optimal solution because of the possibility of multiple local optima, algorithm failure, or numerical difficulties. Although CADOP3 is among the most reliable of the nonlinear mathematical programming codes it cannot be considered absolutely reliable. Thus, it is desirable to use several optimization runs with widely separated starting points to confirm the convergence has been achieved or to determine if local optima are present.

## 2. Problem Formulation

This code treats the problem:

Find those values  $\bar{x}_i$  of the set of "variables"  $x_i$ ,  $i = 1, 2, \dots, IP$  and the constant "parameters"  $p_k$ ,  $k = 1, 2, \dots, KP$ , that result in the minimum value of the objective function  $f(p_k, x_i)$ , that is

$$f(p_k, \bar{x}_i) = \min [f_k(p, x_i)] \quad (1)$$

while also satisfying the "behavior" constraints

$$g_j(p_k x_i) \leq 0 \quad j = 1, 2, \dots, JP \quad (2)$$

and the "regional" constraints

$$x_i^l \leq x_i \leq x_i^u \quad (3)$$

Two types of behavior constraints are recognized, one type are "linear" constraints of the form

$$g_j(x_i) = \sum_{i=1}^{IP} a_{ij} x_i \quad (4)$$

Where  $f(x_i)$  is also of this form one has a linear programming problem. If any of Eqs. (1 and 2) is not of this form the problem is nonlinear. Constraints not of the form of Eq. (4) are nonlinear constraints.

The program in its present form treats problems with

$$2 \leq IP \leq 10, \quad 0 \leq JP \leq 10 \text{ and } \leq KP \leq 100$$

### 3. Specification of Objective and Constraint Functions

The subprogram FUNCTION OBJ(J) is essentially a dummy subprogram created to accept the users FORTRAN IV program statements defining the objective and behavior constraint functions. The normal form of the behavior constraints is

$$B_j^+ (p_k, x_i) \leq U_j^- (p_k, x_i) \quad (4)$$

where  $B_j^+$  can be thought of as the behavior and  $U_j^-$  the upper limit of behavior.

The FUNCTION OBJ(J) is modified so as to define the objective and constraint functions by inserting the following statements immediately after the COMMON statement

```
GO TO (1,2,----jp), J      (omit if no constraints are
used)
```

```
OBJ = expression defining f (p_k, x_i) using arrays
P(k), X(i)
```

```
RETURN
```

```
j B = expression defining B_j^+
```

```
U = expression defining U_j^-
```

```
GO TO 101
```

A constraint statement group is used for every constraint.

The last constraint group need not use the final GO TO statement.

Example

For the problem:

minimize

$$2x_1 x_2 - p_1 x_3 x_4^2$$

such that

$$2x_3^3 - x_1 \leq p_2$$

and

$$-6 \leq x_2^2 - x_4 \leq 0$$

one could define three constraints and insert the statements

GO TO (1,2,3),J

OBJ = 2. \* X (1) \* X (2) - P (1) \* X (3) \* X (4) \*\*2

RETURN

1 B = 2.\* X (3) \*\* 3 - X (1)

U = P(2)

GO TO 101

2 B = X(2) \*\* 2 - X (4)

U = 0.

GO TO 101

3 U = X (2) \*\* 2 - X (4)

B = - 6.

Alternately, noting that both upper and lower limits on the second constraint equation cannot be active simultaneously, one could use two constraint equations and insert the program segment:

```
GO TO (1,2),J
OBJ = 2. * X (1) * (2) - P (1) * X (3) * X (4) ** 2
RETURN
1 B = 2. * X (3) ** 3 - X (1)
U = P
GO TO 101
2 TB = X (2) ** 2 - X (4)
IF(TB.LT.-3.) GO TO 3
B = TB
U = 0.
GO TO 101
3 B = -6.
U = TB
```

The constraints values at the optimum are printed in the form of equation (2) where

$$g_j = \begin{cases} (B_j - U_j)/|U_j| & U_j \neq 0 \text{ and } B_j \neq 0 \\ (B_j - U_j) & U_j = 0 \text{ or } B_j = 0 \end{cases}$$

thus a negative value of  $g_j$  indicates the constraint is satisfied.

#### 4. Data Cards

The program allows multiple optimization runs of a particular problem within a single computer run using different parameters, for parametric studies, different starting points, for optimality confirmation or search for local optima, and different sets of regional constraint limits.

The user must, by use of the data card set, specify the number of parameters, variables and behavior constraints used, identify the linear constraints, specify for the first run if regional constraints are to be used, and for subsequent runs whether the parameters, initial point or regional limits are to be changed. If parameters are used they must be entered. The initial point must be entered. The values for the regional constraints must be entered if such limits are imposed.

The first data set is entered on the "Problem Constants" card which specifies, in order, the number of parameters (KP), variables (IP) and behavior constraints (JP), used (see line 9 of program listing). The data is entered in 3I10 format in the first 3 fields in order and must be right justified.

If constraints are used ( $JP > 0$ ) a second data set is entered on the "Constraint Identification" card which identifies the linear behavior constraints. If a constraint is linear a digit is entered in the column with the same number as the constraint. If no linear behavior constraints are used a blank card is employed (see line 10).

The third data set is entered on the "Data Control" card and is used to specify whether: 1) new parameters are to be used (ICNTR2), 2) a new initial point is to be used (ICNTR3) and, 3) regional limits or new regional limits are to be used (ICNTR4) (see lines 26-43). The entries are made in 3I10 format in the order given above. For the first run no entries in the first two fields are needed. If regional limits are used an entry, any nonzero digit, is made in the third field (col 21-30).

If and only if the number of parameters specified is greater than zero ( $KP > 0$ ) a fourth data set is used to enter the problem parameters  $p_k$ . The entries are made 5 to a card in F15.5 format in order  $i = 1, 2, \dots, KP$ , (see lines 21 and 22).

The fifth data set specifies the starting point and entries are made, in order, 8 to a card in F10.5 format (see lines 23 and 29).

If and only if an entry is made in the third field of the Data Control card ( $ICNTR4 = 0$ ) a sixth data set defining the lower limits is entered, 8 to a card, in F10.5 format followed by an upper limit set (starting with a new card) with similar format. Both complete limit sets must be entered (see lines 30-34).

For every additional run an additional Data Control card is added followed by parameter and/or initial point and/or regional limit sets, where and only where the need for a new set is indicated by an appropriate entry ( $ICNTR2 = 1$  or  $(ICNTR3 = 1$  or  $ICNTR4 = 1$ ) on the Data Control card (see lines 26-34).

A blank Data Control card is used to terminate the run(s) (see lines 26 and 27).

## 5. Conversion To Single Precision

The program as supplied is a double precision form. Experience has shown, however, that the great majority of problems can be treated adequately in single precision and thus it is recommended that single precision be tried first. To convert to single precision remove the IMPLICIT statement at the beginning of MAIN and all subprograms and the function conversion statements such as SQRT(B) = DSQRT(B) etc., near the beginning of most subprograms. Furthermore, the REAL FUNCTION OBJ\*8(J) should be put in single precision form.

## 6. Change of Problem Size

To change the maximum number of variables (IP) or the maximum number of constraints (JP) the program can treat, change the arrays in MAIN and all subprograms as follows:

1. change all D, DZ, X, DL, and DU arrays in common to D(IP), DZ(IP) etc.,
2. change the G array (T in OPT2 and LINPRO) as well as the CK, IA, IC array in COMMON and the BL array in COMMON/OPT/ where they occur to G(JP), CK(JP) ETC.,
3. change the V array (A array in OPT2 and LINPRO) to V(M,N) in COMMON where

$$M = JP + 4IP + 5$$

$$N = JP + 5IP + 6$$

4. change all DUMMY arrays such as DUM, IDUM, etc., to equalize common sizes.
5. In the OPT2 SUBROUTINE DIMENSION statement change DT, SS, IB, IC, to DT(IP) etc., change GG to GG(JP), TN to TN(JP+1), ID to ID(JP), DK to DK(JP+1, IP), DBJ to DBJ (JP+1, IP) and SSS to SSS(M) where M is as above.
6. In the FIND SUBROUTINE DIMENSION statement change DT(10) to DT(IP).
7. In the PATTRN SUBROUTINE DIMENSION statement change all 10's to IP.

## **APPENDIX B**

CADOP 3

## **PROGRAM LISTING**

# MAIN

```

    IMPLICIT REAL*8(A-H,O-Z)
    COMMON D(10),P(100),DZ(10),X(10),G(10),A,AMIN,DL(10),DU(10),
    1V(55,66),CK(10),Q,IP,JP,IA(10),KP,KW,KL,IK,IDUM(21),JX
    COMMON/OPT/BL(10),TM
1   FORMAT(5F15.5)
2   FORMAT(4I10)
3   FORMAT(2F10.5)
101  FORMAT(8F10.5)
     READ2,KP,IP,JP
     IF(JP.NE.0)READ 44,(CK(K),K=1,JP)
     IF(JP.EQ.0)GO TO 45
     DO 46 K=1,JP
     IF(CK(K).NE.0.)CK(K)=0.01
     IF(CK(K).EQ.0.)CK(K)=1.
     IF(CK(K).LT..5)CK(K)=0.
46  CONTINUE
45  DE 21 I=1,IP
     DL(I)=-1.E+48
44  FORMAT(80F1.0)
21  DU(I)=1.E+48
     READ2,ICNTR2,ICNTR3,ICNTR4
     IF(KP.EQ.0) GO TO 12
     READ1,(P(I),I=1,KP)
12  READ101,(D(I),I=1,IP)
18  FORMAT(IHO,' STARTING DESIGN VARIABLE VALUES' )
     GO TO 23
24  READ(5,2,END=33)ICNTR2,ICNTR3,ICNTR4
     IF(ICNTR2+ICNTR3+ICNTR4.EQ.0)STOP
     IF(ICNTR2.NE.0)READ1,(P(I),I=1,KP)
     IF(ICNTR3.NE.0)READ101,(D(I),I=1,IP)
23  IF(ICNTR4.EQ.0)GO TO 22
     READ 101,(DL(I),I=1,IP)
17  FORMAT(IHO,' LOWER LIMITS OF DESIGN VARIABLES' )
     READ 101,(DU(I),I=1,IP)
19  FORMAT(IHO,' UPPER LIMITS OF DESIGN VARIABLES' )
22  IF(KP.LE.0) GO TO 104
     PRINT 6
     PRINT7,(K,P(K),K=1,KP)
104  PRINT 18
     PRINT101,(D(I),I=1,IP)
     PRINT 17
     PRINT101,(DL(I),I=1,IP)
     PRINT 19
     PRINT101,(DU(I),I=1,IP)
     CALL PATTRN
     IK=0
     CALL FIND(TD)
     OBJD=OBJ(0)
     PRINT4,KL
4   FORMAT('NO. OF REDESIGN CYCLES=' ,15)
     PRINT5,DBJD
5   FORMAT('OPTIMUM VALUE=' ,F13.8)
6   FORMAT('DESIGN PARAMETERS' /)
7   FORMAT(' P' ,I2,'=' ,F14.4)
     PRINT8
8   FORMAT('DESIGN VARIABLE VALUES' /)

```

MAIN

```
9 PRINT9,(K,D(K),K=1,IP)
FORMAT(' D',I1,'=',F15.8)
IF(JP-EQ.0) GO TO 24
PRINT10
10 FORMAT('ONEARNESS TO CONSTRAINTS')
PRINT11,(K,G(K),K=1,JP)
11 FORMAT(' G',I2,'=',F15.5)
100 GO TO 24
33 STOP
END
```

## PATTRN

```

SUBROUTINE PATTRN
  IMPLICIT REAL*8(A-H,O-Z)
  COMMON D(10),P(100),DZ(10),X(10),G(10),A,AMIN,DL(10),DU(10),
  IV(55,66),CK(10),Q,IP,JP,IA(10),KP,KW,KL,IK,JDUM(21),JX
  COMMON/DPT/BL(10),TN
  DIMENSION AD(10,10),EI(10,10),ET(10,10),AZ(10)
  TCF=10.E+10
  IPP=IP
  AL=A
  IM=0
  STEST=0.
  KA=0
  KW=0
  KT=0
  KL=0
  KF=0
  KY=0
  KWN=0
  DO 112 K=1,JP
112  BL(K)=-0.1
  DO 111 K=1,IPP
  IF(D(K).LT.DL(K))D(K)=DL(K)
  IF(D(K).GT.DU(K))D(K)=DU(K)
111  X(K)=D(K)
  CALL SIZE
25   KK=1
  DO 1 K=1,IPP
  DO 22 I=1,IPP
  EI(I,K)=0.
  ET(I,K)=0.
22   AD(I,K)=0.
  ET(K,K)=1.
  EI(K,K)=1.
1  CONTINUE
  ST=SB
  IF(KA.EQ.1)GO TO 9
  KA=1
190  IK=0
  JX=0
  CALL FIND(SB)
9   ST=SB
  JX=0
  GO TO 100
19   IK=0
  JX=0
  CALL FIND(ST)
100  KL=KL+1
  IK=1
  KF=KF+1
  DO 2 J=1,IPP
  DE 17 K=1,1PP
  AZ(K)= EI(J,K)*A
  IF(KK.NE.0)AZ(K)=ET(J,K)*A
17   D(K)=D(K)+AZ(K)
  DE 35 K=1,IPP
  IF(D(K).GT.DU(K))GO TO 4

```

## PATTRN

```

    IF(D(K).LT.DL(K))GO TO 4
35  CONTINUE
    CALL FIND(SA)
    IF((ST-SA)/DABS(SB).GT.2.E-16) GO TO 5
4   DO 18 K=1,IPP
18   D(K)=D(K)-AZ(K)-AZ(K)
    DO 36 K=1,IPP
    IF(D(K).LT.DL(K))GO TO 3
    IF(D(K).GT.DU(K))GO TO 3
36  CONTINUE
    CALL FIND(SA)
    IF((ST-SA)/DABS(SB).LT.2.E-16) GO TO 3
5   ST=SA
    GO TO 2
3   DO 21 K=1,IPP
21   D(K)=D(K)+AZ(K)
2   CONTINUE
    IF((SB-ST)/DABS(SB).LT.2.E-16) GO TO 7
26   SUM=0.
    DO 8 K=1,IPP
    DB=2.*D(K)-X(K)
    AD(1,K)=D(K)-X(K)
    IF(DABS(AD(1,K)).LT.+1.E-32) AD(1,K)=0
    SUM=SUM+AD(1,K)*AD(1,K)
    X(K)=D(K)
    SB=ST
    IF(DB.LT.DL(K))DB=DL(K)
    IF(DB.GT.DU(K))DB=DU(K)
8   D(K)=DB
    SUM=DSQRT(SUM)
    IF(SUM.LE.0)GO TO 23
    IF(SUM-DABS(.5*A)) 7,7,950
950 CONTINUE
    KK=0
    SUM=0.
    DO 11 J=1,IPP
    ET(1,J)=EI(1,J)
    EI(1,J)=AD(1,J)
11   SUM=SUM+AD(1,J)**2
    SUM=DSQRT(SUM)
    DO 53 J=1,IPP
53   EI(1,J)=EI(1,J)/SUM
    DO 13 I=2,IPP
    K=I-1
    DO 14 J=I,IPP
14   AD(I,J)=AD(I,J)
    SUMA=0.
    DO 15 J=1,IPP
    IF(DABS(EI(K,J)).LT.+1.E-32) EI(K,J)=0
15   SUMA=SUMA+EI(K,J)*AD(I,J)
    SUM=0.
    DO 16 J=1,IPP
    ET(I,J)=EI(I,J)
    EI(I,J)=AD(I,J)-EI(K,J)*SUMA
    IF(DABS(EI(I,J)).LT.+1.E-32) EI(I,J)=0
16   SUM=SUM+EI(I,J)**2

```

## PATTRN

```

SUM=DSQRT(SUM)
IF(SUM)27,27,28
27  DC 29 J=1,IPP
29  E1(I,J)=ET(I,J)
   GO TO 19
28  DC 13 J=1,IPP
13  E1(I,J)=E1(I,J)/SUM
   GO TO 19
    7 DC 20 K=1,IPP
20  D(K)=X(K)
   IF(KK)25,24,23
24  KK=-1
   IF(KF.GT.200)GO TO 23
   GO TO 9
23  CONTINUE
   TN=SB
   IF(JP.EQ.0)GO TO 32
   IK=0
   CALL FIND(SB)
   DO 30 I=1,JP
   IF(G(I).GT.-.01*BL(I))GO TO 31
30  CONTINUE
   IF(KWW.LE.3) KW=0
   KY=0
   GO TO 32
31  KW=KW+1
   KY=KY+1
   KWW=KWW+1
   IK=0
   IF(KW.NE.0) CALL FIND(SB)
32  IF(KL.GT.KT+4) GO TO 33
   A=A/2.
   DO 34 K=1,JP
34  BL(K)=BL(K)/2.
   33 CALL OPT2(SB,ST,JK,KY)
   KT=KL
   IF(JK.EQ.2) RETURN
311  KF=0
   DO 300 I=1,JP
   IF(G(I).GT.-.01*BL(I))GO TO 301
300  CONTINUE
   IF(KWW.LE.3) KW=0
   KY=0
301  IF(JK.NE.0.) GO TO 46
   IF(A.EQ.AL) GO TO 26
   PRINT 114,A,SB,ST,(D(I),I=1,IP)
114  FORMAT( 0 OPT2BASEA=' ,F11.9,'SB=' ,E16.8,'ST=' ,E16.8,'X=' ,10F10.6)
   AL=A
   DIF=DABS((SB-STEST)/SB)
   IF(DIF.GT..00001) GO TO 45
   IM=IM+1
   IF(IM.LT.2) GO TO 150
   IF(DIF.GT.TDIF) GO TO 45
   RETURN
45  IM=0
149  STEST=SB

```

## PATTRN

```
TDIF=DIF
GO TO 26
148 IM=0
A=.5*A
DO 113 K=1,JP
113 BL(K)=BL(K)/2.
46 D(1)=X(1)
47 D(K)=X(K)
IF(A.LT.AMIN) RETURN
GO TO 190
150 A=A/2
DO 115 K=1,JP
115 BL(K)=BL(K)/2.
IF(A.LE.AMIN) RETURN
GO TO 149
END
```

## FIND

```

SUBROUTINE FIND(T0)
IMPLICIT REAL*8(A-H,O-Z)
COMMON D(10),P(100),DZ(10),X(10),G(10),A,AMIN,DL(10),DU(10),
1V(55,66),CK(10),Q,IP,JP,IA(10),KP,KW,KL,IK,IL,LDUM(10),IC(10),JX
COMMON/OPT/BL(10),TN
DIMENSION DT(10)
I=0
T0=OBJ(0)
TN=T0
IF(JP.EQ.0) RETURN
DO 25 K=1,IP
25 DT(K)=D(K)
IF(IK.EQ.0) GO TO 51
IF(IL.EQ.0) RETURN
DO 52 J=1,IL
IB=IC(J)
G(IB)=OBJ(IB)
IF(G(IB)) 52,52,5
5 I=I+1
IA(I)=IB
CONTINUE
GO TO 56
51 IL=0
DO 3 J=1,JP
G(J)=OBJ(J)
IF(G(J).LT.-2.*BL(J))GO TO 3
6 IL=IL+1
IC(IL)=J
IF(G(J)) 3,3,4
4 I=I+1
IA(I)=J
3 CONTINUE
56 IF(I)41,41,100
100 HTEMP=0.
DO 21 K=1,I
IB=IA(K)
IF(KW.NE.0) GO TO 72
IF(G(IB).LT.-BL(IB))GO TO 21
72 E=10000.*G(IB)
IF(E.GT.HTEMP) HTEMP=E
21 CONTINUE
IF(HTEMP.GT..001) GO TO 19
IF(JX.NE.0) GO TO 42
SUM=0.0
JX=1
DO 1 K=1,IP
DT(K)=D(K)
D(K)=D(K)+.001
DZ(K)=(OBJ(0)-TN)/.001
SUM=SUM+DZ(K)**2
1 D(K)=DT(K)
SUM=DSQRT(SUM)
DO 53 K=1,IP
53 DZ(K)=.001*DZ(K)/SUM
42 DO 50 K=1,IP
50 D(K)=DT(K)+DZ(K)

```

FIND

```
TA=OBJ(0)
DC 18 K=1,1
IB=IA(K)
TT=OBJ(IB)
IF (TT-G(IB)) 37,54,37
37 E=DABS((2.* (T0-TA)/(TT-G(IB)))*G(IB))
IF(G(IB)+BL(IB))55,55,54
54 E=10000.*G(IB)
55 IF(E.GT.HTEMP) HTEMP=E
18 CCONTINUE
19 CCONTINUE
T0=T0+HTEMP
DC 26 K=1,IP
26 D(K)=DT(K)
41 RETURN
END
```

## OPT2

```

SUBROUTINE OPT2(SB,ST,JK,KY)
IMPLICIT REAL*8(A-H,O-Z)
INTEGER Z,E,G,BB,W,B,H
COMMON D(10),F(100),DZ(10),X(10),T(10),V,AMIN,DL(IC),DU(10),
1A(55,66),CK(10),Q,IP,JP,IA(10),KP,KW,KL,IK,IL,M,N,B,L,BB,E,G,Z,H,
2W,IDUM(10),J5
COMMON/OPT/BL(10),TP
DIMENSION DT(10),SS(10),IB(10),IC(10),SSS(55),GG(IC),TN(11),DK(10)
1,20),ID(10),OBJ(11,10)
VT=V/10.
TQ=TP
J6=0
JIN=0
JXN=0
DEL=.00001
IF(V.LT.DEL) DEL=V
IP1=JP+4*IP+4
DO 116 K=1,JP
GG(K)=T(K)
116 ID(K)=0
117 DO 103 I=1,IP1
SSS(I)=0.
JK=1
N=2*IP+1.0
N=IP+1
Z=-1
MM=JP+1
M=MM+2*N+2*IP
L=MM+N+IP
E=0
G=0
NN=6
90 Q=99999
190 B=M+N+G+I
W=M
320 H=1
BB=B+1
J=0
DO 4 I=1,IP
IB(I)=0
IC(I)=0
IF(D(I)-DL(I).LT.V)GO TO 5
IF(DU(I)-D(I).LT.V)GO TO 6
GE TO 4
5 J=J+1
IB(I)=J
GO TO 4
6 J=J+1
IC(I)=J
4 CONTINUE
JX=J
MM=JP+1
IL=0
156 J=0
DO 22 I=1,MM
I1=I-1

```

## OPT2

```

IF(I.EQ.1) TM=TQ
IF(I.EQ.1) GO TO 77
TM=GG(I1)
IF(GG(I1).LE. BL(I1)) GO TO 22
ID(I1)=1

77 J=J+1
TN(J)=TM
22 CONTINUE
IF(JX.EQ.JXN.AND.J.EQ.JIN) GO TO 14
JIN=J
JXN=JX
DO 23 K=1,IP
DT(K)=D(K)
D(K)=D(K)+DEL
J=0
DC 24 I=1,MM
I1=I-1
IF(I.EQ.1) GO TO 777
IF(GG(I1).LE.BL(I1)) GO TO 24
777 J=J+1
IF(J6.NE.0)GO TO 778
DBJ(I,K)=OBJ(I1)
778 DK(J,K)=(DBJ(I,K)-TN(J))/DEL
24 CONTINUE
23 D(K)=DT(K)
J6=1
DO 25 I=1,J
SUM=0.
DC 26 K=1,IP
SUM=SUM+DK(I,K)**2
IF(SUM.EQ.0.) GO TO 25
SUM=DSQRT(SUM)
DO 27 K=1,IP
27 CK(I,K)=DK(I,K)/SUM
25 CONTINUE
JQ=0
330 LX=W+2
IF(J.GT.1)GO TO 775
IF(JX.EQ.0)GO TO 150
775 DO 350 I=1,LX
340 DC 350 J=1,B
350 A(I,J)=0
J=0
DO 2 I=1,MM
I1=I-1
IF(I.EQ.1) GO TO 7
IF(GG(I1).LE. BL(I1)) GO TO 2
7 J=J+1
L=0
DO 3 K=1,IP
IF(IC(K).NE.0) GO TO 3
L=L+1
A(J,L)=DK(J,K)
CONTINUE
DC 10 K=1,IP
IF(IB(K).NE.0) GO TO 10

```

## OPT2

```

L=L+1
A(J,L)=-DK(J,K)
10  CONTINUE
N=L+1
IF(J.EQ.1)GO TO 2
A(J,N)=CK(I1)
IF(KY.NE.0)A(J,N)=10.
2   CONTINUE
A(I,N)=1.
JA=J+1
DC 8 I=1,L
J=J+1
8   A(J,I)=1.0
W=J
A(W+1,N)=1.0
B=J+N+1
BB=B+1
DC 9 I=JA,W
9   A(I,B)=1.0
L=W
DO 510 I=1,N
510  A(W+1,I)=FLOAT(Z)*A(W+1,I)
     M=J
     NN=6
     CALL LINPRO(K2)
     IF(K2.EQ.0) STOP
     LL=M+1
     DC 1460 I=1,LL
     J=A(I,BB)
     SSS(J)=A(I,B)
1460  CONTINUE
     L=0
     DC 19 I=1,IP
     IF(IC(I).NE.0)GO TO 12
     L=L+1
     SS(I)=SSS(L)
     GC TO 19
12   SS(I)=0
19   CONTINUE
     DC 13 I=1,IP
     IF(IB(I).NE.0)GO TO 13
     L=L+1
     SS(I)=SS(I)-SSS(L)
13   CONTINUE
14   SUM=0.
     DG 11 I=1,IP
11   SUM=SUM+SS(I)*SS(I)
     SUM=DSQRT(SUM)
     IF(SUM.EQ.0.)GO TO 187
     DC 16 I=1,IP
     SS(I)=V*SS(I)/SUM
     D(I)=DT(I)+SS(I)
16   CONTINUE
18   IK=0
     CALL FIND(ST)
     IF(JP.EQ.0)GO TO 157

```

## OPT2

```
j=0
DO 185 I=1,JP
IF(T(I).LE.0.) GO TO 185
IF(ID(I).EQ.0) BL(I)=-1.5*DABS(GG(I))
IF(ID(I).EQ.0) J=1
185 CONTINUE
IF(J.GT.0)GO TO 15
157 IF(SB.GT.ST) GO TO 101
187 V=.5*V
DO 113 I=1,JP
113 BL(I)=BL(I)/2.
IF(V.LT.VT) RETURN
IF(V.LT.AMIN) RETURN
15 DE 163 I=1,IP
163 D(I)=DT(I)
GO TO 117
101 JK=0
RETURN
150 SUMM=0.0
DO 1 J1=1,IP
1 SUMM=SUMM+DK(I,J1)**2
SUMM=DSQRT(SUMM)
IF(SUMM.EQ.0) JK=2
IF(SUMM.EQ.0.) RETURN
162 DE 151 I=1,IP
D(I)=DT(I)-DK(I,I)*V/SUMM
IF(D(I).LT.DL(I)) D(I)=DL(I)
151 IF(D(I).GT.DU(I)) D(I)=DU(I)
SUMM=0.
DE 152 I=1,IP
152 SUMM=SUMM+(D(I)-DT(I))**2
SUMM=DSQRT(SUMM)
DO 153 I=1,IP
153 D(I)=DT(I)+(D(I)-DT(I))*V/SUMM
JK=0
CALL FIND(ST)
IF(SB.GT.ST)GO TO 101
V=V/2.
DO 112 I=1,JP
112 BL(I)=BL(I)/2.
IF(V.LT.VT) RETURN
IF(V.GT.AMIN)GO TO 162
RETURN
END
```

## SIZE

```
SUBROUTINE SIZE
  IMPLICIT REAL*8(A-H,O-Z)
  COMMON D(10),P(100),DZ(10),X(10),G(10),A,AMIN,DL(10),DU(10),
  IV(55,66),CK(10),Q,JP,JP,IA(10),KP,KW,KL,IK,IDUM(22)
  TN=OBJ(0)
  DO 1 I=1,IP
    X(I)=D(I)
    D(I)=D(I)+.001
    DZ(I)=(OBJ(0)-TN)/.001
  1  D(I)=X(I)
    DTEMP=0.0
    ITEMP=0
    DO 2 I=1,IP
      IF(DABS(DZ(I)).LT.DTEMP) GO TO 2
      DTEMP=DABS(DZ(I))
      ITEMP=I
  2  CONTINUE
    A=DABS(10.01*TN)/DZ(ITEMP)
    IF(A.LT.0.005)A=0.005
    SUM=0.
    DO 3 I=1,IP
      SUM=SUM+D(I)**2
    SUM=DSQRT(SUM)*.001
    IF(A.LT.SUM) A=SUM
    AMIN=A/100000.
    RETURN
  END
```

## LINPRO

```

SUBROUTINE LINPRO(K2)
  IMPLICIT REAL*8(A-H,D-Z)
  INTEGER Z,E,G,BB,W,B,R,C,H
  COMMON D(10),F(100),DZ(10),X(10),T(10),V,AMIN,DL(10),DU(10),
  1A(55,66),CK(10),Q,IP,JP,IA(10),KP,KW,KL,IK,IL,M,N,B,L,BB,E,G,Z,H,
  2W, IDUM(10),JX
  K2=1
  NA=6
  M=M-1.0
  560  LL=M+2
  DE 580 K=2,LL
  570  A(K-1,N+G+K-1)=1
  580  A(K-1,BB)=K+N+G-1
  600  IF(G.NE.0) GO TO 620
  610  IF(E.EQ.0) GO TO 780
  611  GC TU 650
  620  KK=L+E+2
  LL=M+2
  DE 630 K=KK,LL
  630  A(K-1,K+N-L-E-1)=-1
  650  W=W+1
  660  Q=0
  670  LL=N+G
  DE 760 J=1,LL
  680  S=0
  690  LL1=M-G-E+2
  KK1=M+1
  DE 700 I=LL1,KKI
  700  S=S+A(I,J)
  720  A(W+1,J)=-S
  730  IF(A(W+1,J).GT.Q) GO TO 760
  740  Q=A(W+1,J)
  750  C=J
  760  CONTINUE
  761  S=0
  762  LL=M-G-E+2
  KK=M+1
  DE 763 J=LL,KK
  763  S=S+A(J,B)
  765  A(W+1,B)=-S
  780  CONTINUE
  790  IF(G.EQ.0) GO TO 810
  LL=N+1
  KK=N+G
  810  IF(L.EQ.0) GO TO 830
  LL=N+G+1
  KK=N+G+L
  IC=1
  830  IF(G+E.EQ.0) GO TO 2000
  831  LL=N+G+L+1
  KK=B-1
  IC=1
  860  GO TO 2000
  895  IF(Q.EQ.99999) GO TO 1230
  900  IF(Q.EQ.0) GO TO 1330
  910  GE TU 1400

```

## LINPRO

```

920 H=M+1
930 Q=.IE39
940 R=-1
LL=M+1
950 DE 1000 I=1,LL
960 IF(A(I,C).LE.0) GO TO 1000
970 IF(A(I,B)/A(I,C).GT.Q) GO TO 1000
980 Q=A(I,B)/A(I,C)
990 R=I
1000 CONTINUE
1010 IF(FLOAT(R).GE.-.5) GO TO 1050
1G=2
1030 GE TO 2000
1050 P=A(R,C)
1060 A(R,BB)=C
1070 DE 1080 J=1,B
1080 A(R,J)=A(R,J)/P
LL=W+1
1100 DE 1180 I=1,LL
1110 IF(I.EQ.R) GO TO 1180
1120 DE 1170 J=1,B
1130 IF(J.EQ.C) GO TO 1170
1140 A(I,J)=A(I,J)-A(R,J)*A(I,C)
1150 IF(DABS(A(I,J)).GT..1E-4) GO TO 1170
1160 A(I,J)=0
1170 CONTINUE
1180 CONTINUE
LL=W+1
1190 DO 1200 I=1,LL
1200 A(I,C)=0
1210 CONTINUE
1220 A(R,C)=1
1230 Q=0
1240 LL=N+G+L
DE 1280 J=1,LL
1250 IF(A(W+I,J).GT.Q) GO TO 1280
1260 Q=A(W+1,J)
1270 C=J
1280 CONTINUE
1290 GO TO 900
1330 IF(W.EQ.M+1) GO TO 1360
1340 W=W-1
1350 IF(A(W+2,B).LT..1E-5) GO TO 1353
K2=0
1352 RETURN
1353 LL=M+1
DO 1358 I=1,LL
1354 IF(IDINT(A(I,BB)).LE.N+G+L) GO TO 1358
1355 DE 1356 J=1,B
1356 A(I,J)=0
1358 CONTINUE
1359 GO TO 1230
1360 CONTINUE
1400 IF(Q.EQ.0) GO TO 1420
1420 LL=M+1
1430 DE 1460 I=1,LL

```

## LINPRO

```
1440 IF(IDINT(A(I,BB)).EQ.0) GO TO 1460
1460 CONTINUE
1470 IF(Q.NE.0) GO TO 920
1520 KK=N+1
      LL=B-G-1
      XJJ=-Z*A(W+1,B)
      LL=H-1
      IC=3
1550 GE TO 2000
2000 LL=H-1
      LL=W+1
      GE TO(895,1050,999),1Q
999 RETURN
END
```

OBJ

```
REAL FUNCTION OBJ*8(J)
IMPLICIT REAL*8(A-H,O-Z)
COMMON X(10),P(100),DUM(3693),IDUM(38)
101  OBJ=B-U
      IF(B.EQ.0. .OR.U.EQ.0.) RETURN
      OBJ=(B-U)/DABS(U)
      RETURN
END
```

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