A Note on a Comparison of Starting Points for the Generation of D-optimal Second-order Designs

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Abstract. Usually, line search algorithms begin by an initial guess of the starting point. However, for quick convergence and reduction in computational cost, it is necessary to determine an optimal starting point. In this paper, comparison of three starting points for the generation of second-order $D$-optimal designs using a variance-gradient based exchange algorithm is considered. Quantile plots are used to compare the variances of $D$-optimal designs generated by each of the three starting points. Result obtained reveals that an optimal choice of the starting point improves $D$-efficiency of the generated design points.

Résumé (Abstract in French) Les algorithmes de recherche en ligne commencent par un point donné et il est souhaitable de déterminer ce point de manière optimale afin d’assurer une convergence plus rapide et de réduire le coût de calculs. Dans ce papier, nous comparons trois méthodes de détermination de trois points de départ pour la génération de second ordre (D-) dans un design optimal, utilisant une méthode de variance-gradient. Des études de simulations illustrent notre méthode.

Key words: $D$-efficiency; line search method; quantile plots; second-order designs; starting points

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1. Introduction

In design of experiments, standard statistical designs such as Box Behnken design, central composite design, factorial and fractional factorial designs which are optimal with respect to the $A^{-}$, $D^{-}$ and $G^{-}$ optimality criteria are usually used. However, these designs are inappropriate in some situations involving non-standard modelling, constrained regions, unusual sample sizes, or ‘tailor-made’ designs (designs arising just to satisfy the exact needs of the experimenter); see Montgomery (2013, p.511). In such cases, algorithmic construction of optimal designs are useful. Algorithms based on the line search equation have three common parts: the starting point, $x_0$, the search direction, $d$ and the step length, $\rho$. Such algorithms begin by an initial guess of $x_0$ and performing some sequences of operations using the $d$ and $\rho$ to obtain an iterate, $x^\ast$. A stopping rule is applied and the iterate is tested for convergence depending on the design criterion to be satisfied. The line search equation is given as

$$x^\ast = x_0 + \rho d$$  \hspace{1cm} (1)

At the $j$th step, the line equation is,

$$x^\ast_j = x^\ast_{j-1} + \rho_j d_j$$  \hspace{1cm} (2)


A 'good' choice of the starting point helps to attain quick convergence in few iterations. Gertz, Nocedal and Sartenar (2004) noted the poor performance (or even failure) in linear and quadratic programming when the starting point is unfavourable. Based on that, they presented a starting point strategy in non-linear programming. An appropriate choice of an initial starting point reduces computational cost and improves the speed of convergence, see Torabi and Hosseini (2018). $x_0$ can be chosen in some arbitrary manner but should be a reasonable estimate of the solution (Nocedal and Wright, 1999). Hence, there is need for optimal choice of the starting point.

In this paper, three starting points shall be compared using the gradient-based algorithm proposed by Onukogu and Nsude (2014). Each starting point is applied to the algorithm to generate $D^{-}$-optimal second-order designs in the experimental
space \([-1,1]\). The optimal design points obtained are compared graphically and with \(D\)-efficiency. For \(D\)-efficiency, the optimal points are compared with Box and Draper (1974) minimum point second-order designs. Comparison of the starting points are carried out for a second order model defined in the experimental space \([-1,1]\).

2. \(D\)-optimality Criterion

Kiefer and Wolfowitz (1959) greatly contributed to the development of \(D\)-optimality criterion. This criterion maximizes the determinant of Fisher’s information matrix. It also minimizes the generalized variance of the regression parameter estimates and the volume of the confidence ellipsoid of the parameter; see Atkinson and Donev (1992). Consider the usual linear model,

\[ y = f'(\bar{x})\beta + e; \quad (3) \]

defined in an experimental space \(\{\bar{X}, F_x, \Sigma_x\}\) where \(\bar{X}\) is a \(k\)-dimensional compact, continuous and metric space of all possible experimental trials, \(F_x\) is a finite dimensional linear space of continuous functions and \(\Sigma_x\) is a non-negative continuous space of random error associated with the response at each point \(\bar{x} \in \bar{X}\): see Onukogu (1997). \(f(\bar{x})\) is a \(p\)-parameter polynomial of degree \(m\), \(\beta\) is a \(p\)-vector of unknown parameters to be estimated by the experimenter, \(\bar{x} = (x_1, x_2, \ldots, x_k)\) are non-stochastic independent variables in \(\bar{X}\), \(e \in \Sigma_x\). The model in (3) can be written in matrix form as

\[ Y = X\beta + e; \quad (4) \]

where \(Y\) is an \(n \times 1\) response vector, \(X\) is an \(n \times p\) design matrix of rank \(p\), \(\beta\) is a \(p \times 1\) vector of unknown parameters, \(e \sim N(0, \sigma^2 I)\) random error term. The least square estimate of \(\beta\) is \(\hat{\beta} = (X'X)^{-1}X'Y\).

An optimal design generated based on this criterion for a specified model is called \(D\)-optimal design. Therefore, \(D\)-optimal design maximizes the determinant of \((X'X)\), that is \(\max|x'Xx|\). For a specified \(p\)-parameter linear model defined in an experimental space, an \(n\)-point \(D\)-optimal design is given as \(|M(\bar{x}, \xi_{n})| = \max_{\xi}|X'X|/n^p\), see Heredia-Langner et al(2003). \(D\)-efficiency is one of the properties of \(D\)- optimum designs. It is defined as

\[ D - ef f = \left( \frac{|M(\bar{x}, \xi_{n})|}{\max_{\xi}[M(\bar{x}, \xi_{n})]} \right)^{1/p} \]

where, the denominator is the maximum of all available \(D\)-optimal designs.

3. Method

The three starting points shall be applied to the algorithm using a second order model with two independent variables. The model is shown in (5)

\[ y = \beta_0 + \beta_1 x_1 + \beta_2 x_2 + \beta_{12} x_1 x_2 + \beta_{11} x_1^2 + \beta_{22} x_2^2 + e \quad (5) \]
3.1. The Algorithm

The algorithm as presented by Onukogu and Nsude (2014) is summarized in the following steps:

1. Construct $N$ grid points $(x_1, x_2, \ldots, x_N)$ for the factors $x = (x_1, x_2, \ldots, x_k)$ spread over $\tilde{X}$ such that $X(\xi_N) = \begin{pmatrix} x_1 \\ x_2 \\ \vdots \\ x_N \end{pmatrix}$ is an $N \times p$ extended design matrix and $M(x, \xi_N^{(0)}) = (X' (\xi_N) X(\xi_N))^{-1}$ is a $p \times p$ normalized Fisher’s information matrix.

2. Define the measure $\xi_N^{(0)} = \{ \frac{x_1}{N}, \frac{x_2}{N}, \ldots, \frac{x_N}{N} \}$

3. Take a sample, $n (\leq N)$ support points, thus define the initial design measure $\xi_n^{(0)} = \begin{pmatrix} x_1 w \\ x_2 w \\ \vdots \\ x_n w \end{pmatrix}$; $w = \frac{1}{n}$ such that $\Delta(0) = |M(x, \xi_n^{(0)})| \neq 0$

4. Define the variance function, $V(x, \xi_N^{(0)}) = x' M^{-1}(x, \xi_N^{(0)}) x$, a $2m$ degree polynomial and compute $V(x, \xi_N^{(0)})$.

5. From step 4, obtain the gradient vector, $g_{ik} = \frac{\delta V(x, \xi_N^{(0)})}{\delta x_{ik}}$, a $2m - 1$ degree polynomial evaluated at all $i = 1, 2, \ldots, N$.

6. Define the k-component direction vector $d = \sum_{i=1}^{N} \theta_i g_{ii}$; $\theta_i \geq 0$, $\sum_{i=1}^{N} \theta_i = 1$

7. Obtain $\theta_i$ from the equation $\frac{\delta V(d)}{\delta \theta_i} = 2 \sum_{i=1}^{N} \theta_i V(x) = 0$

8. Obtain the starting point, $x_0$.

9. Compute the step length, $\rho$ from the equation $\frac{\delta V(x, \xi_N^{(0)})}{\delta \rho} = 0$.

10. Move to the point $x_j^* = x_0 + \rho d_j$ and at the $j^{th}$ step move to $x_j^* = x_{j-1}^* + \rho_j d_j$

11. If $V(x_j^*, \xi_N) > V(x_{j(min)}, \xi_N)$, exchange $x_j^*$ with $x_{j(min)}$, else, stop and check for convergence.

Is $\|\Delta(j) - \Delta(j-1)\| < \epsilon \geq 0$?

Yes: stop, $\xi_N^{(j)}$ is $D-$optimal

No: $\xi_N^{(j)} = \xi_N^{(j-1)} + x_j^*$ and continue from step 2.

3.2. Starting Points

The starting points to be considered for the construction of second order $D-$optimal designs are given in equations (6), (7) and (8).

$\bar{x}_0 = \frac{\sum_{i=1}^{N} x_i}{N}$ (6)

$\bar{x}_0 = \sum_{i=1}^{N} \theta_i x_i$ (7)
where $\theta_i$ is the weighting factor for each $x_i$.

$$\Xi_0 = \sum_{i=1}^{N} \alpha_i x_i$$

(8)

where, $\alpha_i = \frac{(\hat{z}_i^T \hat{z}_i)^{-1}}{\sum_{i=1}^{N} (\hat{z}_i^T \hat{z}_i)^{-1}}$

4. Discussion of Results

Eleven support points chosen from $\tilde{X}$ are:

$$\{(−1, −1); (−\frac{1}{2}, −1); (0, −1); (1, −1); (−1, 0); (−\frac{1}{2}, 0); (0, 0); (1, 0); (−1, 1); (0, 1); (1, 1)\}.$$

$V(x, \xi^{(0)}_n), g, g^j, d^j_0$, and $\rho$ are computed from the algorithm presented in section 3.1. The algorithm was implemented using MATLAB, a software that handles matrices efficiently. The three starting points are separately applied to the variance-exchange algorithm to obtain a minimum point design in the experimental space: $\tilde{X} = \{x : x_1, x_2; −1 \leq x_1, x_2 \leq 1\}$ for the model in (5). This implies that from the starting points in (6), a move is made to $\hat{z}_j^* = \hat{z}_j^{(0)} + \rho_j d_j^*$ and $\hat{z}_j^*$ is obtained. The same procedure is carried out using (7) and (8) separately but with the same support points and $\xi^{(0)}_n$. At convergence, the design points and determinants obtained for each of the starting points are shown in Table 1. For clarity, (6), (7) and (8) shall be referred to as starting points I, II, and III respectively. The $|M(\hat{z}, \xi^{*}_n)|$ is calculated for each starting point and the optimal designs generated are compared with Box and Draper (1974) $D-$optimal minimum point designs using $D-$efficiency. Results obtained show that the $D-$optimal designs generated compare well with those of Box and Draper (1974) but starting point III has the highest $D-$efficiency value. In addition, the combined quantile plots are used to compare the performance of designs generated by the starting points. For more details on the combined quantile plots, see, Khuri, Kim and Um (1996), Chigbu and Umelo-Ibemere (2015). Graphical comparison reveals that starting point III generated design points have the lowest quantile values for radius, $r = 0.3, 0.7$ but have equivalent quantile values for $r = (1.1, 1.4)$. Information obtained from the plots reveal that the three starting points perform well. However, starting point III slightly performs better than I and II.

Table 1: Comparison of the $D$–optimal designs for each starting point

| starting point | $x^*$ | $|M(x^*,\xi)|$ | $D_{-\text{eff}}$ |
|---------------|-------|----------------|----------------|
| I             | (-1,-1);(-0.0813,0.0499); (1,-1);(-1,1);(1,0.4);(0.4,1) | $0.55 \times 10^{-2}$ | 0.9929 |
| II            | (-1,-1);(-0.1403,0.0444); (1,-1);(-1,1);(1,0.4);(0.4,1) | $0.55 \times 10^{-2}$ | 0.9929 |
| III           | (-1,-1);(-0.1867,-0.0952); (1,-1);(-1,1);(1,0.4);(0.4,1) | $0.57 \times 10^{-2}$ | 0.9988 |
| Box and Draper | (-1,-1);(-1,1);(-0.1315,-0.1315); (1,0.3945);(0.3945,1) | | |

5. Conclusion

The choice of optimal starting point in line search methods influence the optimal design points generated at convergence. Although the starting points considered are reasonable estimates of the mean of the support points, starting point III has the highest determinant of the information matrix. The algorithm converged using the three starting points. Graphical comparison of the variances of the optimal design points generated also reveal convergence to $D$–optimality with the lowest variation in starting point III generated designs. In comparison with Box and Draper (1974) minimum point design, the design points generated with starting point III are more efficient than those of the other starting points.

References


Fig. 1. quantiles of the variance functions, $V(x, z_n)$ for the $D$–optimal designs using starting points I, II, and III.
