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COMPARISON OF TWO METHODS FOR DESIGN CENTERING

by

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ABSTRACT

Two methods recently proposed to solve the design cantering problem [2,9] are compared. Although the methods are formulated differently, they are shown, under general assumptions, to yield the same solutions. Simplifications required to make the methods efficiently implementable introduce, however, significant differences from an utilization point of view.

1. Introduction

Finding the best nominal design in order to maximize the yield is an important problem in tC design. Because of the unavoidable fluctuations in the manufacturing process, the actual values of the circuit parameters, denoted by the vector p, is characterized by the known joint density distribution <jrtp-p°). where p° represents the nominal values. Oesign centering methods [1.5,31 try to imbed in the so-called region of acceptability R_{Jf} the largest convex domain B_r(p) related to <pip-p°). as shown below. The region of acceptability can for our purposes be defined as the *i*mt

$$R_{a} = \{p \mid f_{i}(p) \le f_{max}, j = 1,...,m\}$$
(1)

where the f, represent the performance functions which characterize the circuit behavior. R_t is assumed to be simply connected. The contours of equal probability of ?<p) can be associated for all the distribution of interest with a norm n(p). See for example [3]. $B_r(p)$, often referred to as a norm body, is defined relative to n(p)

$$B_r(p^\circ) \ll \{p|n(p-p^\circ) \le r\}$$
⁽²⁾

and represents a body centered at p° whose size is proportional to r. The first method we will look into is the approach referred to as (VTP) in [2], and can be formulated as

uch that max max
$$f_i(y) \leq f_{max}$$
 (3)
i $y \in \mathfrak{S}_f$

In this approach a maximally sized body is to be found, such that inside the body none of the performance functions wiH exceed their maximum allowable value. We note from the outset that the main difficulty in solving (3) derives from the maximization subproblem

as it must be solved for all constraints i * 1....m and is likely to have local maxima, making gradient based methods unreliable. That (4) is likery to have local maxima can be inferred from the fact that quite often the $t_i(y)$ are convex functions and $B_r(p^\circ)$ is a parallelepiped.

The second method, referred to as the *{HP*) method, is formulated as:

$$\{UP\}: \max_{r} \min_{i} \min_{m} rtfyp^{0}\})$$

$$: r i W_{r} > .*_{r}$$

$$(5)$$

In this method the points on the boundary of the region of acceptability which are closest to the nominal design are located and then this distance is maximized. Searching for these near points is based uoon the fact that they limit further expansion of the body $B_r(p)$. It is possible to prove that if the performance functions are either quastconvex or quasiconcave, their accumulation points are the points where the largest body touches the boundary of the region of acceptability.

The domain of the innermost minimization is the intersection of the region of acceptability with the surface $f_{,,}(y) * f^{,}$, i.e. in the boundary of R_a , we will assume this domain to be nonempty. Otherwise the constraint ty $f_{,,}(y) = f_{,,}(y) + f_{,,}(y) + f_{,,}(y)$. Otherwise the constraint ty $f_{,,}(y) = f_{,,}(y) + f_{,,}(y)$. Otherwise the constraint ty $f_{,,}(y) = f_{,,}(y) + f_{,,}(y)$. The main sunchanged An algorithm can easily detect this situation by verifying that for the constraint i there is no feasible solution to the minimization problem. The main difficulty that arises in solving (5) is the subproblem

minimize nfj(y)_{<} < f^{\land}, j > 1....m
 $f_{i}(y) \ge f_{max}$
(6)

The difficulty here is that (6) is a constrained minimization problem which is computationally intensive in itself, and one which must be solved repeatedly.

2. Equivalence of Methods

It is illustrative to recognize that if R_a is simply connected, the performance functions are differentiate and their gradients do no vanish at the boundary of R_{a} . bd(R_{a}). then a locally optimal solution to (3) is also locally optimal to (5).

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For convenience we define

andthesats

$$M_{f}(p,r) = 01 |f_{r}(0) .max f_{r}(p)>$$
 (8)

VTncenteUi proved [10] that a solution 0°, ?) to problem (3)* locally optimal if

$$f_{i}(M_{i}(\beta^{\circ}, r^{\bullet})) \leq f_{max}$$
(9) and

$$0 \in CO y_t \{Vfjtp\} \mid pcM.0 ^{\circ}/).fj(p) = I_{max} \}$$
(10)

This last **expression** can be mierprated **geo**metrically to **mean** that the convex hull defined by gradients of the cunauaima at the points where the body touches the boundary of R_t , contains the origin, intuitively, if this situation occurs, there is no direction in which the center could be moved such that the radius of the body COUWincress.

To establish the equivalence between the two problems^{*} it is enough to show that any point $0 \in M_1(f1^\circ, f)$, exists in $bd_1(Ra)$ and is at a nutimal distance from &. Assume ttwe is a point p"ebd₁(Ra) such that $n < p^* S^\circ > < n(fi^\circ \circ)$ then p* exists in the interior of B_r. (f1°) and there is an «>0 such that $B_e(p \circ CB? \ll \circ)$. By assumption $Vf_1(p^*) = m \circ 0$ and there ta therefore a point $P^* \in B_p(p^*)$ such that $ftp^* > Kp^*$). But this contradicts the hypothesis that (0 V) is optimal as for p* we would have p* $\in B_p(p^\circ)$. If $p > \infty$.

3. Relation Between Subproblems (4) and (6)

The relation between problems (4) and (8) can be made dearer by dropping from (6) the requirement that $f-fi)^{A}_{nm}$. i « i.~ « ii,i + i,...,m. if the constraint i is not superfluous, i.^se constraints for most cases are not active and the solution to (6) would remain unchanged. We rewrite therefore (6) in the following form

$$\begin{array}{ll} \text{minimize} & \mathbf{n}(\mathbf{y} - \mathbf{p}^{\circ}) & (11) \\ \text{such that} & \mathbf{I}(\mathbf{p}) \geq \mathbf{I}_{max} \end{array}$$

On the other hand, using (2) subproblem (4) can also be rewritten as

maximize
$$f_i(y)$$
 •
such that $a(y-p^3) \leq r$ (12)

for p° and r constants.

As illustrated in Rg. 1 for the **norm** $n(x) = ||x|| = max ||x, 1^-1,...^n$, in subproblem (10) for a constant sized paraHeHpiped we search for the



point where f_1 is maximized. In (11) the opposite problem is solved: given a fixed value of the function f_p , the minimum value of r is found such that the resulting norm body will have at least a common point with the surfac $f_1(p) - t_{(xmx)}$.

It is interesting to rfote that (11) and (12) are in some sense duals of each other since they have the objective function *and* the constraint interchanged. We saw before that the solution to these two subproblems for $\{\$, ?\}$ was the same. We further notice that a local solution to (11), assuming that n(x) is locally Lipschitz, is given by [4].

$$0 \notin_{o} 3rKy - p^{\circ}) - r_{1} Vf_{\cdot}(y)$$

$$r_{0} r_{1} \ge 0$$

$$r_{1} [t_{i}(y) + t_{max}] = 0$$
(13)

where $3n(y-p^{\circ})$ represents the generalized gradient of $n(y-p^{\circ})$. Similarly, for (12)

$$0 \in -r_0 \nabla f_j(y) + r_1 \partial n(y-p^0)$$

$$r_0 r_1 \ge 0$$

$$r_1 n(y-p^0)-r_1^2 = 0$$
(14)

The similarity of these problems is enchanced if we nouca that at the solution of (3) and (5) we must have r ^ .

4. Implementation of the Methods

The implementation of the two methods is similar in the sense that the subproblems (4) and (6) are solved at each iteration and the function and gradient information gathered during this process is used by the outer maximization. In the first case the problem can be reduced to a constrained nonlinear problem which can be solved by some constrained variable metric methods. Specsficafly, a variation of Poweffs algorithm [7] is used, in the second method, the equivalent information is used to generate a second order approximation to the constraining surfaces and the largest normbody is inscribed in that simplified approximation to the region of **acceptability.**

Serving subprotiveni (4) at each iteration would result in a **computationally** sxpansnw agofiuvn. nwrarora nw nonnoaay $B_f(p)$ is immicad to ttw cas* wtwr* ft* corresponding norm is $I \mid_{OO} L^{Le}$.

$$\mathbf{B}_{r}(\mathbf{p}) = \{\mathbf{y} \mid |\mathbf{y}_{t} \cdot \mathbf{p}_{t}| \leq r\}$$

$$(15)$$

The basis ror irns resmcoon is xnac m most practical **cases** the maximum of (4) with occur at a vertex of $8_r(p)$, therefore reducing the set where the search is to be done to a finite set. This set can however still be very large and in [2] a scheme referred to as splitting is introduced. TWa technique, which is baaed on previous information, predicts where the local maxima of (4) are Hkaty to occur. Reducwgmeiearchfbramaxirnumtothesatofverticaeof $B_r(p)$ can however introduce significant errors in the case where the region of acceptability is not one (finenakraly convex [1], or the region of acceptability has holee in its interior, as iHuatratad in **Fig. 2**.



The (NP) method doea not have thia limitation*

Although methods exist tar immininad groblems when the objective function is imodifterentiaMe. ag> [6], and thia is presentfy an area of active raae*ch, all algorithms for thia general problem tend to require a very large number of function and gradient (generalized gradients) evaluations making them unsuitable to be used in subproblem (6) where repealed solutions are required. In our HiHxemeiHation we limited the algorithm (NP) to the casee where the norm is differentfable, and used alao Powers constrained variable metric algorithm to solve subproblems (6).

From the designer's point of **view**, the (NP) method yields interesting information on howweilthe constraints are formulated. At early stages of a design, it is very often the case that a large number of cmisuaintt *m* tentatively specified, dearly if the solution to (6) corresponds to a value of t M ^ the constraints can be modified. Further, the relative distances to the final nominal?⁰ give an indication on how strongly the* constraint affects the circuits yield.

A limitation of the NP method which could be serious is 9 the initial design is very far from the final solution, the quadratic approximations might have to be updated, thereby significantly increasaig xnecoai of me atgoninni.

5. Examples

We tried both methods on a group of examples to compare the behavior of both methods. For the (VTP) we used the infinite norm, while for the (NP) we used the L-norm, this accounts for the

different solutions obtained with both methods. (Note also the theoretical advantage of the (VTP) methods of not requiring a feasible starting point).

Example 1 In this example we have a single aonve. Quadratic performance function

constrained to be smaller than a1.









Example 2 Thia example is taken from (8]

$$\begin{split} f_1(\mathbf{p}) &= e^{-p_1 + 1} \left((p_2 - 1)^2 + 1 \right) \\ f_2(\mathbf{p}) &= e^{p_1 - 2p_2 + 1} \\ f_n(\mathbf{p}) &= p_1^2 + p_2^2 - 1 \end{split}$$

The constraints are 4≤1.5, i = 1,2.3

Starting Point FtnalPoint Number of F.E. Number of G.E. (VTP) (2.0,2.0) (03a1.01) 18 20 (NP) (07,09) (035,1.01) 15 46 $\frac{1}{100}$ Example 3 This example was taken from [8]

| $f_1(p) = 1.5 - \rho_1(1 - \rho_2)$ | $f_4 = -f_1$ |
|---|------------------|
| $f_2(p) = 2.25 - p_1(1 - p_2^2)$ | $f_{5} = -f_{2}$ |
| $f_3 = 2.625 \cdot \rho_1 (1 \cdot \rho_2^3)$ | $f_6 = -f_3$ |

constrained to $f_i \leq 1.5$, i = 1,...,6



Example 4 Also taken from [8] has two performance function, one convex and the other concave, both constrained to be smaller than 1.15

 $f_{1}(p) = 0.96 + (p_{1}-1)^{2} + (p_{2}-0.8)_{2}$ $f_{2}(p) = 1.16 - (p_{1}-1)^{2} - (p_{2}-0.6)^{2}$ Starting Point Final Point Number of F.E. Number of G.E. $(VTP) \quad (2.0,2.0) \quad (1.000,1.004) \quad 31 \qquad 32$ $(NP) \quad (0.8,0.9) \quad (1.000,1.016) \quad 11 \qquad 22$

For this example and for function f_2 the solution to subproblem (4) is not in a vertex of the parallelipiped and therefore a significant part of the final tolerance box is infessible.



Conclusion

We have shown that two apparently very different problems can be seen as different formulations of the same body imbedding problem, yielding under general conditions an identical result. With reasonable assumptions the problems lend themselves to very different implementations : one (VTP) suitable for the worst case problem, the other (NP) for a general body center problem when the norm is differentiable. It is interesting to note further that when applied to similar problems they seem to have computational requirements of the same order, when measured in terms of the number of function and gradient evaluations.

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