

TOLERANCE SYNTHESIS FOR NONLINEAR  
SYSTEMS BASED ON SENSITIVITY ANALYSIS

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## **Abstract**

Automatic assignment of tolerances to dimensioned mechanical assemblies is studied as an optimization problem: the objective of which is to minimize the (manufacturing) cost, subject to the constraints of (design) functionality and (assembly) interchangeability. By associating a nominal dimension and a tolerance to the variance, the probabilistic approach is taken.

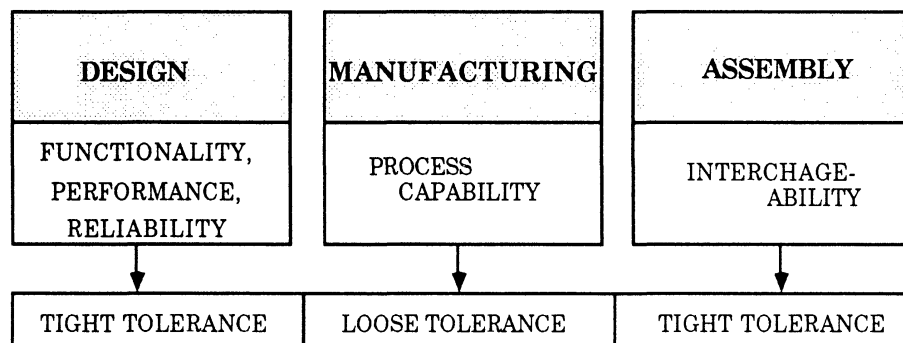
Trigonometric functions relating the component geometries give rise to the nonlinearity in the system. Estimating an n-dimensional nonlinear integral by a polytope converts the probabilistic optimization formulation to a deterministic one. It also allows speedy evaluation of tolerance analysis embedded in tolerance synthesis.

Local optimality is ensured by analysis of convexity and quasi-concavity of the objective function and some of the constraints. Sensitivity analysis is performed to provide search directions for global optimality. An implementation is reported with an example.

# 1. Introduction

Tolerances in an engineering design are intended to capture variations from the ideal, as introduced by the very process of realization such as manufacturing and assembly. Nominal dimensions specify idealized geometries by size, location and form. The range between the upper and lower limits of the variation from the nominal dimension is called *tolerance* [1].

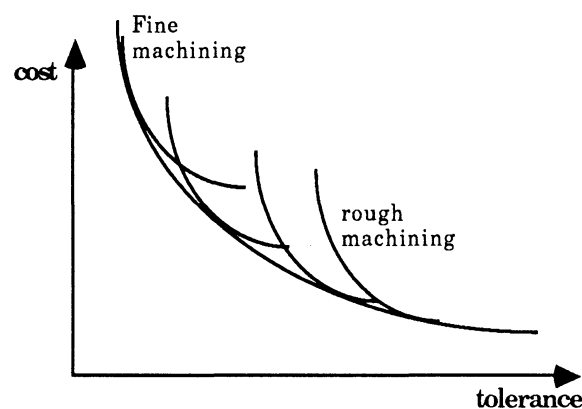
At the design stage, functionality, performance and reliability are the major issues under consideration. Tolerance, or variation from the ideal, should be set to be as close to zero as possible. However, high precision or tight tolerances are usually associated with high costs; looser tolerances are less costly at the manufacturing stage. Yet, at the assembly stage, due to the objective of optimizing the interchangeability of components, tight tolerances are desirable. While design and assembly prefer tight tolerancing, virtual components (from design) must be brought to realization by manufacturing before physical components can be assembled. This requirement results in a three-way trade-off amongst design, manufacturing and assembly as shown in Figure 1.



**Figure 1. Concurrent Consideration of Tolerance**

The concurrent consideration of opposing criteria on tolerancing, between design and manufacturing, and between manufacturing and assembly, can only be resolved by compromise. The result of such a rationalization, if agreeable to all concerns, is effectively a synthesis of tolerances. This paper deals with computational techniques for tolerance synthesis by analysis.

Basic to component manufacturing is cost. To assign cost-effective product tolerances, probabilistic tolerancing is considered in this paper. The probabilistic approach is considered to be advantageous over the deterministic approach, because it is possible to perform trade-off analysis with the probabilistic approach [13,19]. The problem, simply stated, is to convert the designer's "function-oriented" specifications into manufacturable specifications by allocating the tolerances to the ideal dimensions such that the manufacturing cost is minimized. Models of tolerance-cost functions from [10,19,22,23,25] are employed. Figure 2 shows a typical inverse relation of manufacturing cost to tolerances: the tighter the tolerance the higher the cost.



**Figure 2. A Typical Tolerance-Cost Curve**

To make the optimization effective for the probabilistic tolerancing approach requires efficient algorithms to estimate the yield and its sensitivity with respect to tolerances. A major obstacle of probabilistic tolerancing for general nonlinear system is the tremendous amount of CPU time required to recompute the yield at each iteration of the optimization. To reduce the time for the yield computation, Parkinson [20] and the authors [14,15] used the notion of reliability index [9] as an approximation of the yield. However, the computation of the approximation is still intensive. In [14], a local circular search was used to find a global optimal solution. Michael and Siddall [19] decomposed the random variable space into orthogonal  $n$ -dimensional cubes, where  $n$  is the number of dimensions considered; that gives  $O(2^n)$  cells to be tested. Other attempts which impose restrictions on the domain of the problem such as, linearity [2,7,11,24,25,26] and single design constraint [22], have practical limitations.

The purpose of this paper is to provide a general framework for tolerance synthesis for nonlinear systems with multiple, dependent<sup>1</sup> design constraints. Least-cost tolerance synthesis is mathematically formulated as a nonlinear programming (NLP) problem. Algorithms are provided and illustrated by an example. Post-optimality analysis is also considered so as to allow the possibility of modifying some yield constraints in the design process.

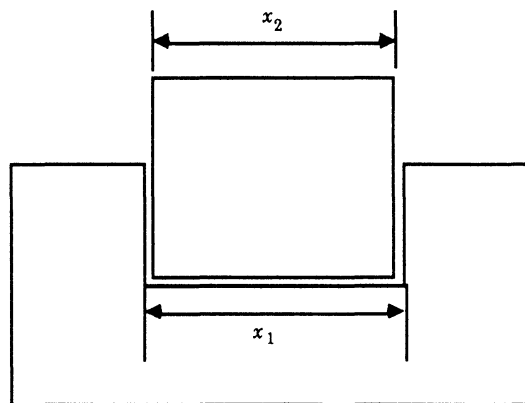
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<sup>1</sup>. The design constraints are termed "dependent" if they share the same dimension variables. For instance, the two design constraints  $x_1 - x_2 - 0.1 \leq 0$  and  $x_1 + x_3 - 0.05 \leq 0$  are dependent because they share the same variable  $x_1$ .

## 2. Basic Concepts

### 2.1 Design Function

Design and assembly are concerning more with inter-component relations. Consider a simple assembly as shown in Figure 3.



**Figure 3. An Assembly**

Suppose the clearance between the components has to be greater than or equal to 0.01 to achieve certain performance and assemblability. This requirement can be expressed mathematically as:

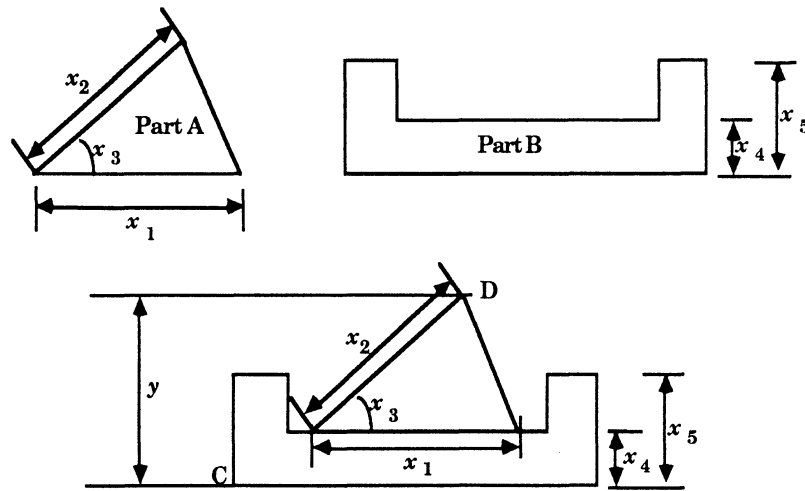
$$x_1 - x_2 \geq 0.01, \quad (1)$$

where  $x_1$  and  $x_2$  denote the dimensions of the hole and the shaft respectively as shown in the figure. Rewriting inequality (1) as a function  $F(\mathbf{x},c)$  gives:

$$F(\mathbf{x},0.01) = -x_1 + x_2 + 0.01 \leq 0. \quad (2)$$

Functions such as (2) are called *design functions*. They describe the inter-component relations, and provide a mathematical basis for

controlling functionality and interchangeability. In (2), 0.01 is referred to as a *design constant*  $c$  of the design function  $F(\mathbf{x},c)$ . A design function is not always linear. For example, if the assembly is not recti-linear, as illustrated in Figure 4, some of the design functions will take on trigonometric terms.



$y$  : vertical distance from C (of part B) to D (of part A) is equal to  $x_4 + x_2 \sin x_3$

**Figure 4. A Non-linear Design Function**

A linear system is one in which an assembly dimension  $y$  is defined by a linear combination of component dimensions  $x_1, x_2, \dots, x_n$ :

$$y = \sum_{j=1}^n a_j x_j \quad (3)$$

where  $a_j$  is a signed binary integer. As variations accumulate, the tolerance analysis (or "stack-up", as it is commonly referred to) for linear systems is useful and has been studied extensively [2,7,11,24,25,26]. The key property that facilitates the analysis is that the variance of the linear sum is

the sum of the variances of component dimensions under the assumption of independence, that is,

$$\sigma_y^2 = \sum_{j=1}^n (a_j)^2 \sigma_{x_j}^2. \quad (4)$$

When tolerance  $t_j$  of dimension  $x_j$  is defined as  $\pm 3\sigma_j$  from the mean  $\mu_j$ , the assembly tolerance  $t_y$  can be then represented by

$$t_y = k_y \sqrt{\sum_{j=1}^n (a_j)^2 \left(\frac{t_j}{6}\right)^2} \quad (5)$$

where  $k_y$  is a constant derived from an allowable percentage  $\lambda_y$  of defect in the assembly. In the case of normal distribution,  $k_y = 2 * \Phi^{-1}(1 - \frac{\lambda_y}{2})$ .

However, the analysis procedure for linear system can not be extended to nonlinear system because of the lack of a general rule for an aggregate such as equation (4).

## 2.2 Yield

The probabilistic approach is performed under the assumption that the dimension vector  $\mathbf{x}$  follows the multivariate normal distribution. (Indeed, Mansoor [18] shows that most manufacturing processes produce dimensions with normal distributions.) Let  $\mu_j$  and  $\sigma_j$  denote the mean and the standard deviation of the normally distributed random variable  $x_j$ . The mean  $\mu_j$  is typically fixed by the designer, whereas the standard deviation  $\sigma_j$  is chosen according to the precision of the controls exercised over the manufacturing process. This parameter  $\sigma_j$  is therefore a function of the



tolerance  $t_j$ , and according to standard practice,  $\sigma_j$  is normally set to  $\frac{t_j}{6}$ . Clearly, if  $t_j$  is given,  $x_j$  is a well defined random variable.

Probabilistic tolerance analysis can be stated more succinctly in a mathematical formulation: Given tolerances  $t_j$  (or standard deviations  $\sigma_j$ ), determine the probability such that the design function  $F(\mathbf{x},c)$  is less than or equal to zero, i.e.,  $Pr(F(\mathbf{x},c) \leq 0)$ . In other words, evaluate the integral,

$$\text{yield} \equiv Y(\mathbf{t}) = \int_{F(\mathbf{x},c) \leq 0} f(\mathbf{x};\boldsymbol{\mu},\mathbf{t},\mathbf{R}) d\mathbf{x} \quad (6)$$

where  $F(\mathbf{x},c)$  is a design function, and  $f(\mathbf{x};\boldsymbol{\mu},\mathbf{t},\mathbf{R})$  is the probability density function of multivariate normal vector  $\mathbf{x}$  for which  $\boldsymbol{\mu}$ ,  $\mathbf{t}$ , and  $\mathbf{R}$  denote the mean vector, the tolerance vector (standard deviation vector), and the correlation matrix of  $\mathbf{x}$ , respectively. (If the random variables  $x_j$  are independent, it will not be necessary to specify covariances.) The integral (6) shall be referred as the *yield*.

It helps the intuition to visualize the interaction between the tolerance and the design function. Consider an assembly of two random variables,  $x_1$  and  $x_2$ . A design function  $F(\mathbf{x},c)$  partitions the two-dimensional space into two regions. The *safe region*  $\mathbf{R}_S$  in Figure 5(a) corresponds to the region in which  $F(\mathbf{x},c) \leq 0$ . (The complement of the safe region is called the *failure region*  $\mathbf{R}_F$ .) Now, the given tolerances also prescribe a region in the same space. In Figure 5(b), the upper and lower limits of a random variable  $x_j$  define a strip. Intersecting the strips for  $x_1$  and  $x_2$  gives the *tolerance region*  $\mathbf{R}_T$ . It is noted that the size of  $\mathbf{R}_T$  varies with tolerances, but  $\mathbf{R}_S$  (or  $\mathbf{R}_F$ ) is independent of tolerances. Combining  $\mathbf{R}_T$  with  $\mathbf{R}_S$  gives the *reliable region*  $\mathbf{R}_R$  as shown in Figure 5(c). For any system to perform reliably,  $\mathbf{R}_R$

must be non-empty. (Such would be the case if the tolerances were not assigned properly or the design function was incorrectly specified.)

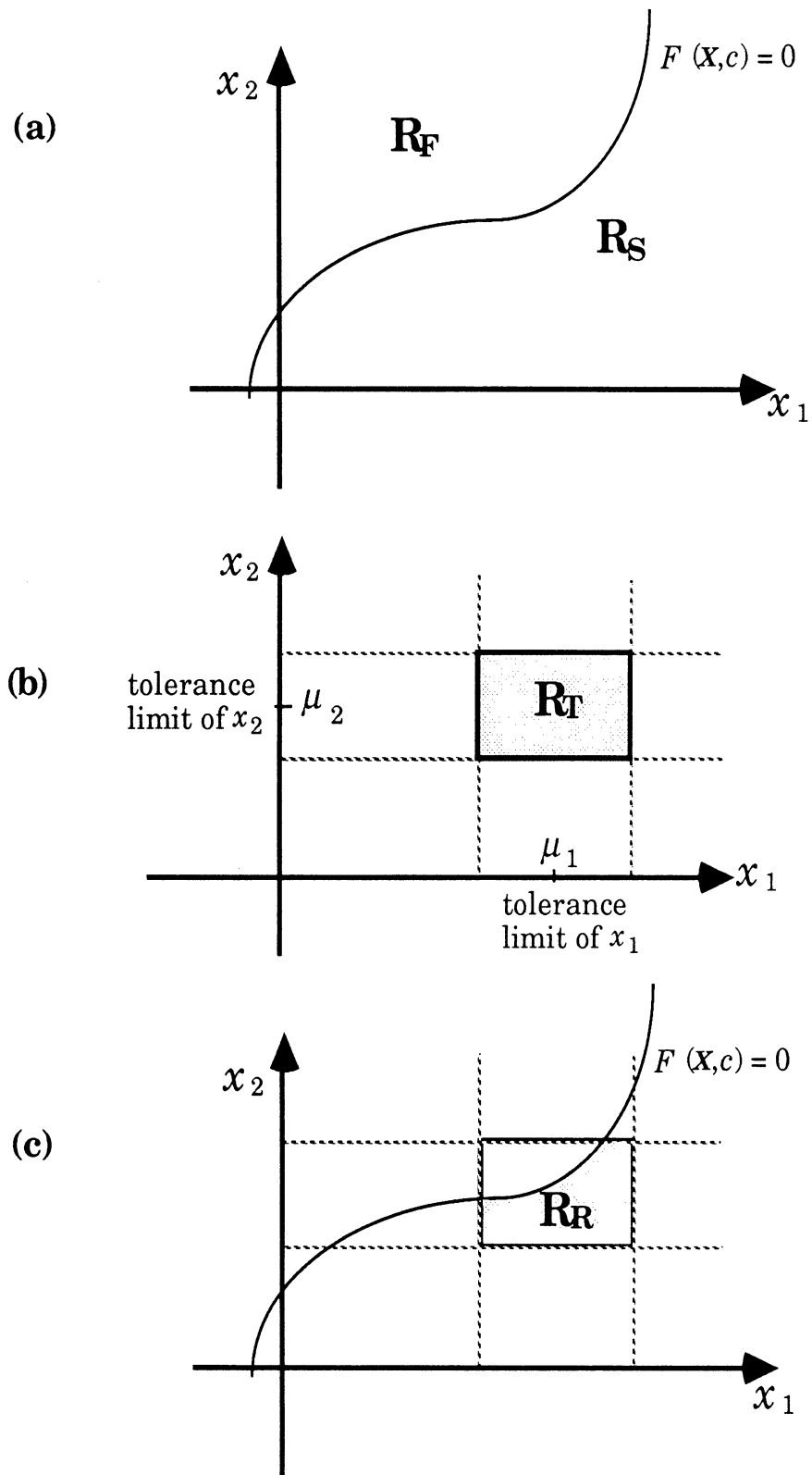


Figure 5. Safe, Tolerance and Reliable Regions

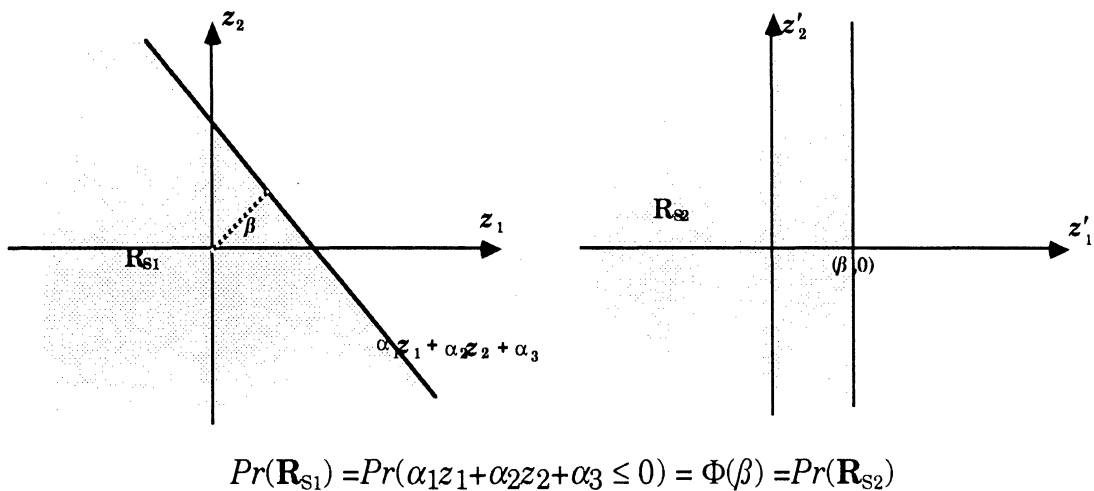
### 3. Tolerance Analysis

#### 3.1 Yield of Linear Design Function

While this paper addresses systems with nonlinear design functions, discussion of the linear case facilitates subsequent development as nonlinear functions will be approximated by hyperplanes. The yield of a linear design function  $F(\mathbf{x},c)$  as computed from (6), involves multiple integration. An approximation method which can be extended to the nonlinear case<sup>2</sup> adopts the notion of the *reliability index*, which was introduced by Hasofer and Lind [9]. Consider Figure 6, in which there are two independent random variables  $z_1$  and  $z_2$  both following the standard normal distribution  $N(0,1)$ . The (standardized) linear design function is given by the form  $\alpha_1 z_1 + \alpha_2 z_2 + \alpha_3$ . The desired yield  $Pr(\alpha_1 z_1 + \alpha_2 z_2 + \alpha_3 \leq 0)$  is equal to  $\Phi(\beta)$ , where  $\beta = \frac{|\alpha_3|}{\sqrt{\alpha_1^2 + \alpha_2^2}}$  is the minimum distance from the design function to the origin. The distance  $\beta$  is referred to as the *reliability index* of the design function. Because of the rotational symmetry of the standard normal coordinate, the yield can be obtained by looking up the value of  $\Phi(\beta)$  in the univariate standard normal distribution table. The above technique is next generalized for computing the yield of a design function with any number of random variables.

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<sup>2</sup> The observation that a linear combination of multivariate normal random variables follow a univariate normal distribution [6, p.56] does not extend readily to the case of nonlinear design functions.



**Figure 6. Rotational Symmetry of Standard Coordinate**

In general, tolerance analysis for linear design functions  $F(\mathbf{x}, c) = \mathbf{a}^T \mathbf{x} + c$  requires the following steps:

- (i) **Standardization** : Transform the dependent normal variables to the independent standard normal variables by using

$$\mathbf{z} = (\mathbf{PD})^{-1} (\mathbf{x} - \boldsymbol{\mu}) \quad (7)$$

where  $\mathbf{z}$  is the transformed standard normal vector and  $\mathbf{P}$  is the orthogonal matrix for diagonalizing a given covariance matrix  $\mathbf{V}$  such that  $\mathbf{P}^T \mathbf{V} \mathbf{P} = \mathbf{D} \mathbf{D}^T$ . (The detailed transformation procedure is given in Appendix A.) The transformed  $\mathbf{z}$  space is referred to as the *standard coordinates*.

- (ii) **Reliability Index Computation** : Compute the minimum distance  $\beta$  from the origin to the transformed design function in the standard coordinates. First, represent the design function in terms of  $\mathbf{z}$ :

$$\mathbf{a}^T \mathbf{x} + c = \mathbf{a}^T \mathbf{P} \mathbf{D} \mathbf{z} + \mathbf{a} \mu + c. \quad (8)$$

The distance from the origin to the right hand side of equation (8) corresponds to  $\beta$  which can be then expressed as

$$\beta = \frac{-\mathbf{a}^T \mu - c}{\sqrt{(\mathbf{a}^T \mathbf{P} \mathbf{D})(\mathbf{a}^T \mathbf{P} \mathbf{D})^T}} = \frac{-\mathbf{a}^T \mu - c}{\sqrt{\mathbf{a}^T \mathbf{V} \mathbf{a}}} \quad (9)$$

since  $\mathbf{P} \mathbf{D} (\mathbf{P} \mathbf{D})^T = \mathbf{P} \mathbf{D} \mathbf{D}^T \mathbf{P}^T = \mathbf{P} (\mathbf{P}^T \mathbf{V} \mathbf{P}) \mathbf{P}^T = \mathbf{V}$ . Note that the numerator of equation (9),  $-\mathbf{a}^T \mu - c$ , is always positive since the given nominal dimensions,  $\mu$ , are assumed to satisfy the design condition.

- (iii) **Table Look-Up** : Look up the univariate standard normal distribution table for  $\Phi(\beta)$ .

### 3.2 Yield of Nonlinear Design Functions

To solve the integral (6) under nonlinear function  $F(\mathbf{x}, c)$  and multivariate normal PDF, two techniques are considered: (i) Monte-Carlo simulation [5,8] and (ii) approximation through the linearization of  $F(\mathbf{x}, c)$  [14,15,20].

Monte-Carlo simulation starts with generating  $N$  sets of random samples  $(x_{11}, x_{21}, \dots, x_{n1}), \dots, (x_{1N}, x_{2N}, \dots, x_{nN})$  from the given multivariate normal PDF, where  $x_{jl}$  denotes the  $l$ -th sample ( $l=1, \dots, N$ ) of the  $j$ -th dimension ( $j=1, \dots, n$ ). Then, each set is substituted into the design function and the sign of the functional value is checked. Suppose  $T$  sets out of  $N$  sets have a nonpositive sign. Then, the estimate of the yield is  $\frac{T}{N}$ . While Monte-Carlo simulation can be applied to the linear or nonlinear  $F(\mathbf{x}, c)$ , it is time intensive because a large number of random samples needs to be taken to

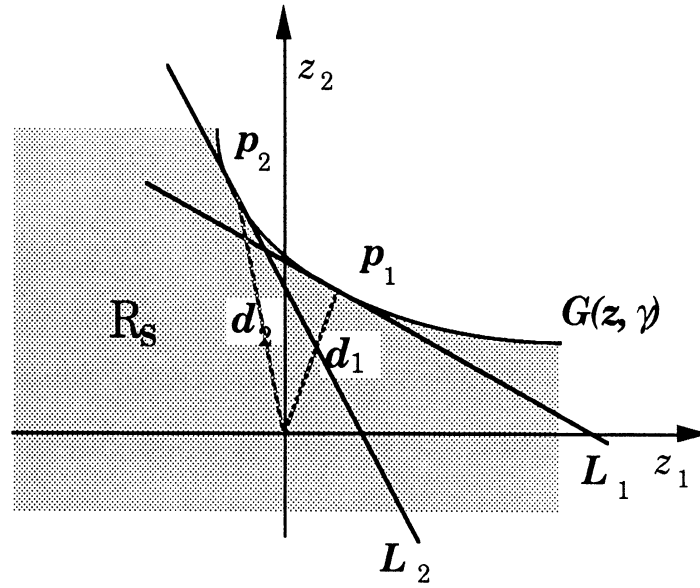
have an accurate result. This intensive consumption of time is compounded in *tolerance synthesis*, which requires iterative tolerance analysis to estimate the yield and the gradient.

In order to reduce the computational time, approximation of yield by linearization of nonlinear function is used. An expansion point of a design function  $F(\mathbf{x},c)$  is selected, then linearization is done through Taylor series expansion<sup>3</sup>. Note that the probability density in the standard coordinate decreases exponentially as the distance from the origin increases. This suggests that for an expansion point to be well-selected the (probabilistically) densest area should be preserved after linearization.

The "dense" area is estimated by the distance  $\beta$ . The "densest" area is estimated by minimizing  $\beta$ . Each design function is standardized by transformation (7), and the point on each standardized design function which has the minimum distance from the origin is selected as the expansion point. Consider two points  $p_1$  and  $p_2$  on  $G(\mathbf{z},\gamma)$ , the standardized  $F(\mathbf{x},c)$ , as candidates for expansion points. In figure 7, lines  $L_1$  and  $L_2$  correspond to linearization of  $G(\mathbf{z},\gamma)$  through  $p_1$  and  $p_2$ . Since  $d_1 < d_2$ ,  $p_1$  is closer to the origin than  $p_2$ .

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<sup>3</sup>. The first order Taylor series expansion gives the tangent hyperplane passing through the expansion point.



**Figure 7. Expansion Point for Linearization**

The expansion point having the minimum distance from the origin is obtained by solving the following single constraint NLP:

$$\text{Min } \beta = \sqrt{\mathbf{z}^T \mathbf{z}} \quad , \quad \text{subject to } G(\mathbf{z}, \gamma) = 0. \quad (10)$$

Based on equation (7), formulation (10) can be rewritten with the original variables:

$$\text{Min } \beta^2 = (\mathbf{x} - \boldsymbol{\mu})^T \mathbf{V}^{-1} (\mathbf{x} - \boldsymbol{\mu}) \quad , \quad \text{subject to } F(\mathbf{x}, \mathbf{c}) = 0. \quad (11)$$

Here, the objective function is expressed as  $\beta^2$  instead of  $\beta$  since the positive definiteness of the covariance matrix  $\mathbf{V}$  always guarantees the same solution. As a solution scheme for (11), the iterative method based on Newton-Raphson method [17] is used:

$$\mathbf{x}^{(k+1)} = \boldsymbol{\mu} + \mathbf{V} \nabla F(\mathbf{x}^{(k)}, \mathbf{c}) \frac{(\mathbf{x}^{(k)} - \boldsymbol{\mu})^T \nabla F(\mathbf{x}^{(k)}, \mathbf{c}) - F(\mathbf{x}^{(k)}, \mathbf{c})}{\nabla F(\mathbf{x}^{(k)}, \mathbf{c})^T \mathbf{V} \nabla F(\mathbf{x}^{(k)}, \mathbf{c})} \quad (12)$$



where  $\mathbf{x}^{(k)}$  denotes the solution after the  $k$ -th iteration and  $\nabla F(\mathbf{x},c)$  is the  $n \times 1$  gradient vector of  $F(\mathbf{x},c)$  at  $\mathbf{x}$ . As an initial solution for (12), the mean nominal vector  $\boldsymbol{\mu}$  is used.

Tolerance analysis for nonlinear design function therefore takes the following steps:

- (i) **Expansion Point Finding** : Find the expansion point  $\mathbf{x}^*$  by iteratively using equation (12).
- (ii) **Reliability Index Computation** : The reliability index  $\beta$  is computed by  $\sqrt{(\mathbf{x}^*-\boldsymbol{\mu})^T \mathbf{V}^{-1}(\mathbf{x}^*-\boldsymbol{\mu})}$
- (iii) **Table Look-Up** : The yield  $Y(\mathbf{t})$  (i.e.,  $Pr(F(\mathbf{x},c) \leq 0)$ ) is approximated by  $\Phi(\beta)$ , which can be looked up in the univariate standard normal distribution table.

#### 4. Least Cost Tolerance Synthesis

The problem in the context of tolerance analysis is: given design functions and tolerances, compute the yield. Suppose, instead, the yield is given, how are the tolerances to be assigned to each dimension? Figure 8 illustrates the impact of two different sets of tolerances on the yield. In the figures, the concentric circles represent the contour of equal probability density. It is noted that the tighter set of tolerances (as in Figure 8(a)) results in a higher yield as compared to the yield with the looser set of tolerances (as in Figure 8(b)).

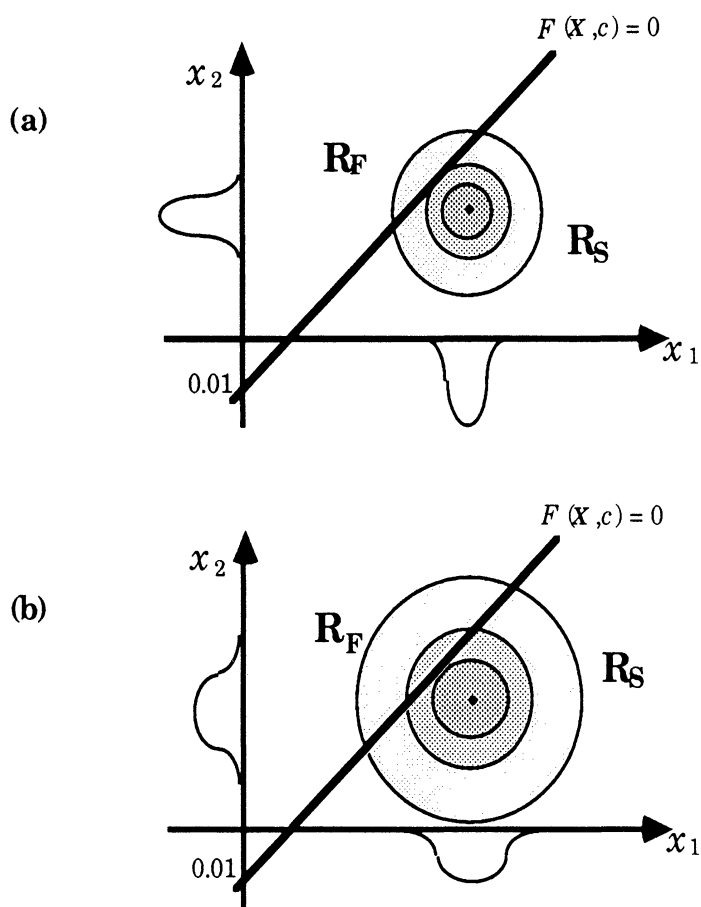


Figure 8. Impact of Different Tolerances on Yield

A higher yield, while desirable, is achieved at the expense of other considerations such as cost and manufacturability. To resolve this conflict, heuristic rules have been proposed [2]: (i) equal tolerances, (ii) tolerances proportional to dimensions, and (iii) tolerances proportional to process deviations. While heuristics are practical, it would be intellectually satisfying to see if assignments of tolerances can be optimized.

#### 4.1 Tolerance-Cost Model

As indicated in the introduction, it is generally accepted that there is an inverse relationship between tolerance and manufacturing cost. A number of cost models have been employed to fit manufacturing tolerance-cost sampled data [10,19,22,23,25], as shown in Table 1.

Model Name	References	Cost Function (*)
Sutherland-Roth Model	[23]	$C(t) = a t^{-b} + f$
Reciprocal Squared Model	[10, 22]	$C(t) = \frac{a}{t^2} + f$
Exponential Model	[25]	$C(t) = a \exp\{-\frac{t}{b}\} + f$
Michael-Siddall Model	[19]	$C(t) = a t^{-b} \exp\{-e t\} + f$

(\*)  $a, b, e$  constants for variable manufacturing cost  
 $f$  : constant for fixed manufacturing cost

**Table 1. Tolerance-Cost Models**

With such tolerance-cost models for the tolerated dimensions, the total manufacturing cost can be obtained by summing the individual manufacturing cost:

$$C(\mathbf{t}) = \sum_{j=1}^n C(t_j). \quad (13)$$

Note that model (13) is based on the "throw-aways" strategy, that is, the repair cost for the defect is not considered in the model. However, if the reworking cost is considered, the model becomes:

$$C(\mathbf{t}) = \sum_{j=1}^n C(t_j) + C_r(p_1, \dots, p_n) \{1 - Y(\mathbf{t})\} \quad (14)$$

where  $C_r(\cdot)$  is the cost function due to reworking and  $p_j$  is the probability of reworking the  $j$ -th dimension in case that the design function is not satisfied. The difficulty of employing model (14) is in procuring the empirical values for  $p_j$ . In this paper, the throw-away cost model (13) is adopted.

## 4.2 Mathematical Formulation

Least-cost tolerance allocation is a procedure for determining an optimal set of tolerances which minimizes the manufacturing cost and satisfies the performance requirement; the decision variables are the tolerances. Therefore, we can formulate the problem with minimizing the manufacturing cost as the objective function and satisfying the performance requirements as the constraints:

$$\text{Min } C(\mathbf{t}), \text{ subject to } Y(\mathbf{t}) \geq 1 - \lambda. \quad (15)$$

where  $C(\mathbf{t})$  is the manufacturing cost function in terms of tolerance  $\mathbf{t}$ , and  $(1-\lambda)$  is the minimal satisfactory yield, i.e., the yield given by tolerances  $\mathbf{t}$  should be greater than or equal to the given level  $1-\lambda$ . Formulation (15) implies that tolerance synthesis includes tolerance analysis.

With cost model (13) and multiple design functions, formulation (15) is rewritten as

$$\begin{aligned} & \text{Min} \quad \sum_{j=1}^n C(t_j) & (16) \\ & \text{subject to} \end{aligned}$$

$$Y_i(\mathbf{t}) \geq 1 - \lambda_i \quad \text{for } i = 1, 2, \dots, m.$$

This probabilistic optimization problem is simplified into a deterministic optimization problem through approximation of the yield by the reliability index.

This conversion into deterministic optimization differs from chance constrained programming due to Charnes and Cooper [3] in that the yield is approximated not at the origin of the standard coordinates but at the point of minimum distance from the origin. Thus,  $Y_i(\mathbf{t})$  is approximated by  $\Phi(\beta_i)$ , as suggested in Section 3.2. The constraints of (16) are rewritten as  $\Phi(\beta_i) \geq 1 - \lambda_i$ . Furthermore, by the monotonic property of the function  $\Phi(\cdot)$  the constraint  $\Phi(\beta_i) \geq 1 - \lambda_i$  can be inverted to the constraint  $\beta_i \geq \Phi^{-1}(1 - \lambda_i)$ , and the formulation becomes

$$\begin{aligned} & \text{Min} \quad \sum_{j=1}^n C(t_j) & (17) \\ & \text{subject to} \end{aligned}$$

$$\beta_i \geq \Phi^{-1}(1 - \lambda_i) \quad \text{for } i = 1, 2, \dots, m.$$

Note that  $\beta_i$  is a function (with respect to  $\sigma_j$ ); and  $\Phi^{-1}(1 - \lambda_i)$  is equal to a value  $q_i$  which can be obtained from the standard normal distribution table.

Recall that an expansion point  $x_j^*$  is obtained from solving formulation (11) with a fixed  $\sigma_j$ . However, the "optimal"  $\sigma_j$  obtained from solving formulation (17) is in general different from the initial  $\sigma_j$ . Therefore, in the solution process, as an  $x_j^*$  is changed, the constraints of (17) are changed as well. Consequently,  $\sigma_j$  may no longer be optimal or even feasible. In order to reflect the changes in  $x_j^*$ , formulation (11) (for all design functions) is added to formulation (17) as constraints, hence, guaranteeing the satisfaction of locally optimal conditions.

The local optima of (11) are the solutions that satisfy the Kuhn-Tucker necessary condition<sup>4</sup> [27]. Since (11) is a minimization problem, its constraints,  $F_i(x,c)=0$ , can be relaxed to  $F_i(x,c) \geq 0$ . With the Kuhn-Tucker necessary conditions, formulation (17) becomes:

$$\text{Min } \sum_{j=1}^n C(t_j) \tag{18}$$

subject to

$$\beta_i \geq \Phi^{-1}(1 - \lambda_i)$$

$$\frac{\partial \beta_i}{\partial x} - u_i \frac{\partial F_i(x^*,c)}{\partial x} = 0 \tag{18.1}$$

$$u_i F_i(x^*,c) = 0 \tag{18.2}$$

$$u_i \geq 0 \tag{18.3}$$

for  $i = 1, 2, \dots, m$ .

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<sup>4</sup>  $\beta_i$  and  $F(x,c)$  being differentiable at  $x^*$  are assumed.

### 4.3 Global Optimality Analysis

In Nonlinear Programming, no existing algorithm guarantees a globally optimal solution unless the objective function and the constraints are of certain forms. The existence check for the special forms is performed based on the following theorem [27, pp.43-44]: that the objective function is convex and the constraint functions are quasi-concave corresponds to a sufficient condition for global optimality. This ensures that the locally optimal solution implied by the Kuhn-Tucker necessary conditions is also a globally optimal solution.

Checking formulation (18) for the satisfaction of the Kuhn-Tucker sufficient condition proceeds as follows. The objective function is the sum of the individual tolerance-cost function (as in (13)). Cost models such as the ones in Table 1 are convex because the derivative of  $C(t_j)$  are monotonically nondecreasing with respect to tolerances. Since the sum of convex functions is also convex, the tolerance-cost function of (13) is always convex.

Similarly, the derivatives of the reliability index function are checked for quasi-concavity.  $\frac{\partial \beta}{\partial \sigma_j}$  is obtained by following the chain rule:

$$\frac{\partial \beta}{\partial \sigma_j} = \frac{\partial \beta}{\partial \beta^2} * \frac{\partial \beta^2}{\partial \sigma_j} = \frac{1}{2\beta} * \frac{\partial \beta^2}{\partial \sigma_j} . \quad (19)$$

For  $\frac{\partial \beta^2}{\partial \sigma_j}$ , rewrite  $\beta^2$  as:

$$\beta^2 = (\mathbf{x}^* - \boldsymbol{\mu})^T \mathbf{V}^{-1} (\mathbf{x}^* - \boldsymbol{\mu}) = (\mathbf{x}^* - \boldsymbol{\mu})^T (\mathbf{DRD})^{-1} (\mathbf{x}^* - \boldsymbol{\mu}) = \mathbf{g}^T \mathbf{R}^{-1} \mathbf{g}$$

where  $\mathbf{D}$  is a diagonal matrix whose the  $k$ -th diagonal element is  $\frac{1}{\sigma_k}$ ,  $\mathbf{R}$  is a  $n \times n$  correlation matrix, and  $\mathbf{g} = (\frac{x_1^* - \mu_1}{\sigma_1}, \dots, \frac{x_n^* - \mu_n}{\sigma_n})^T$ . Then,

$$\begin{aligned} \frac{\partial \beta^2}{\partial \sigma_j} &= \frac{\partial}{\partial \sigma_j} \mathbf{g}^T \mathbf{R}^{-1} \mathbf{g} = \frac{\partial \mathbf{g}^T}{\partial \sigma_j} \mathbf{R}^{-1} \mathbf{g} + \mathbf{g}^T \mathbf{R}^{-1} \frac{\partial \mathbf{g}}{\partial \sigma_j} \\ &= 2 \frac{\partial \mathbf{g}^T}{\partial \sigma_j} \mathbf{R}^{-1} \mathbf{g} = -2 (0, \dots, 0, \frac{x_j^* - \mu_j}{\sigma_j^2}, 0, \dots, 0) \mathbf{R}^{-1} \mathbf{g} \\ &= -2 \frac{x_j^* - \mu_j}{\sigma_j^2} \left( \sum_{i=1}^n \hat{\rho}_{ji} \frac{x_i^* - \mu_i}{\sigma_i} \right). \end{aligned} \quad (20)$$

where  $\hat{\rho}_{ji}$  is the  $(i,j)$ -th element of matrix  $\mathbf{R}^{-1}$ . Substituting (20) into (19) results in

$$\frac{\partial \beta}{\partial \sigma_j} = - \frac{x_j^* - \mu_j}{\beta \sigma_j^2} \left( \sum_{i=1}^n \hat{\rho}_{ji} \frac{x_i^* - \mu_i}{\sigma_i} \right). \quad (21)$$

The quasi-concavity of the reliability index function with respect to standard deviation is checked under the assumption of independence among dimension variables, i.e.,  $\hat{\rho}_{ji} = 1$  if  $j=i$ , and  $\hat{\rho}_{ji} = 0$  otherwise. Therefore, equation (21) can be rewritten as

$$\frac{\partial \beta}{\partial \sigma_j} = - \frac{(x_j^* - \mu_j)^2}{\beta \sigma_j^3}. \quad (22)$$

The quasi-concavity of the reliability index functions is thus shown by the nonincreasing monotonic relationship between the reliability index and the tolerances.

Since there is no restriction on the type of the design functions that can be used, equations (18.1) and (18.2) are not necessarily quasi-concave.



As a result, with arbitrary design functions, the optimal solution of formulation (18) is not guaranteed.

#### 4.4 Algorithmic Analysis

The feasible direction method is commonly used to solve the NLP problems. The constraints of formulation (18) comprise a feasible region. A point in the feasible region corresponds to a feasible assignment of tolerance. The total cost with the assigned tolerances is obtained by evaluating the objective function at that point.

Suppose  $\sigma^{(k)}$  is a point in the feasible region. A direction  $d^{(k)}$  is identified such that, for a sufficiently small  $\lambda > 0$ , the following two properties are true: i)  $\sigma^{(k+1)} = \sigma^{(k)} + \lambda d^{(k)}$  is feasible, and ii) the objective value at  $\sigma^{(k+1)}$  is better than the objective value at  $\sigma^{(k)}$ . In each iteration of the feasible direction method, having determined a feasible direction, a one-dimensional optimization problem is solved to maximize the improvement of the objective value.

If formulation (11) satisfies the Kuhn-Tucker sufficient condition and equations (18.1) and (18.2) are quasi-concave, the optimal solution of (18) is guaranteed by applying the feasible direction method. That is, if the feasible direction algorithm reaches a point that satisfies the Kuhn-Tucker necessary condition, then the corresponding tolerance assignment requires the least cost. The modified version of Zoutendijk's Method due to Topkis and Veinott [16,27] can be applied and is guaranteed to converge.

If either the formulation (11) does not satisfy the Kuhn-Tucker sufficient condition or equations (18.1) and (18.2) are not quasi-concave (or both), the sensitivity of yield is analyzed. The sensitivity of yield provides information about which tolerances are critical and helps determine the search direction in the optimization process for tolerance assignment. With respect to a tolerance  $t_j$ , it is defined as

$$s(\beta, t_j) = \frac{\partial Y(\mathbf{t})}{\partial t_j}. \quad (23)$$

Now that  $Y(\mathbf{t})$  is approximated by  $\Phi(\beta)$  (by step (iii) in tolerance analysis for nonlinear design functions), (23) can be approximated by  $\frac{\partial \Phi(\beta)}{\partial t_j}$ . (For a linear design function,  $\frac{\partial \Phi(\beta)}{\partial t_j}$  is the exact yield sensitivity.)

Expanding  $s(\beta, t_j)$  by the chain rule reveals the search direction:

$$s(\beta, t_j) = \frac{\partial \Phi(\beta)}{\partial \beta} \frac{\partial \beta}{\partial \sigma_j} \frac{\partial \sigma_j}{\partial t_j} = \phi(\beta) * \left\{ - \frac{x_j^* - \mu_j}{\beta \sigma_j^2} \left( \sum_{i=1}^n \hat{\rho}_{ji} \frac{x_i^* - \mu_i}{\sigma_i} \right) \right\} * \frac{1}{6} \quad (24)$$

where  $\hat{\rho}_{ji}$  is the (i,j) element of the inverse matrix of the correlation matrix,

$x_j^*$  is the  $j$ -th coordinate of the expansion point, and

$\phi(\cdot)$  is the PDF of the standard normal distribution, i.e.,

$$\phi(\beta) = \frac{1}{\sqrt{2\pi}} \exp\left\{-\frac{\beta^2}{2}\right\}.$$

Substituting (22) into (24) and by assumption of independence of dimensions, the sensitivity of yield is rewritten as

$$s(\beta, t_j) = \phi(\beta) * \left\{ - \frac{(x_j^* - \mu_j)^2}{\beta \sigma_j^3} \right\} * \frac{1}{6}. \quad (25)$$

Equation (25) shows that the  $s(\beta, t_j)$  are always nonpositive. This means that there is a nonincreasing *monotonicity* between yield and tolerance. This monotonic property demonstrates the trade-off between tolerance and performance: the performance, which is implied by yield, increases as the tolerances are tightened.

#### 4.5. Lagrangean Multiplier

The impact of modifying yields constraints on the manufacturing cost, i.e., the objective value, is considered in this section. This is performed as a post-optimality analysis.

The partial derivative  $\frac{\partial C(\sigma_1, \dots, \sigma_n)}{\partial \lambda_i}$  at the optimal solution  $(\sigma_1^*, \dots, \sigma_n^*)^T$  is decomposed into two parts using the chain rule:

$$\frac{\partial C(\mathbf{t})}{\partial \lambda_i} = \frac{\partial C(\mathbf{t})}{\partial q_i} \frac{\partial q_i}{\partial \lambda_i} \quad (26)$$

The second part of this decomposition,  $\frac{\partial q_i}{\partial \lambda_i}$ , turns out to be:

$$\frac{\partial q_i}{\partial \lambda_i} = - \frac{\sqrt{2\pi}}{\exp\left\{-\frac{q_i^2}{2}\right\}}$$

since  $\lambda_i = 1 - \Phi(q_i)$ . The first part of (26),  $\frac{\partial C(\mathbf{t})}{\partial q_i}$ , called the "Lagrangean multiplier" in NLP has the following characteristic: if the constraint is inactive at the optimal solution, then the corresponding Lagrangean multiplier should be zero. This means that for an inactive constraint at

optimum the corresponding  $q_i$  can be modified slightly without incurring additional manufacturing cost.

In case that a constraint is active at optimum, the corresponding Lagrangean multiplier should be greater than zero. It is computed as:

$$\frac{\partial C(\mathbf{t})}{\partial q_i} = \sum_{j=1}^n \frac{\partial C(\mathbf{t})}{\partial \sigma_j} \frac{\partial \sigma_j}{\partial q_i} = \sum_{j=1}^n \frac{\partial C(\mathbf{t})}{\partial \sigma_j} \frac{\partial \sigma_j}{\partial \beta_i} \quad (27)$$

since the  $i$ -th constraint is assumed to be active, i.e.  $q_i = \beta_i$ . From equation (22),  $\frac{\partial \beta_i}{\partial \sigma_j} = -\frac{(x_{ij}^* - \mu_j)^2}{\beta_i \sigma_j^3}$ , where  $x_{ij}^*$  is the  $j$ -th expansion point for  $\beta_i$ , (27)

becomes

$$\frac{\partial C(\mathbf{t})}{\partial q_i} = \sum_{j=1}^n \frac{\partial C(\mathbf{t})}{\partial \sigma_j} \left\{ -\frac{\beta_i \sigma_j^3}{(x_{ij}^* - \mu_j)^2} \right\}. \quad (28)$$

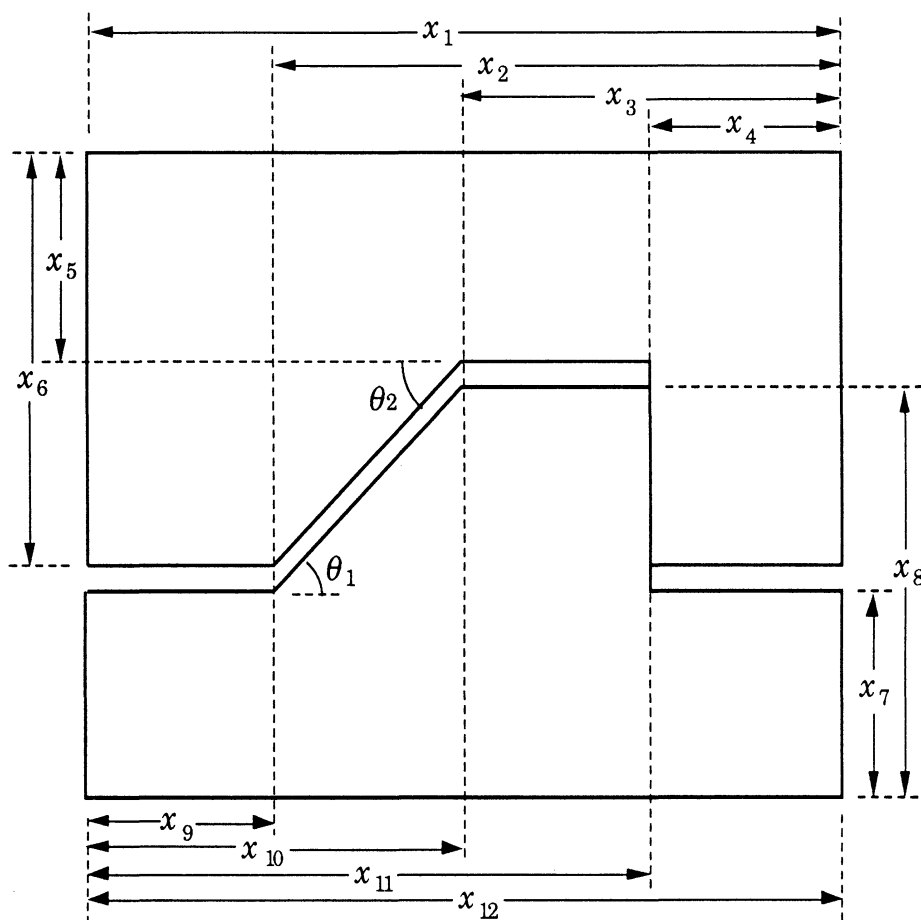
Based on the inverse-squared model, i.e.,  $C(\mathbf{t}) = \sum_{k=1}^n \left\{ \frac{a_k}{(6\sigma_k)^2} + f_k \right\}$ , (28)

becomes

$$\frac{\partial C(\mathbf{t})}{\partial q_i} = \sum_{j=1}^n \frac{-a_j}{18\sigma_j^3} \left\{ -\frac{\beta_i \sigma_j^3}{(x_{ij}^* - \mu_j)^2} \right\} = \sum_{j=1}^n \frac{a_j \beta_i}{18(x_{ij}^* - \mu_j)^2}.$$

## 5. Implementation

An assembly of two parts with twelve dimensions is shown in Figure 9(a). Six design functions, linear and non-linear, are given in Figure 9(b). Design functions  $F_1(X)$  and  $F_2(X)$  represent the vertical and the horizontal clearance conditions of the two parts. Design functions  $F_3(X)$  and  $F_4(X)$  post the restriction on the difference between angles  $\theta_1$  and  $\theta_2$  to ensure feasibility of assembly. Design functions  $F_5(X)$  and  $F_6(X)$  give the requirements for the size difference of those two parts.



(a)

$$F_1(\mathbf{X}) = (x_6 - x_5) - (x_8 - x_7)$$

$$F_2(\mathbf{X}) = (x_3 - x_4) - (x_{11} - x_{10})$$

$$F_3(\mathbf{X}) = (x_8 - x_7)(x_2 - x_3) - (x_6 - x_5)(x_{10} - x_9) \\ + \tan(\pi/180) * \{(x_{10} - x_9)(x_2 - x_3) + (x_8 - x_7)(x_6 - x_5)\}$$

$$F_4(\mathbf{X}) = (x_6 - x_5)(x_{10} - x_9) - (x_8 - x_7)(x_2 - x_3) \\ + \tan(\pi/180) * \{(x_{10} - x_9)(x_2 - x_3) + (x_8 - x_7)(x_6 - x_5)\}$$

$$F_5(\mathbf{X}) = -x_1 + x_{12} + 0.01$$

$$F_6(\mathbf{X}) = x_1 - x_{12} + 0.01$$

(b)

**Figure 9. Examples of Linear and Nonlinear Design Functions.**

The nominal dimensions are given as  $\mathbf{X}^T = (50.0, 40.00125, 20.05, 9.9985, 9.9985, 30.0, 10.0, 30.0, 10.05, 30.0, 40.0, 50.0)$ . The cost function with respect to the tolerance of each dimension is

$$C(\sigma_i) = \frac{a_i \times 10^{-3}}{(6\sigma_i)^{b_i}} .$$

The coefficients of the respective cost functions are:  $a_1 = 0.2$ ,  $a_2 = 1.0$ ,  $a_3 = a_4 = 0.015$ ,  $a_5 = 0.008$ ,  $a_6 = 0.009$ ,  $a_7 = 0.008$ ,  $a_8 = 0.006$ ,  $a_9 = 1.0$ ,  $a_{10} = 0.01$ ,  $a_{11} = 0.015$ , and  $a_{12} = 0.2$ ; and  $b_1 = \dots = b_{12} = 2.0$ .

The algorithm given in this paper is implemented in PASCAL and runs on the IBM PC. For the algorithms to be practical in an interactive design environment, attention is given to speed - in particular, the computation for the Jacobian matrix. The initial tolerances are assigned in accordance to ANSI-Y14.5M for loose fit.

It took 5.25 CPU seconds to obtain the optimal solution, and the resulted tolerances are shown in Table 2; each stack-up condition is satisfied with 95% confidence if the dimensions  $x_1$  to  $x_{12}$  are manufactured within the tolerances obtained. To cross check the algorithm, dimensions  $x_2$  and  $x_9$  are intentionally assigned with higher manufacturing costs, i.e., they are harder to be manufactured than the other dimensions. The tolerances computed are consistent with the manufacturing costs as much looser tolerances are assigned to dimensions  $x_2$  and  $x_9$ .

Dimensions	Cost function coefficients		Tolerances
	$a_i$	$b_i$	
$x_1$	0.2	2.0	0.0093
$x_2$	1.0	2.0	0.6427
$x_3$	0.015	2.0	0.0337
$x_4$	0.015	2.0	0.0337
$x_5$	0.008	2.0	0.0010
$x_6$	0.009	2.0	0.0011
$x_7$	0.008	2.0	0.0010
$x_8$	0.006	2.0	0.0008
$x_9$	1.0	2.0	0.6433
$x_{10}$	0.01	2.0	0.0337
$x_{11}$	0.015	2.0	0.0337
$x_{12}$	0.2	2.0	0.0093

**Table 2. Result of the example.**

## 6. Concluding Remarks

A general framework for tolerance synthesis based on least manufacturing cost has been presented. The algorithm for probabilistic tolerancing has been developed based on the notion of feasible directions. An analytic result of yield sensitivity is used to speed up the computation of the Jacobian matrix inside the optimization loop. Compared to a previously established landmark [14], this new algorithm produces a solution with more than 10 times reduction in CPU time. Also, the post-optimality analysis of the algorithm enables a designer to verify design intention with ease.



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## Appendix : Transformation to the Standard Coordinates

The transformation into the standard coordinates is accomplished by taking the following four steps [6]:

Step 1: Translate into the origin by  $\mathbf{x}^0 = \mathbf{x} - \boldsymbol{\mu}$ ;

Step 2: Diagonalize the given covariance matrix  $\mathbf{V}$  through the operation

$$\mathbf{P}^T \mathbf{V} \mathbf{P} = \mathbf{V}_z, \text{ where } \mathbf{P} \text{ is the orthogonal matrix;}$$

Step 3: Orthogonally transform by  $\mathbf{z}^0 = \mathbf{P}^T \mathbf{x}^0$ ;

Step 4: Standardize by  $\mathbf{z} = \mathbf{D}^{-1} \mathbf{z}^0$ , where  $\mathbf{D} \mathbf{D}^T = \mathbf{V}_z$ .



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