

—NOTES—

A GEOMETRICAL INTERPRETATION OF THE
RELAXATION METHOD*By J. L. SYNGE (*The Ohio State University*)

Let a_{ij}, B_i ($i, j = 1, 2, \dots, n$) be given constants such that $a_{ij} = a_{ji}$ and $\sum_{i,j=1}^n a_{ij}x_i x_j$ is a positive definite form. Consider the equations

$$\sum_{j=1}^n a_{ij}x_j - B_i = 0 \quad (i = 1, 2, \dots, n). \quad (1)$$

The solution is easily expressed as a set of quotients of determinants. However, as n increases, the task of calculating the determinants becomes excessively burdensome. The relaxation method¹ provides a set of easy steps by which the solution of (1) is approached. The method has been compactly described by Temple.²

The purpose of the present note is to give a geometrical description of the relaxation method. For the trivial case $n = 2$ the geometrical description may be displayed accurately in a diagram. For $n = 3$ a model may be visualized. For $n > 3$ we pass beyond the region of simple concrete geometrical representation, but in many ways geometry in an n -space is closely analogous to geometry in 2-space or 3-space, and the geometrical description continues to serve as a general guide to procedure.

Let us regard x_i as rectangular Cartesian coordinates in a Euclidean n -space. Let us define

$$H(x) = \frac{1}{2} \sum_{i,j=1}^n a_{ij}x_i x_j - \sum_{i=1}^n B_i x_i. \quad (2)$$

The equation $H(x) = \text{const.}$ represents a family of ellipsoids E ; these ellipsoids have a common center, common directions for their principal axes, and common values for the ratios of their principal axes. They form, in fact, a family of similar and similarly situated ellipsoids.

The equations (1) represent a set of planes (i.e., flats of $n-1$ dimensions). The point of intersection of these planes is the common center G of E . Thus the problem of solving (1) is the problem of finding the center of an ellipsoid when its equation is given.

It is important to note that $H(x)$ takes a minimum value at G . H is constant over each ellipsoid, and increases steadily as we pass out from G .

It is not possible to define precisely what procedures are to be regarded as permissible. It is a question of ease of computation. Let us follow Southwell and consider an approach to G by steps each of which is parallel to one of the axes of coordinates x_i .

Fig. 1 shows Southwell's procedure. It is a schematic diagram in which the ellip-

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¹ R. V. Southwell, *Relaxation methods in engineering*, Oxford, 1940.

² G. Temple, *Proc. Roy. Soc. London A*, **169**, 476-500 (1939).

soids are represented by circles. (The ellipsoids can of course be transformed into concentric spheres by a linear transformation, which however destroys the orthogonality of perpendicular lines.) We start with an arbitrary point P_0 (the zero approximation). Let E_0 be the ellipsoid which passes through P_0 . Through P_0 we draw a straight line L parallel to one of the coordinate axes. Let Q_1 be the second point in which L cuts E_0 . Let P_1 be the middle point of the chord P_0Q_1 . Then P_1 is the first approximation.

Since the ellipsoid is a convex surface, P_1 lies inside E_0 and so $H(P_1) < H(P_0)$. Moreover it is easy to see that P_0Q_1 is tangent at P_1 to the ellipsoid E_1 which passes through P_1 . Thus, of all points on the chord P_0Q_1 , the point P_1 gives the smallest value of H .

The process is repeated, starting from P_1 . The second approximation P_2 is the middle point of a chord P_1Q_2 of E_1 , drawn parallel to another of the coordinate axes. In this way we get a sequence of points P_0, P_1, \dots . The success of the method depends on the rapidity of the convergence of this sequence to G .

In one important respect the above procedure is incompletely defined. When we have reached P_m , in which of the directions defined by the coordinate axes are we to proceed in order to get P_{m+1} ? There are n coordinate axes. Of these one cannot be used, viz., that which gave the direction of the step $P_{m-1}P_m$. But, of the remaining $n-1$ directions, which should we use?

Gaskell³ has suggested the following plan. Write

$$C_i(x) = \sum_{j=1}^n a_{ij}x_j - B_i. \tag{3}$$

Having reached the point P_m , we calculate the quantities $C_i(P_m)$. Let $C_k(P_m)$ be the greatest of these in absolute value. Then we choose for the step P_mP_{m+1} the direction of the axis of x_k .

This procedure is called the liquidation of the greatest error, since we obtain $C_k(P_{m+1})=0$. It is interesting to see how this result fits into the geometrical discussion. The plane $C_k(x)=0$ is the plane through G conjugate to the direction of the axis x_k . The line P_mP_{m+1} is parallel to this axis and tangent at P_{m+1} to one of the ellipsoids, E_{m+1} . But the point of contact of a line with an ellipsoid lies on the central plane conjugate to the direction of the line. Hence P_{m+1} lies on $C_k(x)=0$, i.e., $C_k(P_{m+1})=0$.

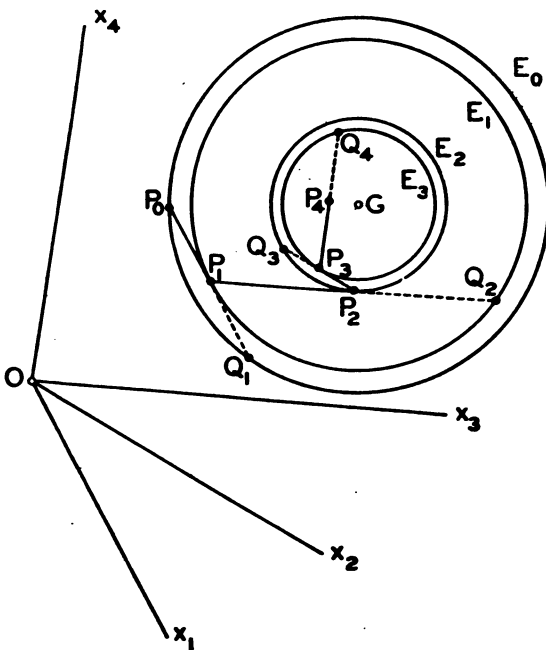


FIG. 1.

³ R. E. Gaskell, Quarterly of Applied Mathematics, 1, 237-249 (1943).

But it may well be asked whether the quantities C_i themselves possess any deep significance. It is true that G satisfies $C_i(x) = 0$, but the quantity $C_i(x)$ for a general point does not represent the perpendicular distance of that point from the plane $C_i(x) = 0$. This perpendicular distance is

$$f_i(x) = \frac{|C_i(x)|}{\left(\sum_{j=1}^n a_{ij}^2\right)^{1/2}}. \tag{4}$$

Should we not liquidate the greatest p_i rather than the greatest C_i ? Or is there a better plan than either?

The following plan is suggested. Having reached the point P_m , we have an option on $n - 1$ next points. Each of these points lies on an ellipsoid of the family E . Choose that point which lies on the innermost ellipsoid. This is equivalent to saying: Choose that point which gives the smallest value to H .

Now⁴ for a step in the direction of the axis x_i the decrease in H is $\frac{1}{2}C_i^2/a_{ii}$. This is to be made as great as possible, and so we should pick the direction of the step $P_m P_{m+1}$ according to the following rule: Proceed in the direction of the axis of x_k where C_k^2/a_{kk} is the greatest of the quantities C_i^2/a_{ii} ($i = 1, 2, \dots, n$).

Thus C_i^2/a_{ii} is made the criterion rather than Gaskell's C_i . The calculation of the former quantities involves slightly more computation, but this may be taken care of by making the initial transformation

$$x'_i = (a_{ii})^{1/2} x_i. \tag{5}$$

Then

$$H = \frac{1}{2} \sum_{i,j=1}^n a'_{ij} x'_i x'_j - \sum_{i=1}^n B'_i x'_i, \tag{6}$$

where

$$a'_{ii} = 1, \quad a'_{ij} = a_{ij}/(a_{ii}a_{jj})^{1/2}, \quad B'_i = B_i/(a_{ii})^{1/2}. \tag{7}$$

Now, with

$$C'_i(x') = \sum_{j=1}^n a'_{ij} x'_j - B'_i, \tag{8}$$

the criterion for the direction of the next displacement is $C_i'^2$ or $|C'_i|$, the same as Gaskell's. Moreover the transformation from P_m to P_{m+1} is now extremely simple. It is⁵

$$\begin{aligned} P_m: & \quad x'_1, \dots, x'_k, \dots, x'_n; \\ P_{m+1}: & \quad x'_1, \dots, x'_k - C'_k(x'), \dots, x'_n. \end{aligned}$$

⁴ R. E. Gaskell, loc. cit., Eq. (23).

⁵ R. E. Gaskell, loc. cit.