

PROBLEM (3). Maximize

$$\begin{aligned} & \sum_{r=1}^R \bar{c}_1^r \lambda_1^r + \sum_{s=1}^S \bar{c}_2^s \lambda_2^s + \cdots + \sum_{t=1}^T \bar{c}_N^t \lambda_N^t \\ & \sum_r \bar{A}_1^r \lambda_1^r + \sum_s \bar{A}_2^s \lambda_2^s + \cdots + \sum_t \bar{A}_N^t \lambda_N^t \leq b_0 \\ & \sum_r \delta_1^r \lambda_1^r = 1 \\ & \sum_s \delta_2^s \lambda_2^s = 1 \\ & \vdots \\ & \sum_t \delta_N^t \lambda_N^t = 1 \end{aligned}$$

where $\bar{c}_j^k = c_j x_j^k$, $\bar{A}_j^k = A_j x_j^k$, $\sigma_j^k = \begin{cases} 1 \\ 0 \end{cases}$ if x_j^k is a $\begin{cases} \text{basic} \\ \text{homogeneous} \end{cases}$ solution of $B_j x_j \leq b_j$, $x_j \geq 0$. The summations are taken over all extreme points and extreme rays of the several polyhedra. The λ_j^k are the system variables and are required to be non-negative; the other symbols are coefficients, though they are generated, not input.

Decomposition algorithms generate coefficient columns for problem (3) as needed by solving subprograms whose objective functions are linear (in the case of the primal algorithm) or quotients of linear functions (in the case of the primal-dual algorithm). Clearly, the efficiency of the algorithms depends largely upon the efficiency in generating the subprogram solutions. We have further comment upon this later in the paper when we contrast the performance of the two algorithms on this problem.

In terms of the system in problem (2) we notice several features which distinguish our blending problem: (a) the b_j vectors are null vectors, with the exception of b_0 , (b) all of the A_j matrices are identity matrices, and (c) the c_j vectors are sum vectors multiplied by constants.

The decomposition algorithms exploit the macro structure; the question now is: can the algorithms be improved by taking explicit account of the micro structure represented by the three features?

(a) When each of the b_j vectors ($j = 1, \dots, N$) is a null vector, the only basic solution of $B_j x_j \leq b_j$ is the trivial solution $x_j = 0$. This means that all of the N rows $\sum_k \delta_j^k \lambda_j^k = 1$ ($j = 1, \dots, N$) in (3) may be dropped, thus further reducing the number of rows and, hence, the basis size in problem (3).³ When N is quite large, as if we were considering a large number of blends, this reduction could be helpful. It will still be necessary to solve the subprograms, however. This reduction will work for either the primal or primal-dual decomposition algorithms.

(b) When the A_j matrices are identity matrices the transformations $A_j x_j^k = \bar{A}_j^k$ are obviously unnecessary and the coefficient columns can be added to the master program just as they are obtained from the subprograms.

³ In such cases all $\delta_j^k = 0$ except that $\delta_j^k = 1$ for the trivial basic solution which implies that its associated $\lambda_j^k = 1$. Hence the constraint may be dropped.

More importantly, in order to extract the optimal solution in terms of the original x_j vectors, it is generally necessary to maintain the solution x_j^k from the subprogram that generated the coefficient column that has just been added to the master program. The reason for this is contained in the relationship of the x_j vectors to the x_j^k and λ_j^k , $x_j = \sum x_j^k \lambda_j^k$ where the only λ_j^k and x_j^k of interest are those in the terminal basic solution of problem (3). Clearly the x_j^k are available if they are stored after being generated, as in the revised Simplex method, before they are transformed by pivoting. This feature also holds for both the primal and primal-dual decomposition algorithms.

(c) Ordinarily, by itself, this feature cannot be exploited, but in conjunction with (a) and (b) it can lead to a substantial improvement in the primal-dual algorithm. As in the general primal-dual algorithm, the decomposition version needs a feasible solution for problem (4), the dual of problem (3).

PROBLEM (4). Minimize

$$\begin{aligned} & \sigma \cdot b_0 + \sum_{j=1}^n \alpha_j \\ & \sigma A_j x_j^k + \beta_j \geq c_j x_j^k \quad j = 1, \dots, N. \\ & \sigma \geq 0. \end{aligned}$$

PROBLEM (5). Maximize

$$\begin{aligned} & \sum_{j=1}^N c_j x_j \\ & \sum_{j=1}^N A_j x_j \leq b_0, \quad x_j \geq 0 \quad j = 1, \dots, N. \end{aligned}$$

The recommended approach for general decomposition problems is to obtain optimal-dual variables associated with an optimal solution of problem (5) and from them obtain an $(m_0 + N)$ vector satisfying the dual constraints in problem (4).⁴

By virtue of (a), the sum rows of (3) are eliminated. Thus, we need only an m_0 -vector σ satisfying:

$$\sigma A_j x_j^k \geq c_j^k x_j^k \quad (6)$$

and when the A_j are identity matrices⁵ this reduces to

$$\sigma x_j^k \geq c_j x_j^k. \quad (7)$$

Clearly, to satisfy (7), it is sufficient to find a vector $\sigma \geq c_j$ for all $j = 1, \dots, N$.⁵ Thus, when (c) prevails, σ may be determined at the time of input by simply determining which c_j vector has the largest constant and setting all the elements of σ at that value initially. Thus, for our blending problem

$$\sigma_1 = \sigma_2 = \sigma_3 = \sigma_4 = 1.494.$$

This feature has two redeeming characteristics beyond avoiding a lot of computation:

(d) It establishes an upper bound on the maximum of

⁴ See [1] for a detailed outline of this procedure.

⁵ When the A_j are identity matrices, it must be the case that $m_0 = n_j$ for all j .

$\sum_{j=1}^N c_j x_j$ at the value of $\sigma \cdot b_0 = 1.494 \cdot 11,833 = 17,678$ (the optimal value is at 15,425).

(e) It furnishes an immediate indication concerning which subprogram to solve in order to obtain candidates for the initial restricted master program (in the blending problem start with the third subprogram).

Point (e) relates to the relative efficiency of the primal-dual and primal algorithms, at least as far as their performance on this problem is concerned. The primal-dual algorithm for general linear programs, and also for special network structures such as the capacitated transportation problem, are known to have a tighter selection criterion than the primal algorithms. This is because the primal-dual algorithm must move the solution towards primal feasibility while simultaneously causing the dual feasible solutions to move towards optimality and maintaining complementary slackness between primal and dual. This parallelism holds true in the decomposition algorithms as well. To illustrate the lower stringency of the primal decomposition algorithm, note that any candidate vector x_j^k from any of the subprograms which satisfies $(\sigma A_j - c_j) \cdot x_j^k < 0$ (where σ is the current vector of multipliers) can be introduced into the master program and cause an increase (under conditions of nondegeneracy) in the function to be maximized. In the blending problem when starting from a full slack basis $\sigma = (0, 0, 0, 0)$ and because all of the c_j vectors have no negative elements,⁶ any $x_j^k \geq 0$ satisfying $B_j x_j \leq 0$ also satisfies $(\sigma A_j - c_j) x_j^k < 0$. Because of the homogeneous nature of the solutions $(\sigma A_j - c_j) x_j^k$ can be made even more negative by multiplying by a positive scalar. Therefore, there is no way of distinguishing in this case, as in the usual primal decomposition algorithm, which candidate is "best" in the sense of the usual Simplex criterion of minimal $(\sigma A_j - c_j) x_j$.

In the blending problem this means that each of the three subprograms could furnish a possible candidate for the first master pivot operation. In anticipation of this "indeterminate" situation we might suspect that candidates drawn from the subprograms with the higher c_j values—that is, subprograms 2 and 3 with values .889 and 1.494—would tend to supply an optimal solution. Post-optimal analysis of this problem in fact confirms this suspicion. It turns out that x_1, x_2, x_3, x_4 can never appear in any optimal solution. This means that any candidate from subprogram 1 which is introduced into the master program must later be rejected.

Thus, there are three major solution paths, depending upon which subprogram furnishes the first candidate. The number of candidates which had to be generated in each case was 11, 7, and 4, respectively. We remark again that the primal decomposition algorithm, by itself, has no means of determining which of these paths is "best." Only

after the fact can we confirm our earlier suspicion. It is now interesting to compare these numbers with the four candidates generated by the primal-dual algorithm, starting with $\sigma = 1.494 (1, 1, 1, 1)$, or alternatively the total of five candidates with the initial $\sigma = (0, 0, 0, 0)$.

Hence the primal-dual algorithm was at least as efficient as the primal, two-phase algorithm in spite of the fact that no Phase I was required because a full slack basis was initially available. This illustrates the point that the apparent disadvantage of using a full artificial basis may be deceptive and that methods employing artificial variables, such as primal-dual methods, may be very efficient when few of the available slacks will appear in optimal solutions.

One final point of computational interest: even for this small problem the total computing time of each decomposition algorithm compared favorably with the time to solve as a standard linear program. This was the case in spite of the fact that both decomposition algorithms were "simulated," not automatically programmed. The Dartmouth College Computation Center is to be commended for its excellent time sharing system which provided the flexibility to carry out the many offline calculations without "turnaround" problems, and thus avoid the necessity of large and complex automated computer codes.

It is interesting to note that in making several simplifying assumptions in order to reduce the basis size of this problem, Garvin could make the statement that "the matrix of our problem has structure . . . characteristic of blending problems. . . . It is natural to inquire whether it is possible to take advantage of the matrix structure. . . ." Ironically, within six months the Decomposition Principle had been published (in 1960).

We hope, indeed, that we have shown that this problem has "structure," both *macro* and *micro*. Attention to both kinds of structure can pay dividends, particularly when dealing with calculations of routine frequency.

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⁶ Notationally $x_j \geq 0$, instead of $x_j \geq 0$, excludes $x_j = 0$.